

Supporting Information

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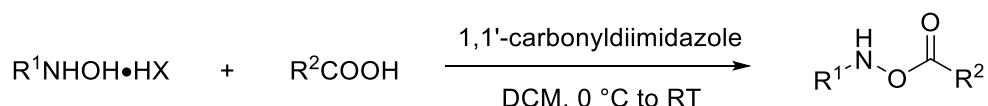
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1. General information

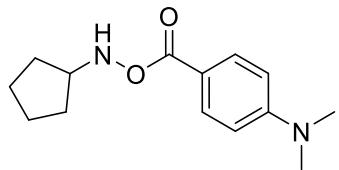
All commercial reagents were purchased from Sigma-Aldrich, Strem, Acros, TCI or Alfa Aesar and used as such unless stated otherwise. Solvents (Anhydrous and under inert atmosphere) were collected from The Solvent purification system by M BRAUN and used under standard schlenk technique. NMR spectra were recorded on Bruker Avance 300 MHz and Bruker ARX 400 MHz spectrometers. Multiplets were assigned as s(singlet), d (doublet), t (triplet), q (quartet), dd (doublet of doublet), m (multiplet) and br. s (broad singlet). Coupling constants reported to 0.5 or 1.0 Hz accuracy. GC-yields were calculated using hexadecane as internal standard. All measurements were carried out at room temperature unless otherwise stated. Electron impact (EI) mass spectra were recorded on AMD 402 mass spectrometer (70 eV). The data are given as mass units per charge (m/z). Gas chromatography analysis was performed on an Agilent HP-7890A instrument with an FID detector and HP-5 capillary column (polydimethylsiloxane with 5% phenyl groups, 30 m, 0.32 mm i.d., 0.25 µm film thickness) using argon as carrier gas. The products were isolated from the reaction mixture by column chromatography on silica gel 60, 0.063-0.2 mm, 70-230 mesh (Merck). For chiral HPLC-analysis a device Agilent 1100 Series was used.

2. General Procedures

2.1 General Procedure I for the Preparation of Amine Electrophiles¹



To a solution of benzoic acid (1.2 equiv) in CH₂Cl₂ (0.3 M) was added CDI (1.2 equiv) at 0 °C. After gas generation was over, hydroxyl amine was added and the mixture was stirred at room temperature for 12 h. The reaction mixture was monitored by TLC. After disappearance of starting material, the reaction mixture was quenched with distilled water and extracted with CH₂Cl₂ (3 times). The combined organic layer was washed with aqueous NaHCO₃ (2 times) and brine (1 time), dried over Na₂SO₄. After removal of solvent, the residue was purified by flash chromatography on silica gel to give a corresponding desired product compound. Hydroxylamine was prepared from estrone by procedure reported by Qiu Wang.²

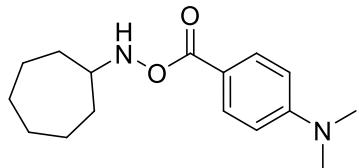


4-(((Cyclopentylamino)oxy)carbonyl)-N,N-dimethylaniline (2f)

¹H NMR (300 MHz, CDCl₃) δ 7.99 – 7.80 (m, 2H), 6.71 – 6.43 (m, 2H), 3.64 (tt, *J* = 6.5, 3.0 Hz, 1H), 3.02 (s, 6H), 1.94 – 1.46 (m, 8H).

¹³C NMR (75 MHz, CDCl₃) δ 167.5, 153.5, 131.1, 114.9, 110.7, 62.2, 40.0, 30.4, 24.4.

HRMS (ESI-TOF) m/z: Calcd. for C₁₄H₂₀N₂O₂ [M+H]⁺: 249.1603, Found: 249.1609.

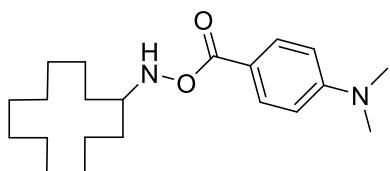


4-(((Cycloheptylamino)oxy)carbonyl)-N,N-dimethylaniline (2g)

¹H NMR (300 MHz, CDCl₃) δ 7.97 – 7.74 (m, 2H), 6.65 – 6.48 (m, 2H), 3.18 – 3.03 (m, 1H), 2.96 (s, 6H), 1.95 – 1.81 (m, 2H), 1.70 – 1.57 (m, 2H), 1.54 – 1.33 (m, 8H).

¹³C NMR (75 MHz, CDCl₃) δ 167.5, 153.5, 131.1, 115.0, 110.8, 61.9, 40.0, 31.6, 28.7, 24.3.

HRMS (ESI-TOF) m/z: Calcd. for C₁₆H₂₄N₂O₂ [M+H]⁺: 277.1916, Found: 277.1912.

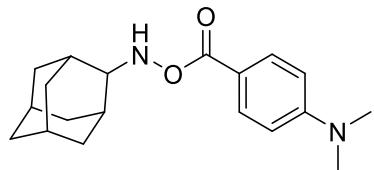


4-(((Cyclododecylamino)oxy)carbonyl)-N,N-dimethylaniline (2h)

¹H NMR (300 MHz, CDCl₃) δ 7.93 – 7.85 (m, 2H), 7.79 (s, 1H), 6.69 – 6.53 (m, 2H), 3.19 (s, 1H), 3.03 (s, 6H), 1.68 – 1.16 (m, 22H).

¹³C NMR (75 MHz, CDCl₃) δ 167.5, 153.5, 131.0, 115.0, 110.8, 58.0, 40.0, 27.5, 24.4, 23.7, 23.5, 23.4, 21.5.

HRMS (ESI-TOF) m/z: Calcd. for C₂₁H₃₄N₂O₂ [M+H]⁺: 347.2698, Found: 347.2702.

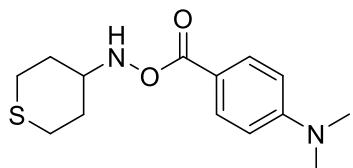


4-(((Adamantan-2-yl)amino)oxy)carbonyl)-N,N-dimethylaniline (2i)

¹H NMR (300 MHz, CDCl₃) δ 8.05 (s, 1H), 7.97 – 7.86 (m, 2H), 6.72 – 6.61 (m, 2H), 3.36 – 3.24 (m, 1H), 3.06 (s, 6H), 2.28 – 2.02 (m, 4H), 1.97 – 1.84 (m, 4H), 1.76 (s, 4H), 1.64 – 1.52 (m, 2H).

¹³C NMR (75 MHz, CDCl₃) δ 167.5, 153.5, 131.1, 115.1, 110.7, 64.6, 40.0, 37.7, 37.2, 31.5, 30.4, 27.7, 27.6.

HRMS (ESI-TOF) m/z: Calcd. for C₁₉H₂₆N₂O₂ [M+H]⁺: 315.2072, Found: 315.2079.

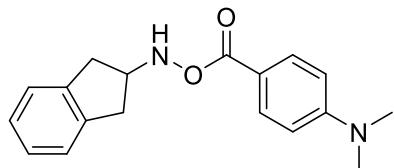


N,N-Dimethyl-4-(((tetrahydro-2H-thiopyran-4-yl)amino)oxy)carbonyl)aniline (2j)

¹H NMR (300 MHz, CDCl₃) δ 7.87 (d, *J* = 9.0 Hz, 2H), 6.64 (d, *J* = 9.0 Hz, 2H), 3.09 – 2.93 (m, 7H), 2.79 – 2.57 (m, 4H), 2.32 – 2.18 (m, 2H), 1.81 – 1.63 (m, 2H).

¹³C NMR (75 MHz, CDCl₃) δ 167.4, 153.6, 131.1, 114.4, 110.8, 59.1, 40.0, 31.6, 27.1.

HRMS (ESI-TOF) m/z: Calcd. for C₁₄H₂₀N₂O₂S [M+H]⁺: 303.1143, Found: 303.1141.

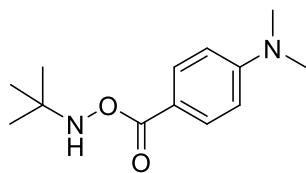


4-(((2,3-Dihydro-1H-inden-2-yl)amino)oxy)carbonyl)-N,N-dimethylaniline (2k)

¹H NMR (300 MHz, CDCl₃) δ 7.99 – 7.82 (m, 2H), 7.33 – 7.10 (m, 4H), 6.80 – 6.60 (m, 2H), 4.20 (tt, *J* = 7.0, 5.0 Hz, 1H), 3.25 (dd, *J* = 16.5, 7.0 Hz, 2H), 3.14 – 2.95 (m, 8H).

¹³C NMR (75 MHz, CDCl₃) δ 167.3, 153.6, 141.0, 131.1, 126.7, 124.8, 114.6, 110.8, 61.5, 40.1, 37.3.

HRMS (ESI-TOF) m/z: Calcd. for C₁₈H₂₀N₂O₂ [M+H]⁺: 297.1603, Found: 297.1607.

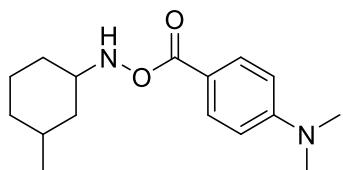


4-(((*tert*-Butylamino)oxy)carbonyl)-*N,N*-dimethylaniline (2l**)**

¹H NMR (300 MHz, CDCl₃) δ 7.82 (d, *J* = 9.0 Hz, 2H), 6.57 (d, *J* = 9.0 Hz, 3H), 2.95 (s, 6H), 1.13 (s, 9H).

¹³C NMR (75 MHz, CDCl₃) δ 167.3, 153.5, 131.0, 114.8, 110.8, 55.9, 40.0, 26.6.

HRMS (ESI-TOF) m/z: Calcd. for C₁₃H₂₀N₂O₂ [M+H]⁺: 237.1603, Found: 237.1605.

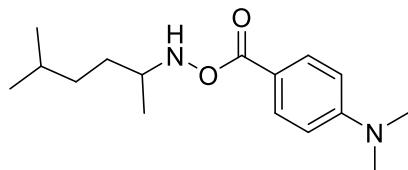


N,N-Dimethyl-4-(((3-methylcyclohexyl)amino)oxy)carbonyl)aniline (2m**)**

¹H NMR (300 MHz, CDCl₃) δ 8.02 – 7.83 (m, 2H), 6.70 – 6.49 (m, 2H), 3.34 (dd, *J* = 6.0, 2.5 Hz, 1H), 3.04 (d, *J* = 1.0 Hz, 6H), 1.98 – 1.04 (m, 8H), 0.93 (dd, *J* = 6.5, 5.5 Hz, 3H).

¹³C NMR (75 MHz, CDCl₃) δ 167.5, 167.5, 153.5, 153.5, 131.1, 131.0, 114.9, 114.9, 110.7, 60.2, 55.8, 40.0, 39.0, 36.6, 34.6, 33.9, 31.4, 30.1, 28.8, 26.9, 24.4, 22.6, 21.5, 20.3, 14.2.

HRMS (ESI-TOF) m/z: Calcd. for C₁₆H₂₄N₂O₂ [M+H]⁺: 277.1916, Found: 277.1919.

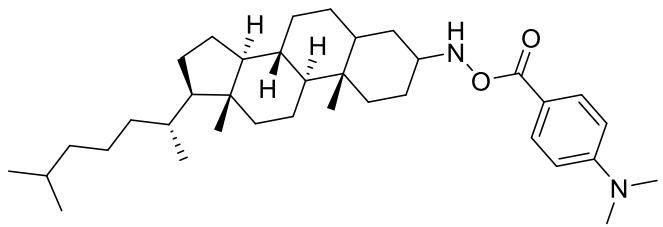


N,N-Dimethyl-4-(((5-methylhexan-2-yl)amino)oxy)carbonyl)aniline (2n**)**

¹H NMR (400 MHz, CDCl₃) δ 7.91 (d, *J* = 9.0 Hz, 2H), 6.67 (d, *J* = 9.0 Hz, 2H), 3.14 (q, *J* = 6.5 Hz, 1H), 3.06 (s, 6H), 1.69 – 1.52 (m, 2H), 1.43 – 1.36 (m, 1H), 1.34 – 1.26 (m, 2H), 1.19 (d, *J* = 6.5 Hz, 3H), 0.91 (dd, *J* = 65, 1.0 Hz, 6H).

¹³C NMR (101 MHz, CDCl₃) δ 167.5, 153.5, 131.0, 114.9, 110.8, 57.0, 40.0, 35.0, 31.9, 28.2, 26.9, 22.6, 22.5, 18.2.

HRMS (ESI-TOF) m/z: Calcd. for C₁₆H₂₆N₂O₂ [M+H]⁺: 279.2072, Found: 279.2072.



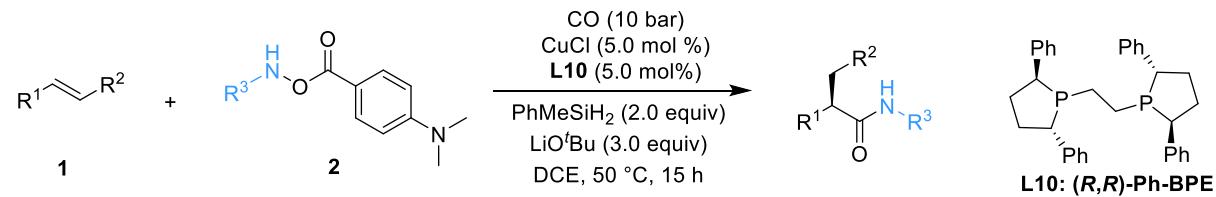
4-((((8*R*,9*S*,10*S*,13*R*,14*S*,17*R*)-10,13-Dimethyl-17-((*R*)-6-methylheptan-2-yl)hexadecahydro-1*H*-cyclopenta[a]phenanthren-3-yl)amino)oxy)carbonyl)-*N,N*-dimethylaniline (2o**)**

¹H NMR (400 MHz, CDCl₃) δ 7.92 (d, *J* = 9.0 Hz, 2H), 6.67 (dd, *J* = 9.0, 1.5 Hz, 2H), 3.06 (d, *J* = 1.0 Hz, 6H), 1.98 (d, *J* = 12.5 Hz, 1H), 1.87 – 1.75 (m, 3H), 1.71 – 1.48 (m, 7H), 1.45 (s, 3H), 1.40 – 1.23 (m, 9H), 1.19 – 1.08 (m, 6H), 1.06 – 0.97 (m, 3H), 0.92 (dd, *J* = 6.5, 2.0 Hz, 3H), 0.88 (dt, *J* = 6.5, 2.0 Hz, 7H), 0.83 (d, *J* = 2.0 Hz, 3H), 0.67 (d, *J* = 1.0 Hz, 3H).

¹³C NMR (101 MHz, CDCl₃) δ 167.5, 167.5, 153.5, 153.5, 131.1, 115.0, 114.9, 110.8, 110.7, 60.5, 56.5, 56.3, 56.2, 55.6, 54.4, 54.2, 45.0, 42.6, 40.0, 39.8, 39.5, 39.5, 37.0, 36.2, 36.0, 35.8, 35.8, 35.5, 32.8, 32.7, 32.1, 31.9, 31.1, 28.8, 28.7, 28.3, 28.0, 26.9, 26.0, 24.2, 23.9, 23.8, 22.8, 22.6, 21.2, 20.8, 18.7, 12.3, 12.1, 12.1, 11.4.

HRMS (ESI-TOF) m/z: Calcd. for C₃₆H₅₈N₂O₂ [M+H]⁺: 551.4576, Found: 551.4580.

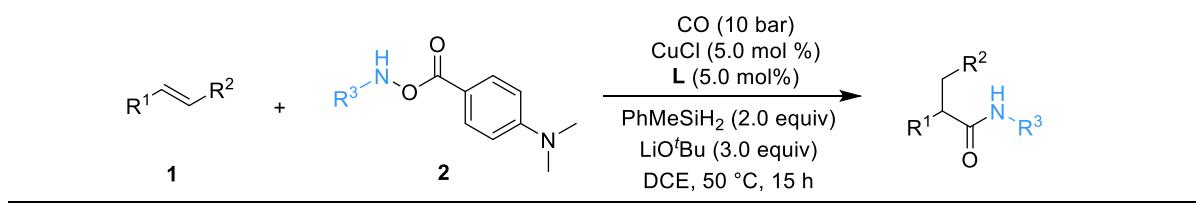
2.2 General Procedure II for Cu-catalyzed Asymmetric Hydroaminocarbonylation Reaction



A vial (4 mL) was charged with CuCl (5.0 mol%), (*R*, *R*)-Ph-BPE (5.0 mol%), LiO'Bu (24 mg, 3.0 equiv), and a stirring bar. The vial was closed by PTFE/white rubber septum (Wheaton 13 mm Septa) and phenolic cap and connected with atmosphere with a needle. The vial was evacuated under vacuum and recharged with argon for three times. Then, DCE (0.25 mL) were injected under argon by using a syringe, and stirred for 5 min at RT. After that, vinylarenes **1** (0.12 mmol, 1.2 equiv), MePhSiH₂ (0.2 mmol, 28 μL, 2.0 equiv) were added, then hydroxylamines **2** (0.1 mmol, 1.0 equiv) dissolved in 0.25 mL DCE was added by using a syringe, the vial (or several vials) was placed in an alloy plate, which was transferred into a 300 mL autoclave of the 4560 series from Parr Instruments. After flushing the autoclave three times

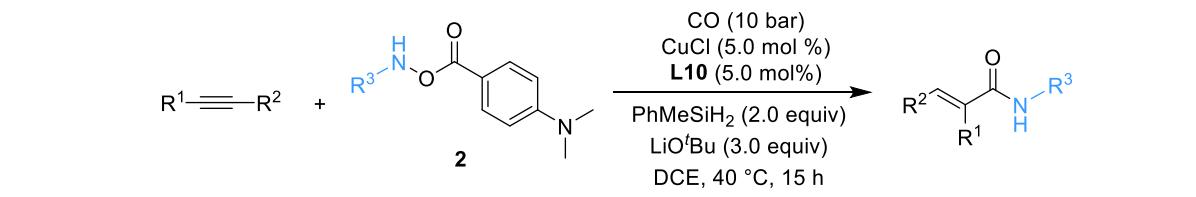
with CO, a pressure of 10 bar of CO was adjusted at ambient temperature. Then, the reaction was performed for 15 h at indicated temperature. After 15 hours, the autoclave was cooled down with ice water to room temperature and the pressure was released carefully. The solution was then filtered through celite and concentrated in *vacuo*. After that the residue was purified by column chromatography to afford the corresponding products. The resulting products was used to determine the enantiomeric excess by HPLC analysis using chiral stationary phases.

2.3 General Procedure III for Cu-catalyzed Racemic Hydroaminocarbonylation Reaction



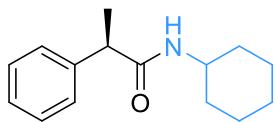
A vial (4 mL) was charged with CuCl (5.0 mol%), Xantphos (5.0 mol%) or racemic Ph-BPE phos (5.0 mol%), LiO*t*Bu (24 mg, 3.0 equiv), and a stirring bar. The vial was closed by PTFE/white rubber septum (Wheaton 13 mm Septa) and phenolic cap and connected with atmosphere with a needle. The vial was evacuated under vacuum and recharged with argon for three times. Then, DCE (0.5 mL) were injected under argon by using a syringe. After that, styrenes **1** (0.12 mmol, 1.2 equiv), hydroxylamines **2** (0.1 mmol, 1.0 equiv) and MePhSiH₂ (0.2 mmol, 28 μL, 2.0 equiv) were added, and the vial (or several vials) was placed in an alloy plate, which was transferred into a 300 mL autoclave of the 4560 series from Parr Instruments. After flushing the autoclave three times with CO, a pressure of 10 bar of CO was adjusted at ambient temperature. Then, the reaction was performed for 15 h at 50 °C. After 15 hours, the autoclave was cooled down with ice water to room temperature and the pressure was released carefully. The solution was then filtered through celite and concentrated *in vacuo*. After that the residue was purified by column chromatography to afford the corresponding racemic products.

2.4 General Procedure IV for Cu-catalyzed Hydroaminocarbonylation of Alkynes



A vial (4 mL) was charged with CuCl (5.0 mol%), Ph-BPE (5.0 mol%), LiO'Bu (24 mg, 3.0 equiv), and a stirring bar. The vial was closed by PTFE/white rubber septum (Wheaton 13 mm Septa) and phenolic cap and connected with atmosphere with a needle. The vial was evacuated under vacuum and recharged with argon for three times. Then, DCE (0.5 mL) were injected under argon by using a syringe. After that, alkynes (0.1 mmol, 1.0 equiv), hydroxylamines **2** (0.12 mmol, 1.2 equiv) and MePhSiH₂ (0.2 mmol, 28 μ L, 2.0 equiv) were added, and the vial (or several vials) was placed in an alloy plate, which was transferred into a 300 mL autoclave of the 4560 series from Parr Instruments. After flushing the autoclave three times with CO, a pressure of 10 bar of CO was adjusted at ambient temperature. Then, the reaction was performed for 15 h at 40 °C. After 15 hours, the autoclave was cooled down with ice water to room temperature and the pressure was released carefully. The solution was then filtered through celite and concentrated *in vacuo*. After that the residue was purified by column chromatography to afford the corresponding products.

3. Characterization Data



(*R*)-*N*-Cyclohexyl-2-phenylpropanamide (**4**)

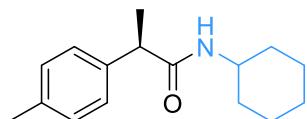
21.0 mg, 91% yield, 98:2 er, white solid, $[\alpha]_D^{24} = -7.9$ ($c = 0.67$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-4/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.36 – 7.09 (m, 5H), 5.12 (s, 1H), 3.77 – 3.56 (m, 1H), 3.44 (q, $J = 7.0$ Hz, 1H), 1.89 – 1.65 (m, 2H), 1.59 – 1.47 (m, 2H), 1.43 (d, $J = 7.0$ Hz, 3H), 1.28 – 1.15 (m, 3H), 1.08 – 0.83 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.2, 141.7, 128.8, 127.6, 127.1, 48.1, 47.2, 32.9, 32.9, 25.5, 24.8, 24.7, 18.6.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{15}\text{H}_{21}\text{NO}$ [$\text{M}+\text{H}]^+$: 232.1701, Found: 232.1704.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 95:5 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (major) = 6.141 min, t_R (minor) = 7.272 min.



(*R*)-*N*-Cyclohexyl-2-(p-tolyl)propanamide (**5**)

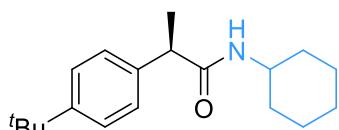
19.2 mg, 78% yield, 99.5:0.5 er, white solid, $[\alpha]_D^{24} = -6.9$ ($c = 0.67$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-4/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.22 – 7.09 (m, 4H), 5.17 (s, 1H), 3.72 (m, 1H), 3.47 (q, $J = 7.2$ Hz, 1H), 2.33 (s, 3H), 1.91 – 1.72 (m, 2H), 1.65 – 1.53 (m, 3H), 1.48 (d, $J = 7.0$ Hz, 3H), 1.39 – 1.20 (m, 2H), 1.12 – 0.91 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.5, 138.6, 136.8, 133.4, 129.6, 127.8, 127.5, 48.1, 46.8, 32.9, 32.9, 25.5, 24.8, 24.7, 21.1, 18.6.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{16}\text{H}_{23}\text{NO}$ [$\text{M}+\text{H}]^+$: 246.1858, Found: 246.1856.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 98:2 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (major) = 12.000 min, t_R (minor) = 14.959 min.



(*R*)-2-(4-(*tert*-Butyl)phenyl)-*N*-cyclohexylpropanamide (**6**)

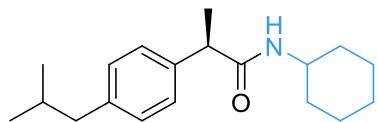
24.1 mg, 84% yield, 98:2 er, white solid, $[\alpha]_D^{24} = -7.9$ ($c = 0.67$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-4/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.39 – 7.31 (m, 2H), 7.20 (d, $J = 8.0$ Hz, 2H), 5.18 (s, 1H), 3.82 – 3.62 (m, 1H), 3.48 (q, $J = 7.0$ Hz, 1H), 1.82 – 1.89 (m, 2H), 1.69 – 1.52 (m, 3H), 1.49 (d, $J = 7.0$ Hz, 2H), 1.31 – 1.38 (m, 1H), 1.14 – 0.90 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.4, 150.0, 138.5, 127.2, 125.7, 48.1, 46.7, 34.5, 33.0, 32.9, 31.3, 25.5, 24.8, 24.7, 18.6.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{19}\text{H}_{29}\text{NO}$ [$\text{M}+\text{H}]^+$: 288.2327, Found: 288.2327.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 98:2 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (major) = 9.830 min, t_R (minor) = 12.114 min.



(*R*)-*N*-Cyclohexyl-2-(4-isobutylphenyl)propanamide (7)

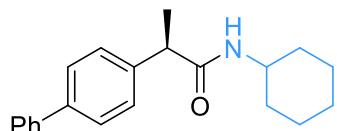
24.4 mg, 85% yield, 98:2 er, white solid, $[\alpha]_D^{24} = -8.6$ ($c = 1.43$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-5/1.

$^1\text{H NMR}$ (400 MHz, CDCl_3) δ 7.21 – 7.00 (m, 4H), 5.14 (s, 1H), 3.89 – 3.65 (m, 1H), 3.49 (q, $J = 7.0$ Hz, 1H), 2.45 (d, $J = 7.0$ Hz, 2H), 1.92 – 1.73 (m, 3H), 1.67 – 1.52 (m, 3H), 1.50 (d, $J = 7.0$ Hz, 3H), 1.39 – 1.24 (m, 2H), 1.15 – 0.94 (m, 3H), 0.89 (d, $J = 6.6$ Hz, 6H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.5, 140.6, 138.8, 129.6, 127.3, 48.0, 46.9, 45.0, 32.9, 32.8, 30.2, 25.5, 24.7, 24.6, 22.4, 18.5.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{19}\text{H}_{29}\text{NO}$ [$\text{M}+\text{H}]^+$: 288.2327, Found: 288.2332.

The enantiomeric excess was determined by Chiralcel OJ column, heptane/EtOH = 99:1 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (major) = 8.540 min, t_R (minor) = 12.889 min.



(*R*)-2-([1,1'-Biphenyl]-4-yl)-*N*-cyclohexylpropanamide (8)

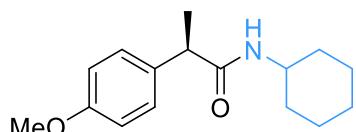
26.1 mg, 85% yield, 93:7 er, white solid, $[\alpha]_D^{24} = -12.7$ ($c = 0.57$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-4/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.64 – 7.53 (m, 4H), 7.50 – 7.38 (m, 2H), 7.41 – 7.30 (m, 3H), 5.25 (d, $J = 7.5$ Hz, 1H), 3.85 – 3.68 (m, 1H), 3.56 (q, $J = 7.0$ Hz, 1H), 1.94 – 1.76 (m, 2H), 1.68 – 1.55 (m, 3H), 1.54 (d, $J = 7.0$ Hz, 3H), 1.42 – 1.21 (m, 2H), 1.18 – 0.95 (m, 3H).

^{13}C NMR (75 MHz, CDCl_3) δ 173.1, 140.7, 140.6, 140.0, 128.8, 128.0, 127.5, 127.3, 127.0, 48.2, 46.9, 33.0, 32.9, 25.5, 24.8, 24.7, 18.7.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{21}\text{H}_{25}\text{NO} [\text{M}+\text{H}]^+$: 308.2014, Found: 308.2016.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 99:1 v/v, $v = 1.0 \text{ mL/min}$, $\lambda = 210 \text{ nm}$, t_R (major) = 35.029 min, t_R (minor) = 41.613 min.



(*R*)-*N*-Cyclohexyl-2-(4-methoxyphenyl)propanamide (9)

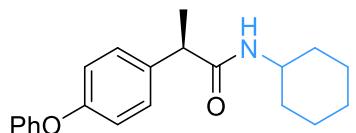
20.1 mg, 77% yield, 97:3 er, white solid, $[\alpha]_D^{24} = -3.9$ ($c = 0.87, \text{CHCl}_3$). Eluent: pentane/ethyl acetate = 5/1-2/1.

^1H NMR (300 MHz, CDCl_3) δ 7.16 – 7.05 (m, 2H), 6.84 – 6.62 (m, 2H), 5.14 (d, $J = 6.5 \text{ Hz}$, 1H), 3.73 (s, 3H), 3.69 – 3.58 (m, 1H), 3.39 (q, $J = 7.0 \text{ Hz}$, 1H), 1.83 – 1.63 (m, 2H), 1.51 (ddd, $J = 10.5, 8.0, 2.5 \text{ Hz}$, 3H), 1.40 (d, $J = 7.0 \text{ Hz}$, 3H), 1.28 – 1.21 (m, 2H), 1.08 – 0.84 (m, 3H).

^{13}C NMR (75 MHz, CDCl_3) δ 173.6, 158.7, 133.7, 128.6, 114.2, 55.3, 48.1, 46.4, 33.0, 32.9, 25.5, 24.8, 24.7, 18.7.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{16}\text{H}_{23}\text{NO}_2 [\text{M}+\text{H}]^+$: 262.1807, Found: 262.1803.

The enantiomeric excess was determined by Chiraldak AD-H column, heptane/EtOH = 80:20 v/v, $v = 0.5 \text{ mL/min}$, $\lambda = 210 \text{ nm}$, t_R (minor) = 8.801 min, t_R (major) = 9.676 min.



(*R*)-*N*-Cyclohexyl-2-(4-phenoxyphenyl)propanamide (10)

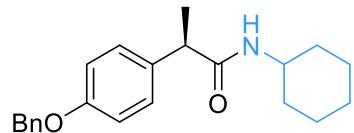
24.6mg, 76% yield, 96:4 er, white solid, $[\alpha]_D^{24} = -5.7$ ($c = 0.8, \text{CHCl}_3$). Eluent: pentane/ethyl acetate = 10/1-3/1.

^1H NMR (300 MHz, CDCl_3) δ 7.41 – 7.31 (m, 2H), 7.31 – 7.23 (m, 2H), 7.18 – 7.09 (m, 1H), 7.07 – 6.97 (m, 4H), 5.25 (d, $J = 7.5 \text{ Hz}$, 1H), 3.91 – 3.68 (m, 1H), 3.51 (q, $J = 7.0 \text{ Hz}$, 1H), 1.95 – 1.76 (m, 2H), 1.72 – 1.56 (m, 3H), 1.52 (d, $J = 7.0 \text{ Hz}$, 3H), 1.44 – 1.27 (m, 2H), 1.22 – 0.92 (m, 3H).

^{13}C NMR (75 MHz, CDCl_3) δ 173.3, 157.1, 156.4, 136.4, 133.4, 129.8, 128.9, 127.8, 123.4, 119.0, 118.9, 48.2, 46.6, 33.0, 32.9, 25.5, 24.7, 24.7, 18.8.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{21}\text{H}_{25}\text{NO}_2 [\text{M}+\text{H}]^+$: 324.1964, Found: 324.1960.

The enantiomeric excess was determined by Chiralpak AD-H column, heptane/EtOH = 80:20 v/v, v = 0.5 mL/min, λ = 210 nm, t_R (minor) = 8.545 min, t_R (major) = 10.925 min.



(R)-2-(4-(Benzylxy)phenyl)-N-cyclohexylpropanamide (11)

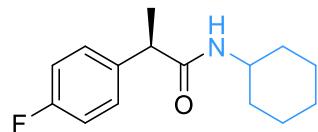
27.9 mg, 83% yield, 98:2 er, slight yellow solid, $[\alpha]_D^{24} = -12.0$ ($c = 0.53$, CHCl₃). Eluent: pentane/ethyl acetate = 10/1-4/1.

¹H NMR (300 MHz, CDCl₃) δ 7.73 – 7.64 (m, 1H), 7.52 – 7.33 (m, 6H), 7.24 – 7.19 (m, 2H), 7.05 – 6.92 (m, 2H), 5.18 (d, $J = 8.1$ Hz, 1H), 5.08 (s, 2H), 3.86 – 3.65 (m, 1H), 3.48 (q, $J = 7.0$ Hz, 1H), 1.93 – 1.75 (m, 2H), 1.72 – 1.54 (m, 3H), 1.50 (d, $J = 7.0$ Hz, 3H), 1.41 – 1.28 (m, 2H), 1.18 – 0.89 (m, 3H).

¹³C NMR (75 MHz, CDCl₃) δ 173.7, 157.9, 136.9, 133.9, 133.4, 130.0, 128.7, 128.6, 128.0, 127.9, 127.5, 115.2, 70.1, 48.1, 46.4, 32.9, 32.9, 25.5, 24.7, 24.7, 18.7.

HRMS (ESI-TOF) m/z: Calcd. for C₂₂H₂₇NO₂ [M+H]⁺: 338.2120, Found: 338.2123.

The enantiomeric excess was determined by Chiralpak AD-H column, heptane/EtOH = 80:20 v/v, v = 0.5 mL/min, λ = 210 nm, t_R (minor) = 11.384 min, t_R (major) = 17.706 min.



(R)-N-Cyclohexyl-2-(4-fluorophenyl)propanamide (12)

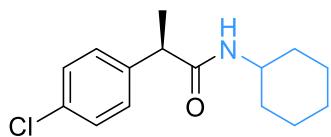
19.0 mg, 77% yield, 97:3 er, white solid, $[\alpha]_D^{24} = -1.5$ ($c = 0.83$, CHCl₃). Eluent: pentane/ethyl acetate = 10/1-5/1.

¹H NMR (300 MHz, CDCl₃) δ 7.36 – 7.22 (m, 2H), 7.04 (t, $J = 8.7$ Hz, 2H), 5.17 (s, 1H), 3.96 – 3.62 (m, 1H), 3.50 (q, $J = 7.0$ Hz, 1H), 1.94 – 1.76 (m, 2H), 1.73 – 1.54 (m, 3H), 1.50 (d, $J = 7.0$ Hz, 3H), 1.39 – 1.28 (m, 2H), 1.17 – 0.96 (m, 3H).

¹³C NMR (75 MHz, CDCl₃) δ 173.04, 160.69 (d, $J = 245.0$ Hz), 135.40 (d, $J = 3.0$ Hz), 129.96, 129.06 (d, $J = 8.0$ Hz), 127.84, 115.65 (d, $J = 21.5$ Hz).

HRMS (ESI-TOF) m/z: Calcd. for C₁₅H₂₀NFO [M+H]⁺: 250.1607, Found: 250.1610.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 98:2 v/v, v = 0.5 mL/min, λ = 210 nm, t_R (major) = 24.800 min, t_R (minor) = 29.149 min.



(R)-2-(4-Chlorophenyl)-N-cyclohexylpropanamide (13)

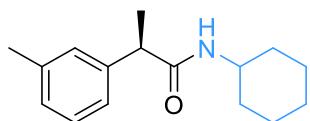
19.3 mg, 73% yield, 92:8 er, white solid, $[\alpha]_D^{24} = -11.8$ ($c = 0.63$, CHCl₃). Eluent: pentane/ethyl acetate = 10/1-5/1.

¹H NMR (300 MHz, CDCl₃) δ 7.37 – 7.02 (m, 4H), 5.11 (s, 1H), 3.79 – 3.56 (m, 1H), 3.39 (q, $J = 7.0$ Hz, 1H), 1.85 – 1.65 (m, 2H), 1.62 – 1.45 (m, 3H), 1.40 (d, $J = 7.0$ Hz, 3H), 1.33 – 1.16 (m, 2H), 1.09 – 0.87 (m, 3H).

¹³C NMR (75 MHz, CDCl₃) δ 172.6, 140.2, 132.9, 128.9, 128.9, 48.3, 46.6, 33.0, 32.9, 25.5, 24.8, 24.7, 18.8.

HRMS (ESI-TOF) m/z: Calcd. for C₁₅H₂₀NOCl [M+H]⁺: 266.1312, Found: 266.1315.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 98:2 v/v, $v = 0.8$ mL/min, $\lambda = 210$ nm, t_R (major) = 12.790 min, t_R (minor) = 14.384 min.



(R)-N-Cyclohexyl-2-(m-tolyl)propanamide (14)

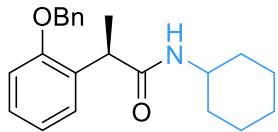
22.8 mg, 93% yield, 97:3 er, slight yellow solid, $[\alpha]_D^{24} = -17.0$ ($c = 0.60$, CHCl₃). Eluent: pentane/ethyl acetate = 10/1-5/1.

¹H NMR (300 MHz, CDCl₃) δ 7.29 – 7.16 (m, 1H), 7.12 – 7.02 (m, 3H), 5.16 (s, 1H), 3.83 – 3.63 (m, 1H), 3.47 (q, $J = 7.0$ Hz, 1H), 2.34 (s, 3H), 1.92 – 1.72 (m, 3H), 1.68 – 1.50 (m, 3H), 1.49 (d, $J = 7.2$ Hz, 3H), 1.38 – 1.24 (m, 2H), 1.14 – 0.93 (m, 3H).

¹³C NMR (75 MHz, CDCl₃) δ 173.3, 141.6, 138.5, 128.7, 128.4, 127.9, 124.6, 48.1, 47.2, 33.0, 32.9, 25.5, 24.8, 24.7, 21.5, 18.7.

HRMS (ESI-TOF) m/z: Calcd. for C₁₆H₂₃NO [M+H]⁺: 246.1858, Found: 246.1859.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 98:2 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (major) = 11.198 min, t_R (minor) = 13.769 min.



(R)-2-(2-(Benzyl)phenyl)-N-cyclohexylpropanamide (15)

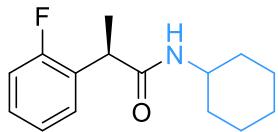
28.0 mg, 83% yield, 99:1 er, white solid, $[\alpha]_D^{24} = -75$ ($c = 0.74$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-4/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.46 – 7.31 (m, 6H), 7.28 – 7.17 (m, 1H), 7.07 – 6.88 (m, 2H), 5.50 (d, $J = 8.1$ Hz, 1H), 5.11 (s, 2H), 3.99 (q, $J = 7.2$ Hz, 1H), 3.75 – 3.56 (m, 1H), 1.85 – 1.61 (m, 2H), 1.48 (d, $J = 7.2$ Hz, 6H), 1.35 – 1.19 (m, 2H), 1.12 – 0.78 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.3, 155.7, 136.8, 130.2, 128.7, 128.2, 128.0, 127.9, 127.4, 121.5, 111.7, 70.2, 47.8, 39.6, 32.9, 32.7, 25.5, 24.7, 24.6, 16.4.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{22}\text{H}_{27}\text{NO}_2$ [$\text{M}+\text{H}]^+$: 338.2120, Found: 338.2120.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 95:5 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (major) = 8.468 min, t_R (minor) = 12.413 min.



(R)-N-Cyclohexyl-2-(2-fluorophenyl)propanamide (16)

15.9 mg, 64% yield, 98:2 er, white solid, $[\alpha]_D^{24} = -19.2$ ($c = 0.33$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-5/1.

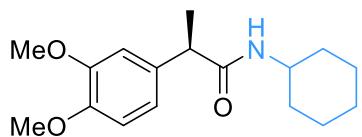
$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.38 (td, $J = 7.5, 2.0$ Hz, 1H), 7.26 – 7.19 (m, 1H), 7.13 (td, $J = 7.5, 1.5$ Hz, 1H), 7.04 (ddd, $J = 10.5, 8.0, 1.5$ Hz, 1H), 5.29 (s, 1H), 3.90 – 3.62 (m, 2H), 1.97 – 1.73 (m, 2H), 1.71 – 1.52 (m, 3H), 1.49 (d, $J = 7.0$ Hz, 3H), 1.40 – 1.23 (m, 2H), 1.17 – 0.93 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 172.1, 160.3 (d, $J = 245.0$ Hz), 128.7, 128.6 (d, $J = 12.5$ Hz), 124.6 (d, $J = 3.5$ Hz), 115.5, 115.2, 48.2, 39.4, 33.0, 32.8, 25.5, 24.8, 24.7, 17.2.

$^{19}\text{F NMR}$ (282 MHz, CDCl_3) δ -118.42.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{15}\text{H}_{20}\text{NFO}$ [$\text{M}+\text{H}]^+$: 250.1607, Found: 250.1611.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 98:2 v/v, $v = 0.5$ mL/min, $\lambda = 210$ nm, t_R (major) = 10.456 min, t_R (minor) = 11.795 min.



(R)-N-Cyclohexyl-2-(3,4-dimethoxyphenyl)propanamide (17)

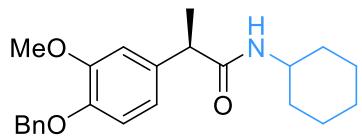
21.5 mg, 74% yield, 96.5:3.5 er, slight yellow solid, $[\alpha]_D^{24} = -10.7$ ($c = 0.47$, CHCl_3). Eluent: pentane/ethyl acetate = 5/1-2/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 6.80 (dd, $J = 4.0, 2.0$ Hz, 3H), 5.19 (d, $J = 7.5$ Hz, 1H), 3.86 (d, $J = 1.5$ Hz, 6H), 3.81 – 3.59 (m, 1H), 3.45 (q, $J = 7.0$ Hz, 1H), 1.90 – 1.71 (m, 2H), 1.67 – 1.49 (m, 3H), 1.48 (d, $J = 7.2$ Hz, 3H), 1.41 – 1.21 (m, 2H), 1.16 – 0.91 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.5, 149.2, 148.1, 134.2, 119.6, 111.3, 110.6, 55.9, 48.1, 46.8, 33.0, 32.9, 25.5, 24.7, 24.7, 18.7.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{17}\text{H}_{25}\text{NO}_3$ [M+H] $^+$: 292.1913, Found: 292.1915.

The enantiomeric excess was determined by Chiraldak AD-H column, heptane/EtOH = 80:20 v/v, $v = 0.5$ mL/min, $\lambda = 210$ nm, t_R (minor) = 9.009 min, t_R (major) = 10.294 min.



(R)-2-(4-(Benzylxy)-3-methoxyphenyl)-N-cyclohexylpropanamide (18)

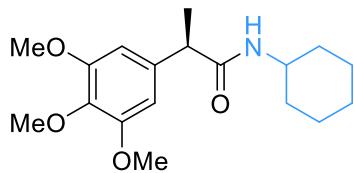
26.4 mg, 72% yield, 95:5 er, white solid, $[\alpha]_D^{24} = -6.3$ ($c = 0.7$, CHCl_3). Eluent: pentane/ethyl acetate = 5/1-3/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.48 – 7.41 (m, 2H), 7.39 – 7.29 (m, 3H), 6.87 – 6.79 (m, 2H), 6.74 (dd, $J = 8.5, 2.0$ Hz, 1H), 5.14 (s, 3H), 3.88 (s, 3H), 3.78 – 3.65 (m, 1H), 3.44 (q, $J = 7.0$ Hz, 1H), 1.90 – 1.72 (m, 2H), 1.67 – 1.52 (m, 3H), 1.48 (d, $J = 7.0$ Hz, 3H), 1.41 – 1.26 (m, 2H), 1.16 – 0.89 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.4, 149.9, 147.3, 137.1, 134.8, 128.6, 127.9, 127.3, 119.6, 114.1, 111.2, 71.1, 56.0, 48.1, 46.8, 32.9, 32.9, 25.5, 24.7, 18.7.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{23}\text{H}_{29}\text{NO}_3$ [M+H] $^+$: 368.2226, Found: 368.2224.

The enantiomeric excess was determined by Chiraldak AD-H column, heptane/EtOH = 80:20 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (minor) = 5.680 min, t_R (major) = 7.218 min.



(*R*)-*N*-Cyclohexyl-2-(3,4,5-trimethoxyphenyl)propanamide (19)

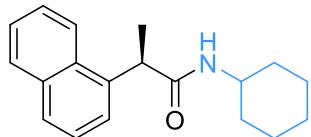
24.1 mg, 75% yield, 99.5:0.5 er, slight yellow solid, $[\alpha]_D^{24} = -12.5$ ($c = 0.57$, CHCl_3). Eluent: DCM/ethyl acetate = 10/1-5/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 6.49 (s, 2H), 5.21 (s, 1H), 3.85 (s, 6H), 3.83 (s, 3H), 3.80 – 3.65 (m, 1H), 3.43 (q, $J = 7.0$ Hz, 1H), 1.90 – 1.73 (m, 2H), 1.69 – 1.53 (m, 3H), 1.48 (d, $J = 7.0$ Hz, 3H), 1.38 – 1.26 (m, 2H), 1.19 – 0.96 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.0, 153.4, 137.4, 104.5, 60.9, 56.2, 48.2, 47.5, 33.0, 32.9, 25.5, 24.7, 18.8.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{18}\text{H}_{27}\text{NO}_4$ [$\text{M}+\text{H}]^+$: 322.2018, Found: 322.2020.

The enantiomeric excess was determined by Chiralpak AD-H column, heptane/EtOH = 80:20 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (minor) = 4.359 min, t_R (major) = 7.000 min.



(*R*)-*N*-Cyclohexyl-2-(naphthalen-1-yl)propanamide (20)

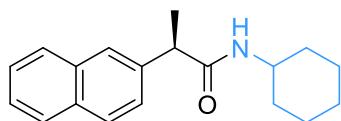
21.4 mg, 56% yield, 90:10 er, white solid, $[\alpha]_D^{24} = -43.4$ ($c = 0.43$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-5/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 8.07 – 7.98 (m, 1H), 7.91 – 7.85 (m, 1H), 7.80 (dd, $J = 6.7, 2.3$ Hz, 1H), 7.56 – 7.44 (m, 4H), 5.05 (d, $J = 7.4$ Hz, 1H), 4.27 (q, $J = 7.2$ Hz, 1H), 3.71 – 3.75 (m, 1H), 1.65 – 1.75 (m, 5H), 1.54 – 1.41 (m, 3H), 1.32 – 1.20 (m, 2H), 1.04 – 0.75 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.7, 137.3, 134.0, 131.5, 129.0, 128.1, 126.5, 125.9, 125.6, 124.9, 123.4, 48.1, 43.9, 32.7, 25.4, 24.6, 18.0.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{19}\text{H}_{23}\text{NO}$ [$\text{M}+\text{H}]^+$: 282.1858, Found: 282.1853.

The enantiomeric excess was determined by Chiralpak AD-H column, heptane/EtOH = 98:2 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (minor) = 19.189 min, t_R (major) = 20.675 min.



(R)-N-Cyclohexyl-2-(naphthalen-2-yl)propanamide (21)

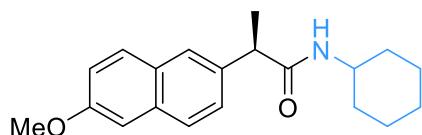
18.0 mg, 64% yield, 92.5:7.5 er, slight yellow solid, $[\alpha]_D^{24} = -21.3$ ($c = 0.37$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-5/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.87 – 7.79 (m, 3H), 7.73 (d, $J = 1.5$ Hz, 1H), 7.52 – 7.45 (m, 2H), 7.41 (dd, $J = 8.5, 1.8$ Hz, 1H), 5.18 (s, 1H), 3.84 – 3.61 (m, 2H), 1.97 – 1.71 (m, 2H), 1.71 – 1.48 (m, 6H), 1.42 – 1.17 (m, 2H), 1.09 – 0.79 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.1, 139.2, 133.5, 132.6, 128.7, 127.8, 127.7, 126.3, 126.2, 125.9, 125.8, 48.3, 47.4, 33.0, 32.9, 25.5, 24.8, 24.7, 18.7.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{19}\text{H}_{23}\text{NO}$ [$\text{M}+\text{H}]^+$: 282.1858, Found: 282.1862.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 98:2 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (minor) = 17.505 min, t_R (major) = 19.761 min.



(R)-N-Cyclohexyl-2-(6-methoxynaphthalen-2-yl)propanamide (22)

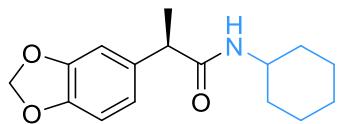
22.7 mg, 73% yield, 95:5 er, white solid, $[\alpha]_D^{24} = -9.7$ ($c = 1.23$, CHCl_3). Eluent: pentane/ethyl acetate = 8/1-3/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.76 – 7.63 (m, 3H), 7.37 (dd, $J = 8.5, 2.0$ Hz, 1H), 7.20 – 7.10 (m, 2H), 5.21 (d, $J = 7.5$ Hz, 1H), 3.92 (s, 3H), 3.83 – 3.59 (m, 2H), 1.90 – 1.71 (m, 2H), 1.67 – 1.46 (m, 6H), 1.37 – 1.23 (m, 2H), 1.13 – 0.82 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.4, 157.7, 136.8, 133.7, 129.2, 129.0, 127.5, 126.3, 126.1, 119.1, 105.7, 55.3, 48.2, 47.2, 33.0, 32.9, 25.5, 24.8, 24.7, 18.6.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{20}\text{H}_{25}\text{NO}_2$ [$\text{M}+\text{H}]^+$: 312.1964, Found: 312.1967.

The enantiomeric excess was determined by Chiralcel OJ-H column, heptane/EtOH = 95:5 v/v, $v = 2.0$ mL/min, $\lambda = 210$ nm, t_R (major) = 13.075 min, t_R (minor) = 23.115 min.



(R)-2-(Benzo[d][1,3]dioxol-5-yl)-N-cyclohexylpropanamide (23)

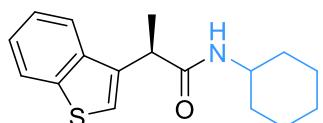
21.2 mg, 77% yield, 95:5 er, pale yellow solid, $[\alpha]_D^{24} = -4.9$ ($c = 0.5$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-3/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 6.90 – 6.62 (m, 3H), 5.95 (s, 2H), 5.18 (s, 1H), 3.89 – 3.61 (m, 1H), 3.42 (q, $J = 7.0$ Hz, 1H), 1.92 – 1.73 (m, 2H), 1.68 – 1.52 (m, 3H), 1.45 (d, $J = 7.0$ Hz, 3H), 1.38 – 1.25 (m, 2H), 1.14 – 0.92 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.3, 148.0, 146.7, 135.4, 133.4, 130.0, 127.8, 120.8, 108.4, 107.9, 101.1, 48.2, 46.9, 33.0, 32.9, 25.5, 24.8, 24.7, 18.8.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{16}\text{H}_{21}\text{NO}_3$ [M+H] $^+$: 276.1599, Found: 276.1601.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 98:2 v/v, $v = 2.0$ mL/min, $\lambda = 210$ nm, t_R (minor) = 23.301 min, t_R (major) = 24.670 min.



(R)-2-(Benzo[b]thiophen-3-yl)-N-cyclohexylpropanamide (24)

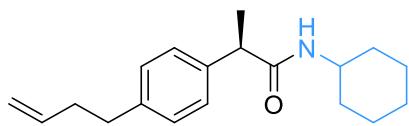
20.1 mg, 70% yield, 94:6 er, white solid, $[\alpha]_D^{24} = -28.9$ ($c = 0.47$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-4/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.93 – 7.82 (m, 1H), 7.79 – 7.73 (m, 1H), 7.40 – 7.34 (m, 3H), 5.21 (d, $J = 7.5$ Hz, 1H), 3.93 (qd, $J = 7.0, 1.0$ Hz, 1H), 3.84 – 3.63 (m, 1H), 1.81 – 1.69 (m, 2H), 1.66 (d, $J = 7.0$ Hz, 3H), 1.57 – 1.45 (m, 3H), 1.33 – 1.20 (m, 2H), 1.12 – 0.81 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 172.6, 140.6, 137.9, 136.0, 124.7, 124.3, 123.0, 122.9, 121.9, 48.1, 41.7, 32.8, 25.4, 24.7, 17.7.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{17}\text{H}_{21}\text{NOS}$ [M+H] $^+$: 288.1422, Found: 288.1426.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 95:5 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (major) = 7.686 min, t_R (minor) = 9.330 min.



(*R*)-2-(4-(but-3-en-1-yl)phenyl)-*N*-cyclohexylpropanamide (25)

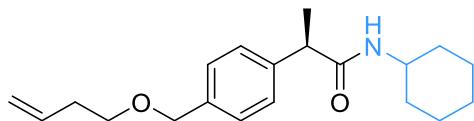
27.1 mg, 95% yield, 97:3 er, white solid, $[\alpha]_D^{24} = -9.5$ ($c = 0.63$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-5/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.24 – 7.10 (m, 4H), 5.97 – 5.74 (m, 1H), 5.14 (s, 1H), 5.10 – 4.92 (m, 2H), 3.80 – 3.63 (m, 1H), 3.48 (q, $J = 7.2$ Hz, 1H), 2.75 – 2.63 (m, 2H), 2.45 – 2.31 (m, 2H), 1.89 – 1.72 (m, 2H), 1.67 – 1.53 (m, 3H), 1.49 (d, $J = 7.2$ Hz, 3H), 1.39 – 1.22 (m, 2H), 1.14 – 0.90 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.4, 140.7, 139.1, 138.0, 128.9, 127.5, 115.0, 48.1, 46.9, 35.4, 34.9, 32.9, 32.9, 25.5, 24.7, 24.7, 18.6.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{19}\text{H}_{27}\text{NO}$ [$\text{M}+\text{H}]^+$: 286.2171, Found: 286.2172.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 98:2 v/v, $v = 1.5$ mL/min, $\lambda = 210$ nm, t_R (major) = 7.188 min, t_R (minor) = 8.811 min.



(*R*)-2-(4-((But-3-en-1-yloxy)methyl)phenyl)-*N*-cyclohexylpropanamide (26)

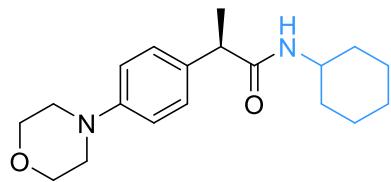
22.1 mg, 70% yield, 96:4 er, pale yellow solid, $[\alpha]_D^{24} = -9.4$ ($c = 0.57$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-4/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.31 (d, $J = 8.5$ Hz, 2H), 7.29 – 7.21 (m, 2H), 5.89 – 5.80 (m, 1H), 5.21 – 4.99 (m, 3H), 4.50 (s, 2H), 3.80 – 3.64 (m, 1H), 3.60 – 3.42 (m, 3H), 2.47 – 2.32 (m, 2H), 1.90 – 1.72 (m, 2H), 1.67 – 1.53 (m, 3H), 1.49 (d, $J = 7.0$ Hz, 3H), 1.41 – 1.22 (m, 2H), 1.14 – 0.91 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.3, 140.9, 137.4, 135.2, 133.4, 130.0, 128.2, 127.9, 127.7, 116.4, 72.6, 69.8, 48.2, 47.0, 34.2, 33.0, 32.9, 25.5, 24.8, 24.7, 18.7.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{20}\text{H}_{29}\text{NO}_2$ [$\text{M}+\text{H}]^+$: 316.2276, Found: 316.2279.

The enantiomeric excess was determined by Chiralpak AD-H column, heptane/iPrOH = 95:5 v/v, $v = 0.5$ mL/min, $\lambda = 210$ nm, t_R (minor) = 19.626 min, t_R (major) = 23.788 min.



(*R*)-*N*-Cyclohexyl-2-(4-morpholinophenyl)propanamide (27)

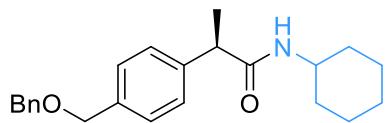
26.9 mg, 81% yield, 97.5:2.5 er, white solid, $[\alpha]_D^{24} = -7.2$ ($c = 0.5$, CHCl_3). Eluent: DCM/ethyl acetate = 10/1-5/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.23 – 7.13 (m, 2H), 6.93 – 6.84 (m, 2H), 5.15 (d, $J = 7.5$ Hz, 1H), 3.91 – 3.82 (m, 4H), 3.79 – 3.65 (m, 1H), 3.44 (q, $J = 7.0$ Hz, 1H), 3.20 – 3.10 (m, 4H), 1.90 – 1.74 (m, 2H), 1.70 – 1.51 (m, 3H), 1.47 (d, $J = 7.0$ Hz, 3H), 1.37 – 1.23 (m, 2H), 1.15 – 0.87 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.7, 150.3, 132.9, 128.4, 115.9, 66.9, 49.3, 48.1, 46.4, 33.1, 33.0, 25.5, 24.8, 24.7, 18.6.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{19}\text{H}_{28}\text{N}_2\text{O}_2$ [$\text{M}+\text{H}]^+$: 317.2229, Found: 317.2229.

The enantiomeric excess was determined by Chiraldak AD-H column, heptane/EtOH = 80:20 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (minor) = 6.840 min, t_R (major) = 7.591 min.



(*R*)-2-(4-((Benzyl)oxy)methyl)phenyl-*N*-cyclohexylpropanamide (28)

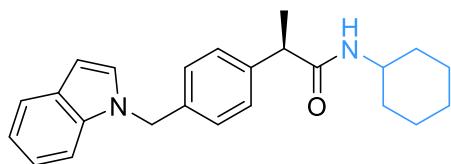
25.3 mg, 72% yield, 96:3 er, pale yellow solid, $[\alpha]_D^{24} = -11.4$ ($c = 0.53$, CHCl_3). Eluent: pentane/ethyl acetate = 8/1-4/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.44 – 7.22 (m, 9H), 5.17 (d, $J = 7.5$ Hz, 1H), 4.58 (s, 2H), 4.54 (s, 2H), 3.83 – 3.65 (m, 1H), 3.51 (q, $J = 7.0$ Hz, 1H), 1.92 – 1.73 (m, 2H), 1.69 – 1.53 (m, 3H), 1.50 (d, $J = 7.0$ Hz, 3H), 1.37 – 1.27 (m, 2H), 1.14 – 0.87 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.2, 141.1, 138.2, 137.2, 133.4, 130.0, 128.5, 128.3, 127.8, 127.74, 127.72, 72.3, 71.8, 48.2, 47.0, 33.0, 32.9, 25.5, 24.8, 24.7, 18.7.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{23}\text{H}_{29}\text{NO}_2$ [$\text{M}+\text{H}]^+$: 352.2276, Found: 352.2271.

The enantiomeric excess was determined by Cellulose4 column, heptane/iPrOH = 95:5 v/v, $v = 0.5$ mL/min, $\lambda = 210$ nm, t_R (major) = 29.114 min, t_R (minor) = 33.019 min.



(*R*)-2-(4-((1*H*-Indol-1-yl)methyl)phenyl)-*N*-cyclohexylpropanamide (29)

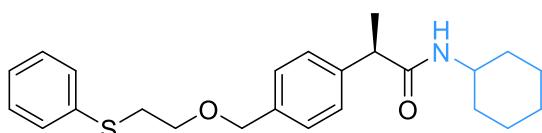
29.2 mg, 81% yield, 93:7 er, white solid, $[\alpha]_D^{24} = -24.2$ ($c = 0.5$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-3/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.71 – 7.61 (m, 1H), 7.32 – 7.04 (m, 8H), 6.56 (dd, $J = 3.0, 1.0$ Hz, 1H), 5.31 (s, 2H), 5.16 (d, $J = 7.5$ Hz, 1H), 3.82 – 3.64 (m, 1H), 3.46 (q, $J = 7.0$ Hz, 1H), 1.91 – 1.71 (m, 2H), 1.65 – 1.52 (m, 3H), 1.46 (d, $J = 7.0$ Hz, 3H), 1.37 – 1.26 (m, 2H), 1.17 – 0.93 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.0, 141.1, 136.5, 128.7, 128.2, 128.0, 127.3, 121.7, 121.0, 119.6, 109.7, 101.7, 49.8, 48.2, 46.9, 33.0, 32.9, 25.5, 24.8, 24.7, 18.7.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{24}\text{H}_{28}\text{N}_2\text{O} [\text{M}+\text{H}]^+$: 361.2280, Found: 361.2277.

The enantiomeric excess was determined by Chiralcel OJ-H column, heptane/EtOH = 95:5 v/v, $v = 2.0$ mL/min, $\lambda = 210$ nm, t_R (major) = 29.265 min, t_R (minor) = 37.592 min.



(*R*)-*N*-Cyclohexyl-2-(4-((2-(phenylthio)ethoxy)methyl)phenyl)propanamide (30)

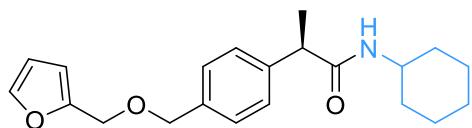
24.6 mg, 62% yield, 93:7 er, white solid, $[\alpha]_D^{24} = -8.8$ ($c = 0.97$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-4/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.41 – 7.18 (m, 9H), 5.21 (d, $J = 7.5$ Hz, 1H), 4.53 (s, 2H), 3.82 – 3.66 (m, 3H), 3.52 (q, $J = 7.0$ Hz, 1H), 3.17 (t, $J = 7.0$ Hz, 2H), 1.94 – 1.74 (m, 2H), 1.69 – 1.55 (m, 3H), 1.51 (d, $J = 7.0$ Hz, 3H), 1.40 – 1.27 (m, 2H), 1.18 – 0.92 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.1, 141.2, 136.9, 136.0, 129.4, 129.0, 128.2, 127.7, 126.2, 72.8, 69.0, 48.2, 47.0, 33.3, 33.0, 32.9, 25.5, 24.8, 24.7, 18.8.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{24}\text{H}_{31}\text{NO}_2\text{S} [\text{M}+\text{H}]^+$: 398.2154, Found: 398.2159.

The enantiomeric excess was determined by Chiraldak AD-H column, heptane/EtOH = 80:20 v/v, $v = 0.5$ mL/min, $\lambda = 210$ nm, t_R (minor) = 11.764 min, t_R (major) = 13.987 min.



(R)-N-Cyclohexyl-2-(4-((furan-2-ylmethoxy)methyl)phenyl)propanamide (31)

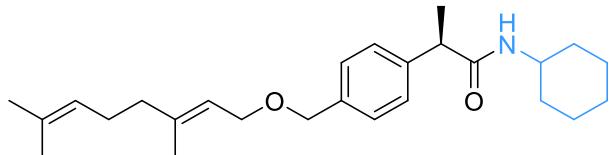
30.0 mg, 88% yield, 93:7 er, slight yellow oil, $[\alpha]_D^{24} = -10.3$ ($c = 0.83$, CHCl₃). Eluent: pentane/ethyl acetate = 5/1-3/1.

¹H NMR (300 MHz, CDCl₃) δ 7.42 (dd, $J = 2.0, 1.0$ Hz, 1H), 7.34 – 7.29 (m, 2H), 7.28 – 7.23 (m, 2H), 6.41 – 6.31 (m, 2H), 5.16 (d, $J = 7.0$ Hz, 1H), 4.51 (d, $J = 7.5$ Hz, 4H), 3.78 – 3.65 (m, 1H), 3.50 (q, $J = 7.0$ Hz, 1H), 1.89 – 1.72 (m, 2H), 1.69 – 1.51 (m, 3H), 1.49 (d, $J = 7.0$ Hz, 3H), 1.36 – 1.22 (m, 2H), 1.15 – 0.89 (m, 3H).

¹³C NMR (75 MHz, CDCl₃) δ 173.1, 151.7, 142.9, 141.2, 136.8, 128.4, 127.7, 110.3, 109.5, 71.6, 64.0, 48.2, 47.0, 33.0, 32.9, 25.5, 24.8, 24.7, 18.7.

HRMS (ESI-TOF) m/z: Calcd. for C₂₁H₂₇NO₃ [M+H]⁺: 342.2069, Found: 342.2066.

The enantiomeric excess was determined by Chiralpak AD-H column, heptane/EtOH = 80:20 v/v, $\nu = 0.5$ mL/min, $\lambda = 210$ nm, t_R (minor) = 10.396 min, t_R (major) = 12.918 min.



(R)-N-Cyclohexyl-2-(4-(((3,7-dimethylocta-2,6-dien-1-yl)oxy)methyl)phenyl)propanamide (32)

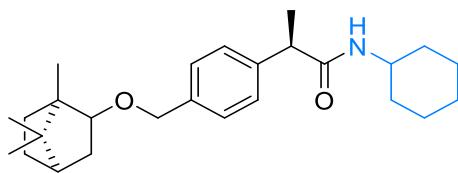
20.2 mg, 51% yield, 95:5 er, colorless oil, $[\alpha]_D^{24} = -12.0$ ($c = 0.4$, CHCl₃). Eluent: pentane/ethyl acetate = 10/1-4/1.

¹H NMR (300 MHz, CDCl₃) δ 7.32 (d, $J = 8.5$ Hz, 2H), 7.25 (d, $J = 8.5$ Hz, 2H), 5.41 (td, $J = 7.0, 1.5$ Hz, 1H), 5.16 – 5.02 (m, 2H), 4.47 (s, 2H), 4.02 (dd, $J = 7.0, 1.0$ Hz, 2H), 3.80 – 3.64 (m, 1H), 3.50 (q, $J = 7.0$ Hz, 1H), 2.07 (d, $J = 3.0$ Hz, 4H), 1.89 – 1.73 (m, 5H), 1.70 – 1.66 (m, 3H), 1.65 – 1.53 (m, 6H), 1.49 (d, $J = 7.5$ Hz, 3H), 1.35 – 1.25 (m, 2H), 1.13 – 0.89 (m, 3H).

¹³C NMR (75 MHz, CDCl₃) δ 173.2, 140.9, 140.7, 137.5, 132.0, 128.4, 127.6, 123.8, 121.8, 71.8, 66.6, 48.2, 47.0, 33.0, 32.9, 32.3, 26.7, 25.7, 25.5, 24.8, 24.7, 23.5, 18.7, 17.7.

HRMS (ESI-TOF) m/z: Calcd. for C₂₆H₃₉NO₂ [M+H]⁺: 398.3059, Found: 398.3063.

The enantiomeric excess was determined by Chiralpak AD-H column, heptane/EtOH = 95:5 v/v, $\nu = 1.0$ mL/min, $\lambda = 210$ nm, t_R (minor) = 6.378 min, t_R (major) = 7.423 min.



(2R)-N-cyclohexyl-2-(4-(((1*S*,4*S*)-1,7,7-trimethylbicyclo[2.2.1]heptan-2-yl)oxy)methyl)phenylpropanamide (33)

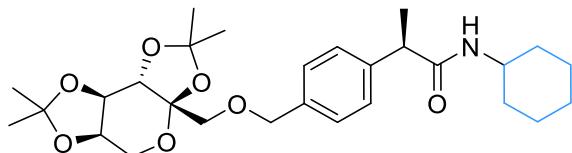
28.1 mg, 71% yield, 97:3 er, pale yellow oil, $[\alpha]_D^{24} = -40.0$ ($c = 0.8$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-4/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.15 – 6.98 (m, 4H), 4.94 (d, $J = 7.5$ Hz, 1H), 4.44 – 4.16 (m, 2H), 3.62 – 3.44 (m, 2H), 3.29 (q, $J = 7.0$ Hz, 1H), 1.99 – 1.81 (m, 2H), 1.67 – 1.33 (m, 6H), 1.28 (d, $J = 7.0$ Hz, 3H), 1.14 – 0.99 (m, 4H), 0.95 – 0.71 (m, 3H), 0.69 (s, 3H), 0.63 (d, $J = 6.5$ Hz, 6H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.3, 140.5, 138.5, 127.7, 127.5, 84.5, 71.3, 49.3, 48.2, 47.9, 47.0, 45.0, 36.2, 33.0, 32.9, 28.3, 26.8, 25.5, 24.8, 24.7, 19.8, 18.9, 18.7, 14.1.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{26}\text{H}_{39}\text{NO}_2$ [M+H] $^+$: 398.3059, Found: 398.3059.

The enantiomeric excess was determined by Cellulose4 column, heptane/*i*PrOH = 95:5 v/v, $v = 0.5$ mL/min, $\lambda = 210$ nm, t_R (major) = 11.179 min, t_R (minor) = 13.100 min.



(2R)-N-Cyclohexyl-2-(4-(((5a*R*,8a*R*,8b*S*)-2,2,7,7-tetramethyltetrahydro-3a*H*-bis([1,3]dioxolo)[4,5-*b*:4',5'-*d*]pyran-3*a*-yl)methoxy)methyl)phenylpropanamide (34)

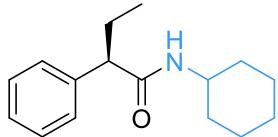
35.8 mg, 71% yield, 99.5:0.5 er, pale yellow oil, $[\alpha]_D^{24} = -16.7$ ($c = 0.87$, CHCl_3). Eluent: pentane/ethyl acetate = 5/1-2/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.43 – 7.12 (m, 4H), 5.14 (d, $J = 7.5$ Hz, 1H), 4.70 – 4.51 (m, 3H), 4.43 (dd, $J = 2.5, 1.0$ Hz, 1H), 4.27 – 4.18 (m, 1H), 3.91 (dd, $J = 13.0, 2.0$ Hz, 1H), 3.79 – 3.67 (m, 2H), 3.68 – 3.54 (m, 1H), 3.49 (q, $J = 7.0$ Hz, 1H), 1.90 – 1.71 (m, 2H), 1.66 – 1.51 (m, 6H), 1.48 (d, $J = 7.0$ Hz, 3H), 1.41 (s, 6H), 1.34 – 1.23 (m, 5H), 1.10 – 0.80 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.1, 140.9, 137.1, 128.0, 127.5, 108.9, 108.6, 102.7, 73.4, 71.7, 71.0, 70.2, 70.2, 61.0, 48.1, 47.0, 32.9, 32.9, 26.6, 25.9, 25.5, 25.4, 24.8, 24.7, 24.0, 18.7.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{28}\text{H}_{41}\text{NO}_7\text{Na}$ [M+Na] $^+$: 526.2786, Found: 526.2789.

The enantiomeric excess was determined by Chiralcel OJ column, heptane/EtOH = 95:5 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (major) = 11.469 min.



(*R*)-*N*-Cyclohexyl-2-phenylbutanamide (35)

From *trans*- β -methylstyrene: 14.5 mg, 59% yield, 98.5:1.5 er, white solid, $[\alpha]_D^{24} = -26.2$ ($c = 0.33$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-6/1.

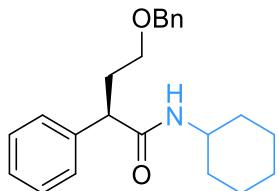
From *cis*- β -methylstyrene: 13.3 mg, 50% yield, >99.5:0.5 er, white solid, $[\alpha]_D^{24} = -34.5$ ($c = 0.33$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-6/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.14 – 6.97 (m, 5H), 5.02 (d, $J = 7.0$ Hz, 1H), 3.50 – 3.55 (m, 1H), 2.95 (t, $J = 7.5$ Hz, 1H), 1.96 (dt, $J = 13.5, 7.5$ Hz, 1H), 1.70 – 1.49 (m, 3H), 1.45 – 1.28 (m, 3H), 1.16 – 1.01 (m, 3H), 0.93 – 0.74 (m, 2H), 0.66 (t, $J = 7.5$ Hz, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 172.5, 140.3, 128.7, 128.0, 127.1, 55.4, 48.1, 33.1, 32.9, 26.5, 25.5, 24.8, 24.7, 12.4.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{16}\text{H}_{23}\text{NO}$ [$\text{M}+\text{H}]^+$: 246.1858, Found: 246.1859.

The enantiomeric excess was determined by Amylose2 column, heptane/EtOH = 90:10 v/v, $v = 0.5$ mL/min, $\lambda = 210$ nm, t_R (minor) = 7.949 min, t_R (major) = 9.778 min.



(*R*)-4-(Benzyl)-*N*-cyclohexyl-2-phenylbutanamide (36)

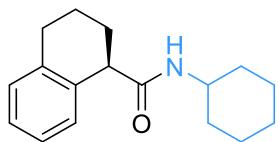
21.1 mg, 60% yield, white solid, $[\alpha]_D^{24} = -39.0$ ($c = 0.67$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-4/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.35 – 7.18 (m, 10H), 5.27 (d, $J = 8.5$ Hz, 1H), 4.40 (s, 2H), 3.76 – 3.60 (m, 1H), 3.52 (t, $J = 7.5$ Hz, 1H), 3.51 – 3.38 (m, 1H), 3.39 – 3.22 (m, 1H), 2.52 – 2.30 (m, 1H), 2.07 – 1.91 (m, 1H), 1.87 – 1.66 (m, 2H), 1.65 – 1.46 (m, 3H), 1.37 – 1.17 (m, 2H), 1.10 – 0.82 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 172.2, 140.0, 138.4, 128.7, 128.4, 128.0, 127.8, 127.6, 127.1, 73.0, 67.8, 49.6, 48.2, 33.5, 33.1, 32.8, 25.5, 24.8, 24.7.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{23}\text{H}_{29}\text{NO}_2$ [$\text{M}+\text{H}]^+$: 352.2276, Found: 352.2274.

The enantiomeric excess was determined by Amylose2 column, heptane/EtOH = 90:10 v/v, $v = 0.5$ mL/min, $\lambda = 210$ nm, t_R (major) = 6.181 min, t_R (minor) = 9.581 min.



(R)-N-Cyclohexyl-1,2,3,4-tetrahydronaphthalene-1-carboxamide (37)

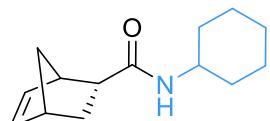
16.0 mg, 62% yield, 93:7 er, white solid, $[\alpha]_D^{24} = -6.3$ ($c = 0.37$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-5/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.23 – 7.06 (m, 4H), 5.13 (d, $J = 7.0$ Hz, 1H), 3.87 – 3.69 (m, 1H), 3.64 (t, $J = 5.5$ Hz, 1H), 2.80 (q, $J = 7.5, 6.5$ Hz, 2H), 2.35 – 2.21 (m, 1H), 2.06 – 1.89 (m, 1H), 1.78 (ddd, $J = 7.0, 6.0, 3.5$ Hz, 4H), 1.65 – 1.47 (m, 3H), 1.42 – 1.22 (m, 2H), 1.16 – 0.92 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 174.0, 137.9, 134.2, 130.0, 129.8, 127.3, 126.1, 48.1, 47.2, 33.1, 32.7, 29.3, 27.4, 25.5, 24.6, 20.3.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{17}\text{H}_{23}\text{NO}$ [$\text{M}+\text{H}]^+$: 258.1858, Found: 258.1865.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 95:5 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (minor) = 6.121 min, t_R (major) = 7.617 min.



(4R)-N-Cyclohexylbicyclo[2.2.1]hept-5-ene-2-carboxamide (38)

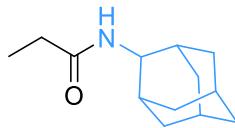
13.8 mg, 63% yield, 78.5:21.5 er, white solid, $[\alpha]_D^{24} = -21.6$ ($c = 0.5$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-6/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 6.11 (qd, $J = 5.5, 3.5$ Hz, 2H), 5.34 (s, 1H), 3.90 – 3.62 (m, 1H), 2.89 (d, $J = 1.5$ Hz, 2H), 1.97 – 1.85 (m, 4H), 1.80 – 1.56 (m, 4H), 1.46 – 1.03 (m, 8H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 174.6, 138.2, 136.1, 48.1, 47.3, 46.3, 44.9, 41.6, 33.4, 33.3, 30.5, 25.6, 24.9.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{14}\text{H}_{21}\text{NO}$ [$\text{M}+\text{H}]^+$: 220.1701, Found: 220.1700.

The enantiomeric excess was determined by Amylose2 column, heptane/EtOH = 90:10 v/v, $v = 0.5$ mL/min, $\lambda = 210$ nm, t_R (minor) = 6.202 min, t_R (major) = 6.589 min.



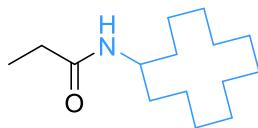
N-(Adamantan-2-yl)propionamide (39)

16.8 mg, 81% yield, white solid. Eluent: pentane/ethyl acetate = 5/1-2/1.

¹H NMR (300 MHz, CDCl₃) δ 5.75 (s, 1H), 4.09 – 4.00 (m, 1H), 2.22 (q, *J* = 7.5 Hz, 2H), 1.99 – 1.58 (m, 14H), 1.16 (t, *J* = 7.5 Hz, 3H).

¹³C NMR (75 MHz, CDCl₃) δ 172.8, 53.0, 37.5, 37.1, 32.0, 31.9, 30.1, 27.2, 27.1, 10.1.

HRMS (ESI-TOF) m/z: Calcd. for C₁₃H₂₁NO [M+H]⁺: 208.1701, Found: 208.1704.



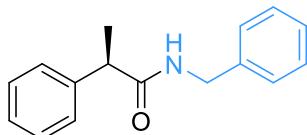
N-Cyclododecylpropionamide (40)

20.6 mg, 86% yield, white solid. Eluent: pentane/ethyl acetate = 5/1-2/1.

¹H NMR (300 MHz, CDCl₃) δ 5.20 (s, 1H), 4.16 – 3.88 (m, 1H), 2.16 (q, *J* = 7.5 Hz, 2H), 1.61 – 1.23 (m, 22H), 1.14 (t, *J* = 7.5 Hz, 3H).

¹³C NMR (75 MHz, CDCl₃) δ 173.0, 45.9, 30.3, 29.9, 24.0, 23.7, 23.5, 23.3, 21.4, 9.9.

HRMS (ESI-TOF) m/z: Calcd. for C₁₅H₂₉NO [M+H]⁺: 240.2327, Found: 240.2329.



(R)-N-benzyl-2-phenylpropanamide (3)

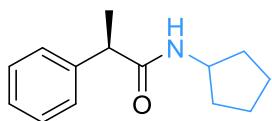
14.8 mg, 62% yield, 99.5:0.5 er, white solid, $[\alpha]_D^{24} = -8.6$ (*c* = 0.5, CHCl₃). Eluent: pentane/ethyl acetate = 10/1-5/1.

¹H NMR (300 MHz, CDCl₃) δ 7.39 – 7.25 (m, 8H), 7.17 (m, 2H), 5.73 (s, 1H), 4.41 (dd, *J* = 6.0, 4.0 Hz, 2H), 3.63 (q, *J* = 7.0 Hz, 1H), 1.58 (d, *J* = 7.0 Hz, 3H).

¹³C NMR (75 MHz, CDCl₃) δ 174.1, 141.3, 138.3, 129.0, 128.6, 127.7, 127.5, 127.4, 127.3, 47.2, 43.6, 18.6.

HRMS (ESI-TOF) m/z: Calcd. for C₁₆H₁₈NO [M+H]⁺: 240.1383, Found: 240.1380.

The enantiomeric excess was determined by Chiralpak AD-H column, heptane/EtOH = 80:20 v/v, *v* = 0.5 mL/min, λ = 210 nm, t_R (minor) = 13.230 min, t_R (major) = 17.448 min.



(R)-N-Cyclopentyl-2-phenylpropanamide (41)

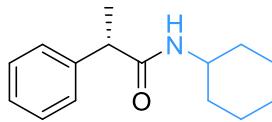
16.1 mg, 74% yield, 94:6 er, white solid, $[\alpha]_D^{24} = -12.2$ ($c = 0.5$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-5/1.

$^1\text{H NMR}$ (400 MHz, CDCl_3) δ 7.31 – 7.17 (m, 5H), 5.15 (s, 1H), 4.08 (h, $J = 7.0$ Hz, 1H), 3.43 (q, $J = 7.0$ Hz, 1H), 1.95 – 1.76 (m, 2H), 1.52 – 1.42 (m, 4H), 1.43 (d, $J = 7.0$ Hz, 3H), 1.22 – 1.04 (m, 2H).

$^{13}\text{C NMR}$ (101 MHz, CDCl_3) δ 173.7, 141.6, 128.9, 127.6, 127.2, 51.3, 47.2, 33.1, 33.0, 23.7, 23.6, 18.6.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{14}\text{H}_{19}\text{NO}$ [$\text{M}+\text{H}]^+$: 218.1545, Found: 218.1542.

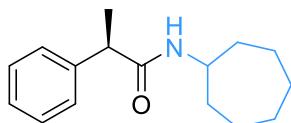
The enantiomeric excess was determined by Chiraldak AD-H column, heptane/EtOH = 80:20 v/v, $v = 0.5$ mL/min, $\lambda = 210$ nm, t_R (minor) = 7.364 min, t_R (major) = 7.822 min.



(S)-N-Cyclohexyl-2-phenylpropanamide (42)

19.2 mg, 83% yield, 95:5 er, white solid, $[\alpha]_D^{24} = +3.75$ ($c = 0.4$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-4/1.

The enantiomeric excess was determined by Cellulose4 column, heptane/iPrOH = 95:5 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (minor) = 9.008 min, t_R (major) = 10.194 min.



(R)-N-Cycloheptyl-2-phenylpropanamide (43)

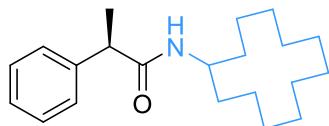
17.2 mg, 70% yield, 95:5 er, white solid, $[\alpha]_D^{24} = -13.9$ ($c = 0.63$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-4/1.

$^1\text{H NMR}$ (400 MHz, CDCl_3) δ 7.30 – 7.24 (m, 2H), 7.23 – 7.17 (m, 3H), 5.16 (s, 1H), 3.94 – 3.76 (m, 1H), 3.43 (q, $J = 7.0$ Hz, 1H), 1.83 – 1.64 (m, 2H), 1.52 – 1.31 (m, 11H), 1.28 – 1.14 (m, 2H).

$^{13}\text{C NMR}$ (101 MHz, CDCl_3) δ 172.9, 141.7, 128.9, 127.6, 127.2, 50.4, 47.2, 34.9, 34.8, 27.8, 24.1, 24.0, 18.6.

HRMS (ESI-TOF) m/z: Calcd. for C₁₆H₂₃NO [M+H]⁺: 246.1858, Found: 246.1857.

The enantiomeric excess was determined by Chiralcel OJ-H column, heptane/EtOH = 95:5 v/v, ν = 1.0 mL/min, λ = 210 nm, t_R (major) = 8.087 min, t_R (minor) = 9.675 min.



(R)-N-Cyclododecyl-2-phenylpropanamide (44)

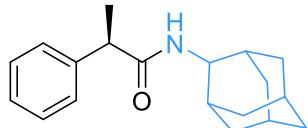
29.0 mg, 73% yield, 96.5:3.5 er, white solid, $[\alpha]_D^{24} = -19.1$ (c = 0.8, CHCl₃). Eluent: pentane/ethyl acetate = 10/1-5/1.

¹H NMR (400 MHz, CDCl₃) δ 7.30 – 7.17 (m, 5H), 5.02 (d, J = 8.0 Hz, 1H), 3.91 – 3.98 (m, 1H), 3.44 (q, J = 7.0 Hz, 1H), 1.44 (d, J = 7.0 Hz, 3H), 1.41 – 1.09 (m, 23H).

¹³C NMR (101 MHz, CDCl₃) δ 173.4, 141.7, 128.9, 127.6, 127.1, 47.2, 46.0, 30.2, 30.0, 23.93, 23.91, 23.7, 23.5, 23.5, 23.4, 23.3, 21.4, 21.2, 18.6.

HRMS (ESI-TOF) m/z: Calcd. for C₂₁H₃₃NO [M+H]⁺: 316.2640, Found: 316.2643.

The enantiomeric excess was determined by Chiralcel OJ-H column, heptane/EtOH = 95:5 v/v, ν = 0.6 mL/min, λ = 210 nm, t_R (major) = 8.077 min, t_R (minor) = 8.495 min.



(R)-N-(Adamantan-2-yl)-2-phenylpropanamide (45)

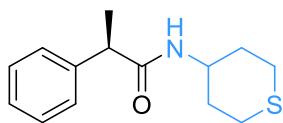
27.2 mg, 96% yield, 99:1 er, white solid, $[\alpha]_D^{24} = -10.7$ (c = 0.97, CHCl₃). Eluent: pentane/ethyl acetate = 10/1-5/1.

¹H NMR (400 MHz, CDCl₃) δ 7.43 – 7.26 (m, 5H), 5.65 (s, 1H), 4.00 (d, J = 8.0 Hz, 1H), 3.61 (q, J = 7.0 Hz, 1H), 1.87 – 1.77 (m, 7H), 1.75 – 1.66 (m, 3H), 1.57 (d, J = 7.0 Hz, 3H), 1.54 – 1.32 (m, 4H).

¹³C NMR (101 MHz, CDCl₃) δ 173.2, 141.6, 128.9, 127.7, 127.3, 53.1, 47.4, 37.5, 37.1, 37.0, 31.83, 31.79, 31.7, 27.1, 27.0, 18.3.

HRMS (ESI-TOF) m/z: Calcd. for C₁₉H₂₅NO [M+Na]⁺: 284.2014, Found: 284.2017.

The enantiomeric excess was determined by Chiralcel OJ-H column, heptane/EtOH = 95:5 v/v, ν = 1.0 mL/min, λ = 210 nm, t_R (major) = 7.079 min, t_R (minor) = 8.602 min.



(R)-2-phenyl-N-(tetrahydro-2H-thiopyran-4-yl)propanamide (46)

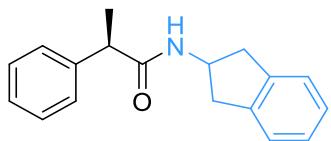
22.4 mg, 90% yield, 93.5:6.5 er, white solid, $[\alpha]_D^{24} = -15.0$ ($c = 0.47$, CHCl_3). Eluent: pentane/ethyl acetate = 6/1-2/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.40 – 7.10 (m, 5H), 5.20 (d, $J = 7.0$ Hz, 1H), 3.85 – 3.68 (m, 1H), 3.51 (q, $J = 7.0$ Hz, 1H), 2.78 – 2.45 (m, 4H), 2.27 – 1.98 (m, 2H), 1.50 (d, $J = 7.0$ Hz, 3H), 1.45 – 1.34 (m, 2H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.2, 141.4, 129.0, 127.5, 127.3, 47.5, 47.2, 34.0, 33.9, 27.7, 27.6, 18.5.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{14}\text{H}_{19}\text{NOS}$ [$\text{M}+\text{H}]^+$: 250.1265, Found: 250.1263.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 95:5 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (major) = 9.014 min, t_R (minor) = 10.639 min.



(R)-N-(2,3-Dihydro-1H-inden-2-yl)-2-phenylpropanamide (47)

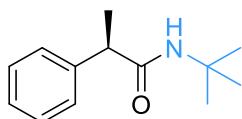
16.4 mg, 62% yield, 91.5:8.5 er, pale yellow solid, $[\alpha]_D^{24} = -9.0$ ($c = 0.4$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-4/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.35 – 7.11 (m, 9H), 5.56 (d, $J = 6.5$ Hz, 1H), 4.77 – 4.65 (m, 1H), 3.48 (q, $J = 7.0$ Hz, 1H), 3.28 (ddd, $J = 16.5, 9.5, 7.5$ Hz, 2H), 2.66 (ddd, $J = 36.0, 16.0, 5.0$ Hz, 2H), 1.50 (d, $J = 7.0$ Hz, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.9, 141.4, 140.8, 140.8, 128.9, 127.5, 127.2, 126.7, 124.7, 50.6, 47.1, 40.1, 40.0, 18.8.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{18}\text{H}_{19}\text{NO}$ [$\text{M}+\text{H}]^+$: 266.1545, Found: 266.1544.

The enantiomeric excess was determined by Cellulose4 column, heptane/EtOH = 95:5 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (minor) = 21.228 min, t_R (major) = 26.005 min.



(*R*)-*N*-(*tert*-Butyl)-2-phenylpropanamide (48)

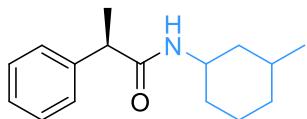
15.6 mg, 76% yield, 95.5:4.5 er, white solid, $[\alpha]_D^{24} = -23.2$ ($c = 0.67$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-5/1.

$^1\text{H NMR}$ (400 MHz, CDCl_3) δ 7.15 – 7.09 (m, 2H), 7.08 – 7.01 (m, 3H), 4.92 (s, 1H), 3.23 (q, $J = 7.0$ Hz, 1H), 1.26 (d, $J = 7.0$ Hz, 3H), 1.05 (s, 9H).

$^{13}\text{C NMR}$ (101 MHz, CDCl_3) δ 173.4, 142.0, 128.8, 127.5, 127.0, 51.1, 47.8, 28.7, 18.7.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{13}\text{H}_{19}\text{NO}$ [$\text{M}+\text{H}]^+$: 206.1545, Found: 206.1550.

The enantiomeric excess was determined by Chiralcel OJ-H column, heptane/EtOH = 95:5 v/v, $v = 1.0$ mL/min, $\lambda = 210$ nm, t_R (major) = 8.003 min, t_R (minor) = 10.066 min.



(2*R*)-*N*-(3-Methylcyclohexyl)-2-phenylpropanamide (49)

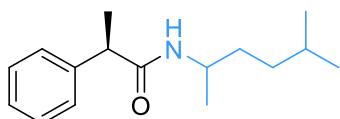
18.2 mg, 74% yield, *dr* 1.5:1, colorless oil, $[\alpha]_D^{24} = -10.9$ ($c = 0.36$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-5/1.

Mixture of two stereoisomers.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.33 – 7.16 (m, 5H), 5.31 (s, 0.6H), 5.03 (s, 0.4H), 4.09 – 3.93 (m, 0.6H), 3.74 – 3.56 (m, 0.4H), 3.52 – 3.37 (m, 1H), 1.88 – 1.54 (m, 2H), 1.52 – 1.36 (m, 6H), 1.27 – 1.09 (m, 3H), 0.85 – 0.73 (m, 3H), 0.72 – 0.45 (m, 1H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 173.2, 141.7, 128.9, 128.9, 127.6, 127.2, 127.1, 48.5, 47.3, 47.2, 44.7, 41.7, 38.5, 38.5, 34.2, 33.6, 32.8, 31.8, 30.4, 27.4, 27.4, 24.8, 22.4, 21.7, 20.63, 20.61, 18.7, 18.4.

HRMS (ESI-TOF) m/z: Calcd. for $\text{C}_{16}\text{H}_{23}\text{NO}$ [$\text{M}+\text{H}]^+$: 246.1858, Found: 246.1857.



(2*R*)-*N*-(5-Methylhexan-2-yl)-2-phenylpropanamide (50)

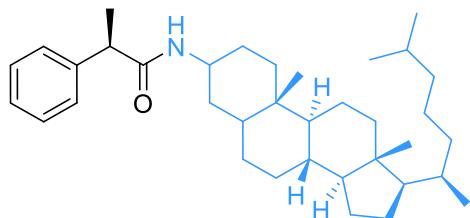
18.8 mg, 76% yield, *dr* 1:1, colorless oil, $[\alpha]_D^{24} = -13.5$ ($c = 0.73$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-5/1.

Mixture of two stereoisomers.

¹H NMR (400 MHz, CDCl₃) δ 7.39 – 7.10 (m, 5H), 5.01 (s, 1H), 3.92 – 3.70 (m, 1H), 3.45 (dq, *J* = 12.0, 7.0 Hz, 1H), 1.44 (dd, *J* = 7.0, 2.5 Hz, 3H), 1.42 – 1.11 (m, 3H), 1.06 – 0.85 (m, 5H), 0.76 (dd, *J* = 6.5, 3.5 Hz, 3H), 0.70 (dd, *J* = 6.5, 4.5 Hz, 3H).

¹³C NMR (101 MHz, CDCl₃) δ 173.5, 173.3, 141.8, 141.6, 128.9, 127.64, 127.61, 127.2, 47.3, 47.2, 45.5, 45.3, 35.0, 34.8, 34.6, 27.9, 27.7, 22.6, 22.54, 22.53, 22.4, 21.0, 20.8, 18.6, 18.5.

HRMS (ESI-TOF) m/z: Calcd. for C₁₆H₂₅NO [M+H]⁺: 248.2014, Found: 248.2018.



(2*R*)-N-((8*R*,9*S*,10*S*,13*R*,14*S*,17*R*)-10,13-dimethyl-17-((*R*)-6-methylheptan-2-yl)hexadecahydro-1*H*-cyclopenta[a]phenanthren-3-yl)-2-phenylpropanamide (51)

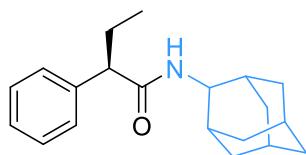
43.7 mg, 84% yield, *dr* 1:1, white solid, $[\alpha]_D^{24} = +14.3$ (c = 1.4, CHCl₃). Eluent: pentane/ethyl acetate = 10/1-4/1.

Mixture of two stereoisomers.

¹H NMR (400 MHz, CDCl₃) δ 7.29 – 7.01 (m, 5H), 5.39 (d, *J* = 7.5 Hz, 1H), 3.89 (d, *J* = 3.5 Hz, 1H), 3.42 (q, *J* = 7.0 Hz, 1H), 1.82 (dt, *J* = 12.5, 3.0 Hz, 1H), 1.75 – 1.61 (m, 1H), 1.53 – 1.34 (m, 8H), 1.31 – 0.82 (m, 20H), 0.78 – 0.72 (m, 9H), 0.64 – 0.45 (m, 9H), 0.34 – 0.23 (m, 1H).

¹³C NMR (75 MHz, CDCl₃) δ 173.3, 173.2, 141.9, 141.7, 133.4, 128.94, 128.91, 127.64, 127.62, 127.3, 127.1, 56.6, 56.4, 56.34, 56.32, 54.6, 54.2, 49.0, 47.3, 45.3, 44.7, 42.6, 40.8, 40.1, 40.0, 39.5, 37.3, 36.2, 35.9, 35.8, 35.5, 35.4, 35.4, 35.3, 33.2, 32.5, 31.9, 28.6, 28.6, 28.4, 28.2, 28.0, 25.8, 24.2, 23.9, 22.9, 22.6, 21.1, 20.7, 18.7, 18.1, 12.3, 12.1, 11.4.

HRMS (ESI-TOF) m/z: Calcd. for C₃₆H₅₇NO [M+H]⁺: 520.4518, Found: 520.4510.



(R)-N-adamantan-2-yl-2-phenylbutanamide (52)

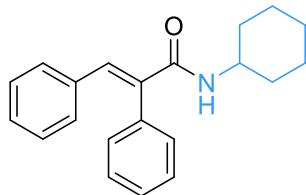
20.5 mg, 69% yield, 98:2 er, white solid, $[\alpha]_D^{24} = -21.4$ ($c = 0.27$, CHCl_3). Eluent: pentane/ethyl acetate = 10/1-5/1.

$^1\text{H NMR}$ (400 MHz, CDCl_3) δ 7.42 – 7.20 (m, 5H), 5.68 (s, 1H), 4.01 (d, $J = 7.5$ Hz, 1H), 3.41 – 3.20 (m, 1H), 2.25 (dt, $J = 14.0, 7.0$ Hz, 1H), 1.96 – 1.77 (m, 6H), 1.72 (d, $J = 16.0$ Hz, 3H), 1.60 – 1.43 (m, 3H), 1.38 – 1.23 (m, 3H), 0.92 (t, $J = 7.5$ Hz, 3H).

$^{13}\text{C NMR}$ (101 MHz, CDCl_3) δ 172.6, 140.2, 128.8, 128.0, 127.2, 55.5, 53.1, 37.5, 37.0, 36.9, 31.83, 31.81, 31.78, 31.7, 27.2, 27.1, 26.2, 12.5.

The enantiomeric excess was determined by Amylose2 column, heptane/EtOH = 90:10 v/v, $v = 0.5$ mL/min, $\lambda = 210$ nm, t_R (minor) = 7.343 min, t_R (major) = 9.026 min.

HRMS (EI) m/z: Calcd. for $\text{C}_{20}\text{H}_{27}\text{NO}$ [M]: 297.20293, Found: 297.20276.



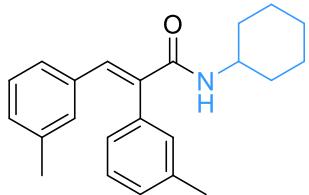
(E)-N-Cyclohexyl-2,3-diphenylacrylamide (53)

12.5 mg, 41% yield, white solid. Eluent: pentane/ethyl acetate = 10/1-5/1.

$^1\text{H NMR}$ (300 MHz, CDCl_3) δ 7.76 (s, 1H), 7.41 – 7.32 (m, 3H), 7.24 – 7.13 (m, 2H), 7.11 – 7.00 (m, 3H), 6.96 – 6.87 (m, 2H), 5.27 (d, $J = 8.0$ Hz, 1H), 3.98 – 3.64 (m, 1H), 1.86 – 1.74 (m, 2H), 1.60 – 1.43 (m, 3H), 1.37 – 1.22 (m, 2H), 1.09 – 0.89 (m, 3H).

$^{13}\text{C NMR}$ (75 MHz, CDCl_3) δ 166.1, 136.8, 136.4, 135.1, 134.7, 130.3, 129.9, 129.5, 128.5, 128.4, 128.1, 48.5, 32.8, 25.5, 24.6.

HRMS (EI) m/z: Calcd. for $\text{C}_{21}\text{H}_{23}\text{NO}$ [M]: 305.17742, Found: 305.17728.



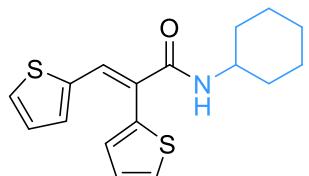
(E)-N-Cyclohexyl-2,3-di-m-tolylacrylamide (54)

14.3 mg, 43% yield, colorless oil solid. Eluent: pentane/ethyl acetate = 10/1-5/1.

¹H NMR (300 MHz, CDCl₃) δ 7.78 (s, 1H), 7.37 – 7.28 (m, 1H), 7.24 – 7.17 (m, 1H), 7.05 – 6.97 (m, 4H), 6.87 – 6.83 (m, 1H), 6.80 – 6.69 (m, 1H), 5.38 (d, *J* = 8.0 Hz, 1H), 4.09 – 3.65 (m, 1H), 2.35 (s, 3H), 2.17 (s, 3H), 1.93 – 1.80 (m, 3H), 1.57 (td, *J* = 9.0, 5.0 Hz, 3H), 1.40 – 1.28 (m, 2H), 1.17 – 0.95 (m, 2H).

¹³C NMR (75 MHz, CDCl₃) δ 166.2, 139.2, 137.6, 136.6, 136.4, 135.1, 134.6, 131.4, 130.3, 129.3, 129.1, 129.1, 127.9, 127.2, 126.8, 48.5, 32.8, 25.5, 24.6, 21.4, 21.3.

HRMS (EI) m/z: Calcd. for C₂₃H₂₇NO [M]: 333.20872, Found: 333.20832.



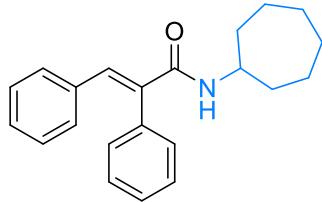
(Z)-N-Cyclohexyl-2,3-di(thiophen-2-yl)acrylamide (55)

15.2 mg, 48% yield, slight yellow oil. Eluent: pentane/ethyl acetate = 10/1-5/1.

¹H NMR (300 MHz, CDCl₃) δ 8.19 (s, 1H), 7.59 (dd, *J* = 5.0, 1.0 Hz, 1H), 7.30 – 7.25 (m, 1H), 7.23 – 7.15 (m, 2H), 7.03 (dd, *J* = 3.5, 1.0 Hz, 1H), 6.95 (dd, *J* = 5.0, 3.5 Hz, 1H), 5.58 (d, *J* = 8.0 Hz, 1H), 4.05 – 3.67 (m, 1H), 1.96 – 1.79 (m, 2H), 1.71 – 1.50 (m, 3H), 1.40 – 1.31 (m, 2H), 1.20 – 0.98 (m, 3H).

¹³C NMR (75 MHz, CDCl₃) δ 165.1, 138.7, 135.2, 134.1, 133.4, 130.5, 129.7, 129.1, 128.4, 126.7, 48.5, 32.8, 25.5, 24.6.

HRMS (EI) m/z: Calcd. for C₁₇H₁₉NOS₂ [M]: 317.09026, Found: 317.08984.



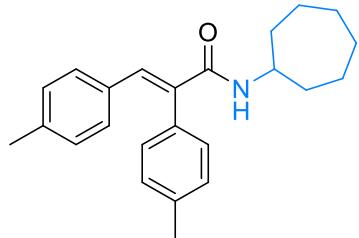
(E)-N-Cycloheptyl-2,3-diphenylacrylamide (56)

13.1 mg, 41% yield, colorless oil. Eluent: pentane/ethyl acetate = 10/1-5/1.

¹H NMR (300 MHz, CDCl₃) δ 7.62 (s, 1H), 7.27 – 7.17 (m, 3H), 7.07 – 6.99 (m, 2H), 6.96 – 6.86 (m, 3H), 6.81 – 6.72 (m, 2H), 5.20 (d, *J* = 8.0 Hz, 1H), 3.95 – 3.70 (m, 1H), 1.71 – 1.54 (m, 2H), 1.41 – 1.08 (m, 10H).

¹³C NMR (75 MHz, CDCl₃) δ 165.9, 136.7, 136.4, 135.1, 134.7, 130.3, 129.8, 129.5, 128.5, 128.4, 128.1, 50.7, 34.7, 27.8, 24.0.

HRMS (EI) m/z: Calcd. for C₂₂H₂₅NO [M]: 319.19307, Found: 319.19276.



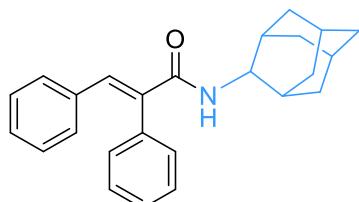
(E)-N-Cycloheptyl-2,3-di-p-tolylacrylamide (57)

16.0 mg, 46% yield, colorless oil. Eluent: pentane/ethyl acetate = 10/1-5/1.

¹H NMR (300 MHz, CDCl₃) δ 7.80 (s, 1H), 7.30 – 7.23 (m, 2H), 7.18 – 7.10 (m, 2H), 7.00 – 6.87 (m, 4H), 5.45 (d, *J* = 8.0 Hz, 1H), 4.24 – 3.91 (m, 1H), 2.44 (s, 3H), 2.28 (s, 3H), 1.95 – 1.78 (m, 2H), 1.63 – 1.27 (m, 10H).

¹³C NMR (75 MHz, CDCl₃) δ 166.2, 138.4, 138.1, 136.5, 133.7, 133.4, 132.4, 130.3, 130.2, 129.7, 128.9, 50.7, 34.8, 27.8, 24.1, 21.4, 21.3.

HRMS (EI) m/z: Calcd. for C₂₄H₂₉NO [M]: 347.22437, Found: 347.22358.



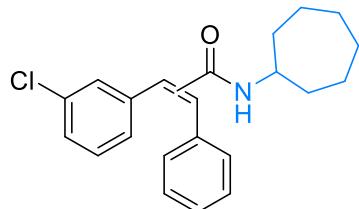
(E)-N-(Adamantan-2-yl)-2,3-diphenylacrylamide (58)

13.5 mg, 38% yield, white solid. Eluent: pentane/ethyl acetate = 10/1-5/1.

¹H NMR (300 MHz, CDCl₃) δ 7.85 (s, 1H), 7.49 – 7.40 (m, 3H), 7.33 – 7.27 (m, 2H), 7.19 – 7.12 (m, 3H), 7.07 – 7.00 (m, 2H), 5.85 (d, *J* = 8.0 Hz, 1H), 4.12 (d, *J* = 8.5 Hz, 1H), 1.84 (d, *J* = 13.5 Hz, 7H), 1.67 (s, 3H), 1.50 (d, *J* = 13.0 Hz, 2H), 1.38 – 1.20 (m, 2H).

^{13}C NMR (75 MHz, CDCl_3) δ 165.9, 136.6, 136.5, 135.1, 134.8, 130.4, 129.8, 129.6, 128.5, 128.4, 128.1, 53.6, 37.5, 37.0, 31.9, 31.8, 27.1, 27.0.

HRMS (EI) m/z: Calcd. for $\text{C}_{25}\text{H}_{27}\text{NO}$ [M]: 357.20872, Found: 357.20778.



(E)-3-(3-Chlorophenyl)-N-cycloheptyl-2-phenylacrylamide (59)

15.9 mg, 45% yield, mixture (1.5:1), colorless oil. Eluent: pentane/ethyl acetate = 10/1-5/1.

^1H NMR (300 MHz, CDCl_3) δ 7.79 (d, $J = 14.0$ Hz, 1H), 7.49 – 7.38 (m, 2H), 7.27 – 7.06 (m, 5H), 7.01 – 6.84 (m, 2H), 5.37 (dd, $J = 22.5, 8.0$ Hz, 1H), 4.35 – 3.89 (m, 1H), 1.95 – 1.77 (m, 2H), 1.64 – 1.31 (m, 10H).

^{13}C NMR (75 MHz, CDCl_3) δ 165.6, 165.5, 137.4, 136.0, 135.4, 135.3, 134.81, 134.77, 134.5, 134.2, 133.7, 133.6, 131.43, 131.35, 130.2, 129.8, 129.70, 129.67, 128.7, 128.6, 128.4, 128.3, 50.9, 50.8, 34.8, 34.7, 27.8, 27.8, 24.1, 24.0.

HRMS (EI) m/z: Calcd. for $\text{C}_{22}\text{H}_{24}\text{NOCl}$ [M]: 353.15409, Found: 353.15315.

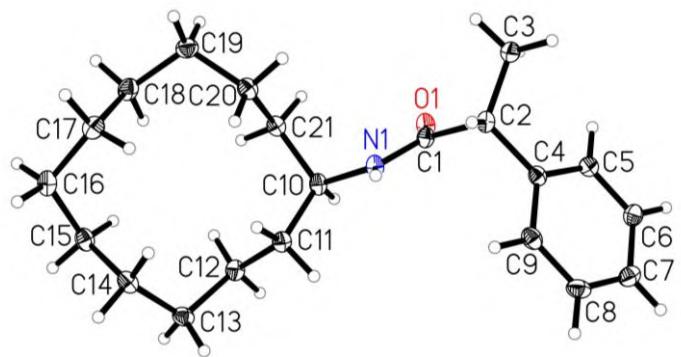
4. X-Ray Crystal Structures of 44

Data were collected on a Bruker Kappa APEX II Duo diffractometer. The structure was solved by direct methods (SHELXS-97: Sheldrick, G. M. *Acta Cryst.* **2008**, *A64*, 112.) and refined by full-matrix least-squares procedures on F^2 (SHELXL-2018: Sheldrick, G. M. *Acta Cryst.* **2015**, *C71*, 3.). XP (Bruker AXS) was used for graphical representations.

CCDC 2084474 contains the supplementary crystallographic data for this paper. These data are provided free of charge by the joint Cambridge Crystallographic Data Centre and Fachinformationszentrum Karlsruhe Access Structures service www.ccdc.cam.ac.uk/structures.

Crystal data of **44**: $\text{C}_{21}\text{H}_{33}\text{NO}$, $M = 315.48$, monoclinic, space group $P2_1$, $a = 5.0832(1)$, $b = 22.6617(7)$, $c = 8.1122(2)$ Å, $\beta = 103.2538(17)$ °, $V = 909.59(4)$ Å³, $T = 150(2)$ K, $Z = 2$, 9911 reflections measured, 3203 independent reflections ($R_{int} = 0.0316$), final R values ($I > 2\sigma(I)$): $R_1 = 0.0329$, $wR_2 = 0.0831$, final R values (all data): $R_1 = 0.0346$, $wR_2 = 0.0847$, 213 parameters, largest diff. peak/hole: 0.13/-0.16 eÅ⁻³.

The absolute configuration was assigned as **R** which is consistent with the determined Flack parameter [0.01(12)].



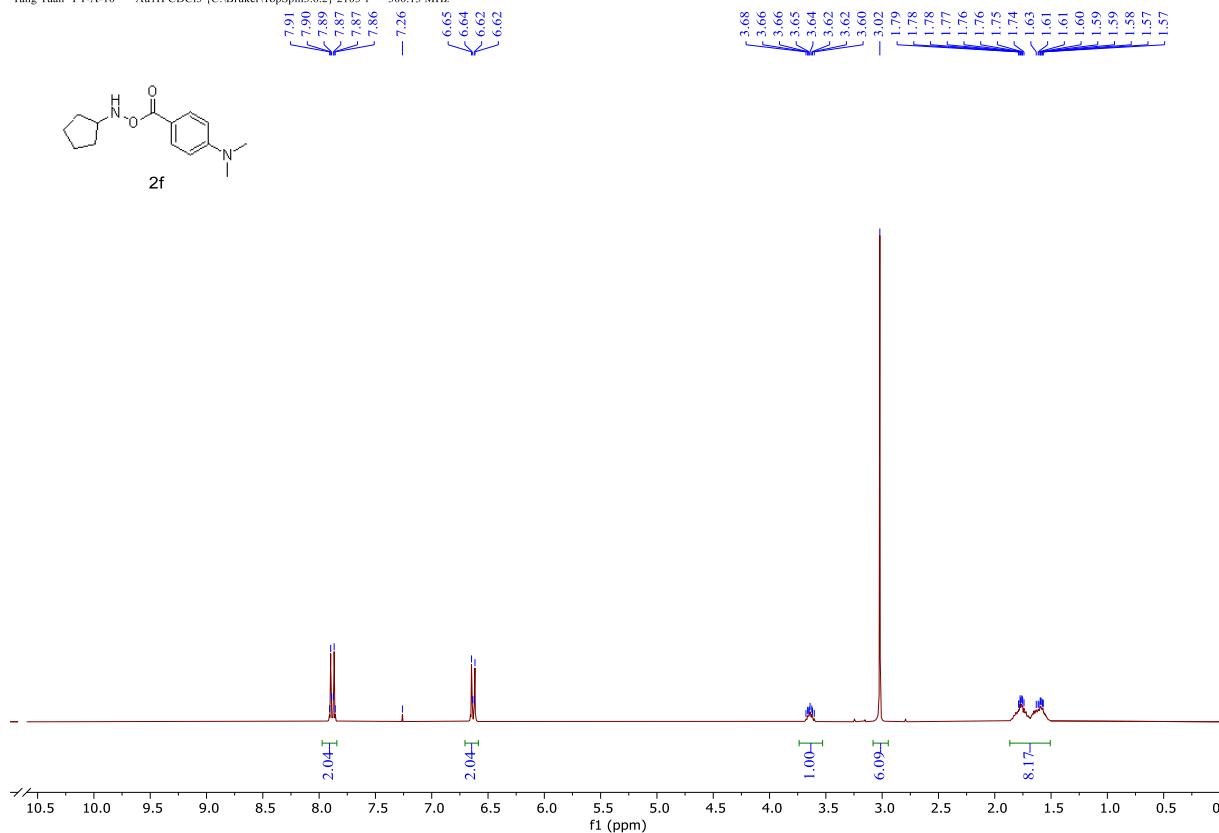
ORTEP representation of **44**. Displacement ellipsoids correspond to 30% probability.

5. References

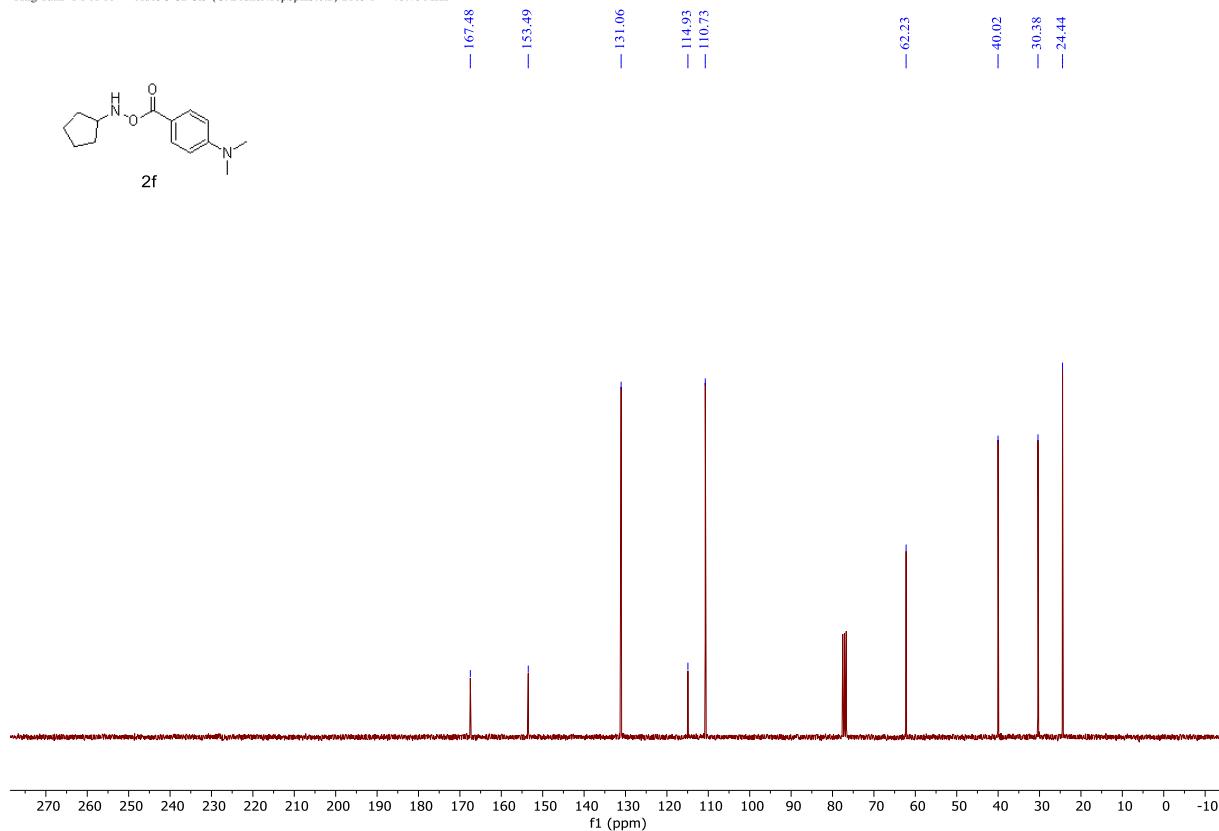
- (1) D. Niu, S. L. Buchwald, *J. Am. Chem. Soc.* **2015**, *137*, 9716-9721.
- (1) G. X. Ortiz, B. N. Hemric, Q. Wang, *Org. Lett.* **2017**, *19*, 1314-1317.

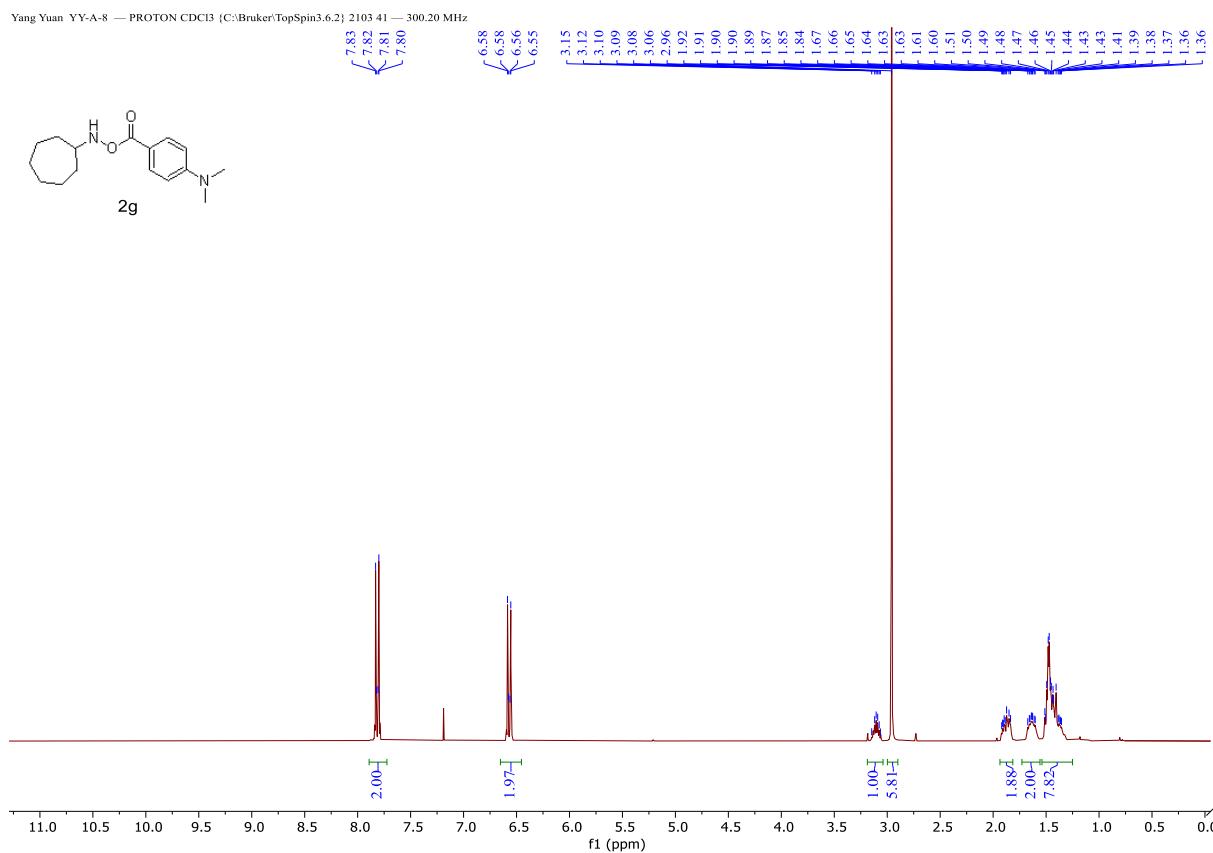
6. NMR Spectra

Yang Yuan YY-A-10 — Au1H CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 1 — 300.13 MHz

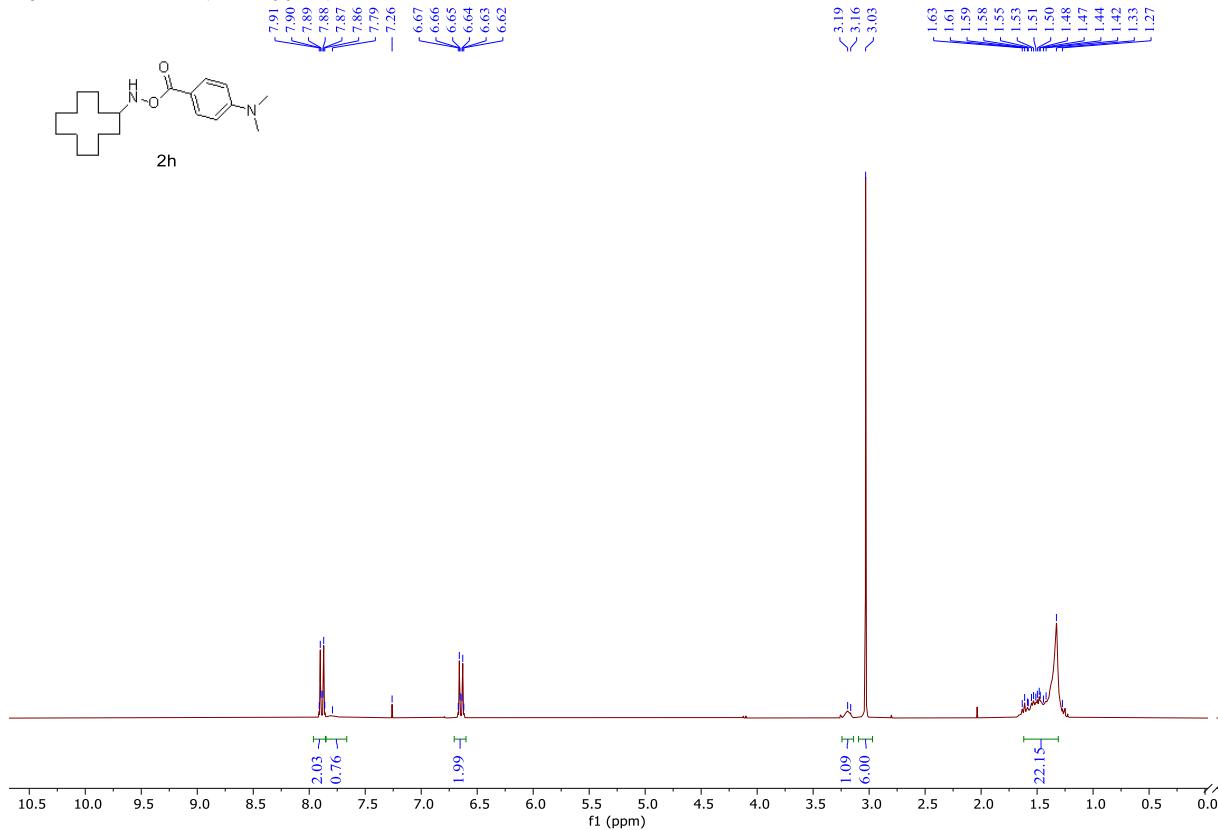


Yang Yuan YY-A-10 — Au13C CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 1 — 75.48 MHz

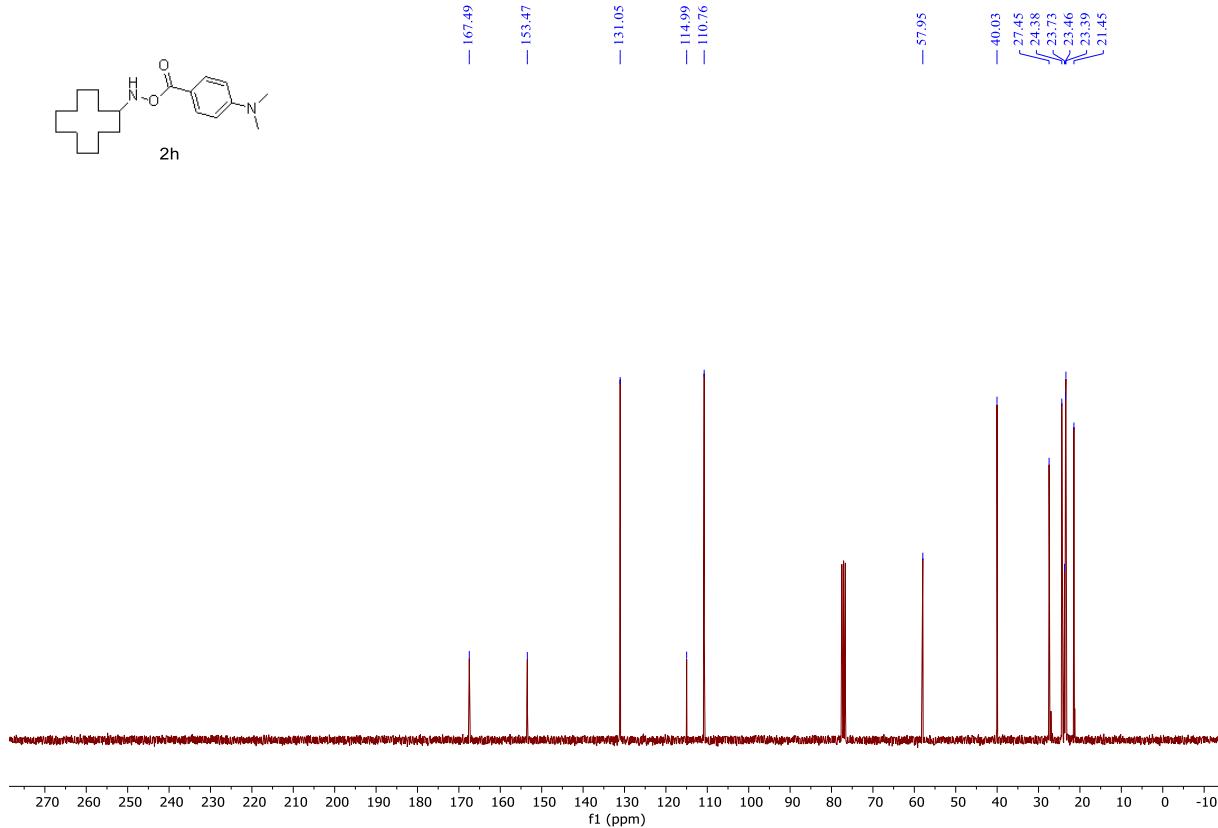


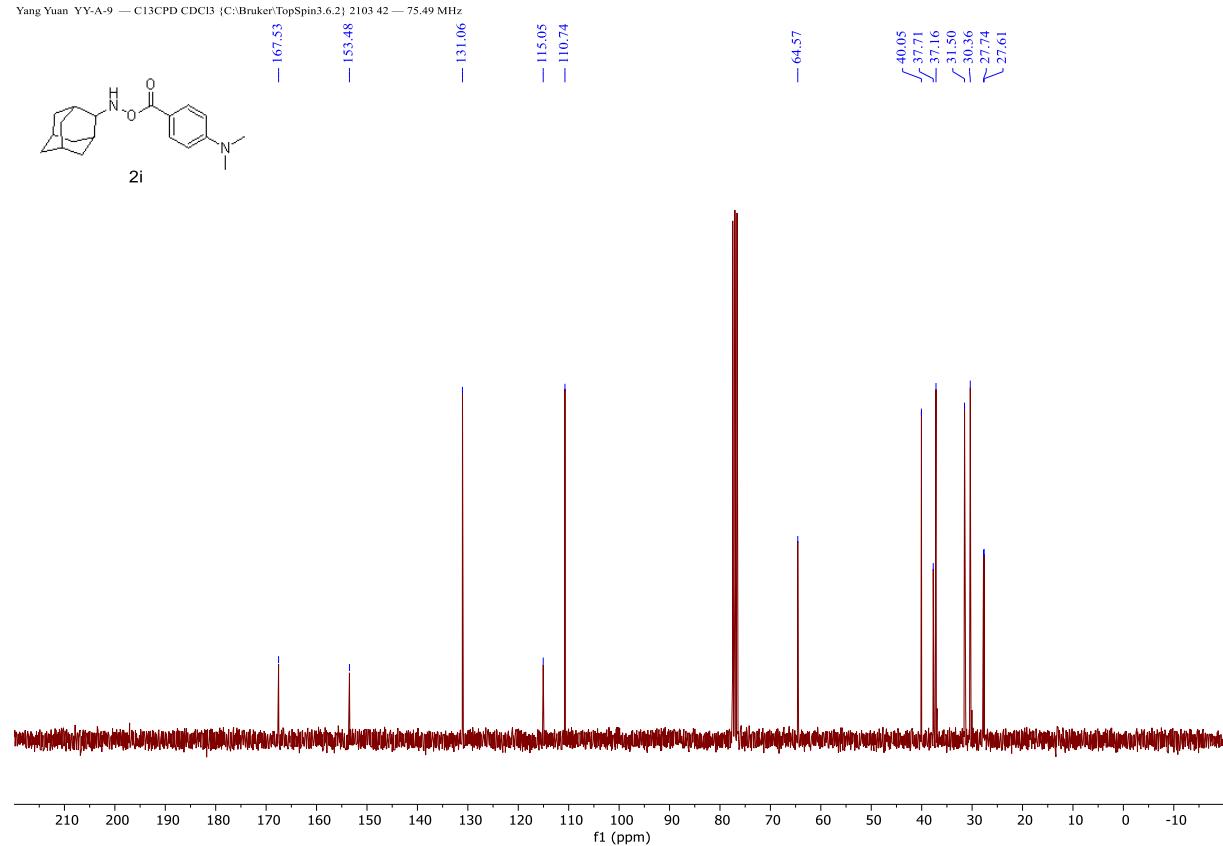
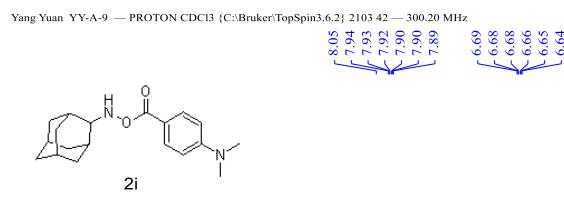


Yang Yuan YY-A-11 — Au1H CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103.2 — 300.13 MHz

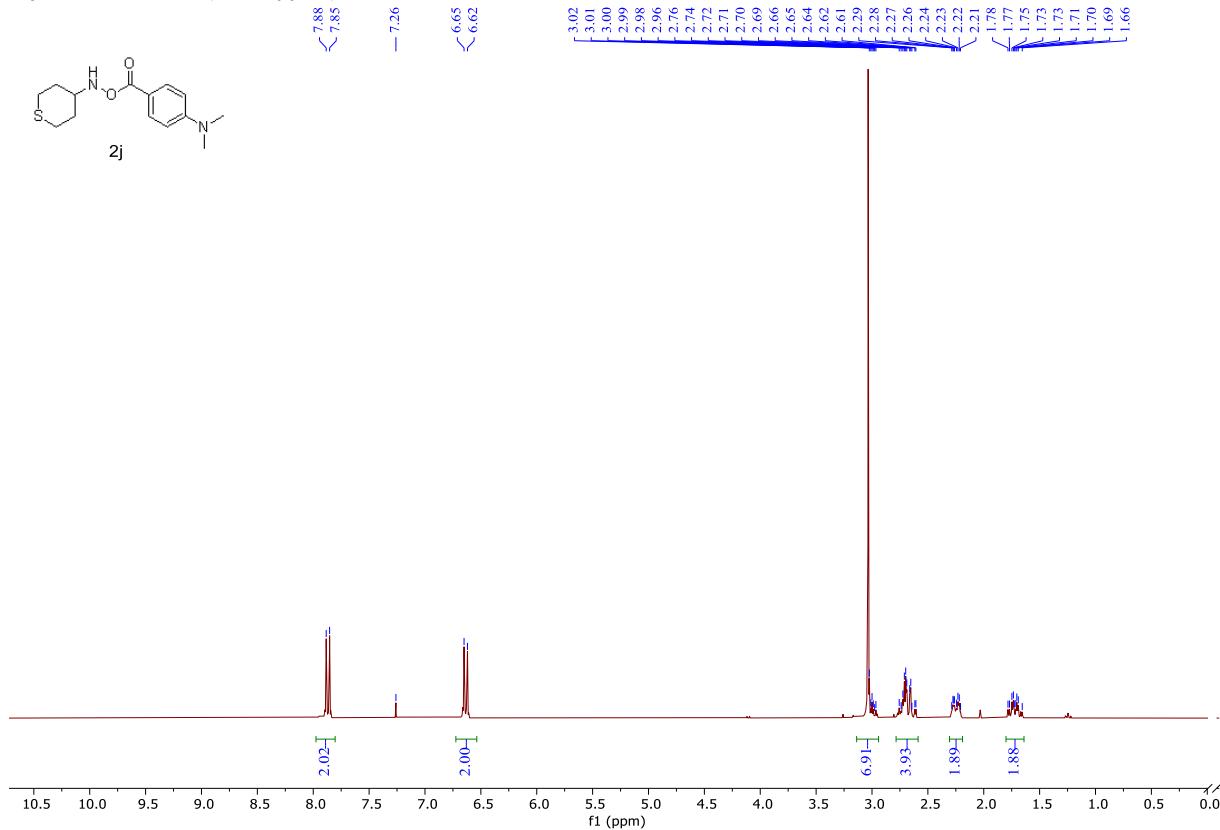


Yang Yuan YY-A-11 — Au13C CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103.2 — 75.48 MHz

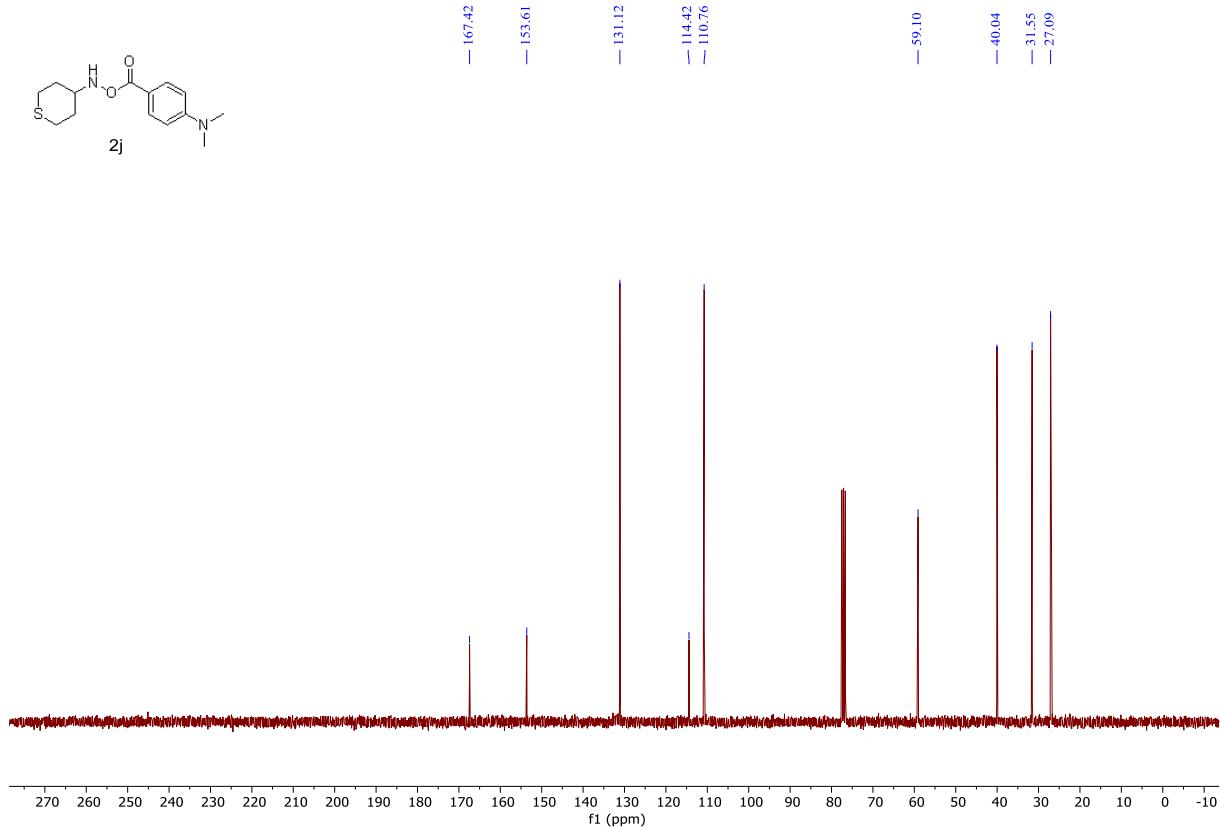




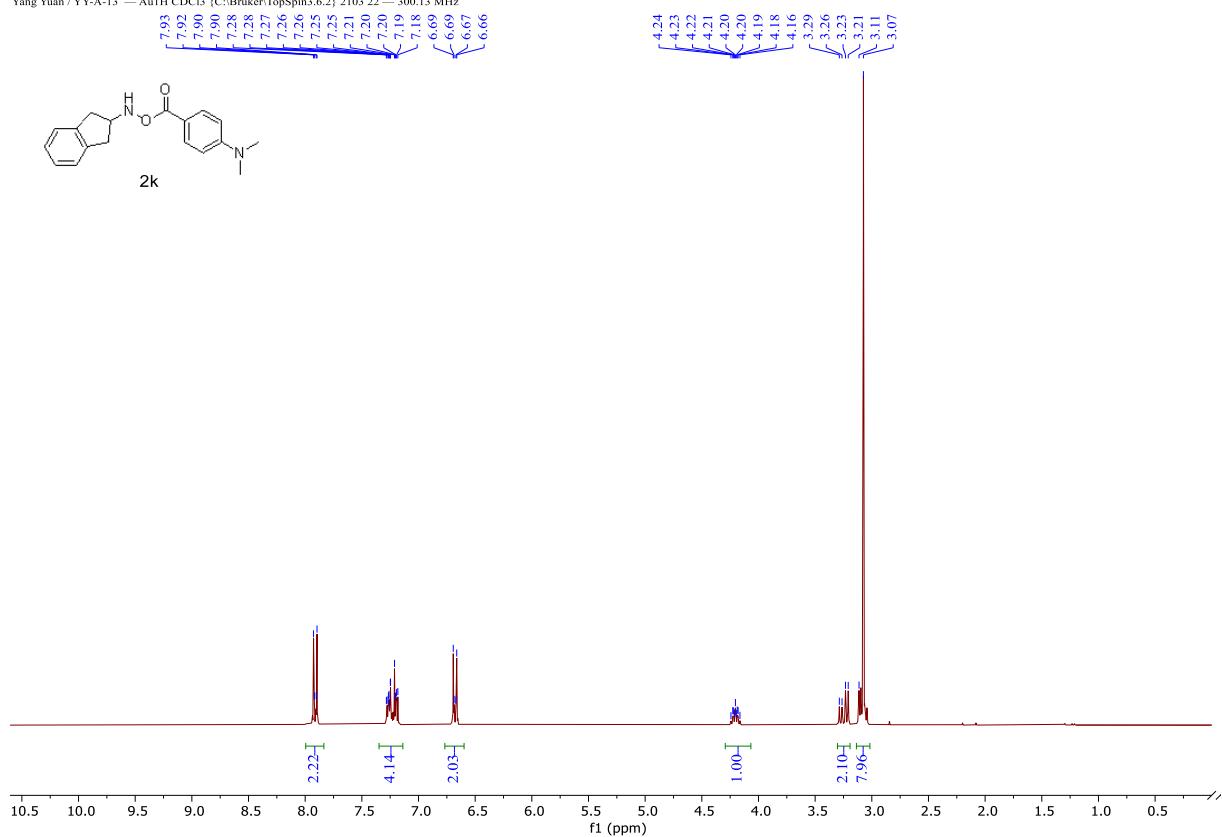
Yang Yuan / YY-A-15 — Au1H CDCl₃ {C:\Bruker\TopSpin3.6.2\} 2103 23 — 300.13 MHz



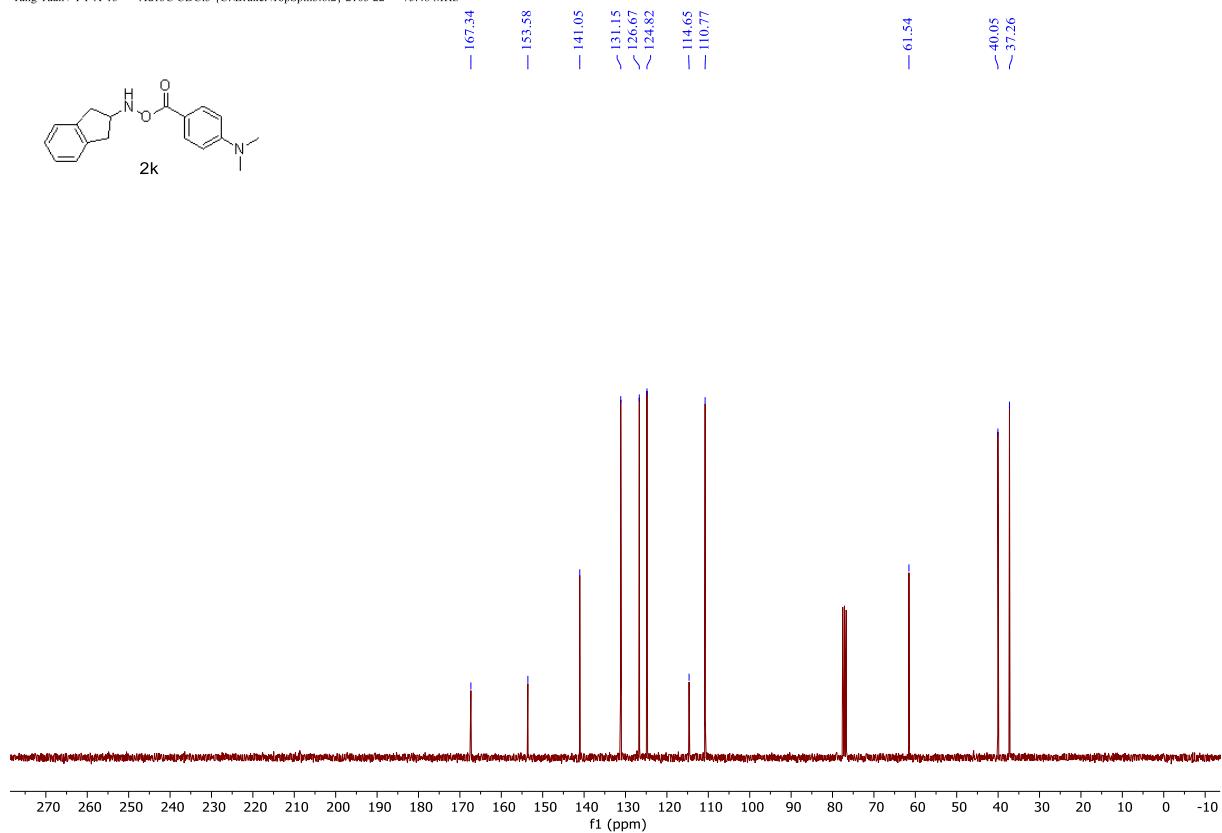
Yang Yuan / YY-A-15 — Au13C CDCl₃ {C:\Bruker\TopSpin3.6.2\} 2103 23 — 75.48 MHz



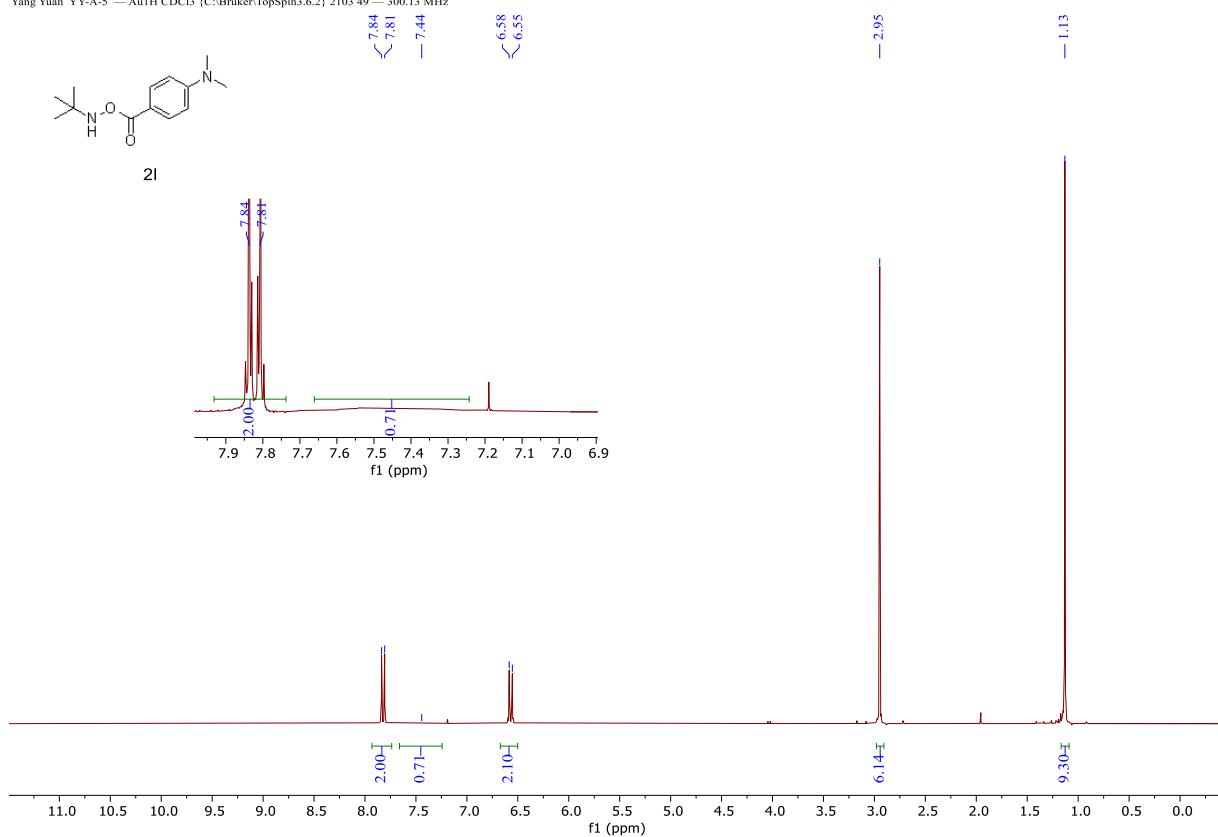
Yang Yuan / YY-A-13 — Au1H CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 22 — 300.13 MHz



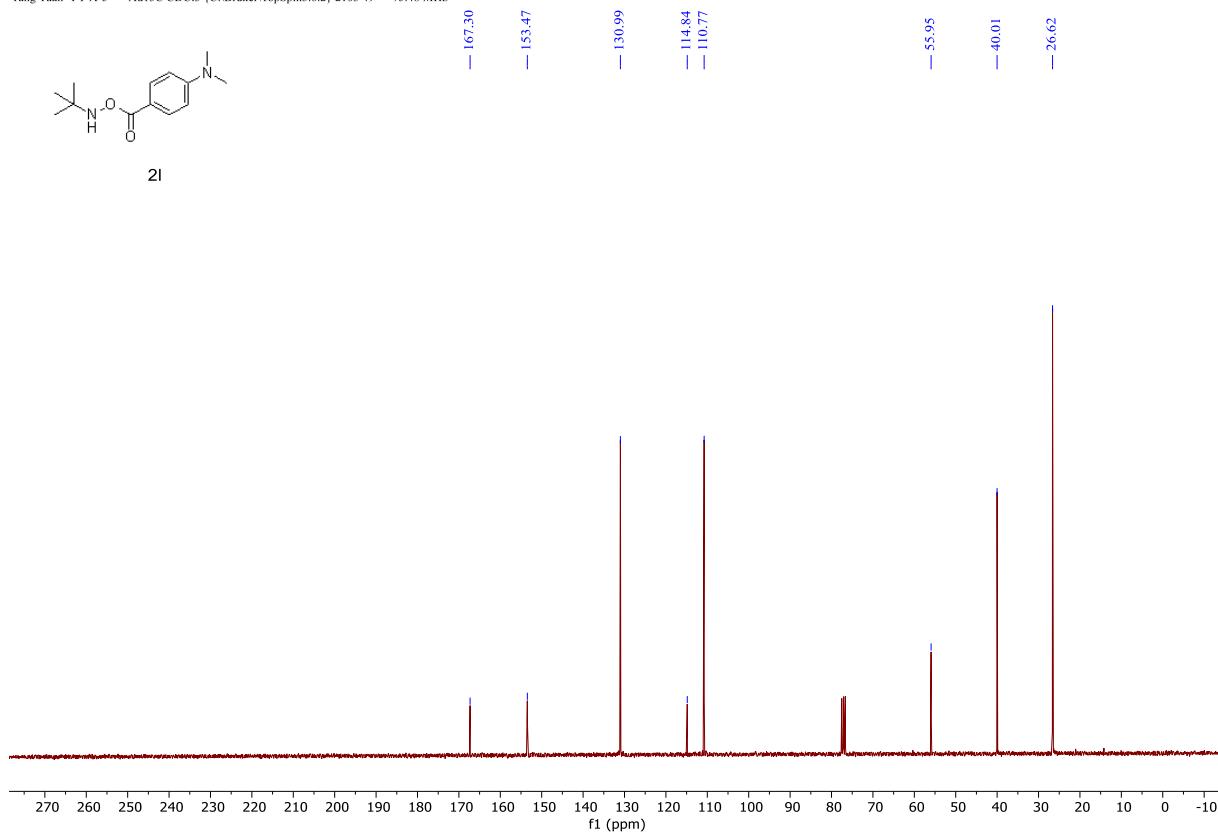
Yang Yuan / YY-A-13 — Au13C CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 22 — 75.48 MHz

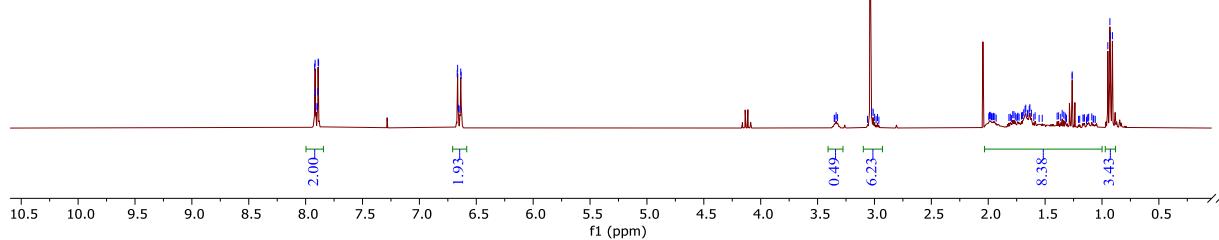
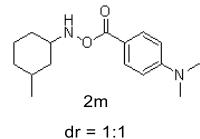


Yang Yuan YY-A-5 — Au1H CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 49 — 300.13 MHz

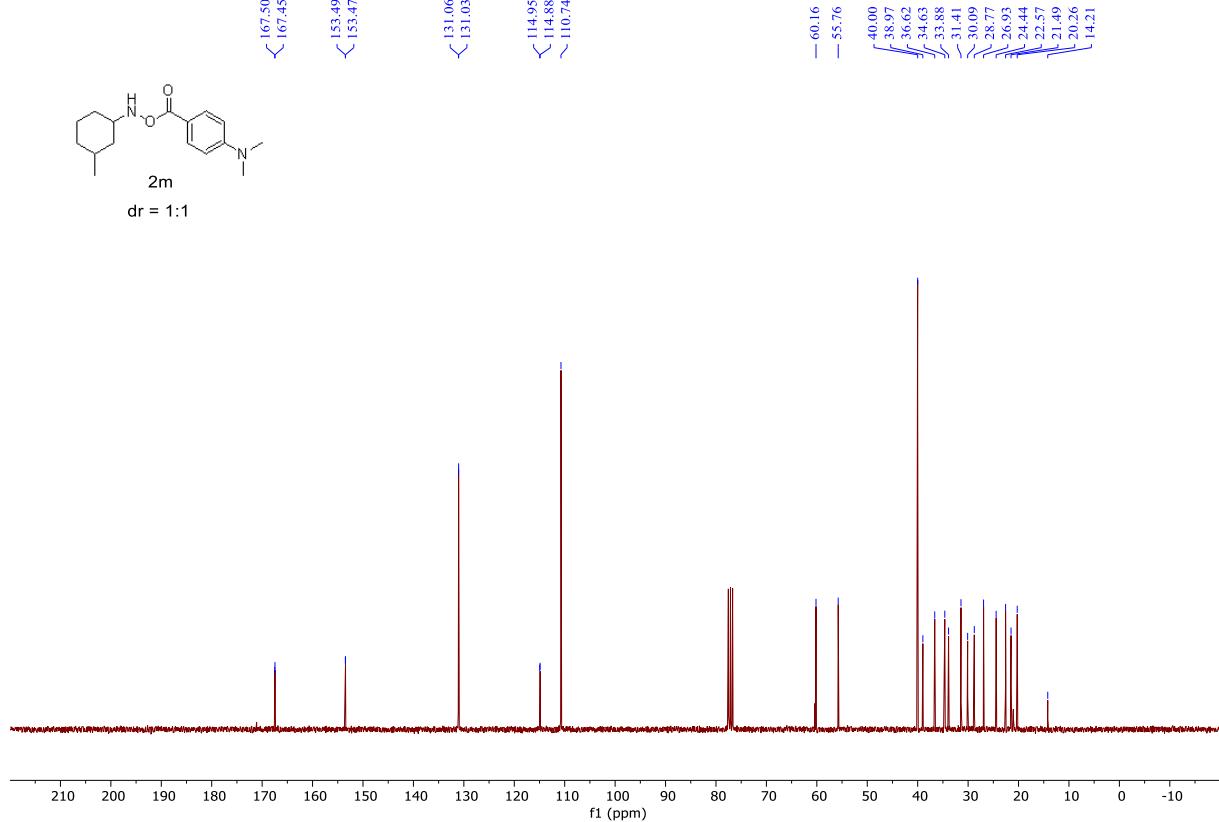
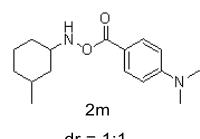


Yang Yuan YY-A-5 — Au13C CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 49 — 75.48 MHz

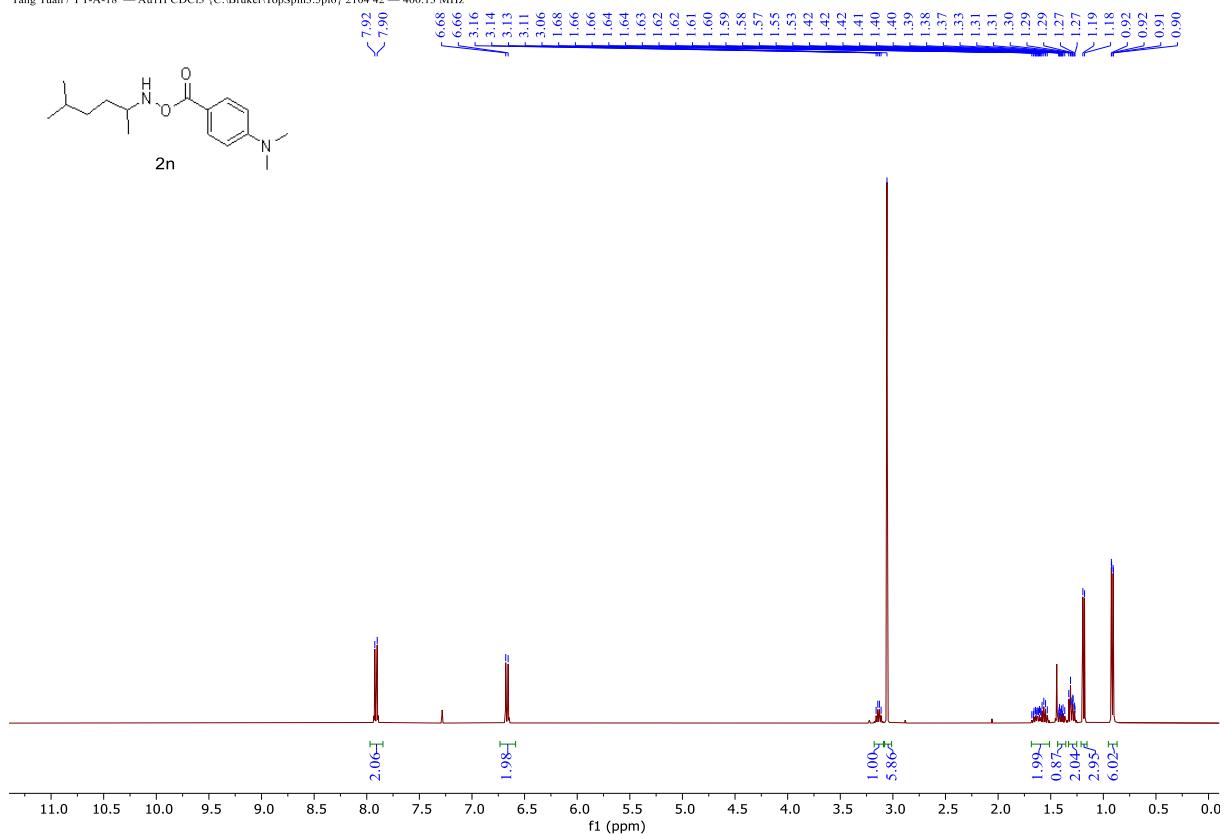




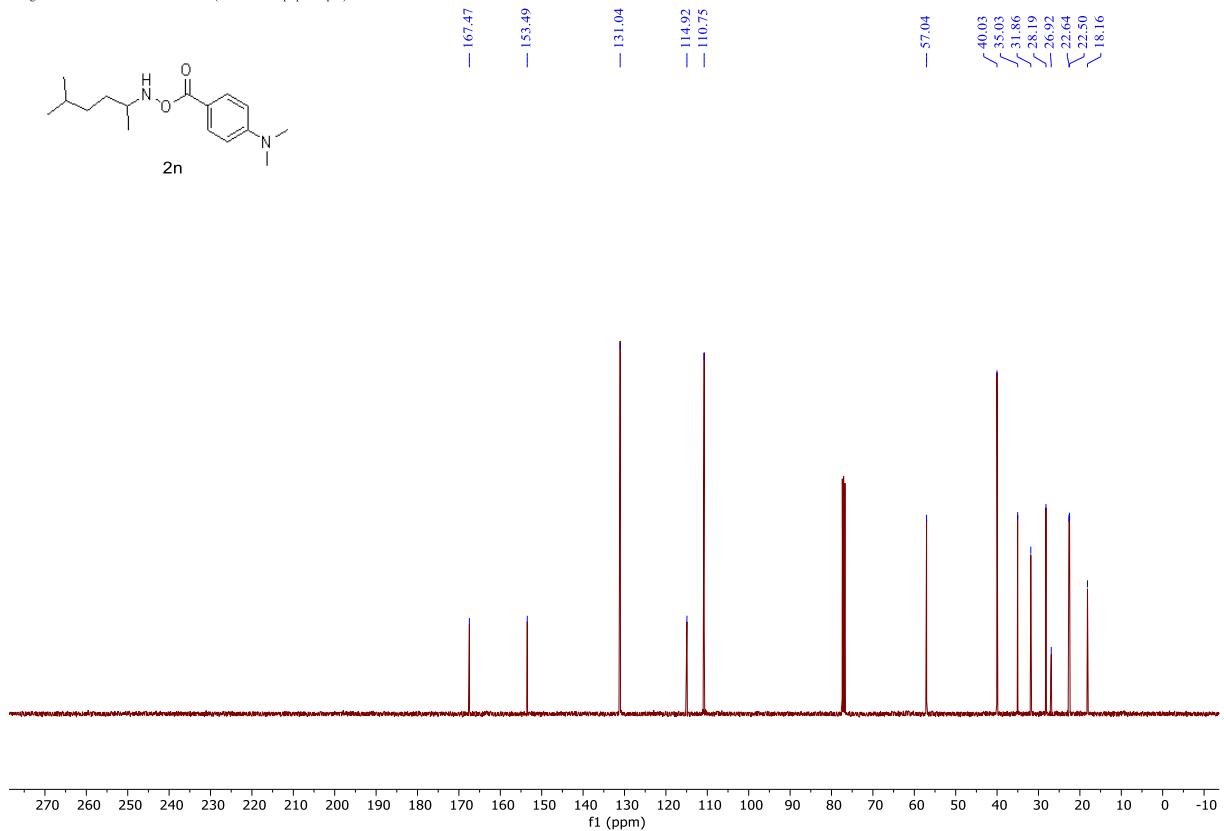
Yuan/YY-A-12 — C13CPD CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 32 — 75.49 MHz



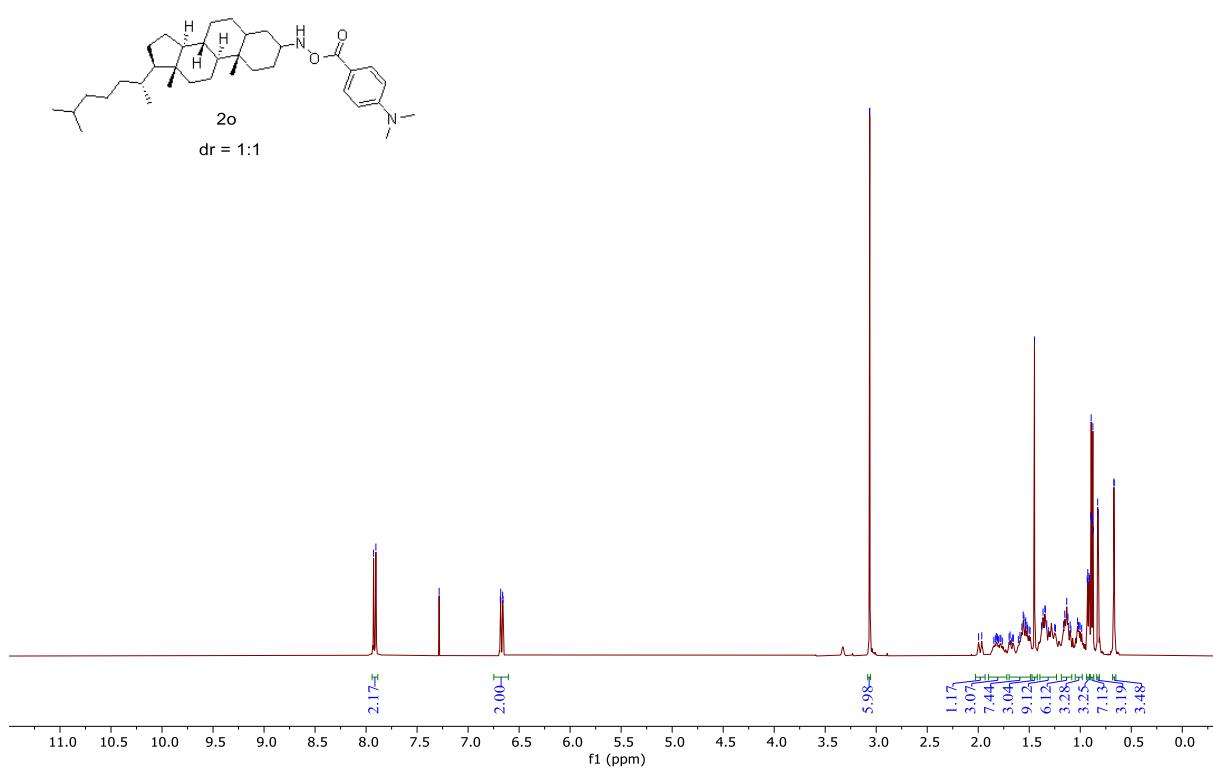
Yang Yuan / YY-A-18 — Au1H CDCl₃ {C:\Bruker\TopSpin3.5\pl6} 2104 42 — 400.13 MHz



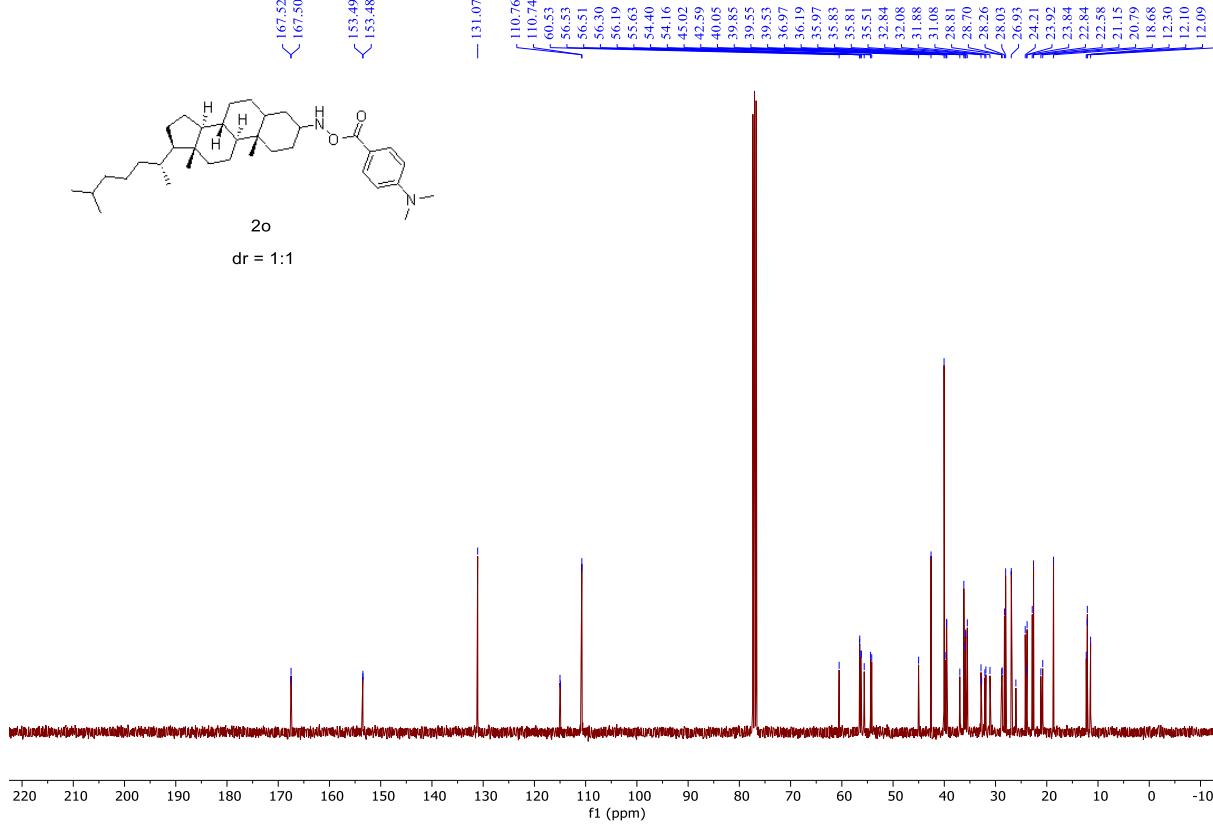
Yang Yuan / YY-A-18 — Au13C CDCl₃ {C:\Bruker\TopSpin3.5\pl6} 2104 42 — 100.63 MHz



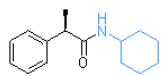
Yang Yuan / YY-A-20 — Au1H CDCl₃ {C:\Bruker\TopSpin3.5\pl6} 2104 45 — 400.13 MHz



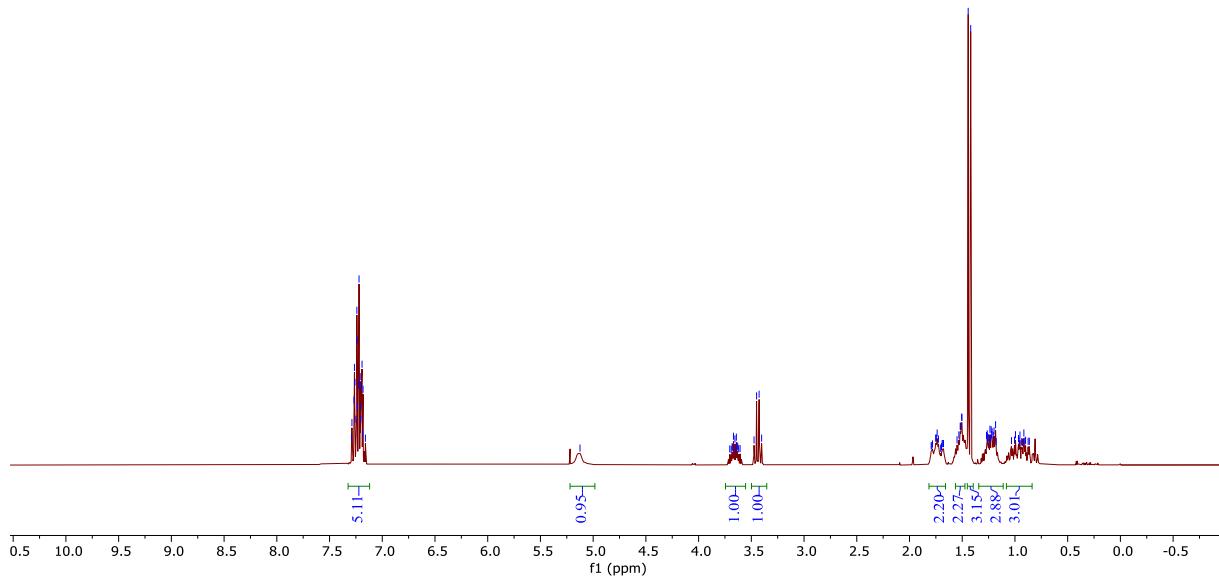
Yang Yuan / YY-A-20 — Au13C CDCl₃ {C:\Bruker\TopSpin3.5\pl6} 2104 45 — 100.63 MHz



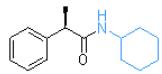
Yang Yuan YY-X-60 — PROTON CDCl₃ {C:\Bruker\TopSpin3.6.2\} 2102 40 — 300.20 MHz



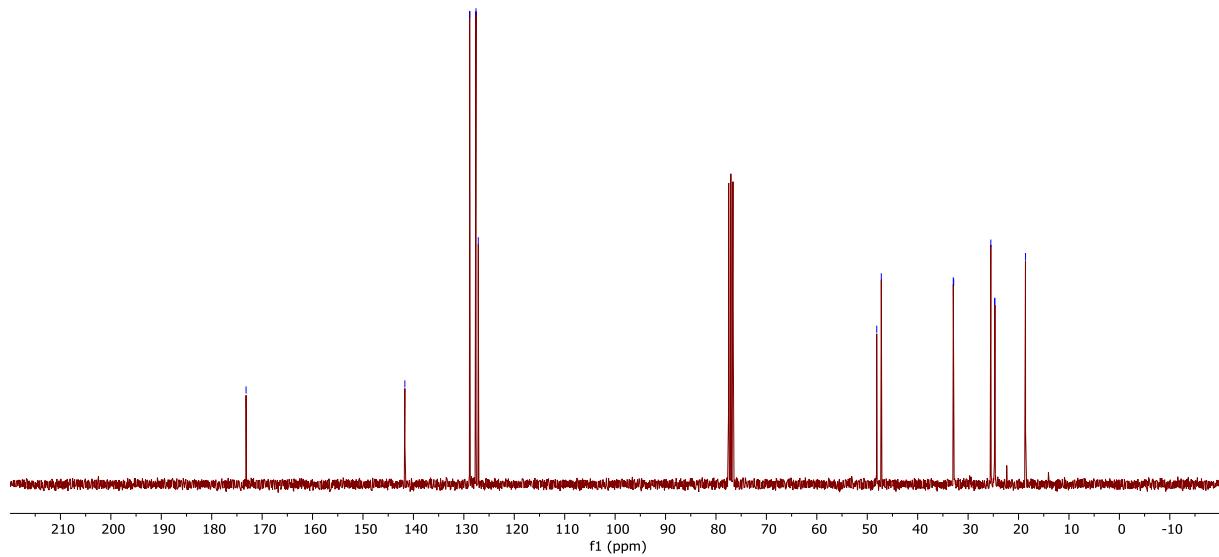
4



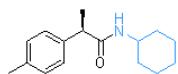
Yang Yuan YY-X-60 — C13CPD CDCl₃ {C:\Bruker\TopSpin3.6.2\} 2102 40 — 75.49 MHz



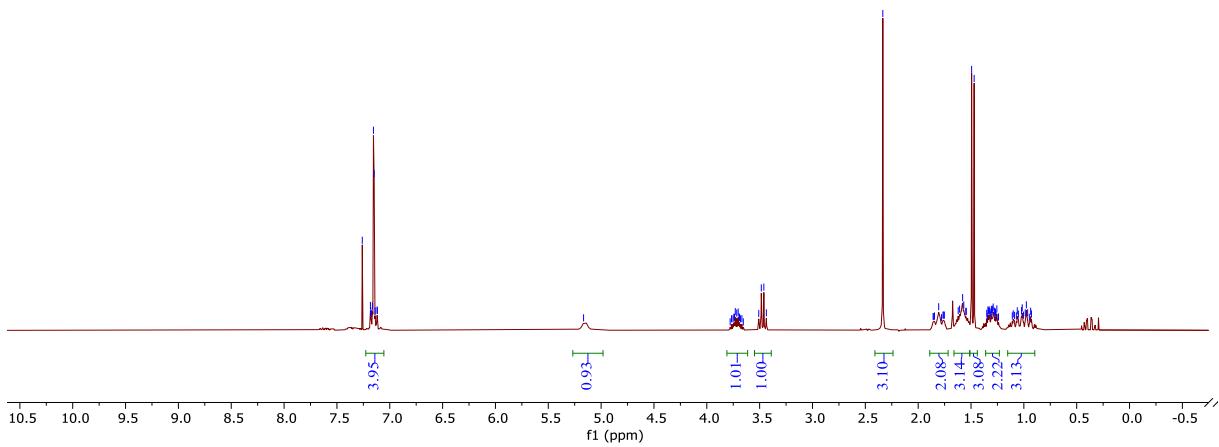
4



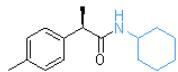
Yang Yuan YY-X-65 — Au1H CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 39 — 300.13 MHz



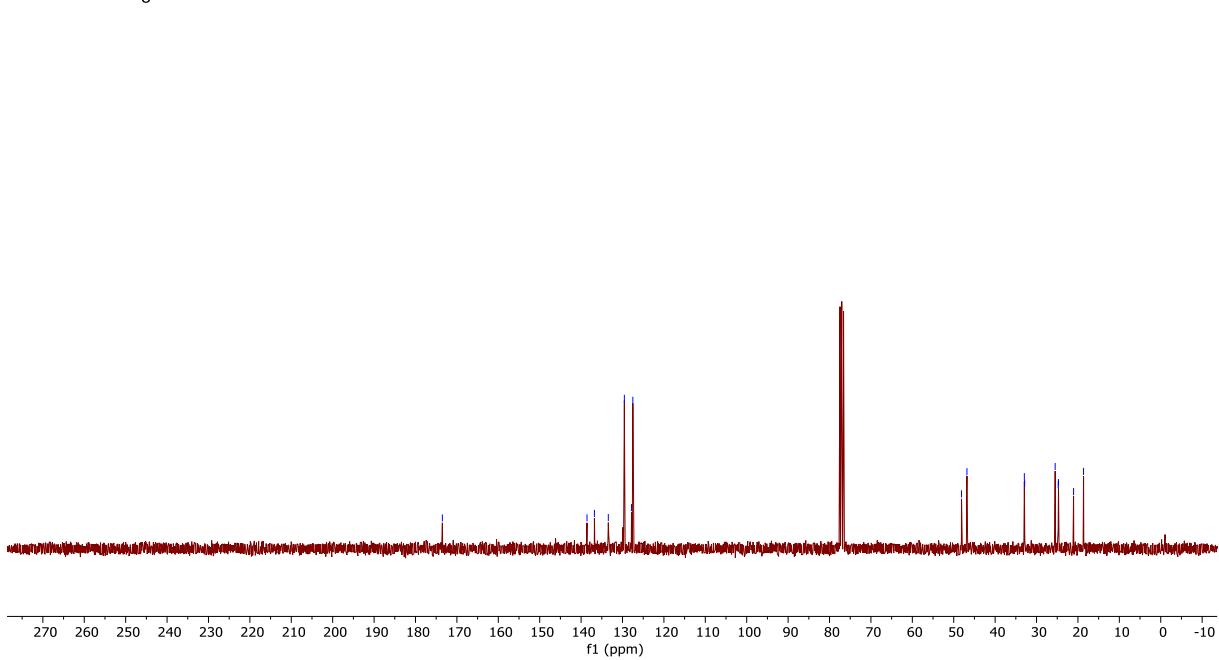
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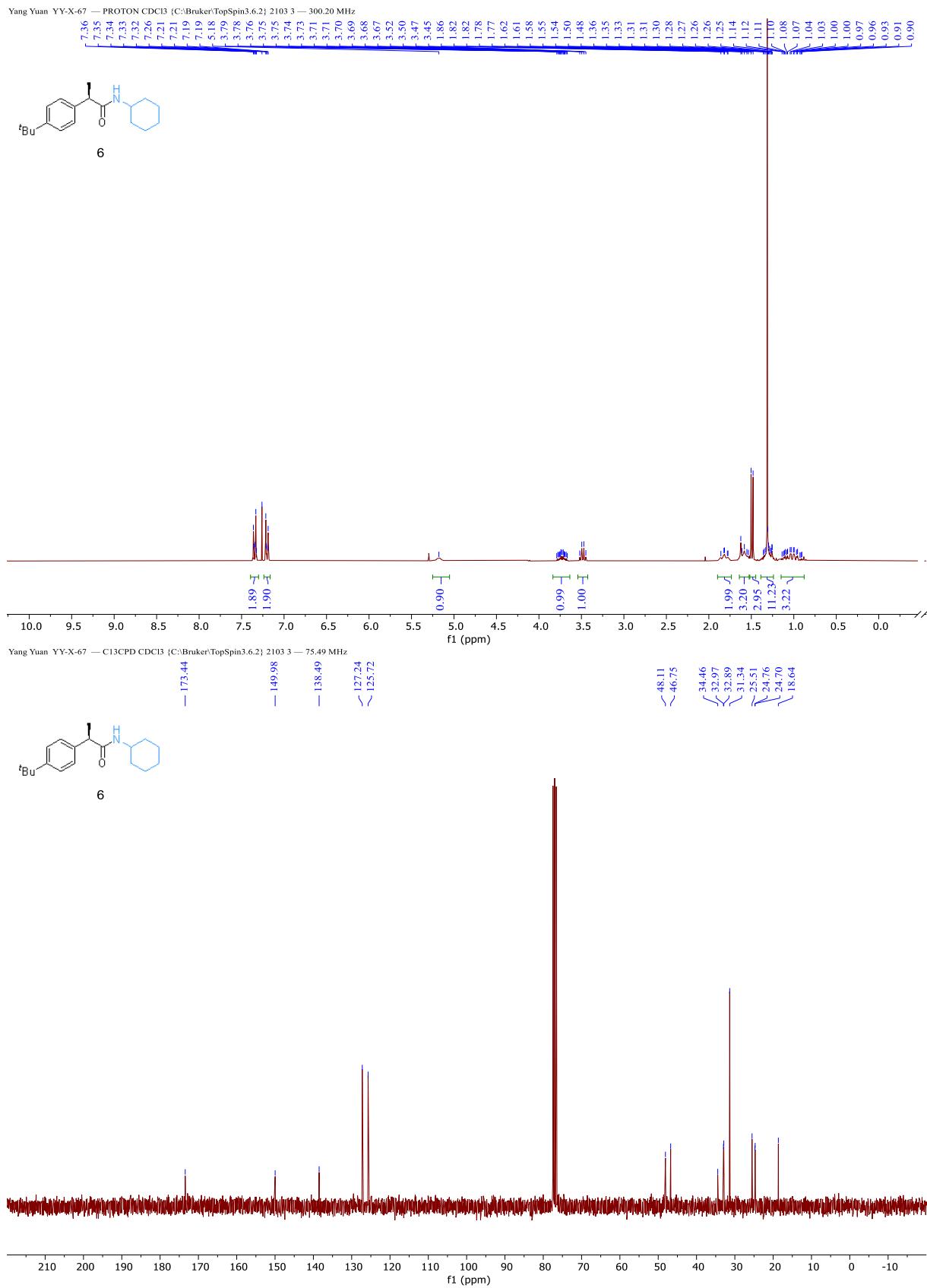


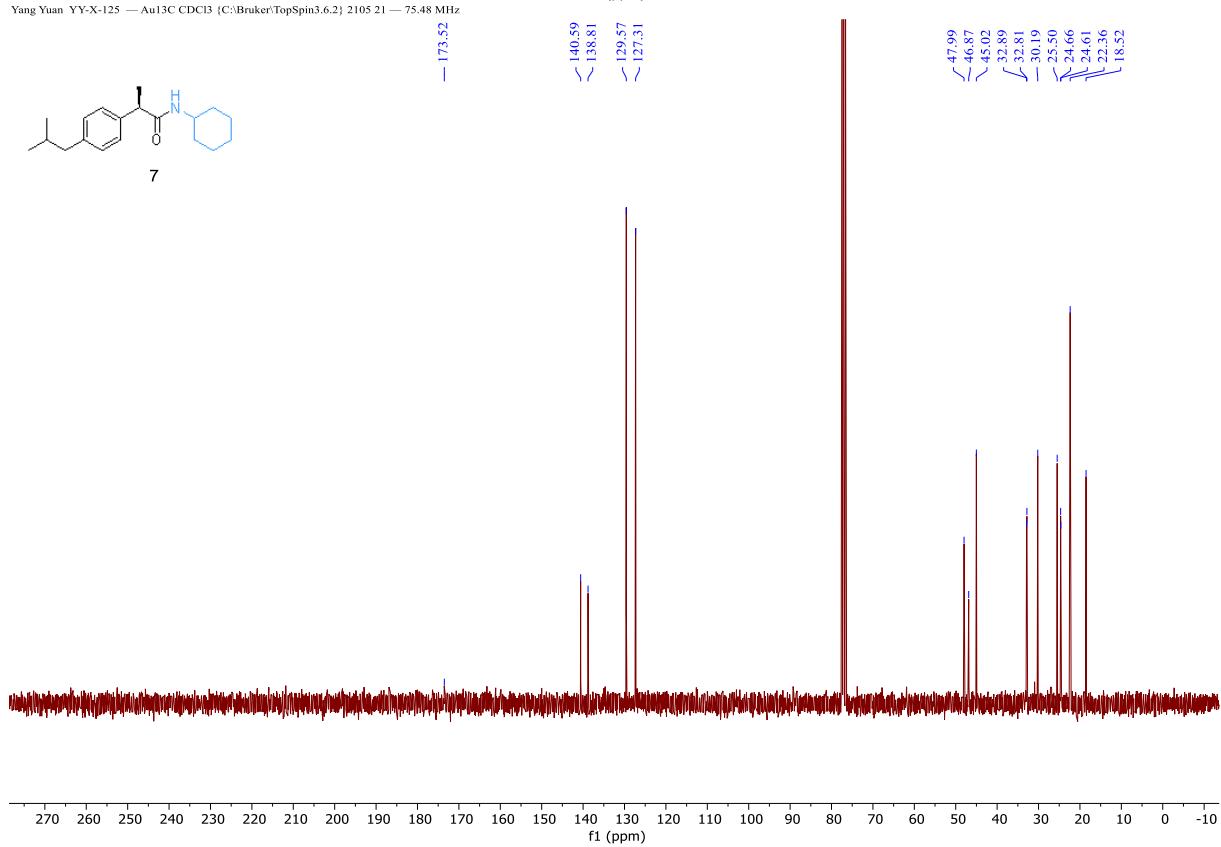
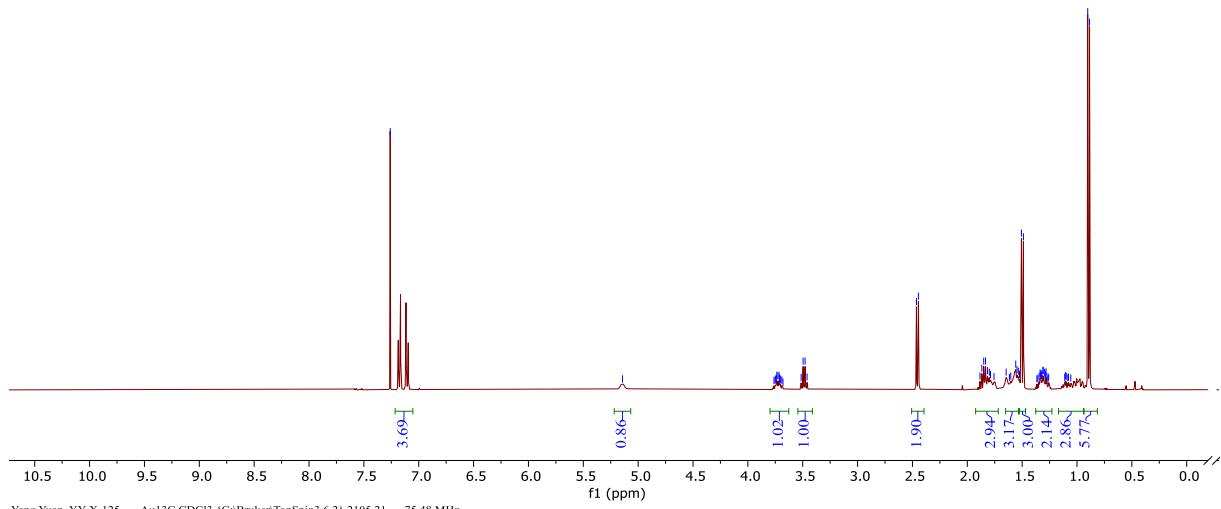
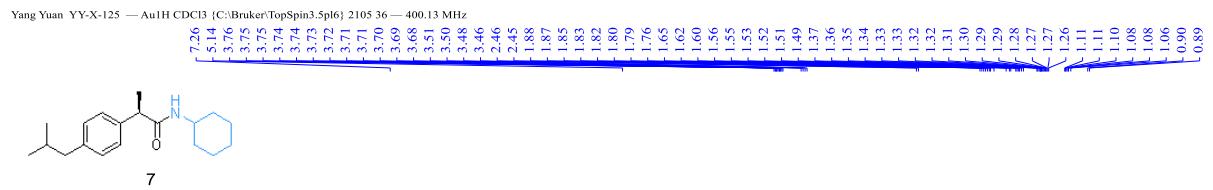
Yang Yuan, YY-X-65 — Au13C CDCl₃ {C:\Bruker\TopSpin3.6.2} 2102 56 — 75.48 MHz

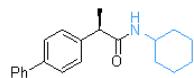
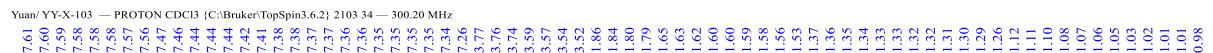


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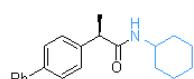
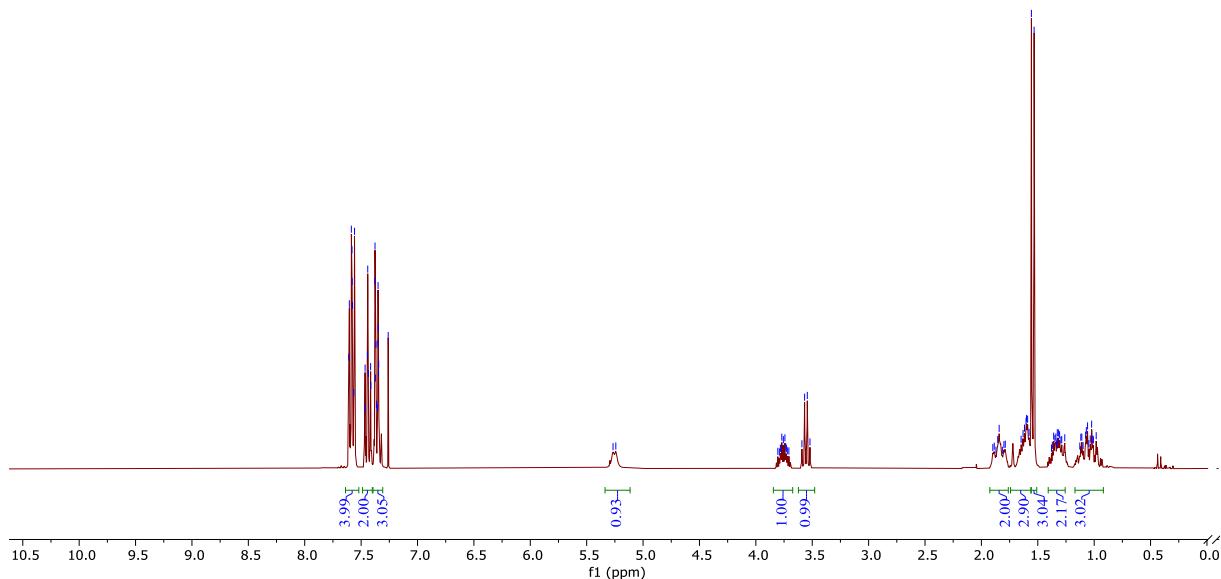




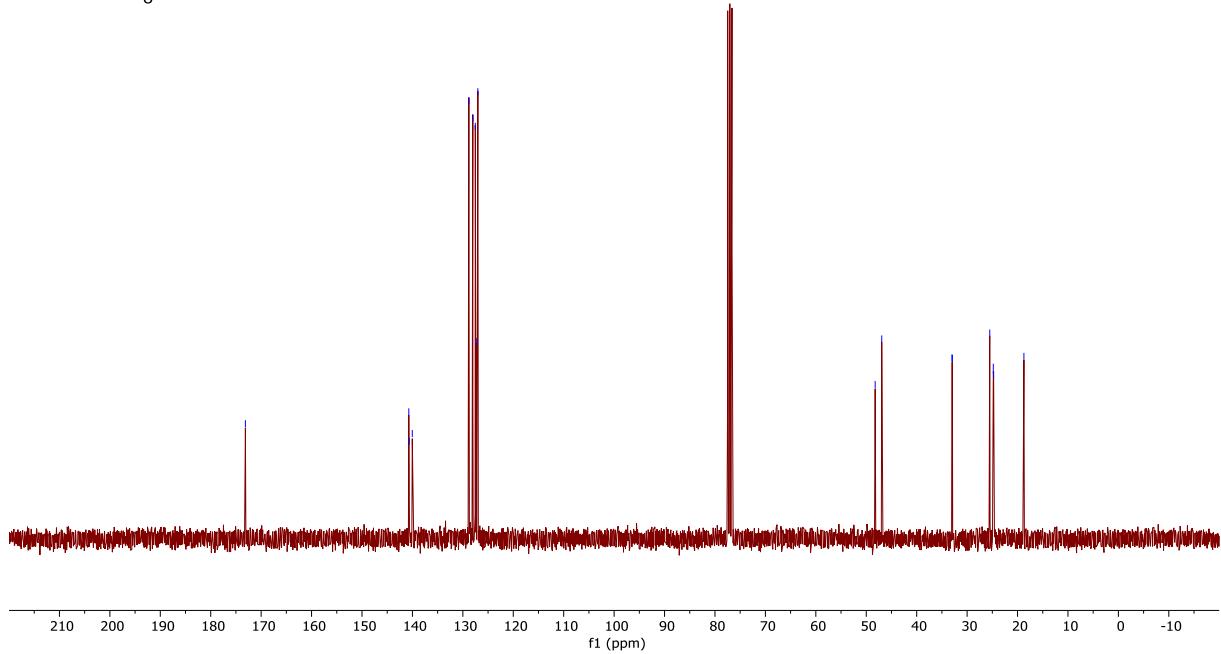


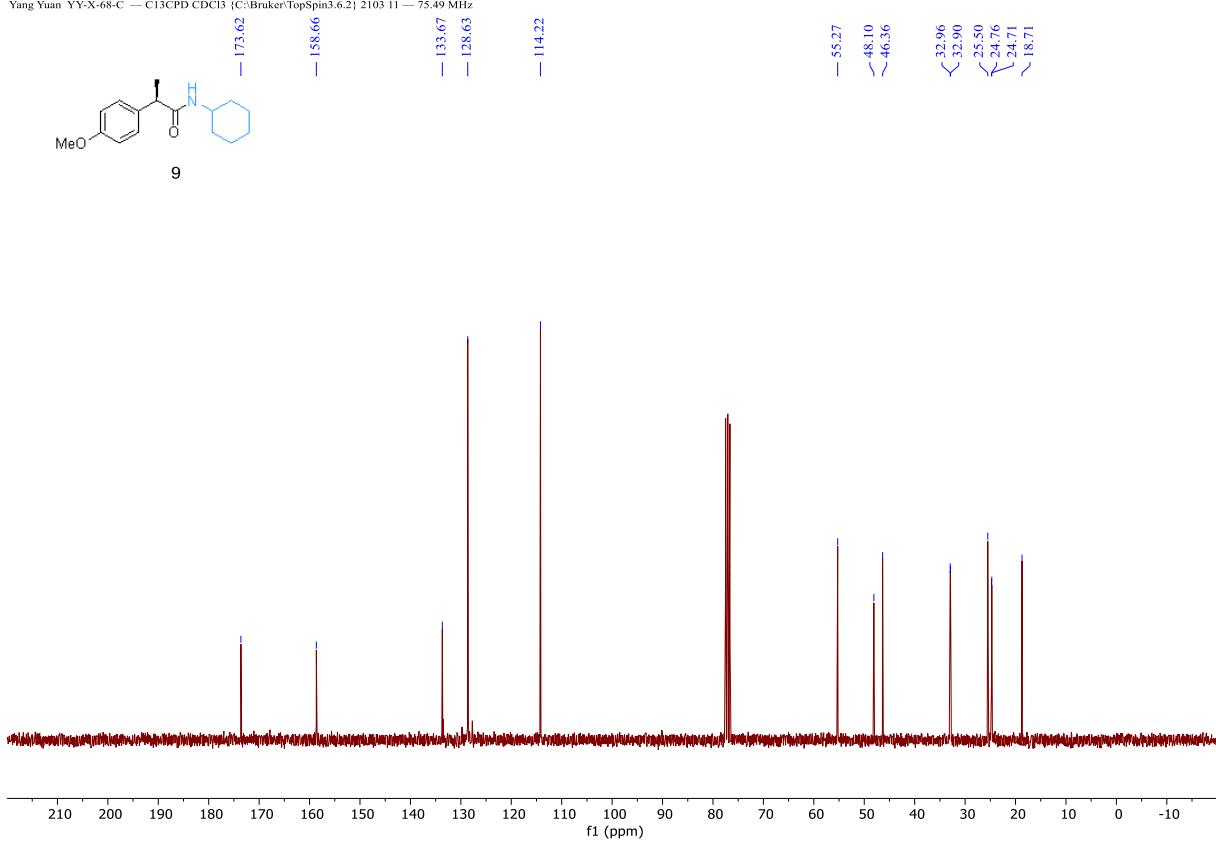
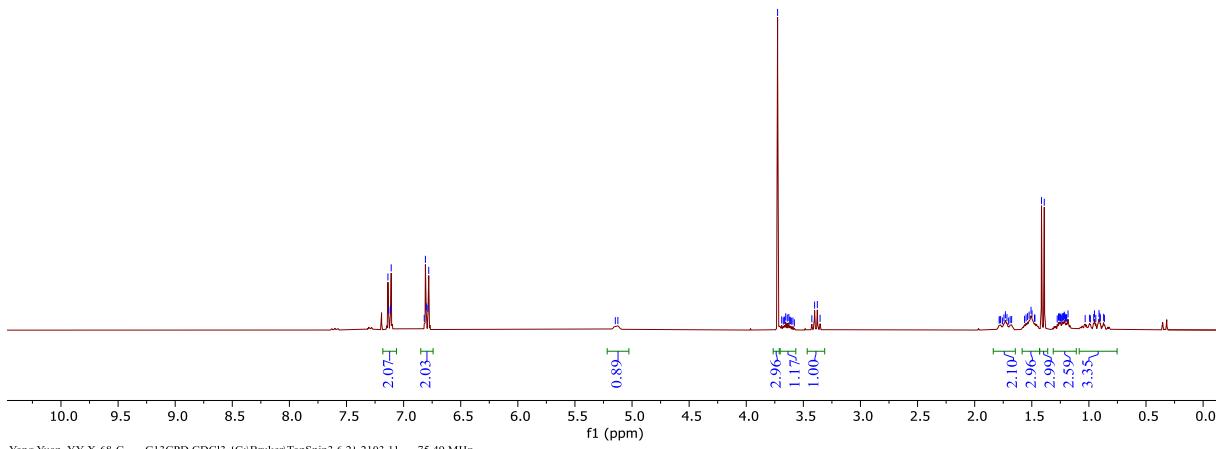
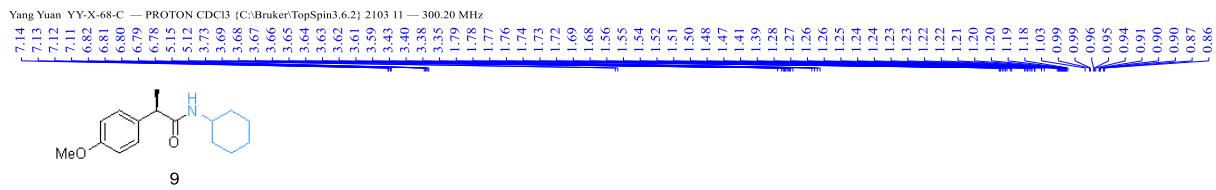


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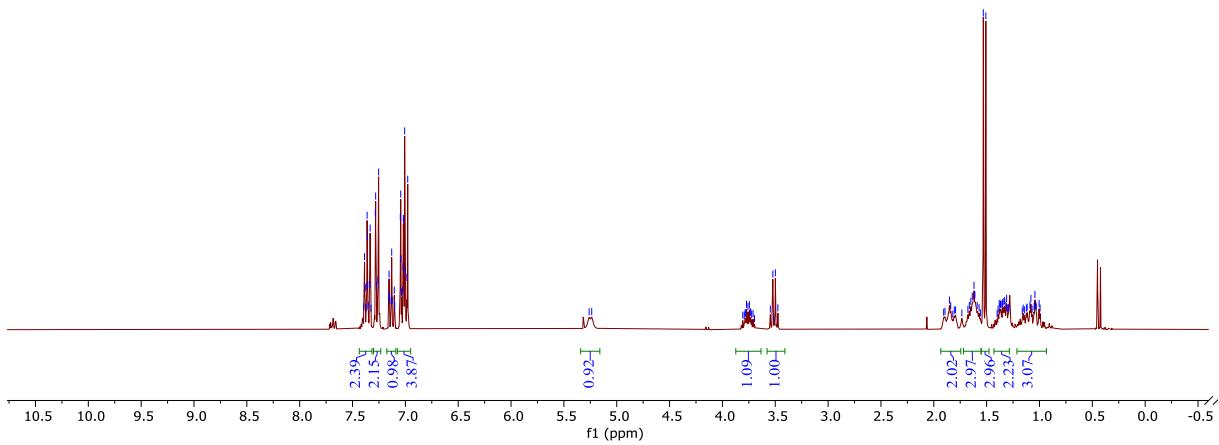
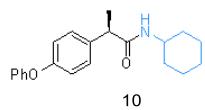


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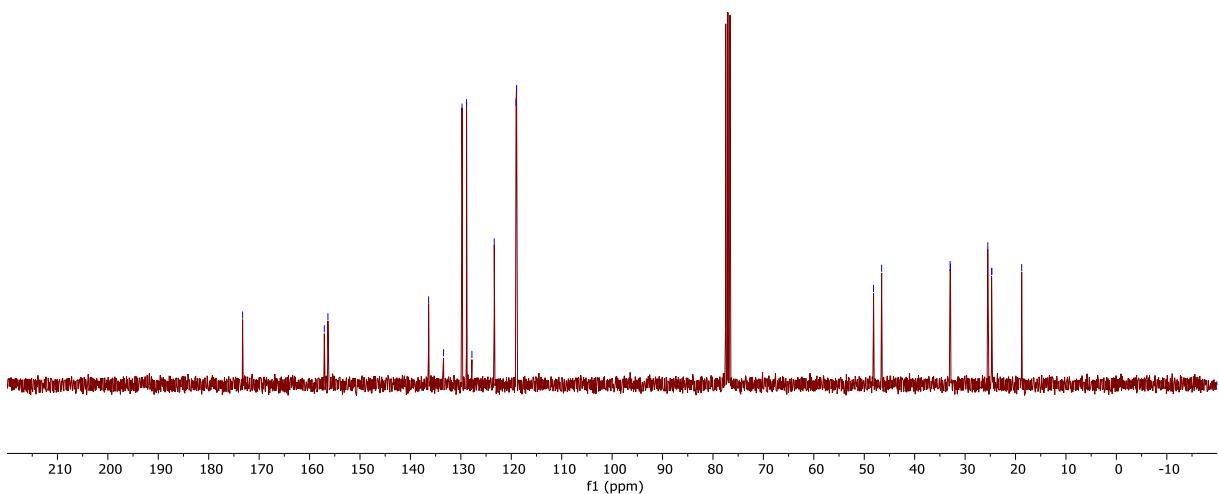
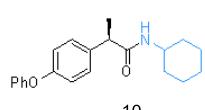


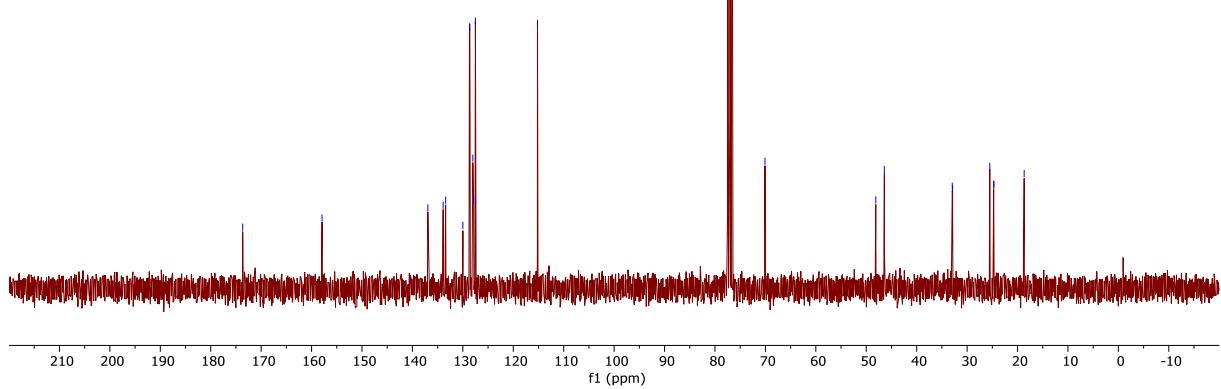
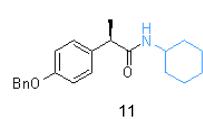
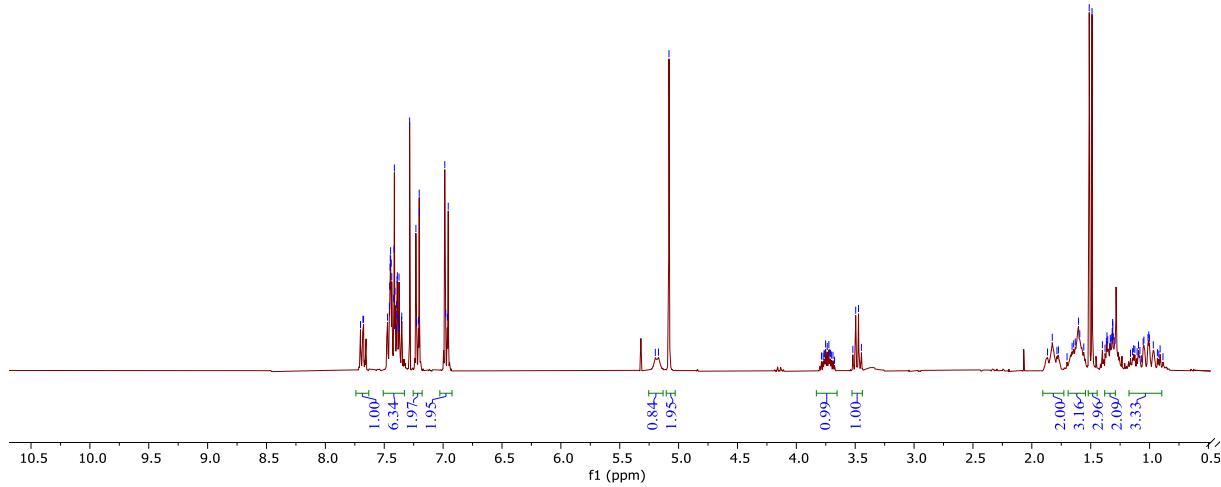
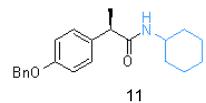


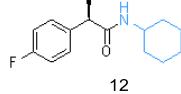
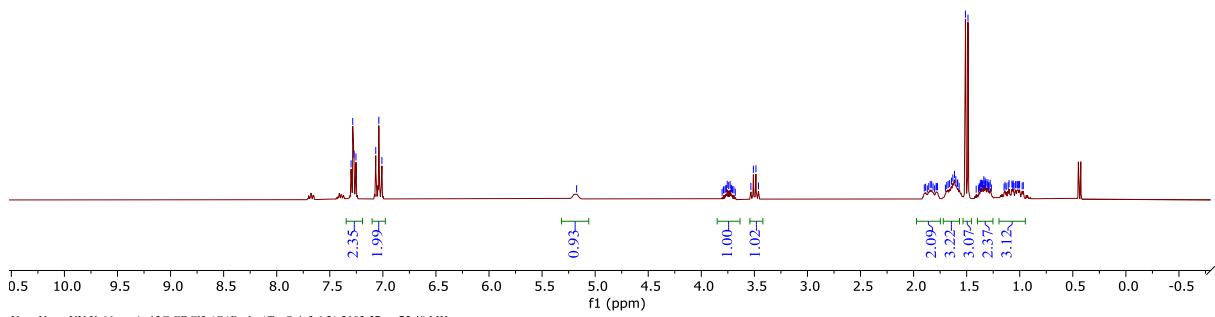
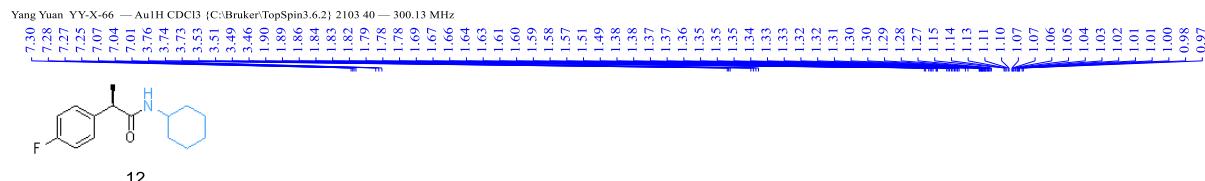
Yang Yuan YY-X-72 — PROTON CDCl₃ {C:\Bruker\TopSpin3.6.2\} 2103 6 — 300.20 MHz



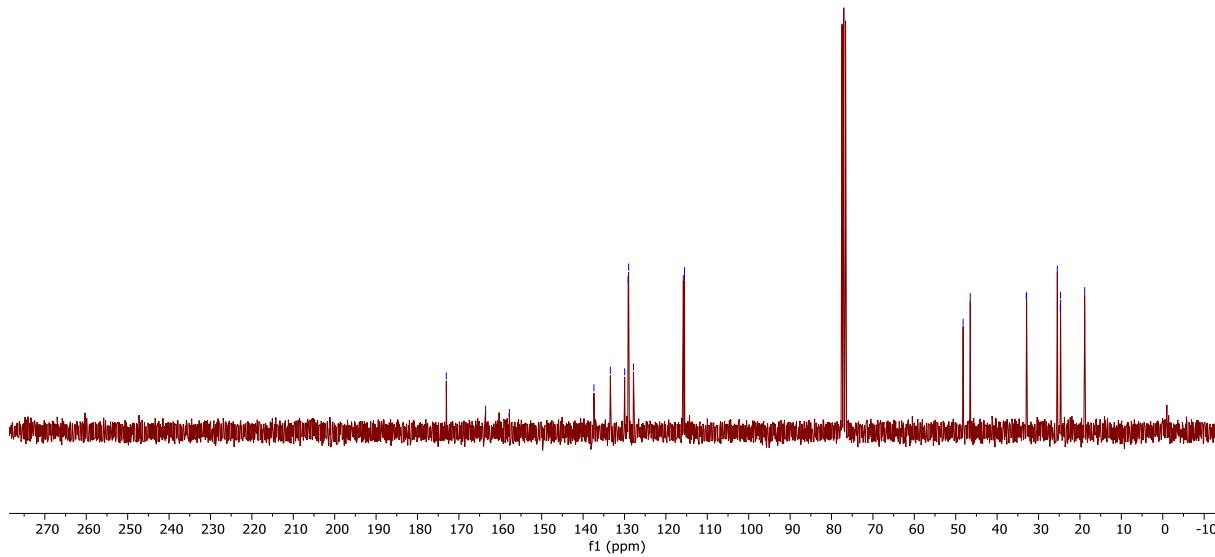
Yang Yuan YY-X-72 — C13CPD CDCl₃ {C:\Bruker\TopSpin3.6.2\} 2103 6 — 75.49 MHz

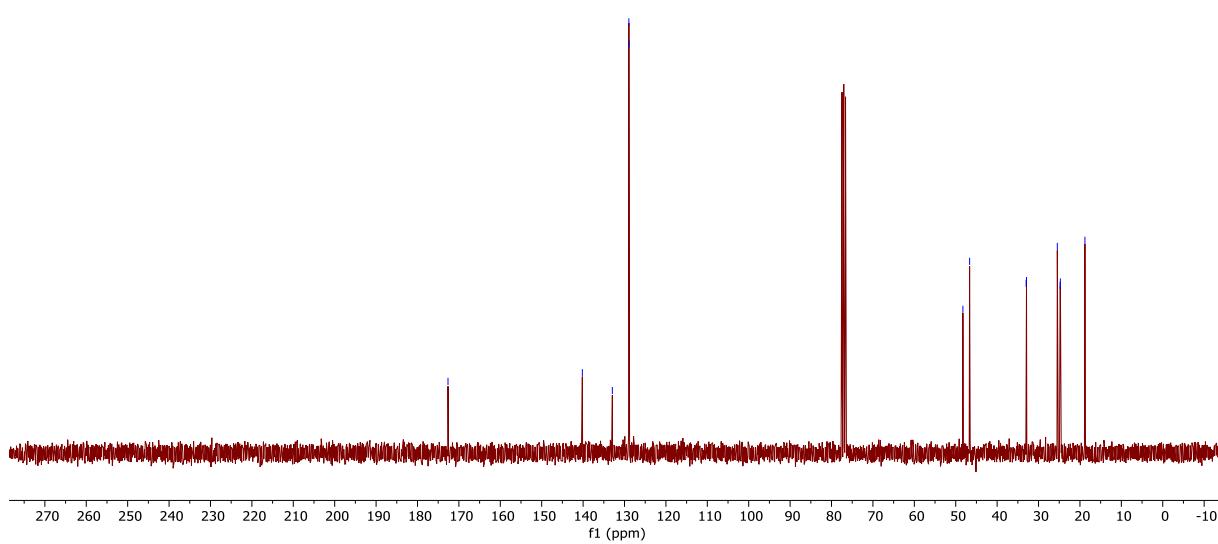
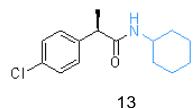
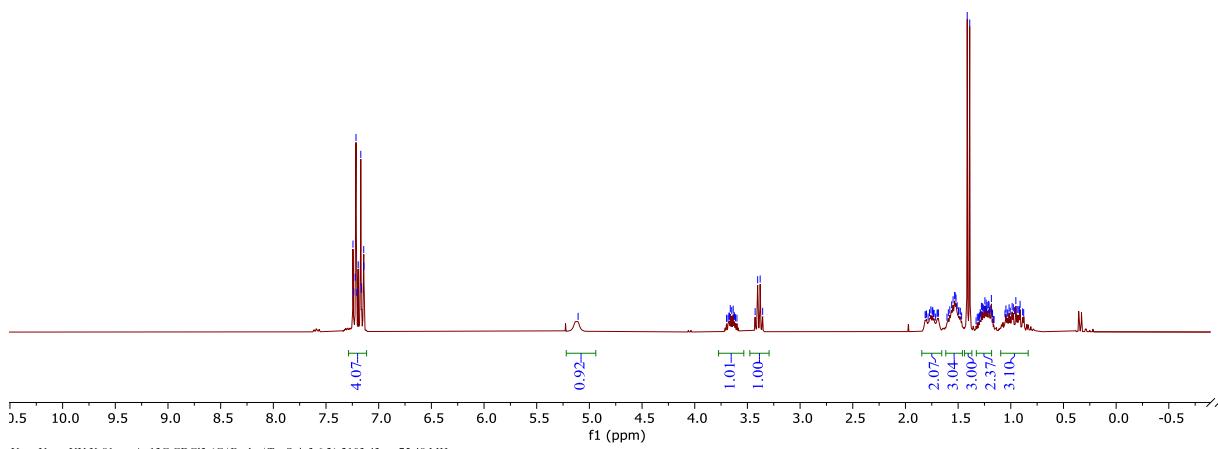
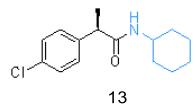




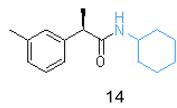


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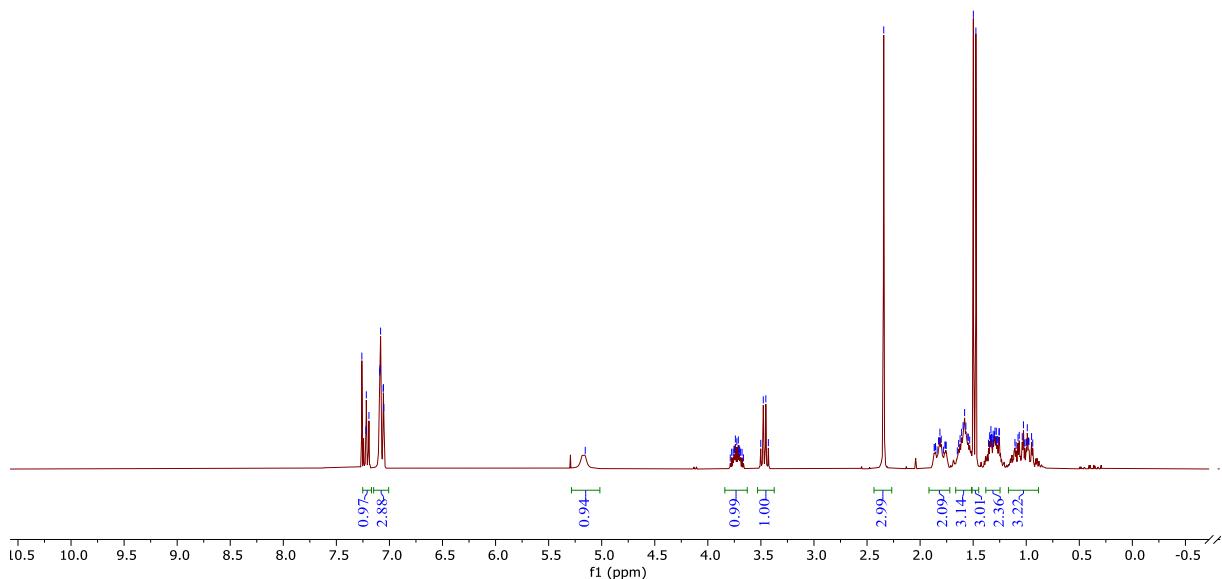




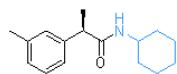
Yang Yuan, YY-X-79 — Au1H CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 42 — 300.13 MHz



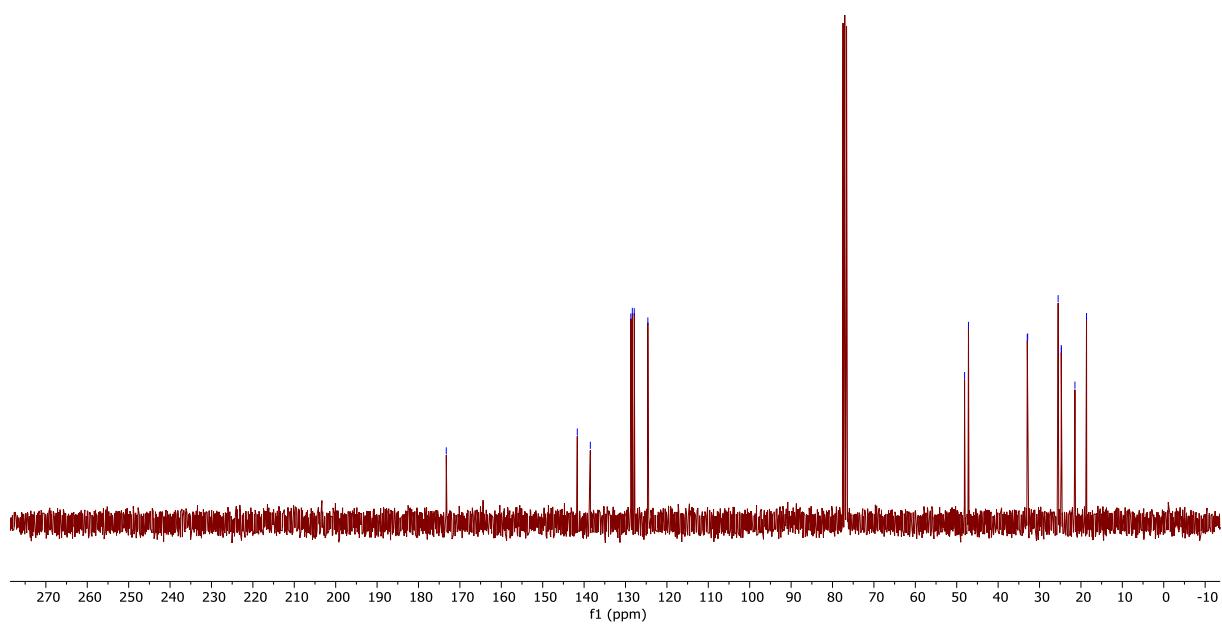
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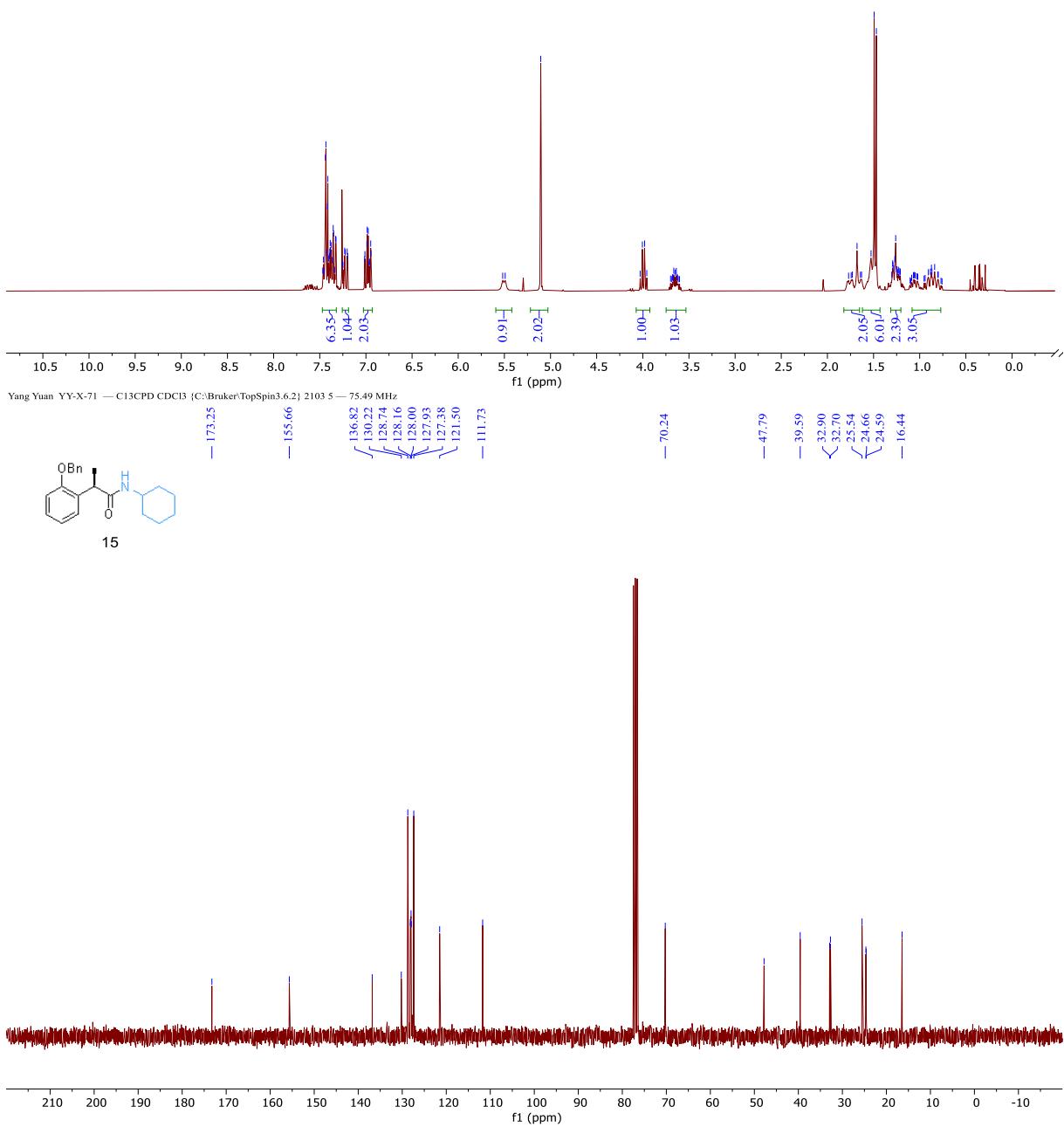
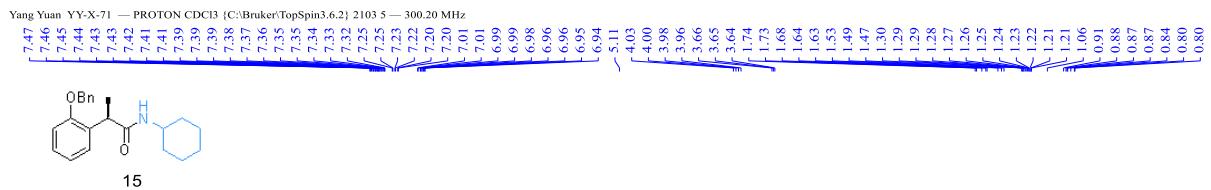


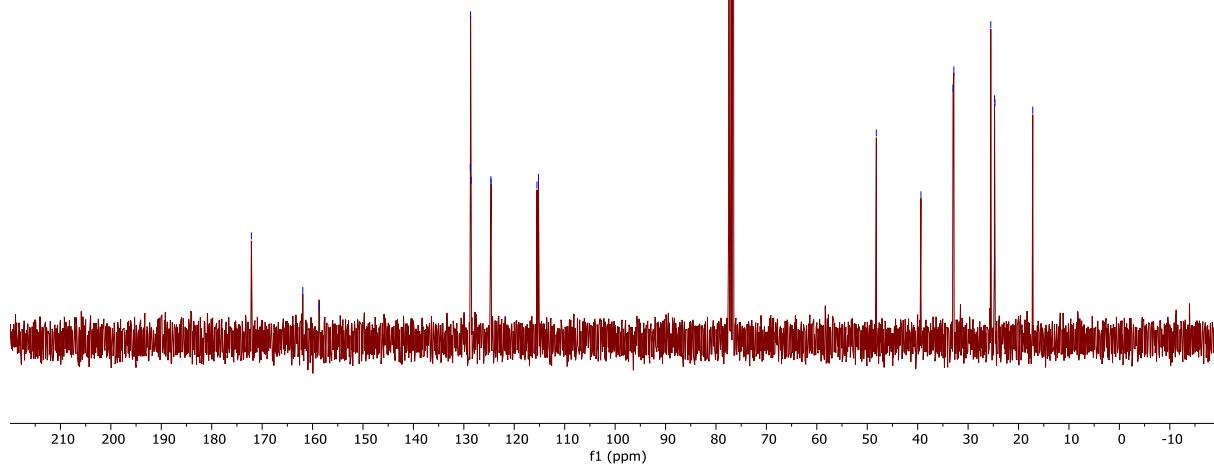
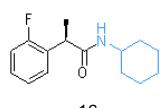
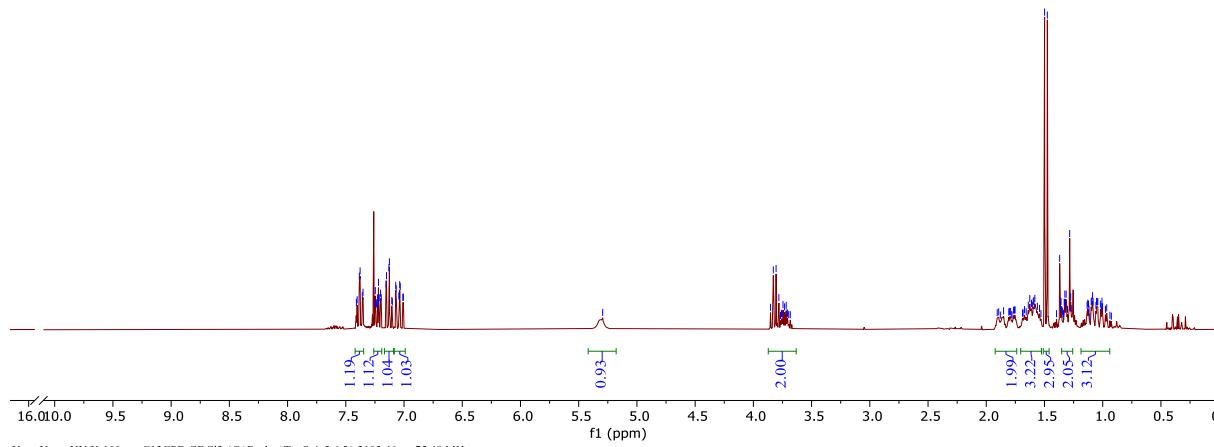
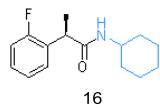
Yang Yuan, YY-X-79 — Au13C CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 42 — 75.48 MHz



14

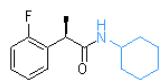




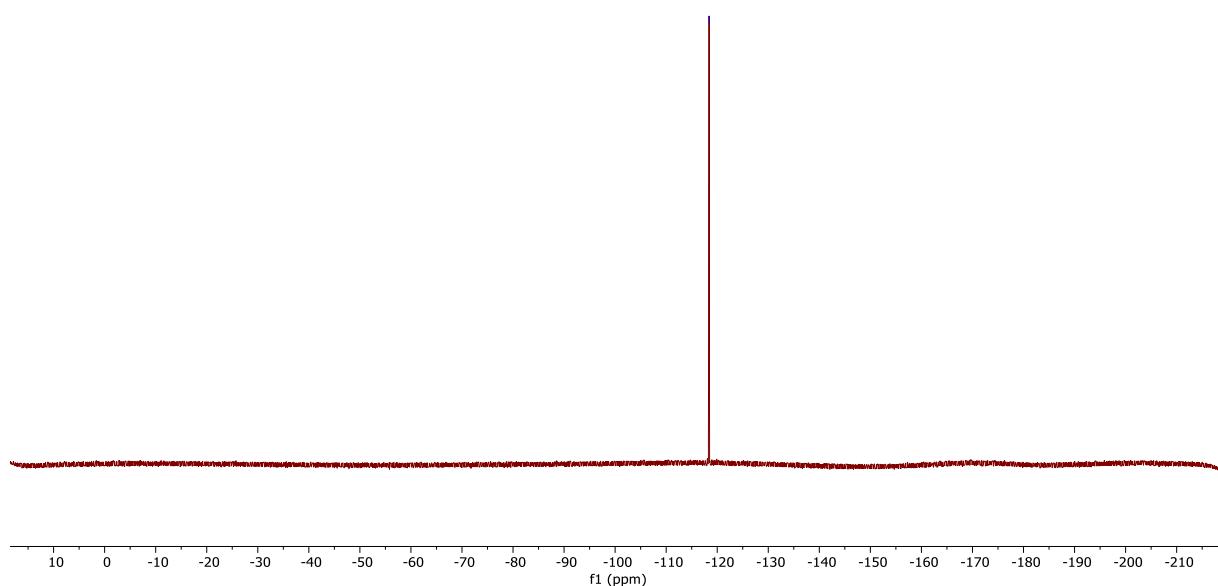


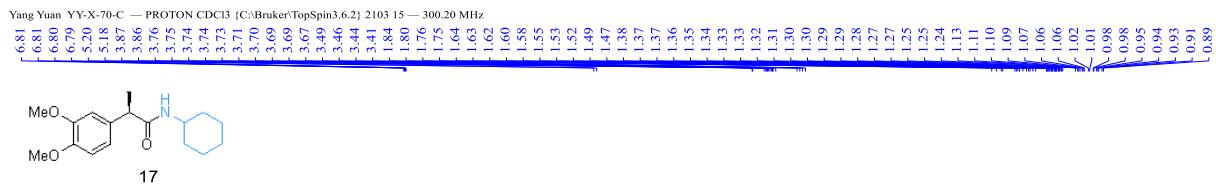
Yang Yuan YY-X-100 — F19 CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103.46 — 282.44 MHz

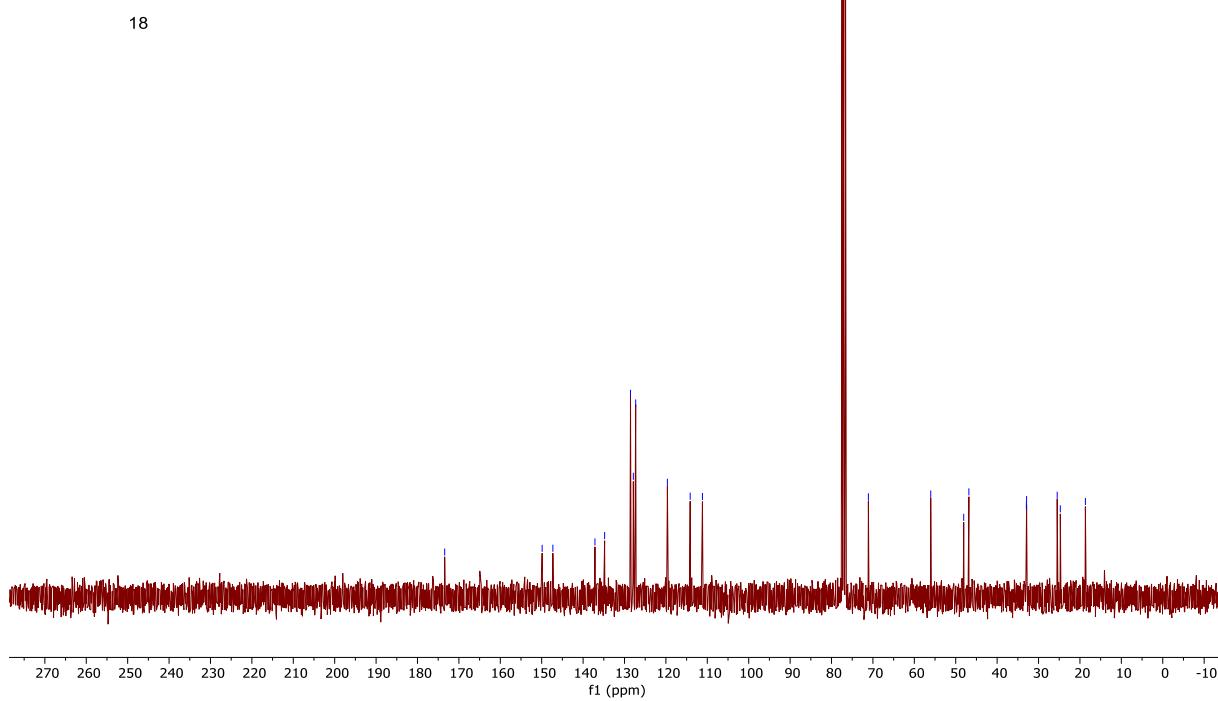
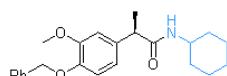
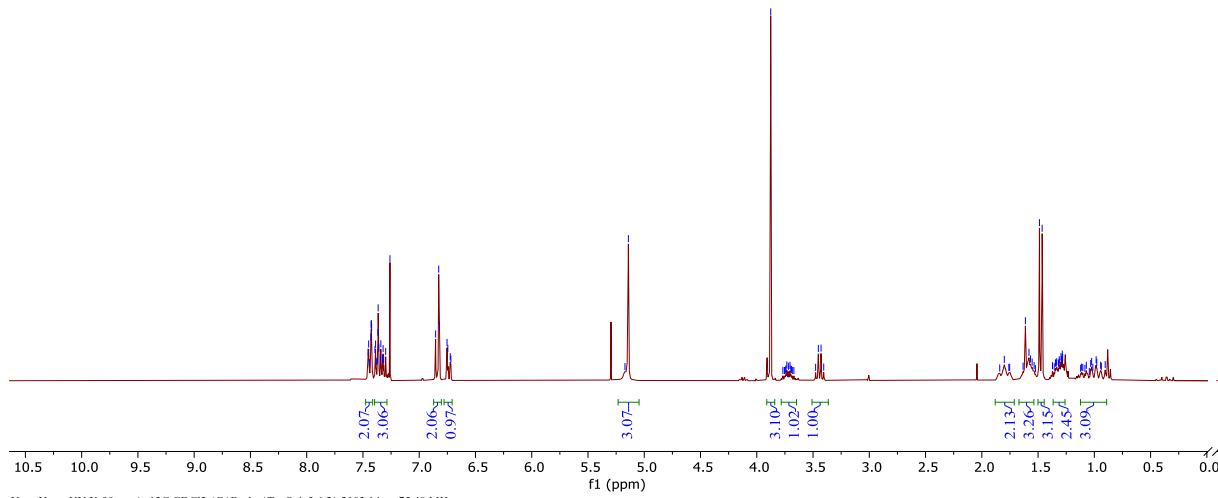
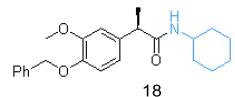
— 118.42

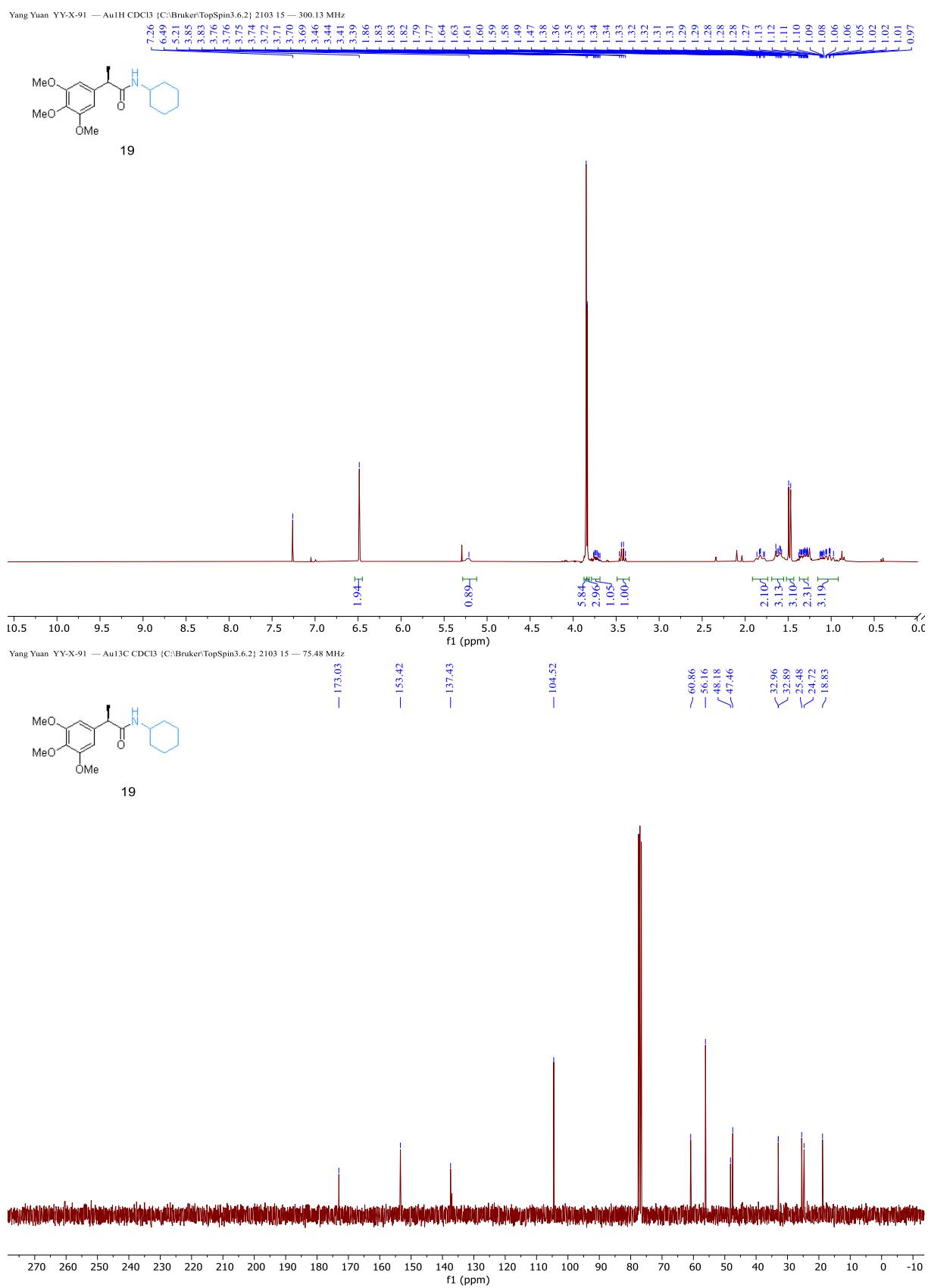


16

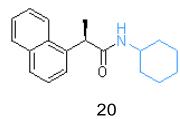




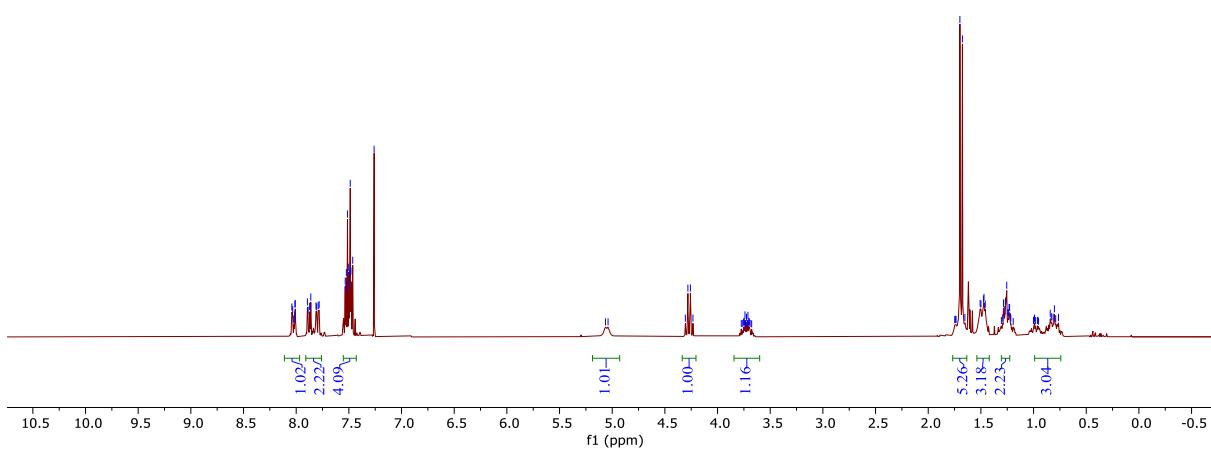




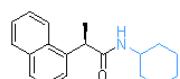
Yang Yuan YY-X-76 — Au1H CDCl₃ {C:\Bruker\TopSpin3.6.2\} 2103.41 — 300.13 MHz



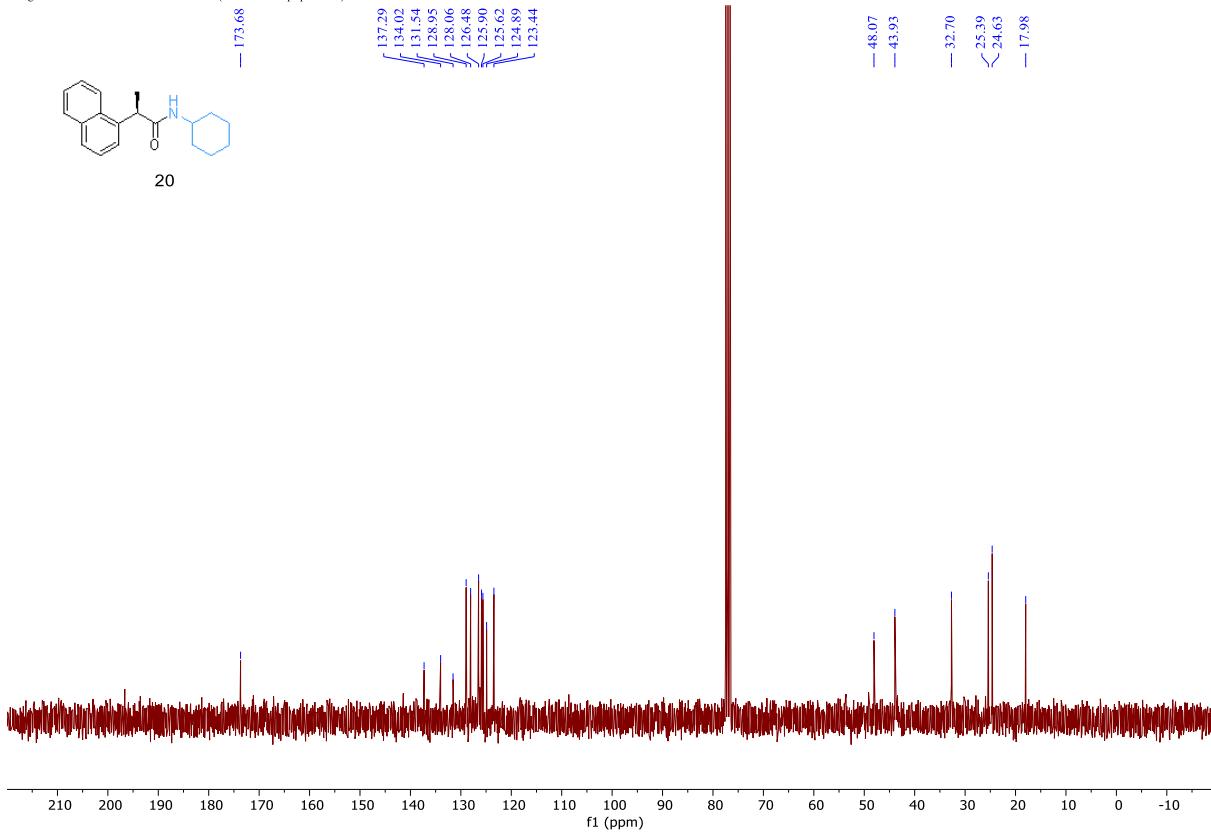
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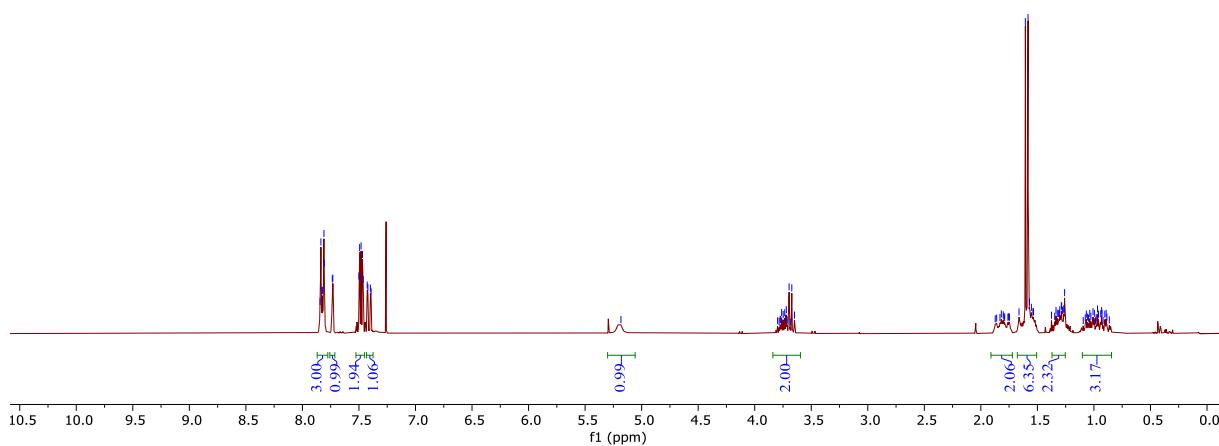
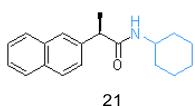
Yang Yuan YY-X-76 — C13CPD CDCl₃ {C:\Bruker\TopSpin3.6.2\} 2103.10 — 75.49 MHz



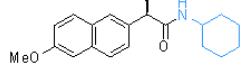
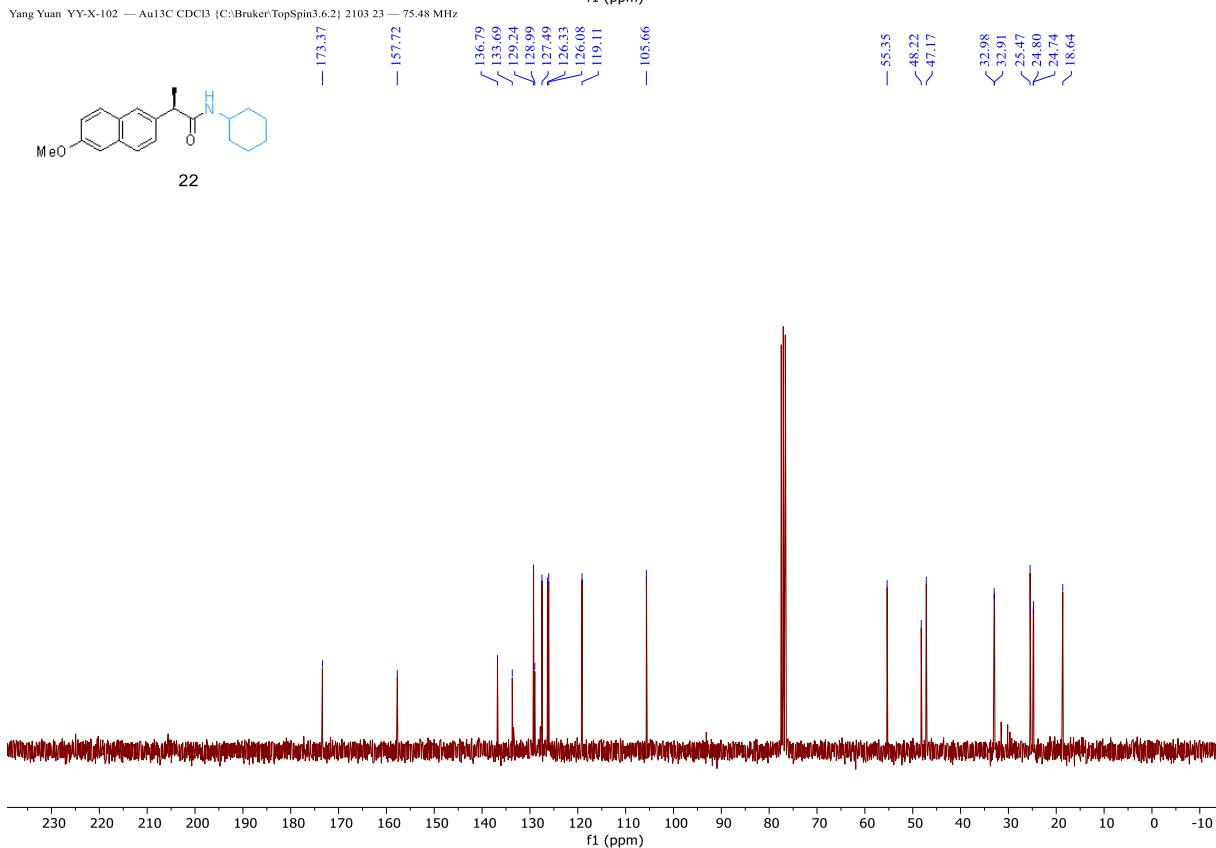
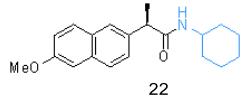
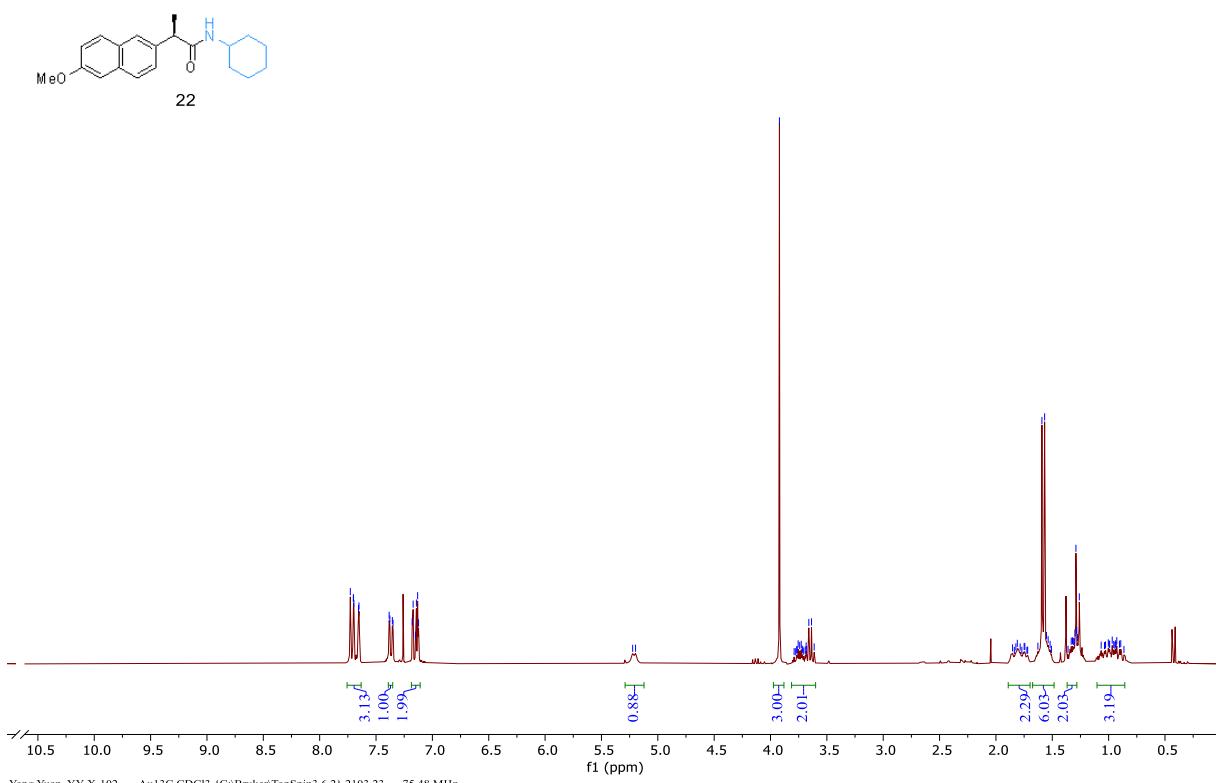
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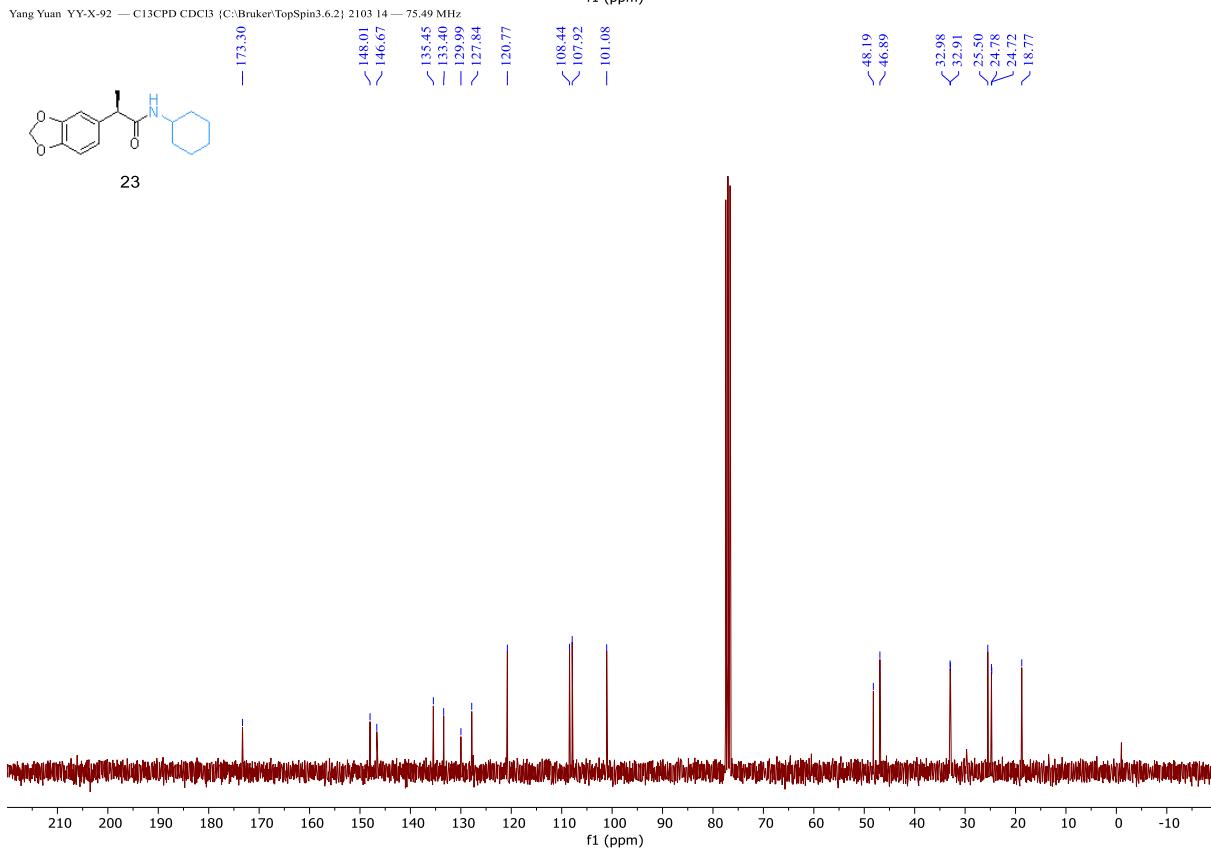
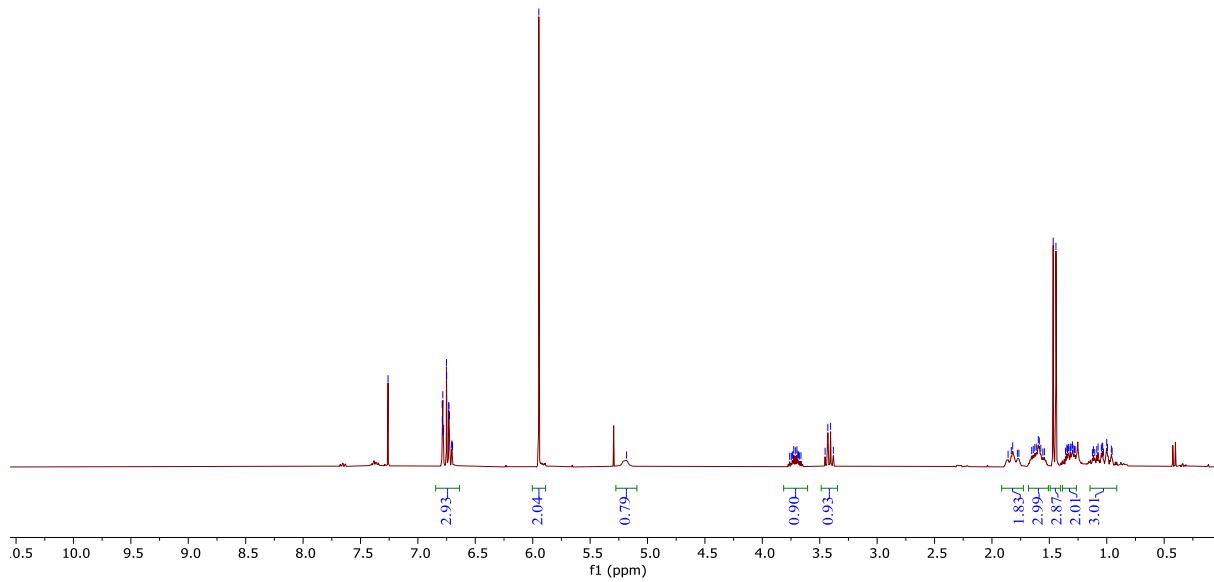
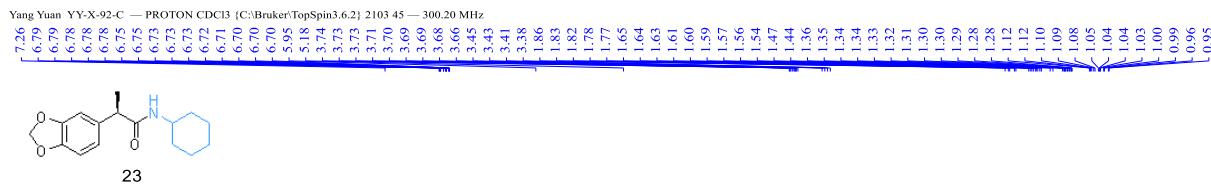


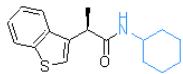
Yang Yuan YY-X-69 — PROTON CDCl₃ {C:\Bruker\TopSpin3.6.2\} 2103 4 — 300.20 MHz



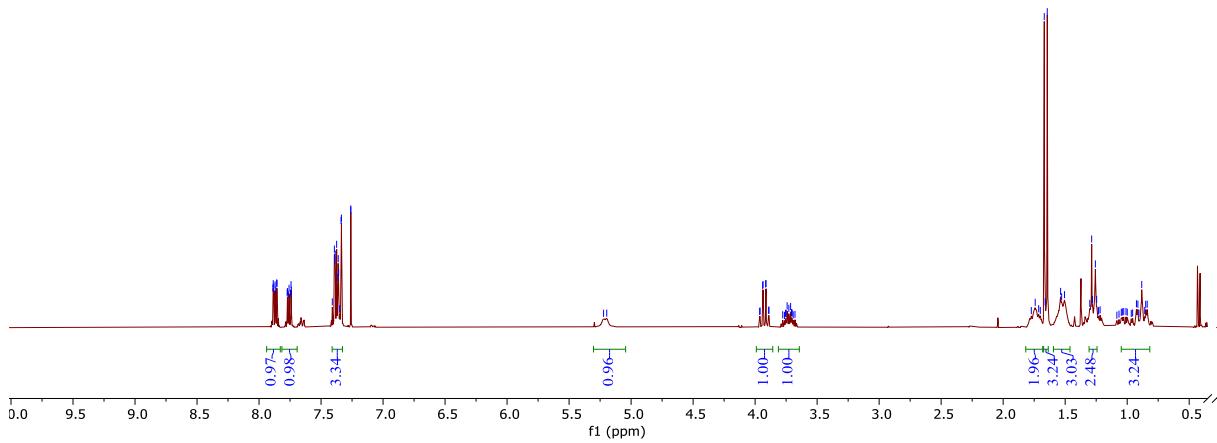
Yang Yuan YY-X-102 — Au1H CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 23 — 300.13 MHz



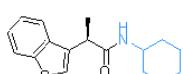




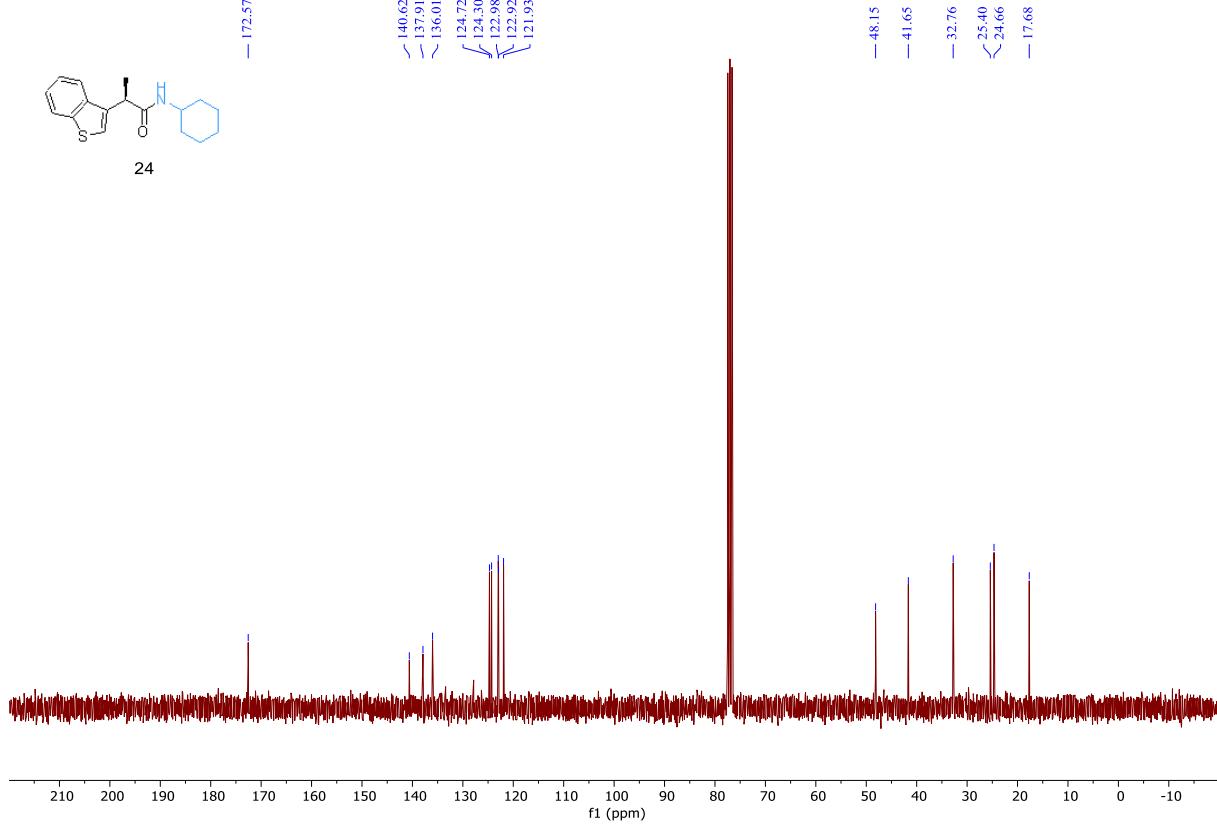
24



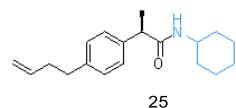
Yang Yuan YY-X-97 — C13CPD CDCI3 {C:\Bruker\TopSpin3.6.2} 2103 43 — 75.49 MHz



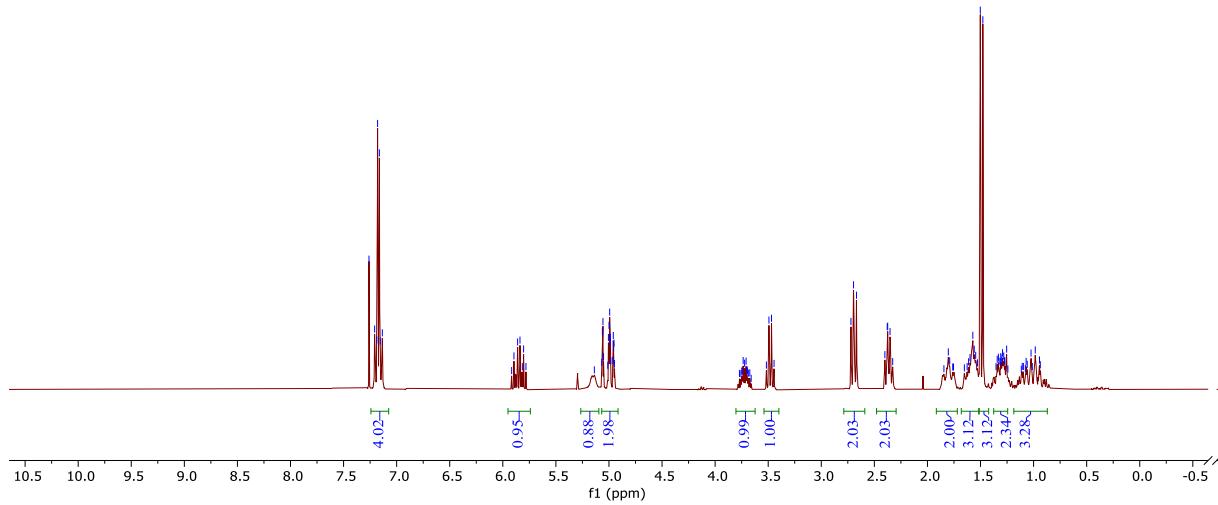
34



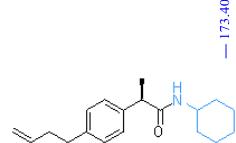
Yang Yuan YY-X-74 — PROTON CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 8 — 300.20 MHz



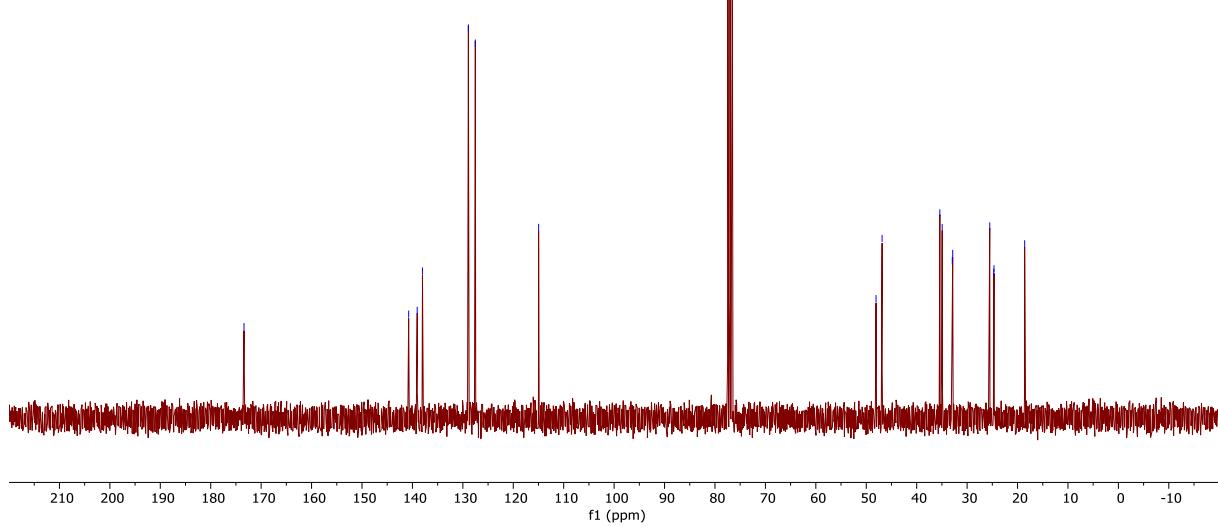
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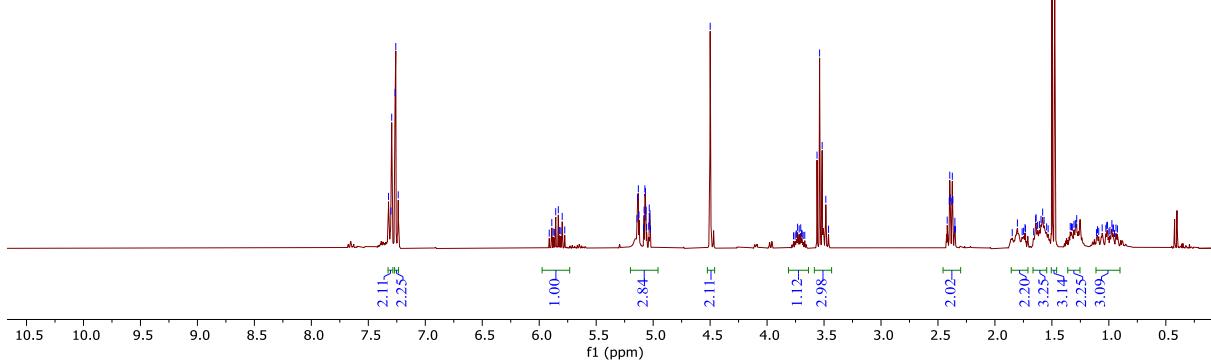
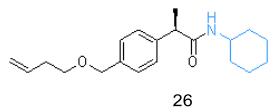
Yang Yuan YY-X-74 — C13CPD CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 8 — 75.49 MHz



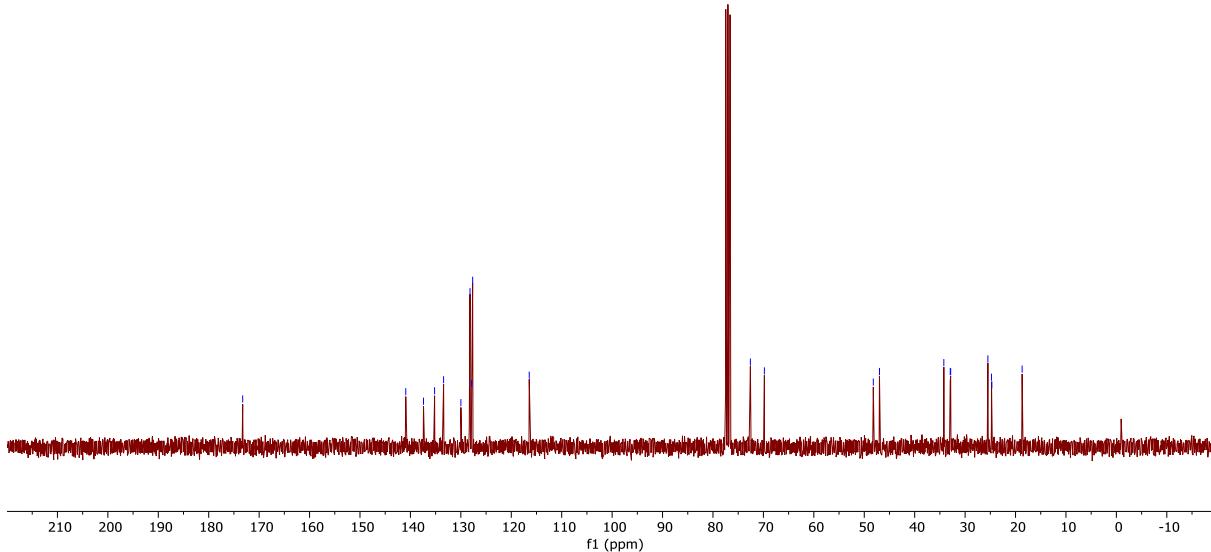
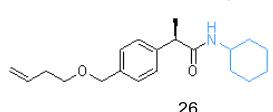
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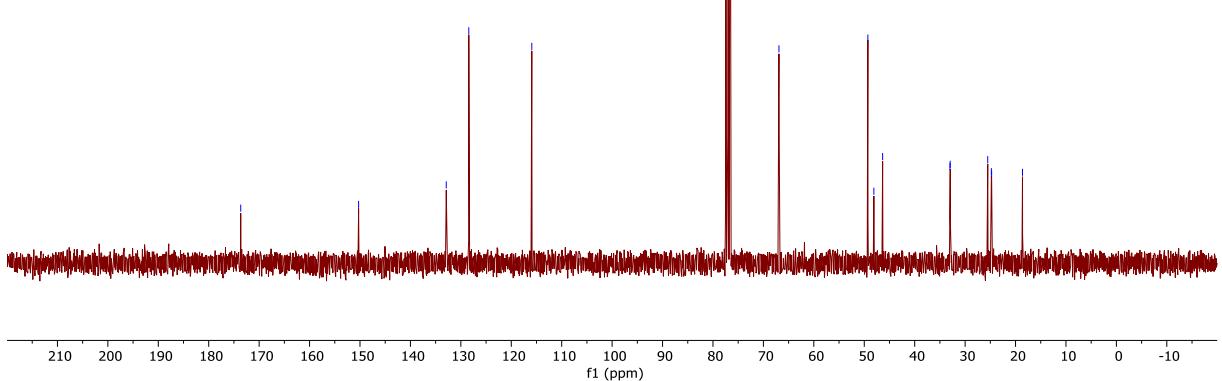
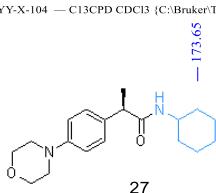
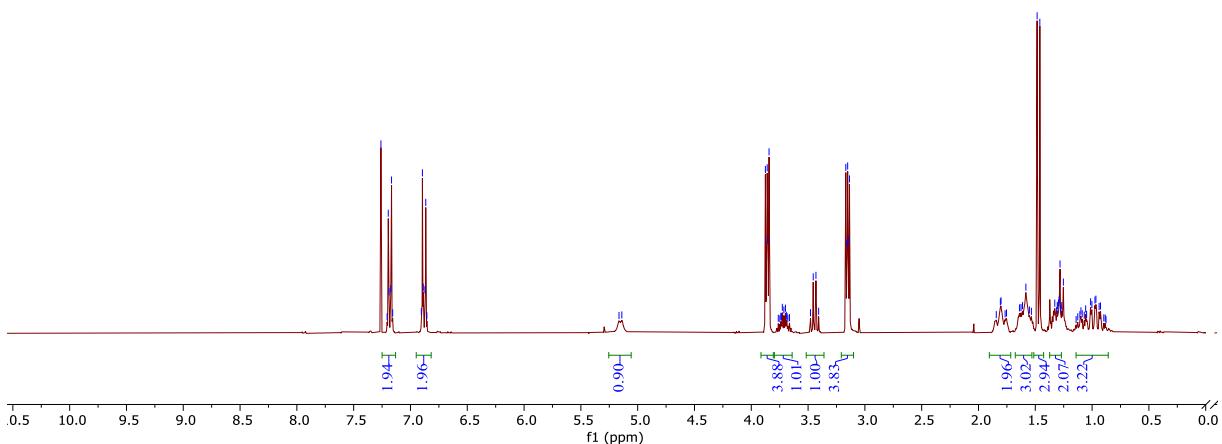
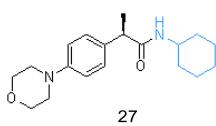
Yang Yuan YY-X-88-C — PROTON CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 44 — 300.20 MHz



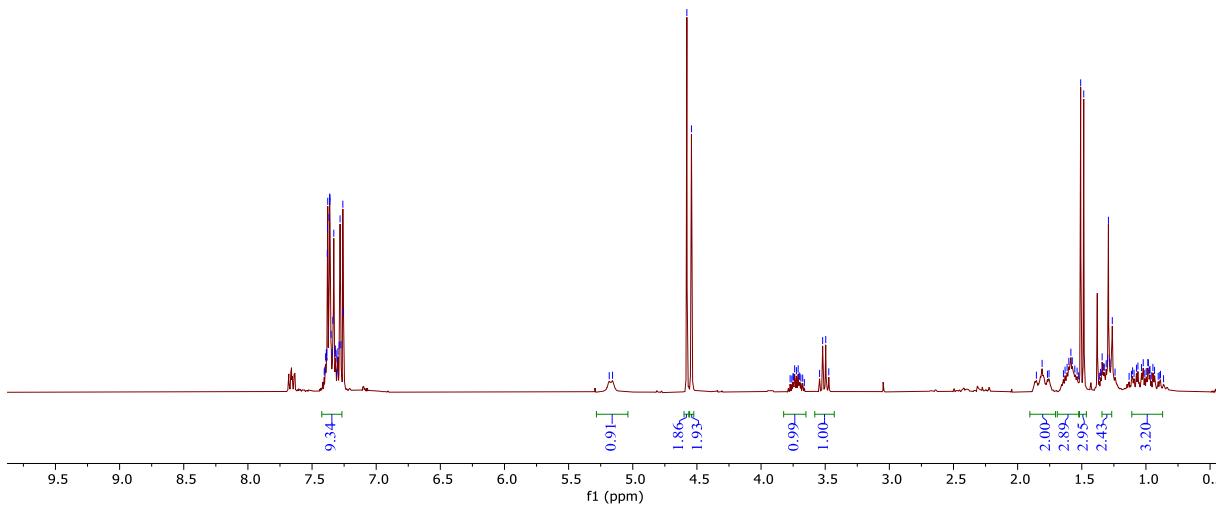
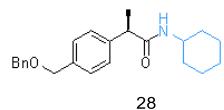
Yang Yuan YY-X-88 — C13CPD CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 13 — 75.49 MHz



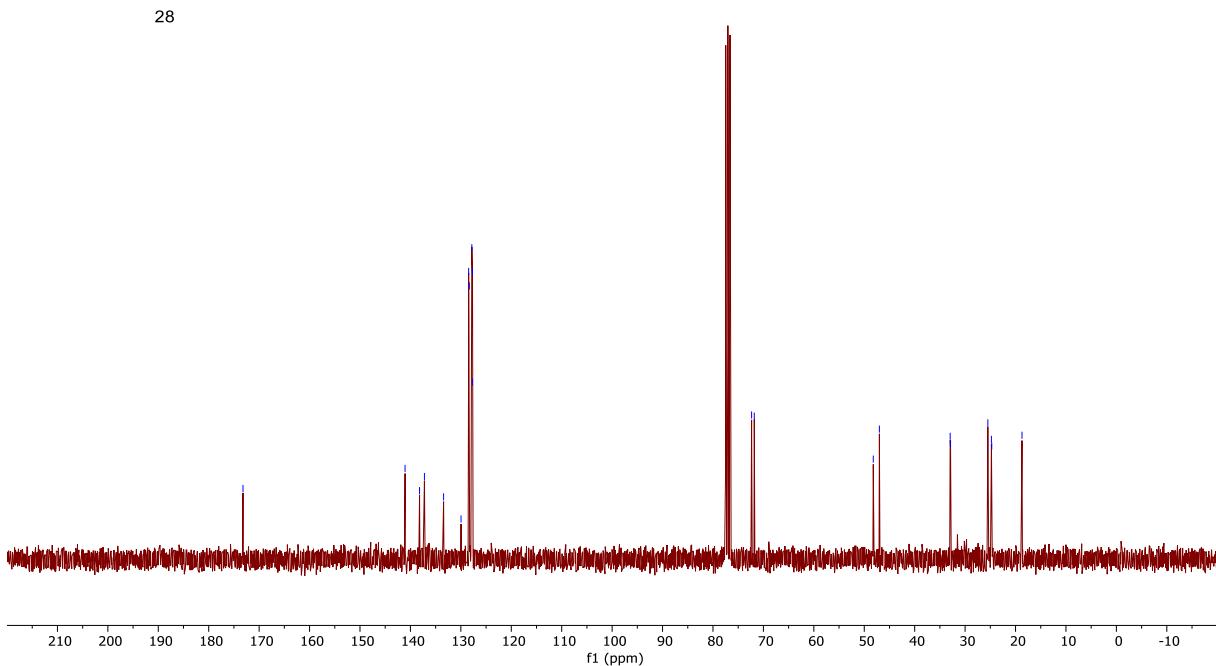
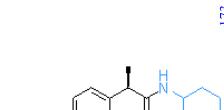
Yuan/YY-X-104 — PROTON CDCl₃ {C(Bruker)TopSpin3.6.2} 2103 33 — 300.20 MHz



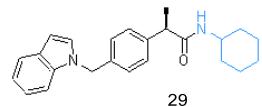
Yang Yuan YY-X-98 — PROTON CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 44 — 300.20 MHz



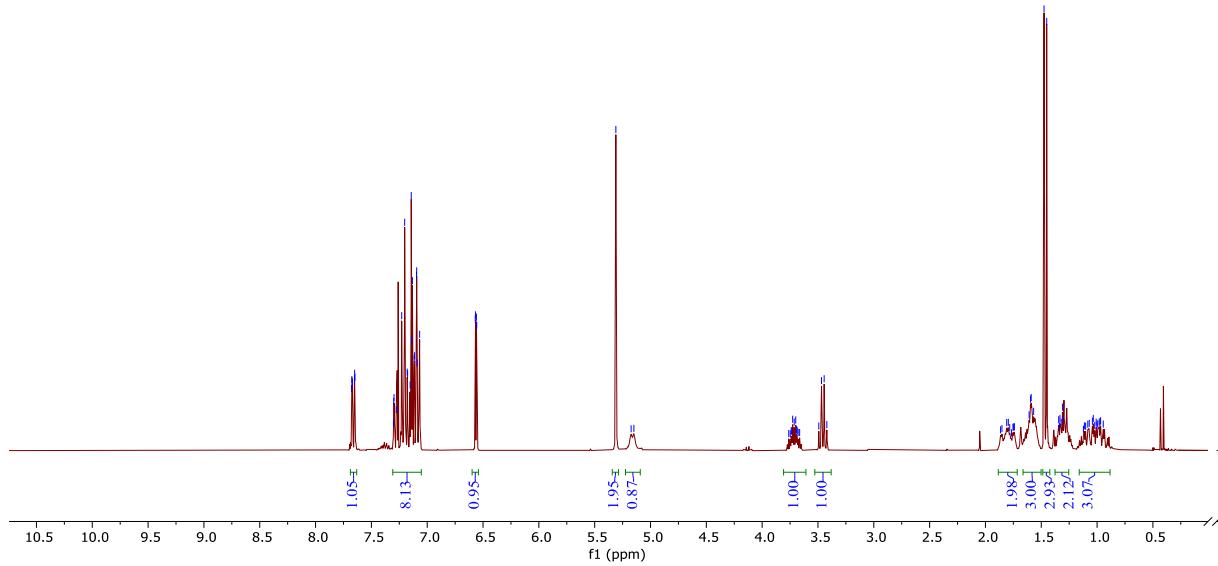
Yang Yuan YY-X-98 — C13CPD CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 44 — 75.49 MHz



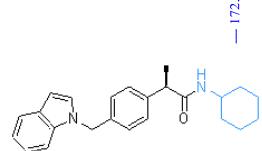
Yang Yuan / YY-X-106	— PROTON C6CD3 {CABruker}TopSpin3.6.2	2103	50 — 300.20 MHz
7.68	7.65	7.28	
7.68	7.65	7.23	
7.67	7.65	7.20	
7.67	7.30	7.18	
7.67	7.30	7.18	
7.68	7.30	7.15	
7.68	7.30	7.14	
7.68	7.30	7.14	
7.68	7.30	7.12	
7.68	7.30	7.11	
7.68	7.30	7.10	
7.68	7.30	7.09	
7.68	7.30	7.09	
7.68	7.30	7.07	
7.68	7.30	7.07	



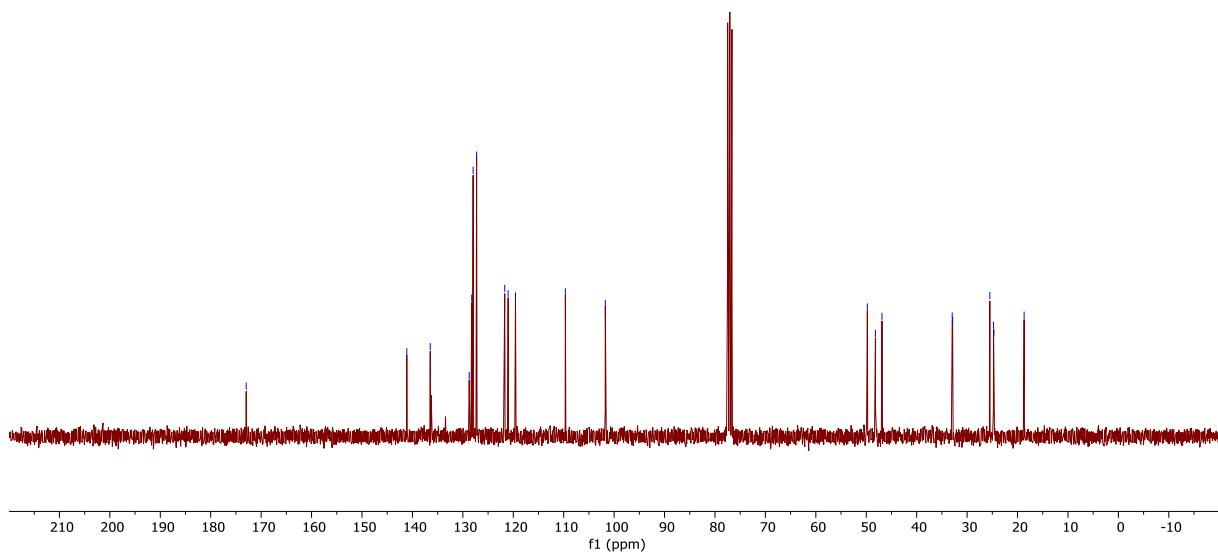
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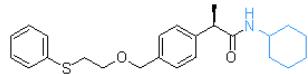


Yang Yuan / YY-X-106 — C13CPD CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 50 — 75.49 MHz

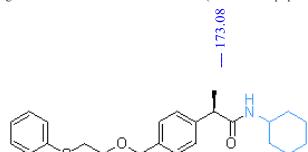
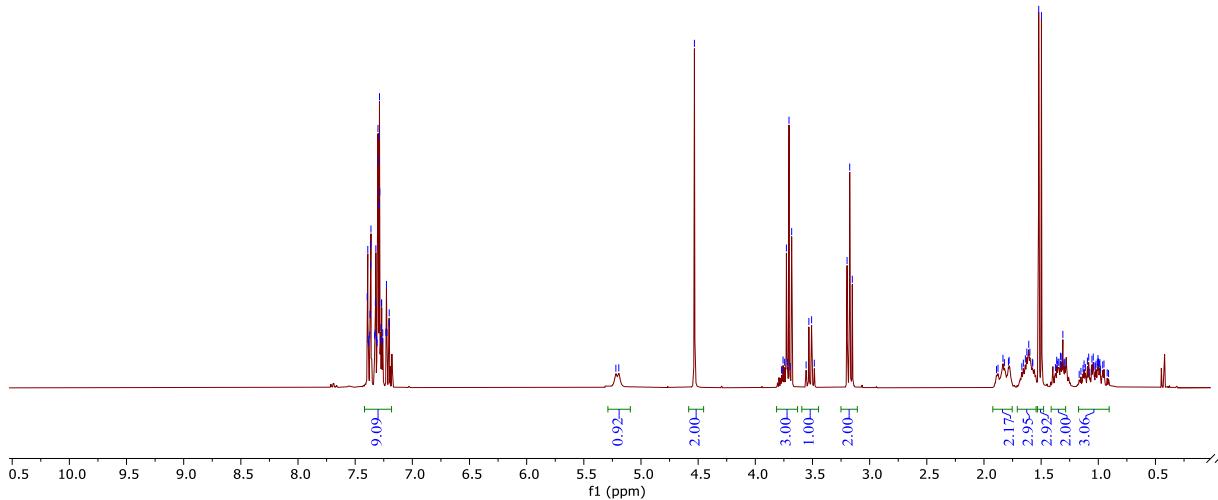


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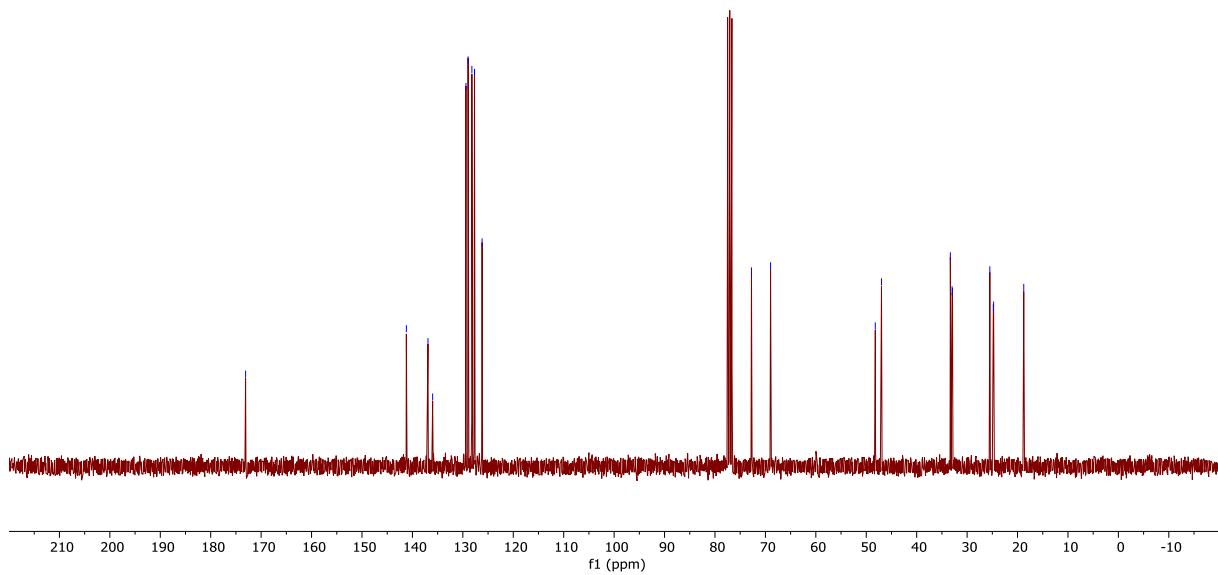


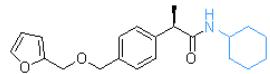
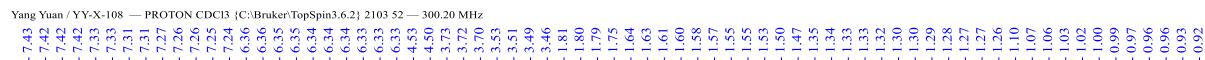


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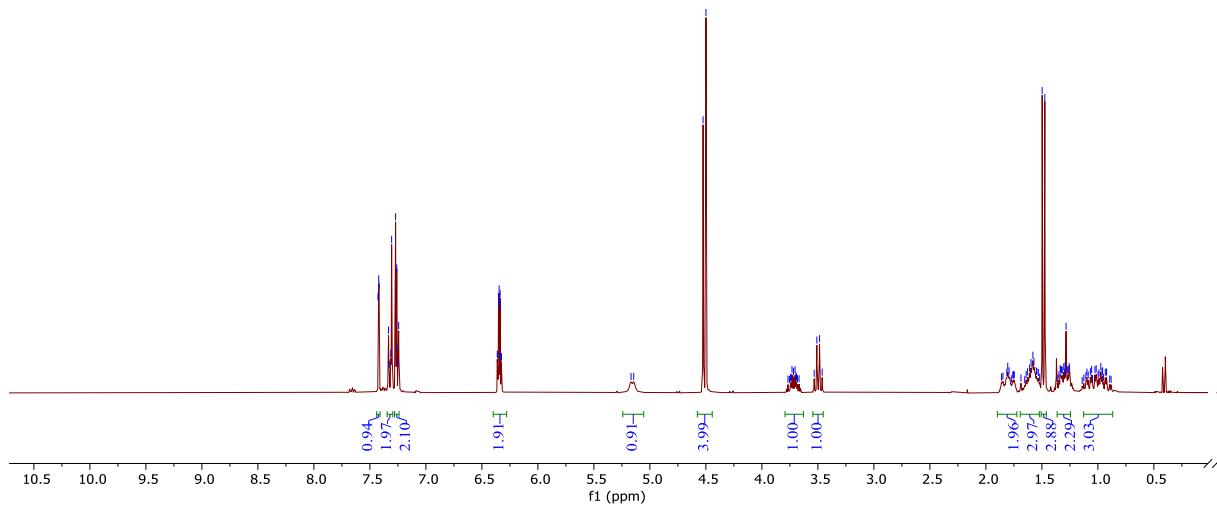


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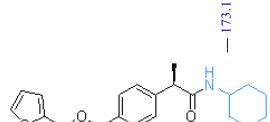




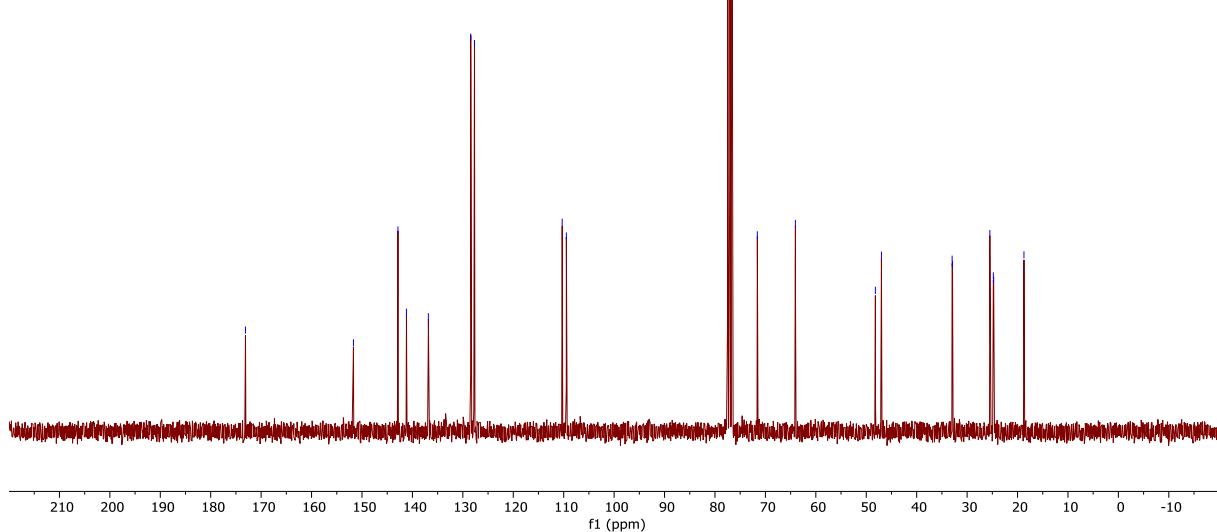
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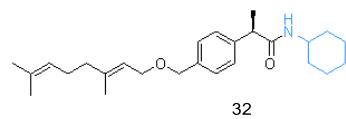
Yang Yuan / YY-X-108 — C13CPD CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 52 — 75.49 MHz



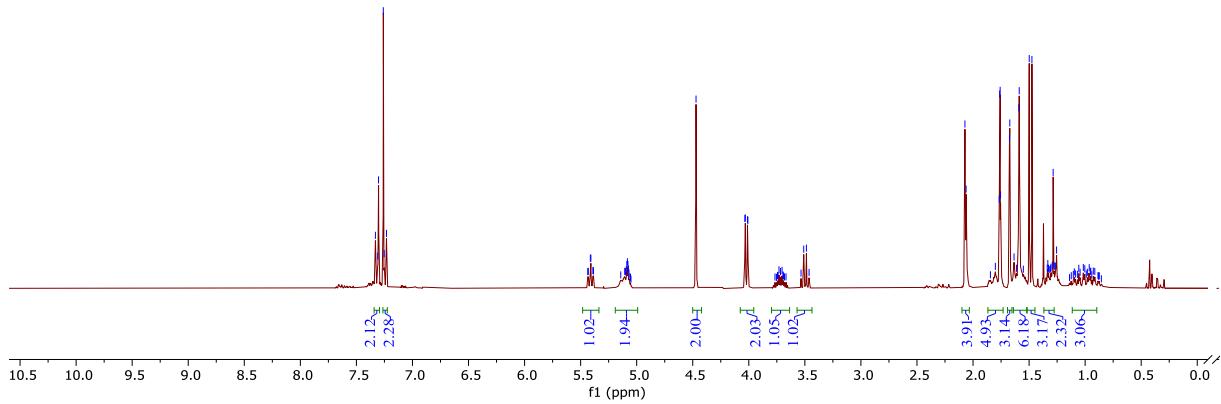
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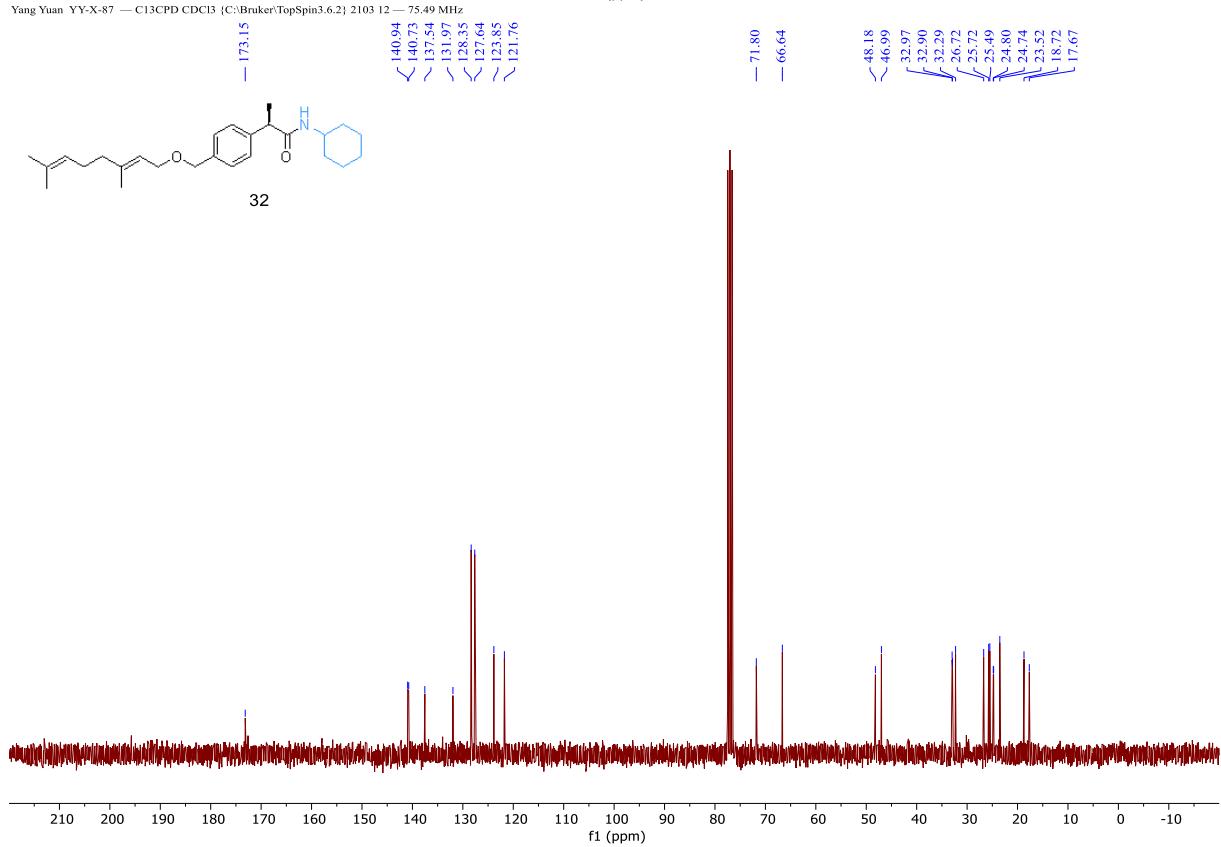
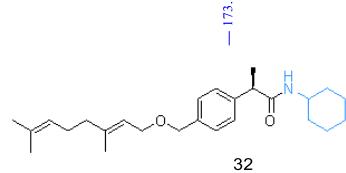
Yang Yuan YY-X-87 — PROTON CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 12 — 300.20 MHz



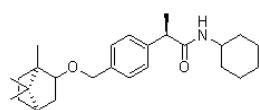
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7.25
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5.43
5.41
5.39
5.38
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5.10
5.09
5.08
5.07
5.07
5.07
4.47
4.03
4.03
4.01
4.01
3.73
3.72
3.70
3.53
3.51
3.49
3.46
2.07
2.06
1.80
1.77
1.76
1.75
1.68
1.67
1.63
1.61
1.59
1.59
1.55
1.50
1.47
1.34
1.33
1.33
1.32
1.31
1.30
1.29
1.29
1.28
1.27
1.25
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0.98
0.96
0.95
0.95
0.93
0.93
0.91



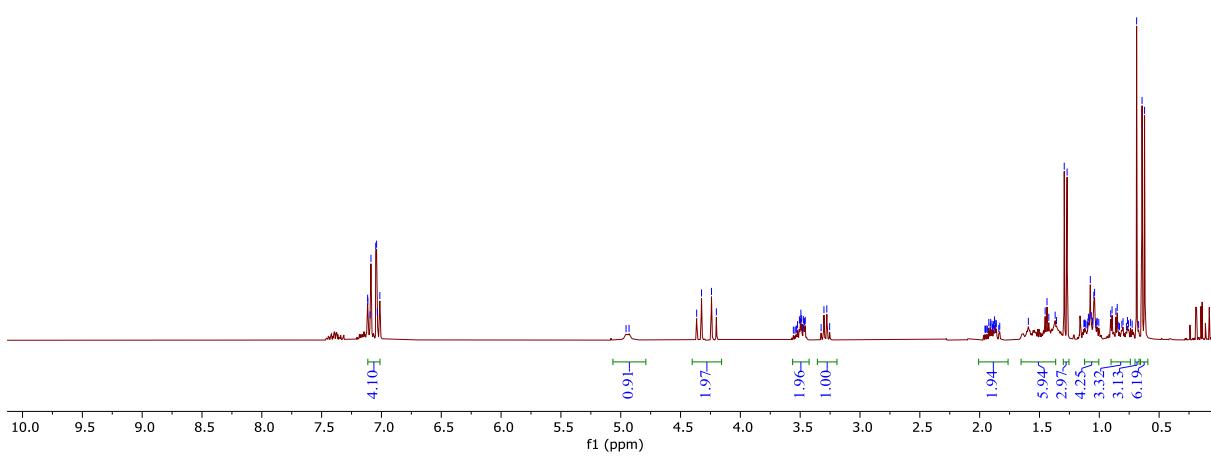
Yang Yuan YY-X-87 — C13CPD CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 12 — 75.49 MHz



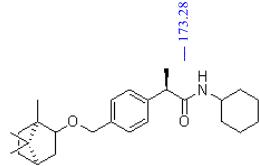
Yang Yuan YY-X-99 — PROTON CDCl₃ [C:\Bruker\TopSpin3.6.2] 2103.45 — 300.20 MHz



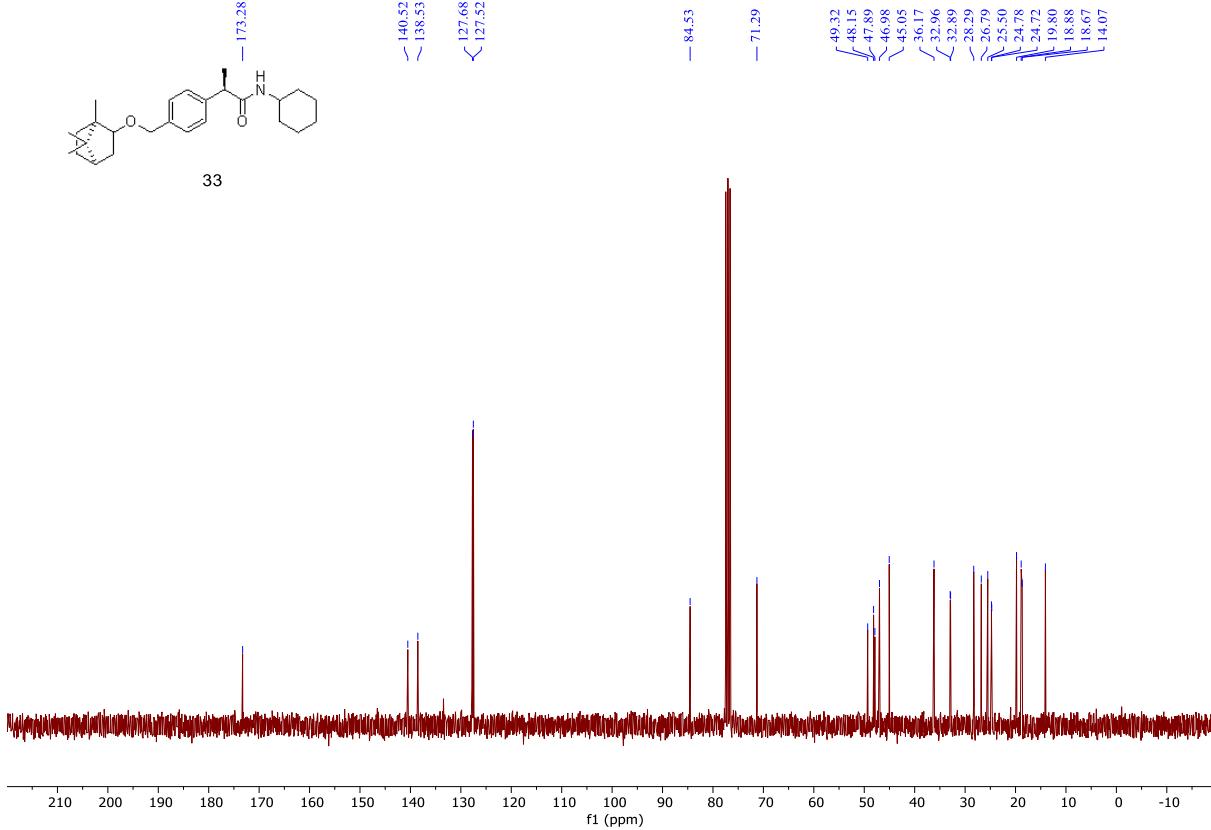
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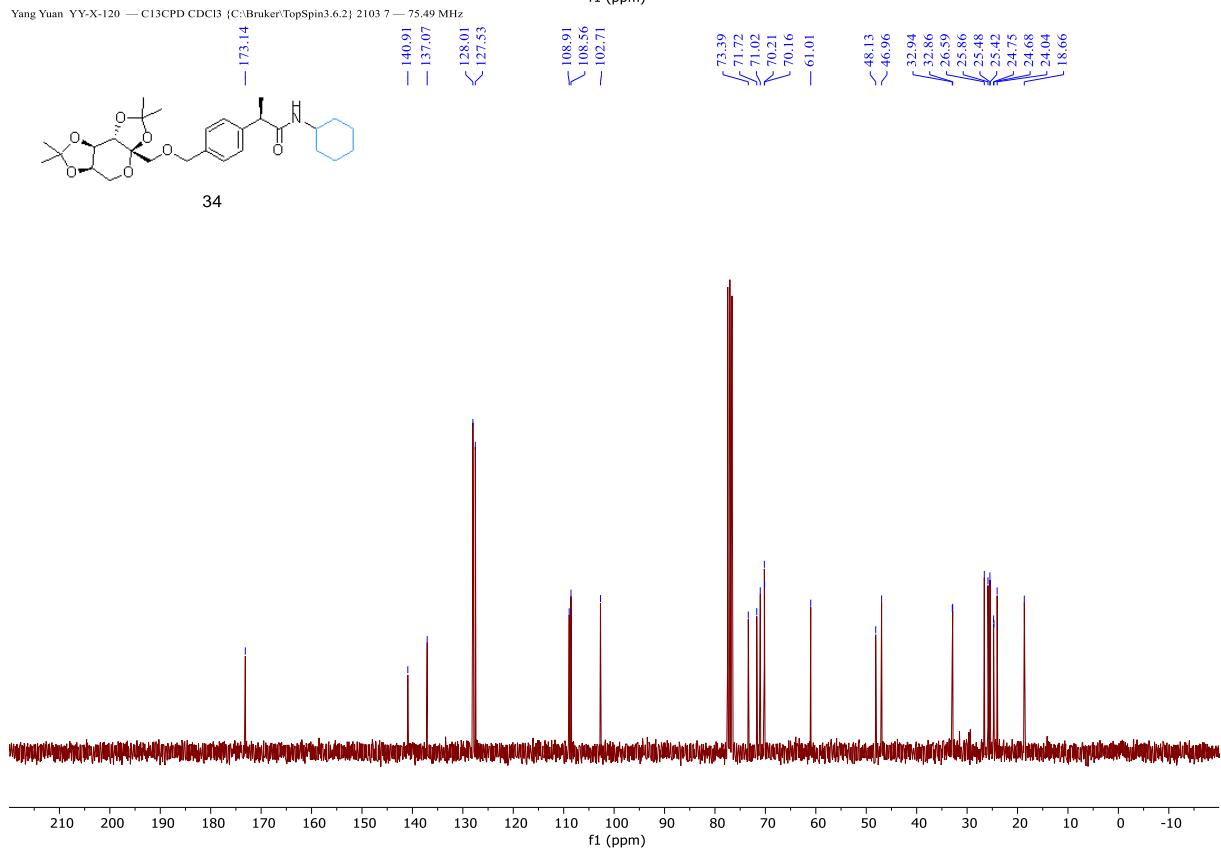
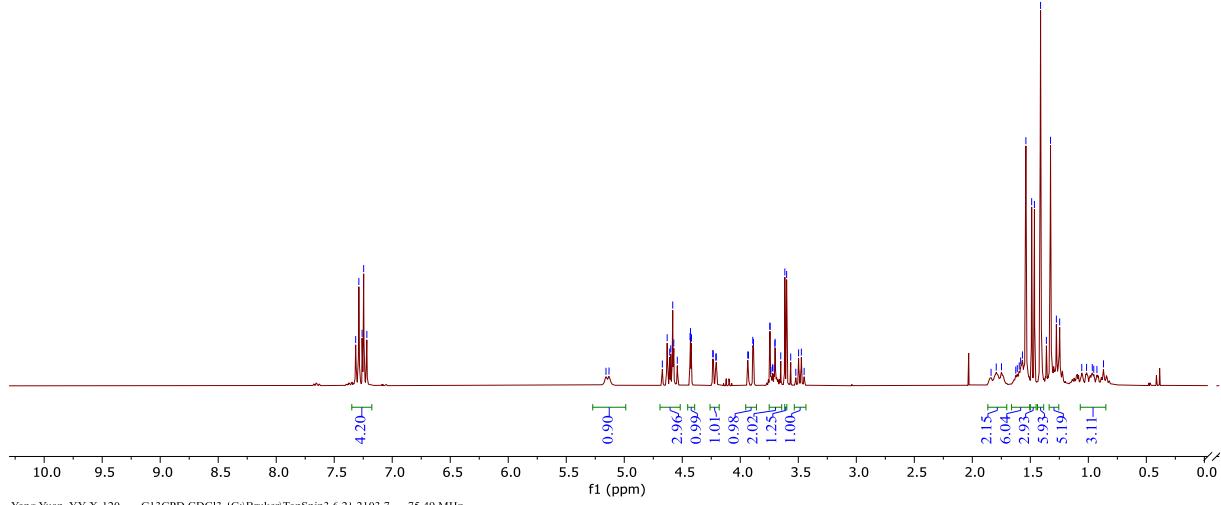
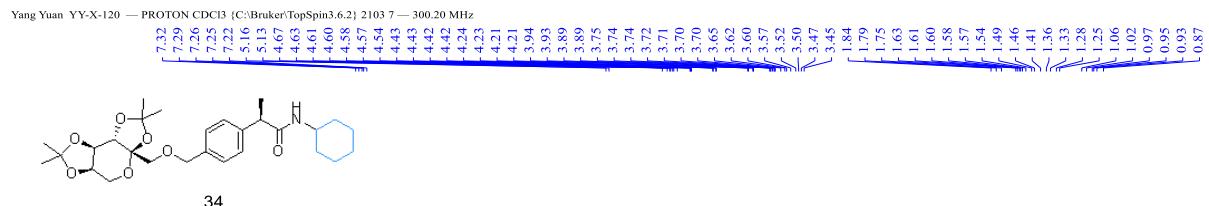


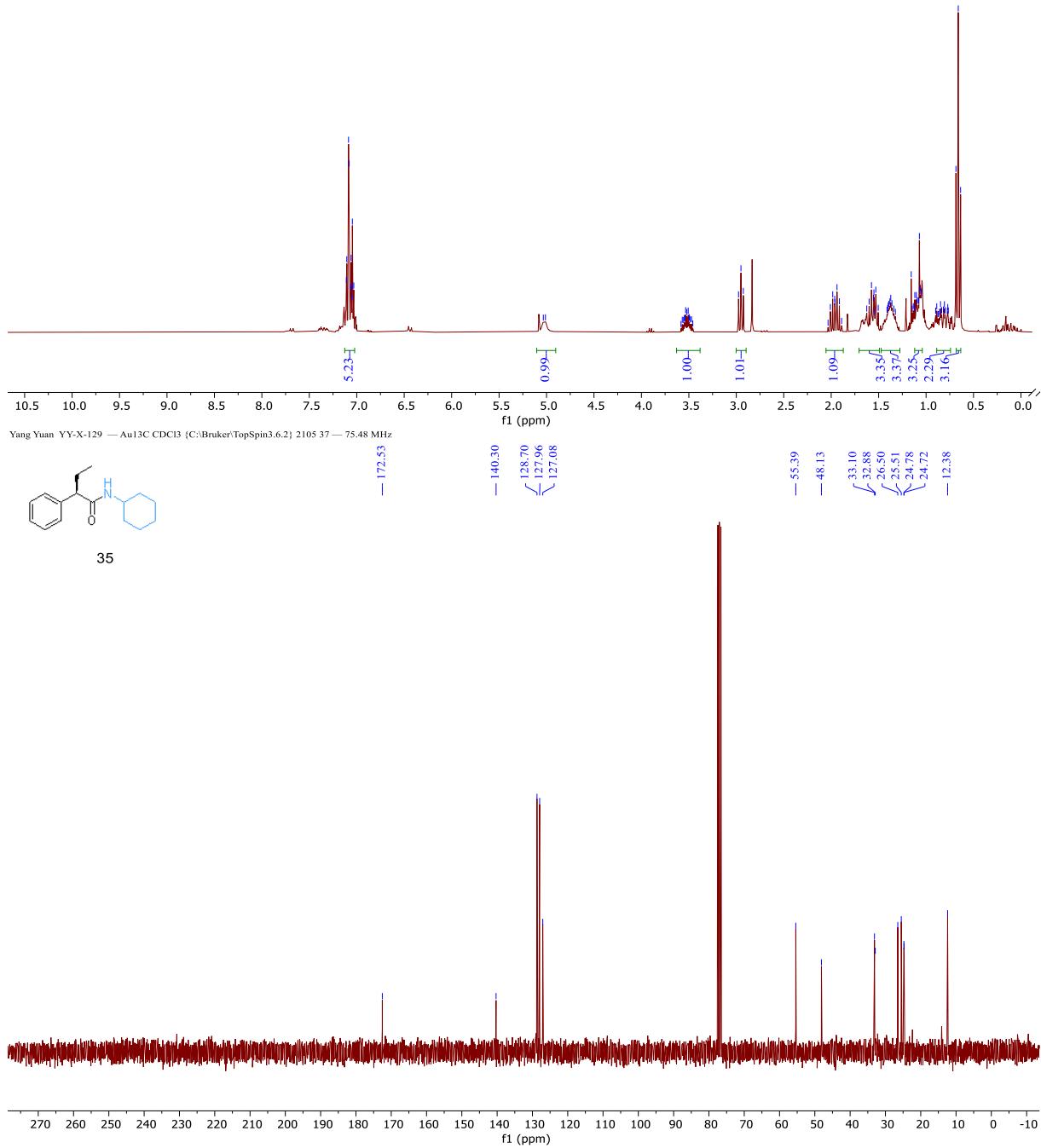
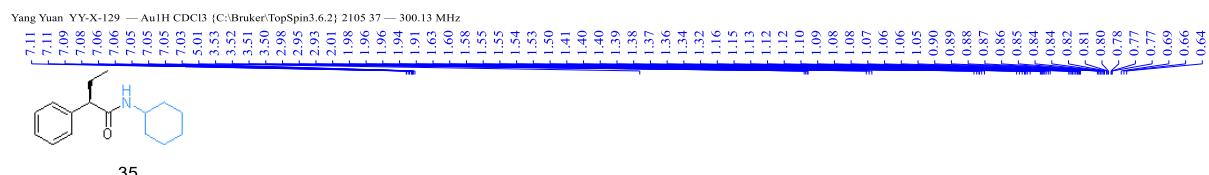
Yang Yuan YY-X-99 — C13CPD CDCl₃ [C:\Bruker\TopSpin3.6.2] 2103.45 — 75.49 MHz

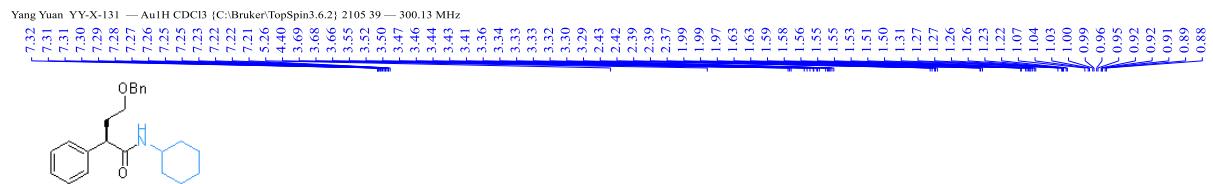


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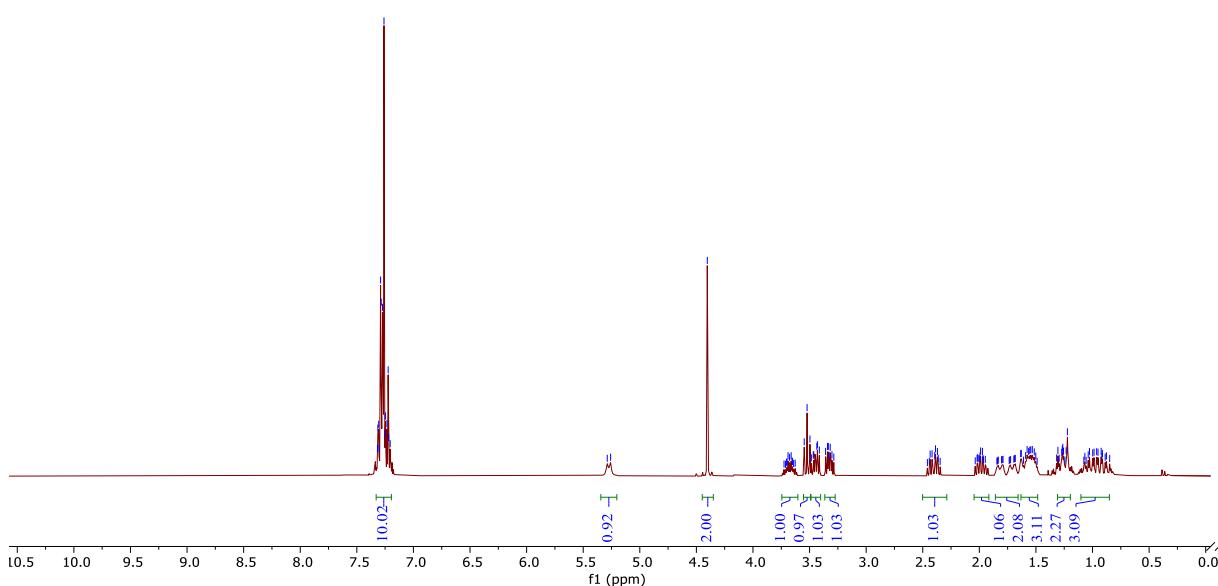




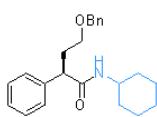




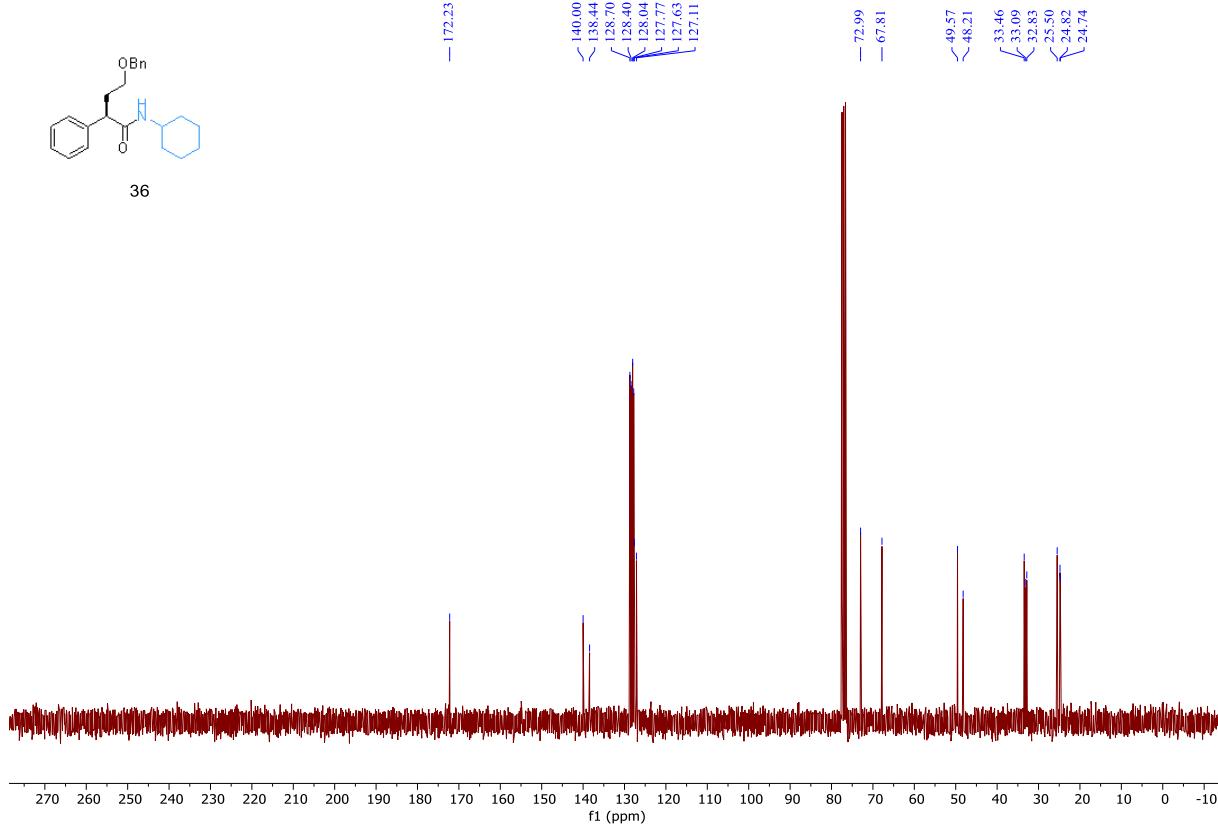
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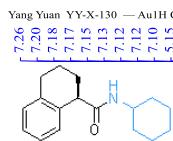


Yang Yuan YY-X-131 — Au13C CDCl₃ {C:\Bruker\TopSpin3.6.2} 2105 39 — 75.48 MHz

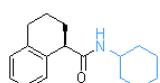
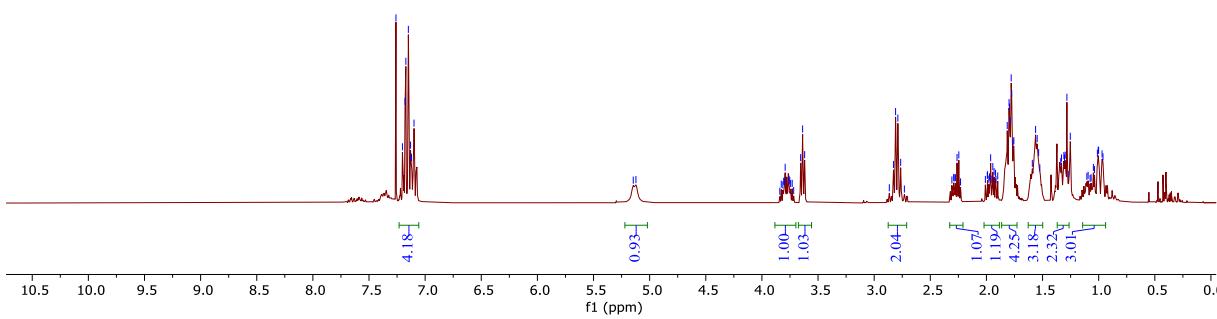


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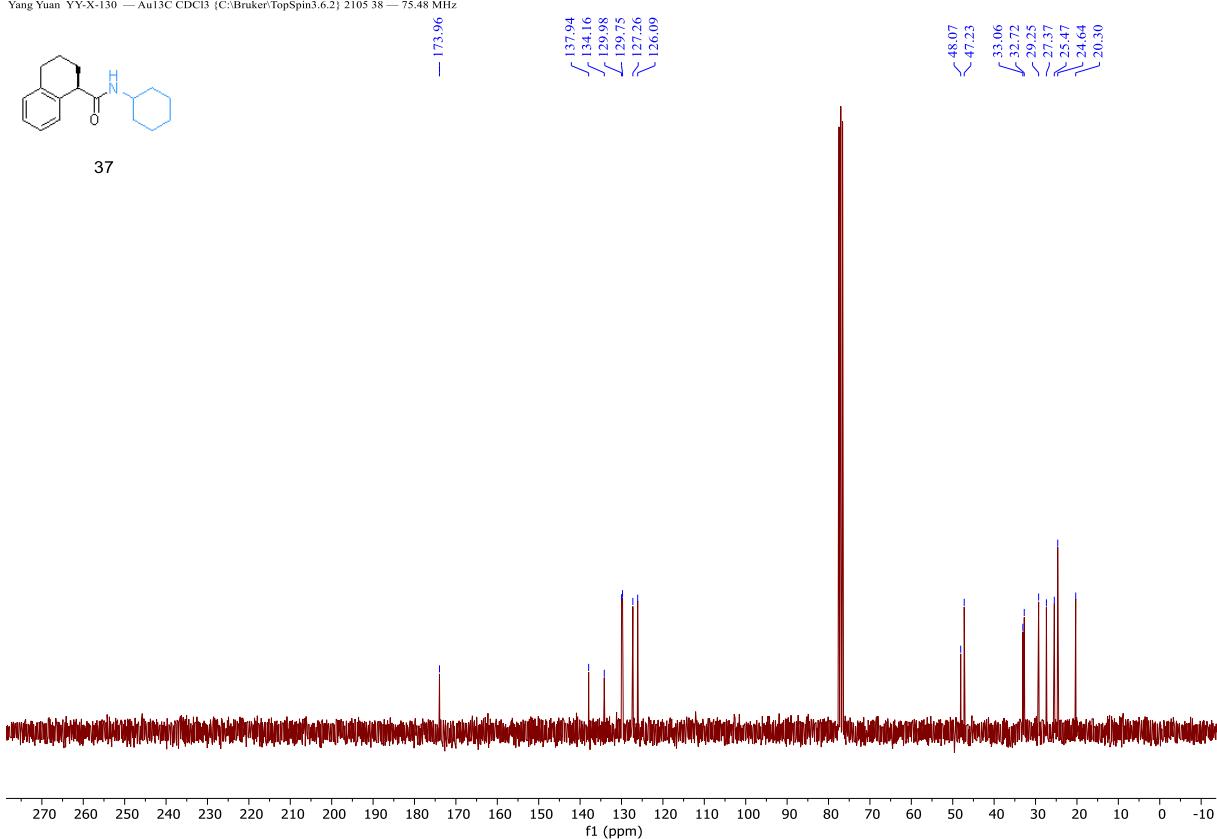


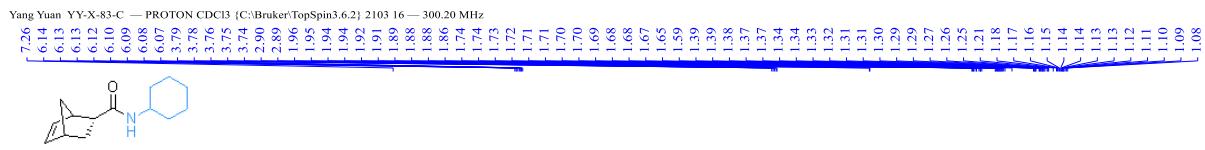


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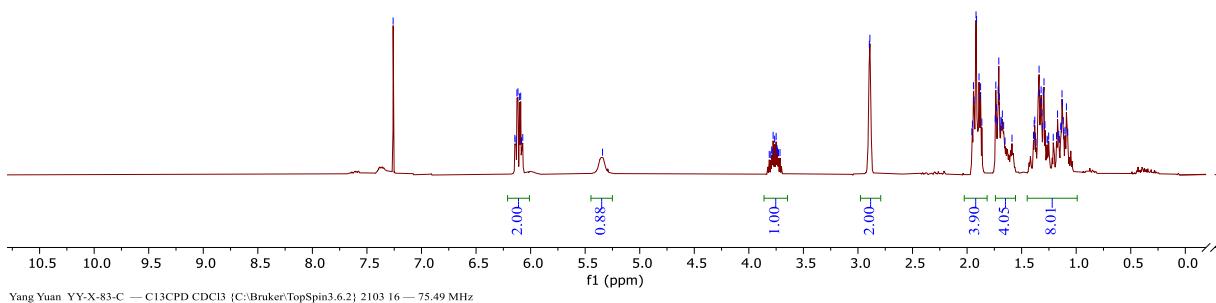


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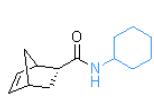




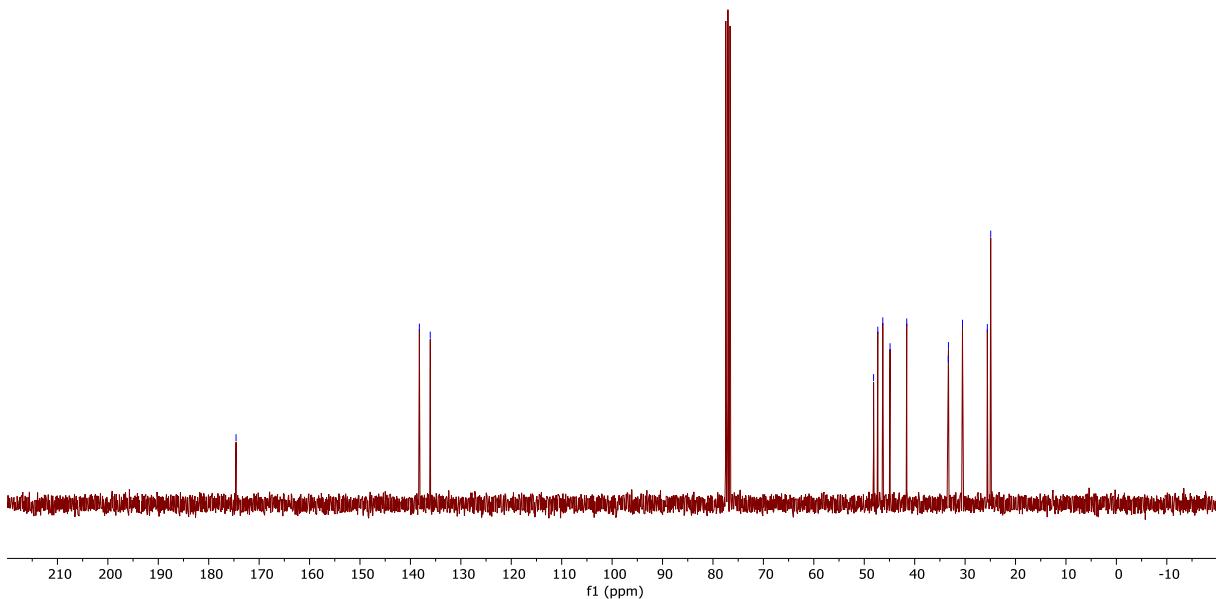
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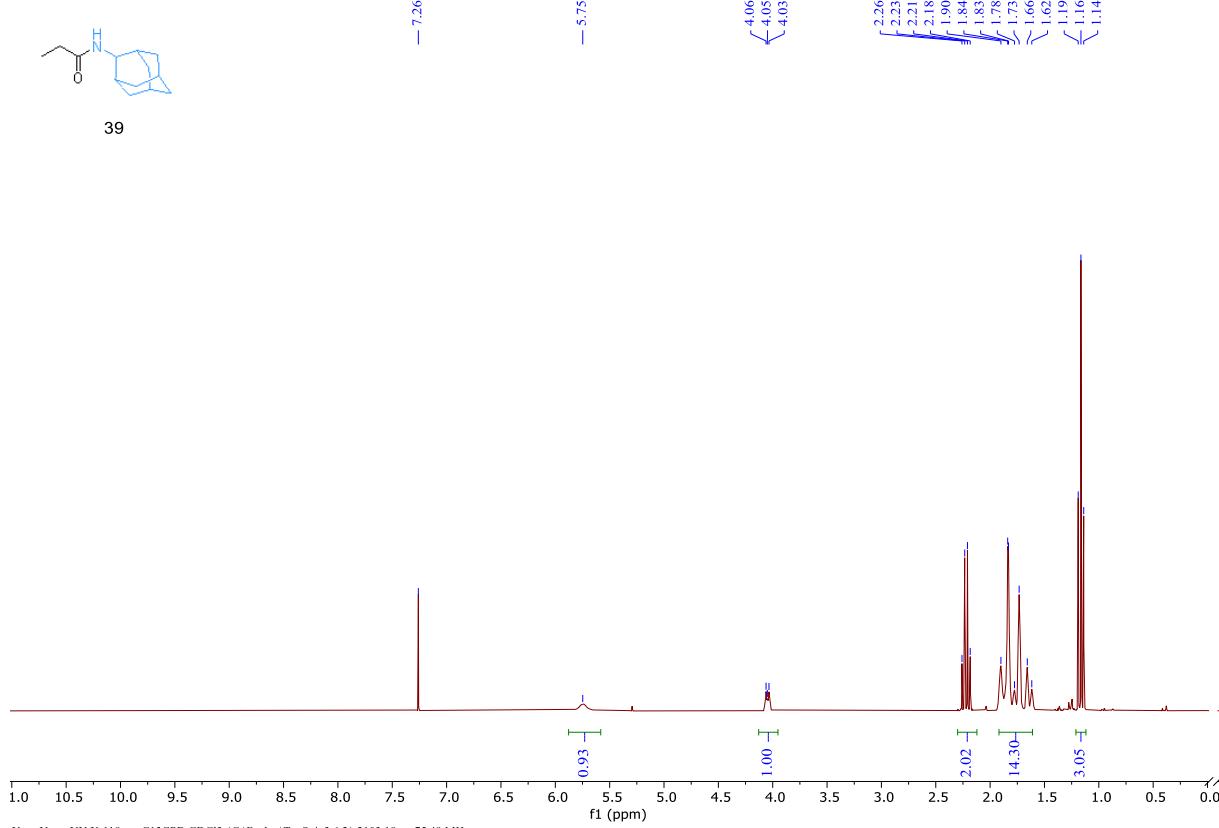
Yang Yuan YY-X-83-C — C13CPD CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103.16 — 75.49 MHz



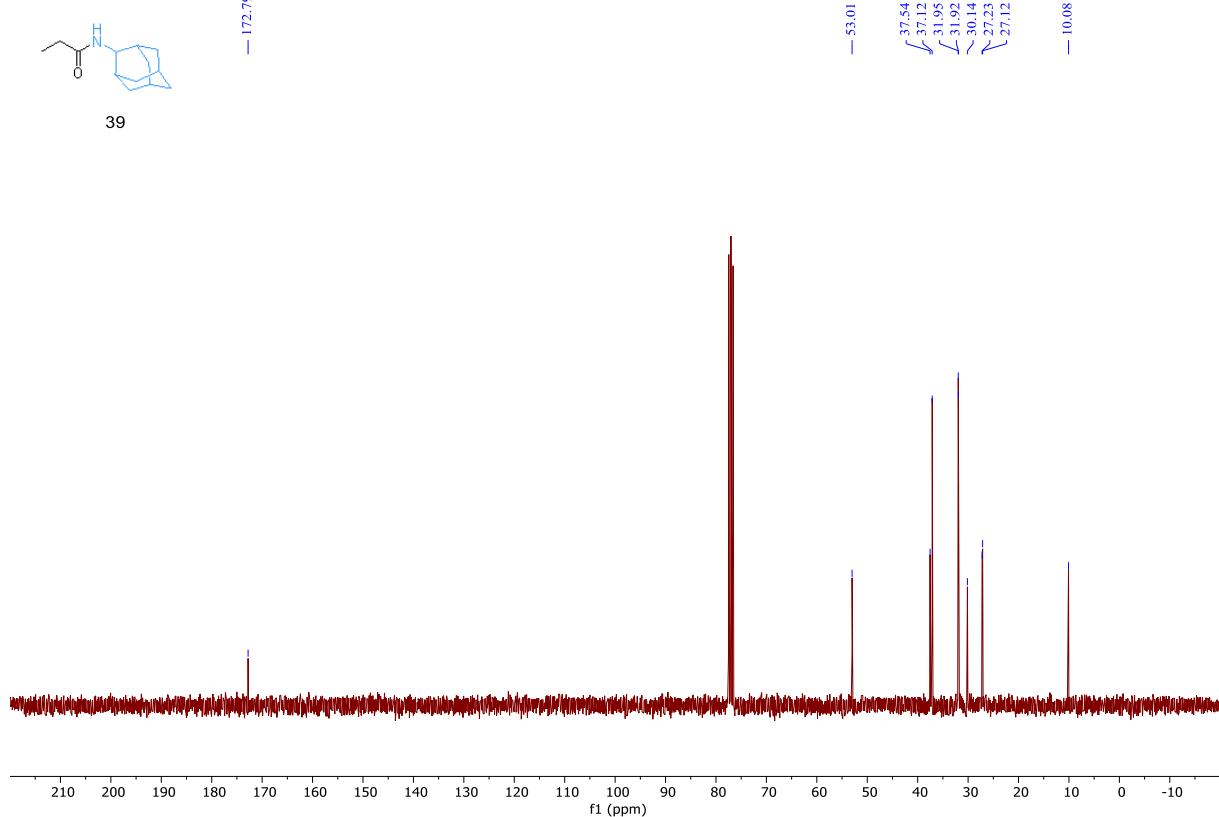
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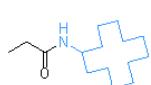
Yang Yuan YY-X-118 — PROTON CDCl₃ {C:\Bruker\TopSpin3.6.2\} 2103 18 — 300.20 MHz



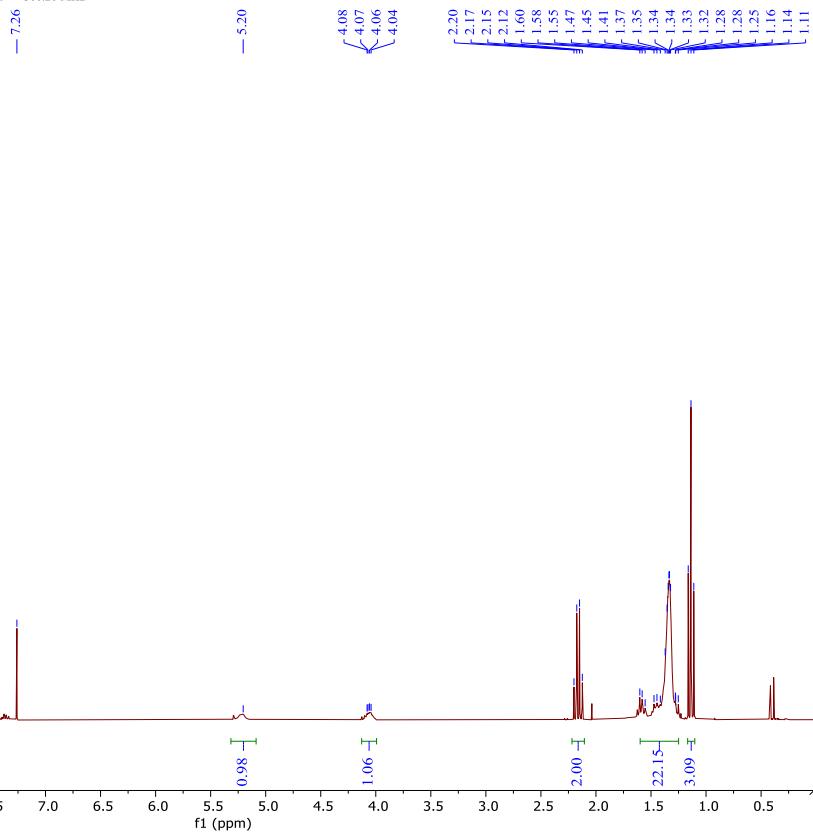
Yang Yuan YY-X-118 — C13CPD CDCl₃ {C:\Bruker\TopSpin3.6.2\} 2103 18 — 75.49 MHz



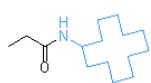
Yang Yuan YY-X-119 — PROTON CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 19 — 300.20 MHz



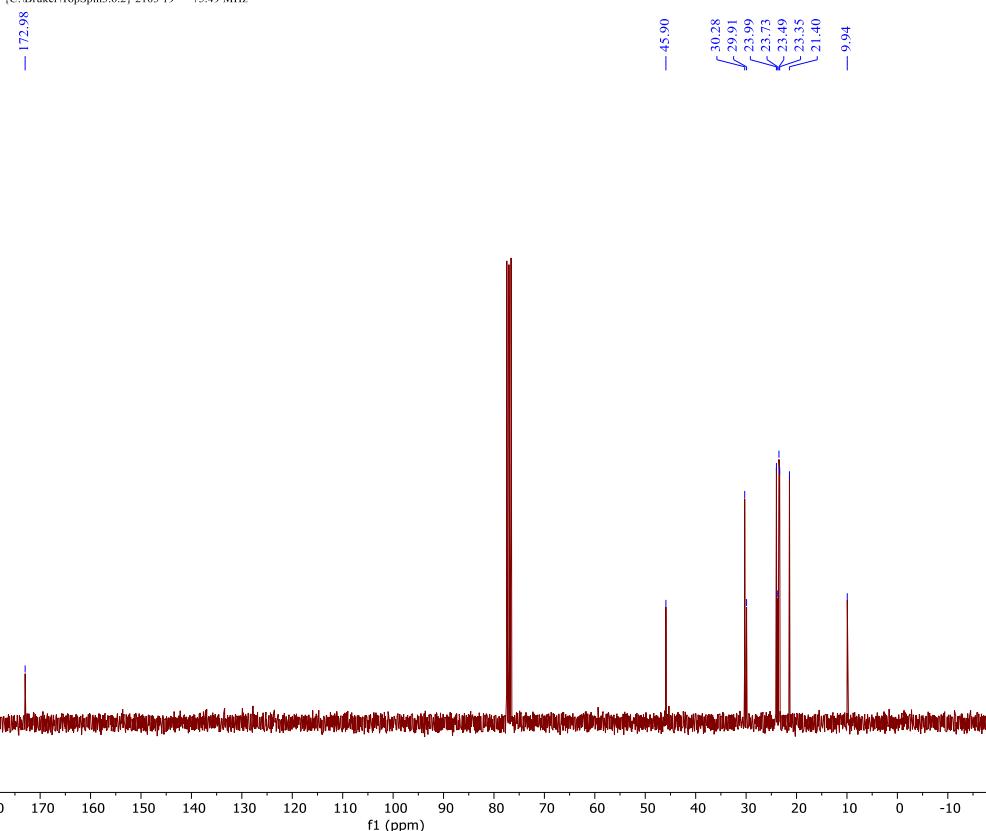
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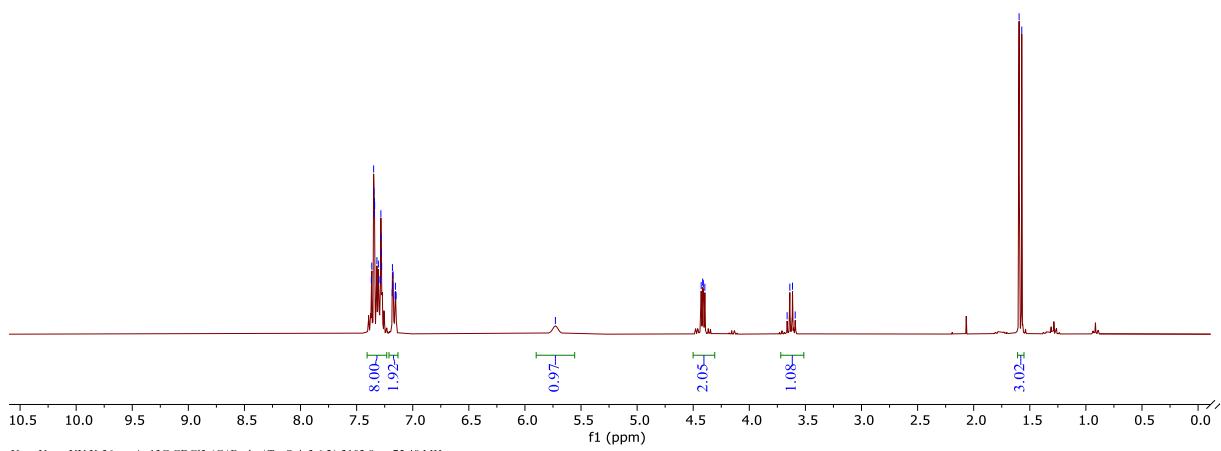
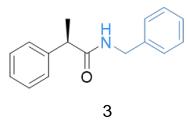
Yang Yuan YY-X-119 — C13CPD CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 19 — 75.49 MHz



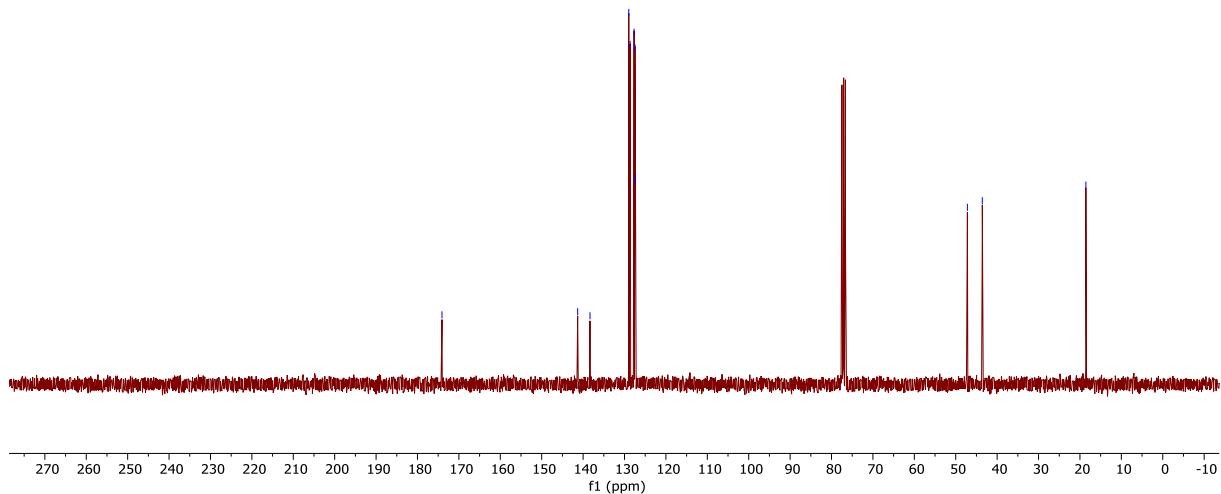
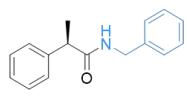
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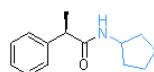
Yang Yuan YY-X-36 — Au1H CDCl₃ {C:\Bruker\TopSpin3.6.2} 2102.8 — 300.13 MHz



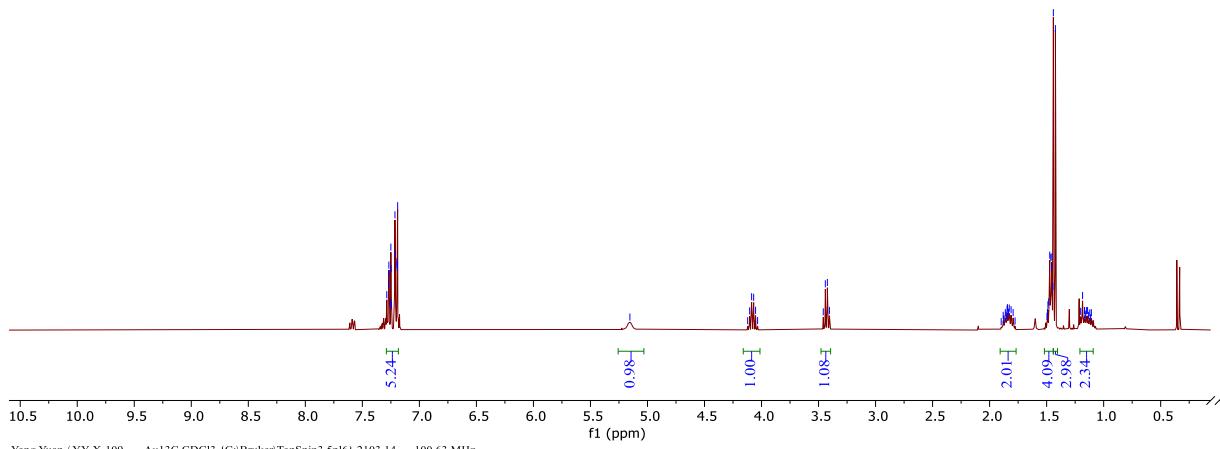
Yang Yuan YY-X-36 — Au13C CDCl₃ {C:\Bruker\TopSpin3.6.2} 2102.8 — 75.48 MHz



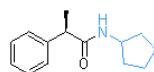
Yang Yuan / YY-X-109 — Au1H CDCl₃ {C:\Bruker\TopSpin3.5pl6} 2103 14 — 400.13 MHz



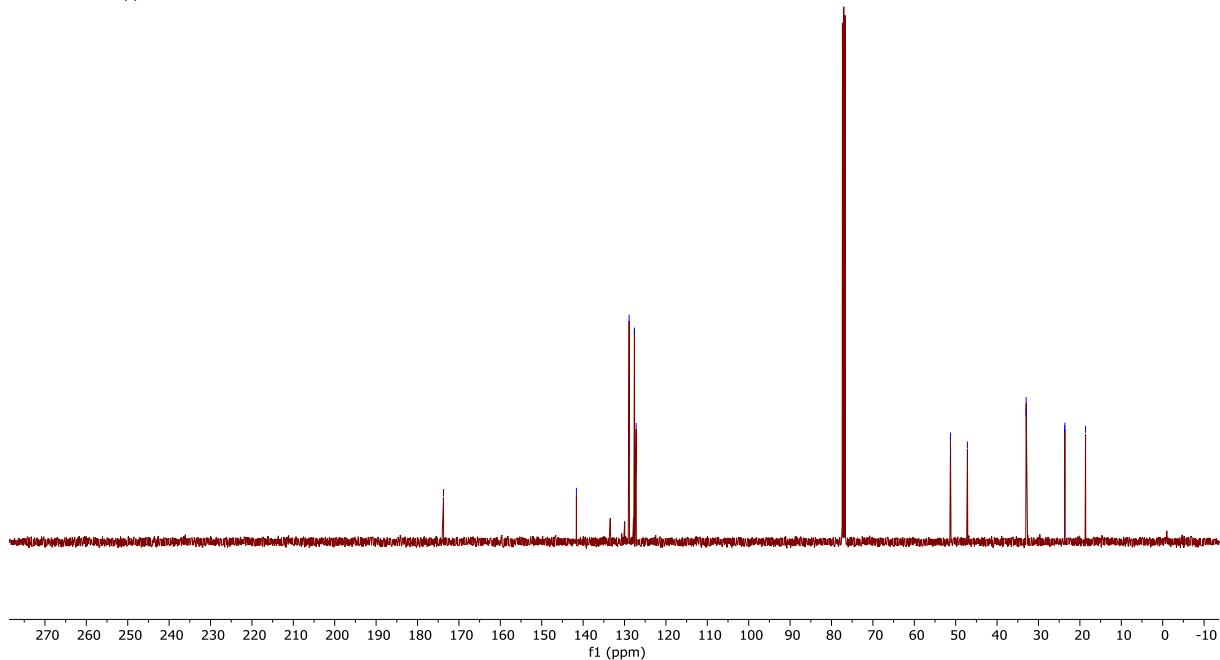
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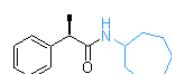
Yang Yuan / YY-X-109 — Au13C CDCl₃ {C:\Bruker\TopSpin3.5pl6} 2103 14 — 100.63 MHz



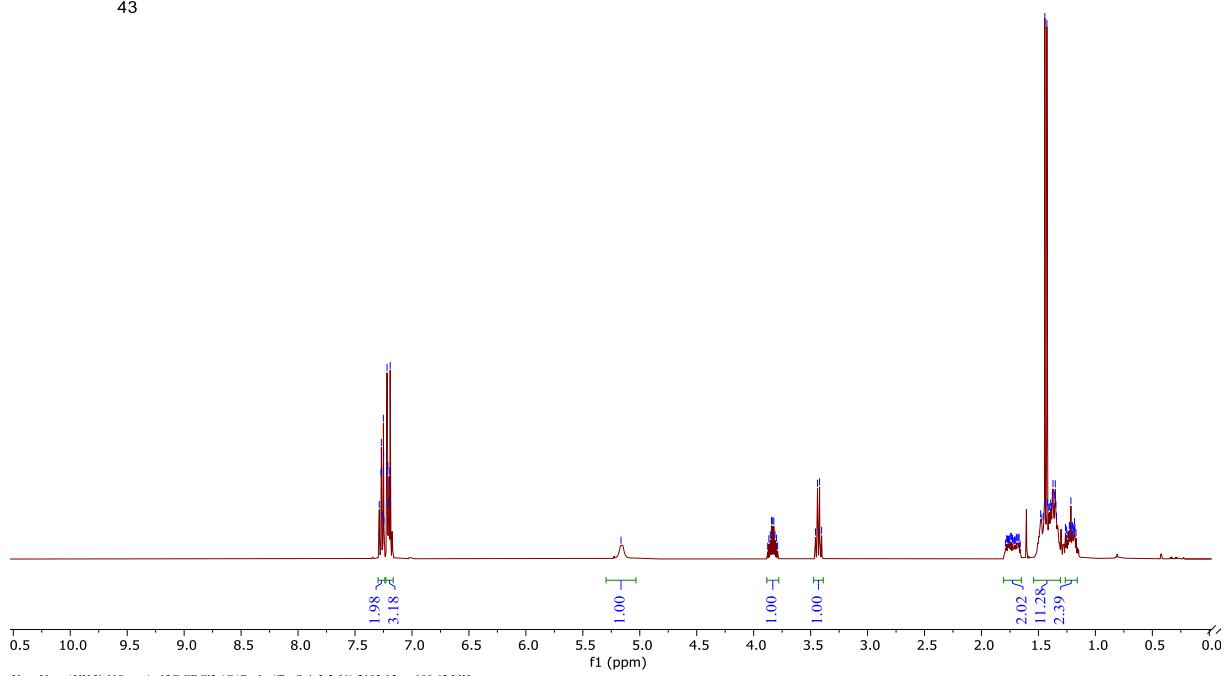
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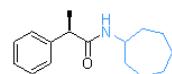
Yang Yuan / YY-X-110 — Au1H CDCl₃ {C:\Bruker\TopSpin3.5\pl6} 2103 15 — 400.13 MHz



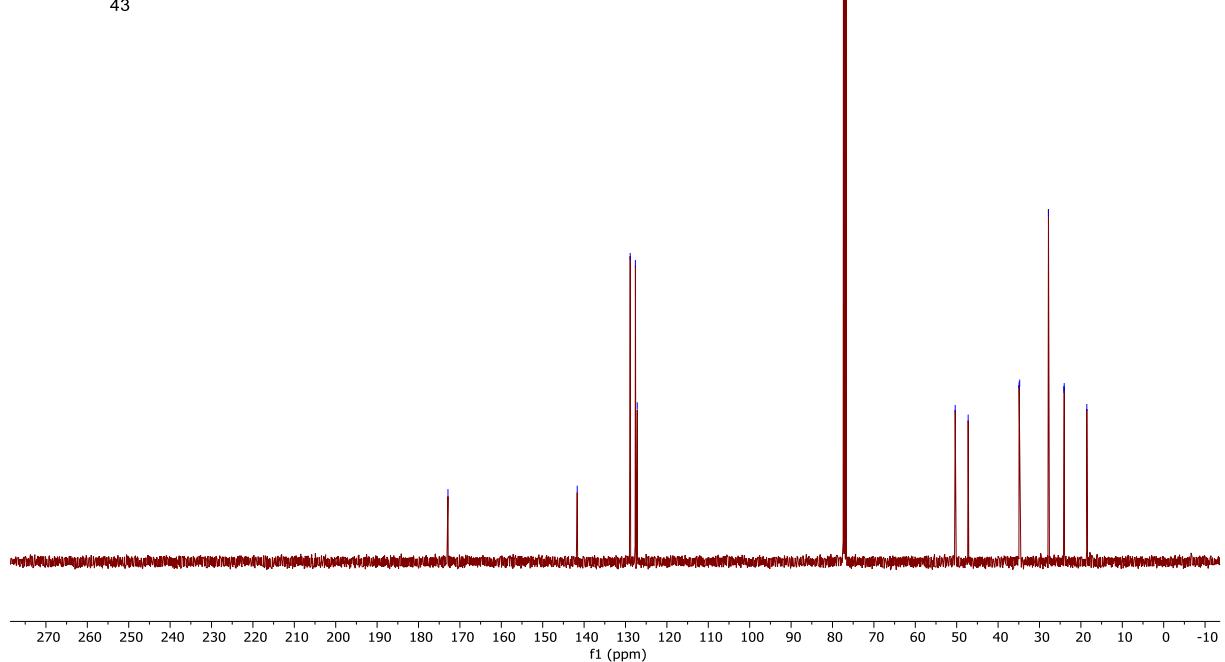
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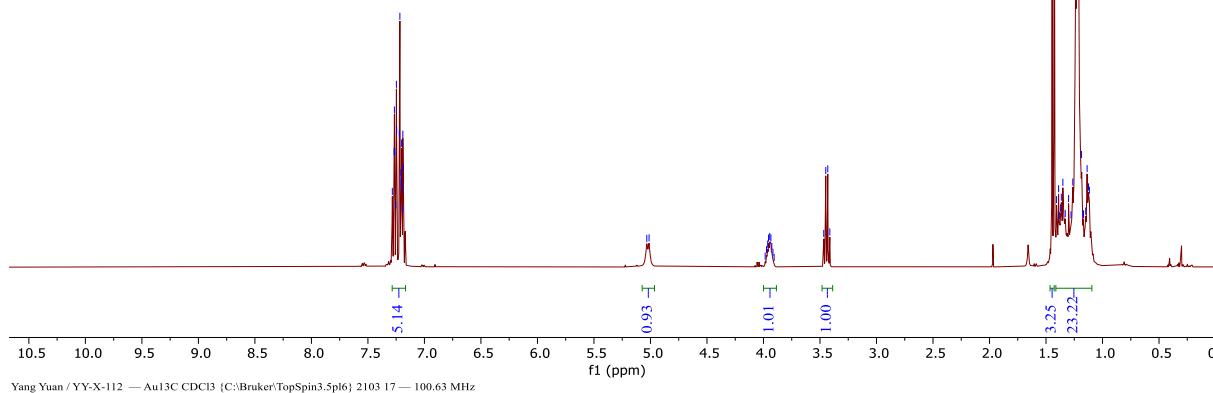
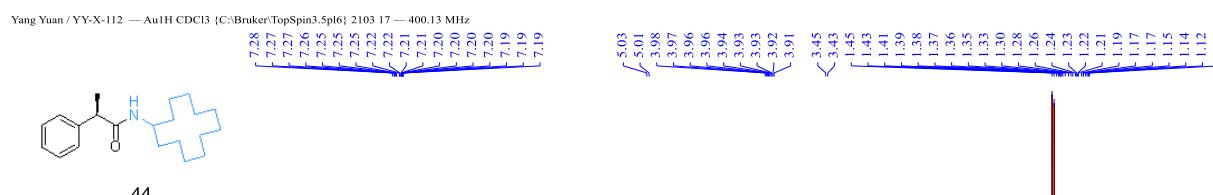


Yang Yuan / YY-X-110 — Au13C CDCl₃ {C:\Bruker\TopSpin3.5\pl6} 2103 15 — 100.63 MHz

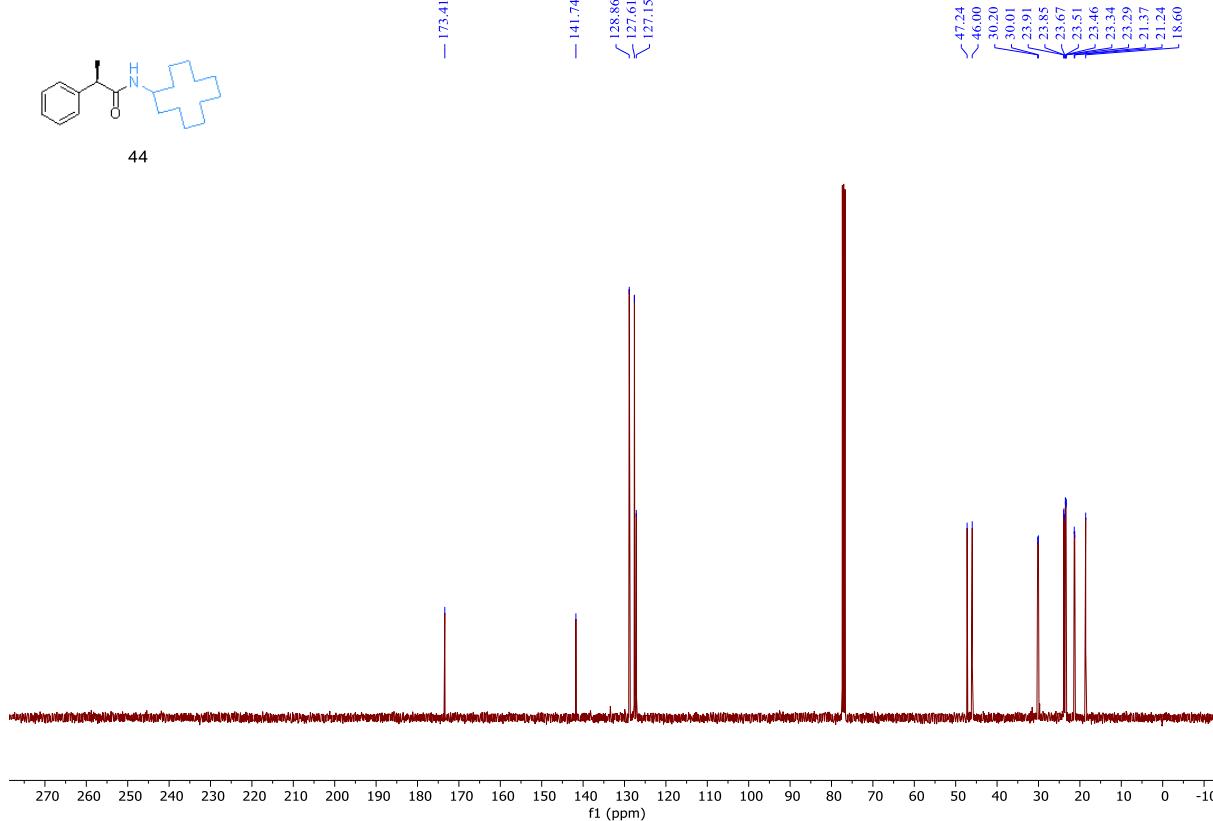


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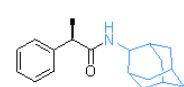




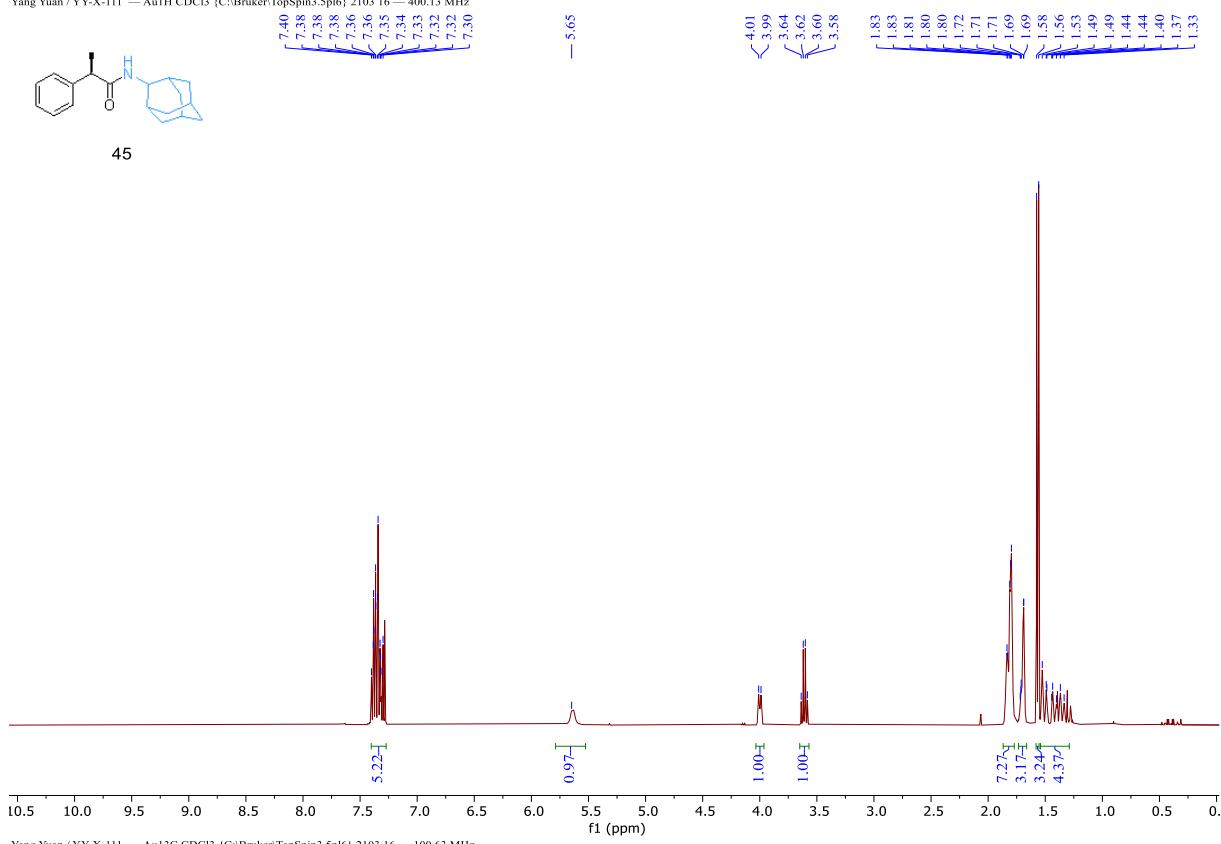
Yang Yuan / YY-X-112 — Au13C CDCl₃ {C:\Bruker\TopSpin3.5\pl6} 2103 17 — 100.63 MHz



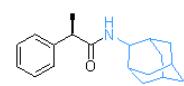
Yang Yuan / YY-X-111 — Au1H CDCl₃ {C:\Bruker\TopSpin3.5pl6} 2103 16 — 400.13 MHz



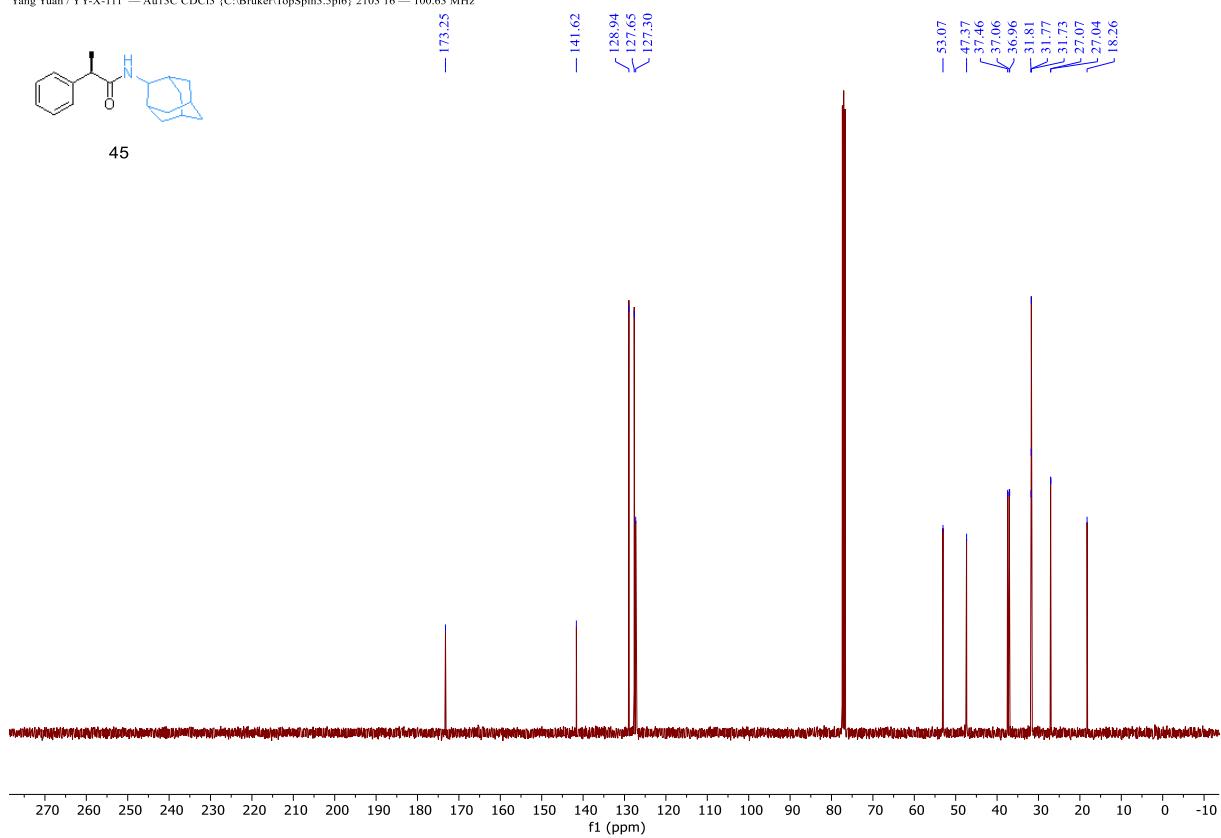
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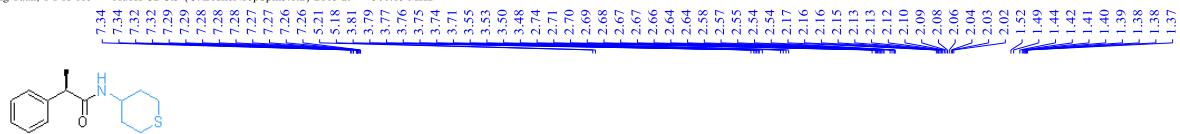
Yang Yuan / YY-X-111 — Au13C CDCl₃ {C:\Bruker\TopSpin3.5pl6} 2103 16 — 100.63 MHz



45

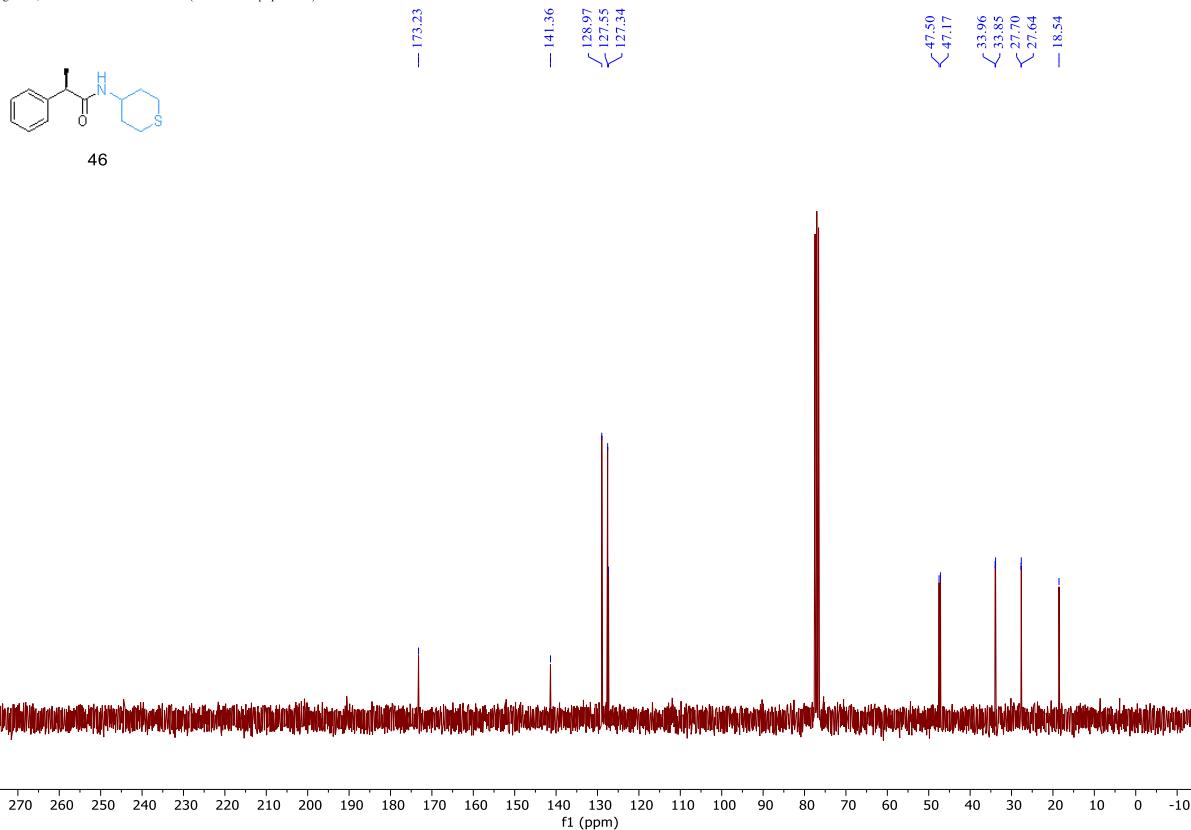


Yang Yuan, YY-X-115 — Au1H CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 29 — 300.13 MHz

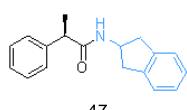


46

Yang Yuan, YY-X-115 — Au13C CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 29 — 75.48 MHz



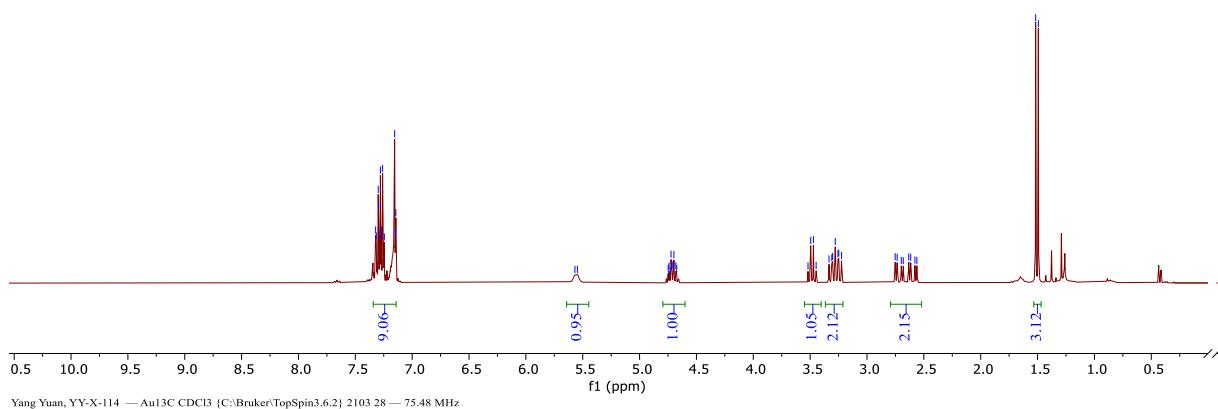
Yang Yuan, YY-X-114 — Au1H CDCl3 {C:\Bruker\TopSpin3.6.2} 2103 28 — 300.13 MHz



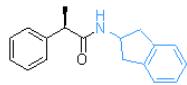
47

7.32
7.32
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7.30
7.28
7.27
7.27
7.26
7.24
7.15
7.15
7.14

5.57
5.55
4.75
4.74
4.73
4.72
4.72
4.71
4.70
4.68
4.67
3.52
3.49
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3.45
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3.31
3.30
3.28
3.26
3.25
3.22
2.75
2.73
2.70
2.68
2.63
2.62
2.58
2.56
1.52
1.49
1.49



Yang Yuan, YY-X-114 — Au13C CDCl3 {C:\Bruker\TopSpin3.6.2} 2103 28 — 75.48 MHz

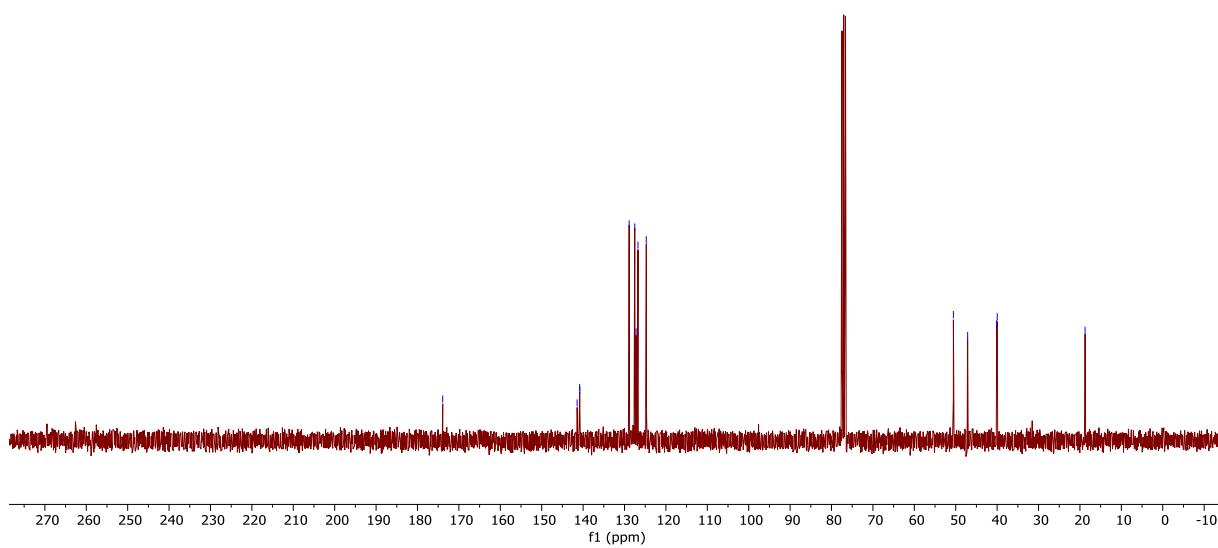


47

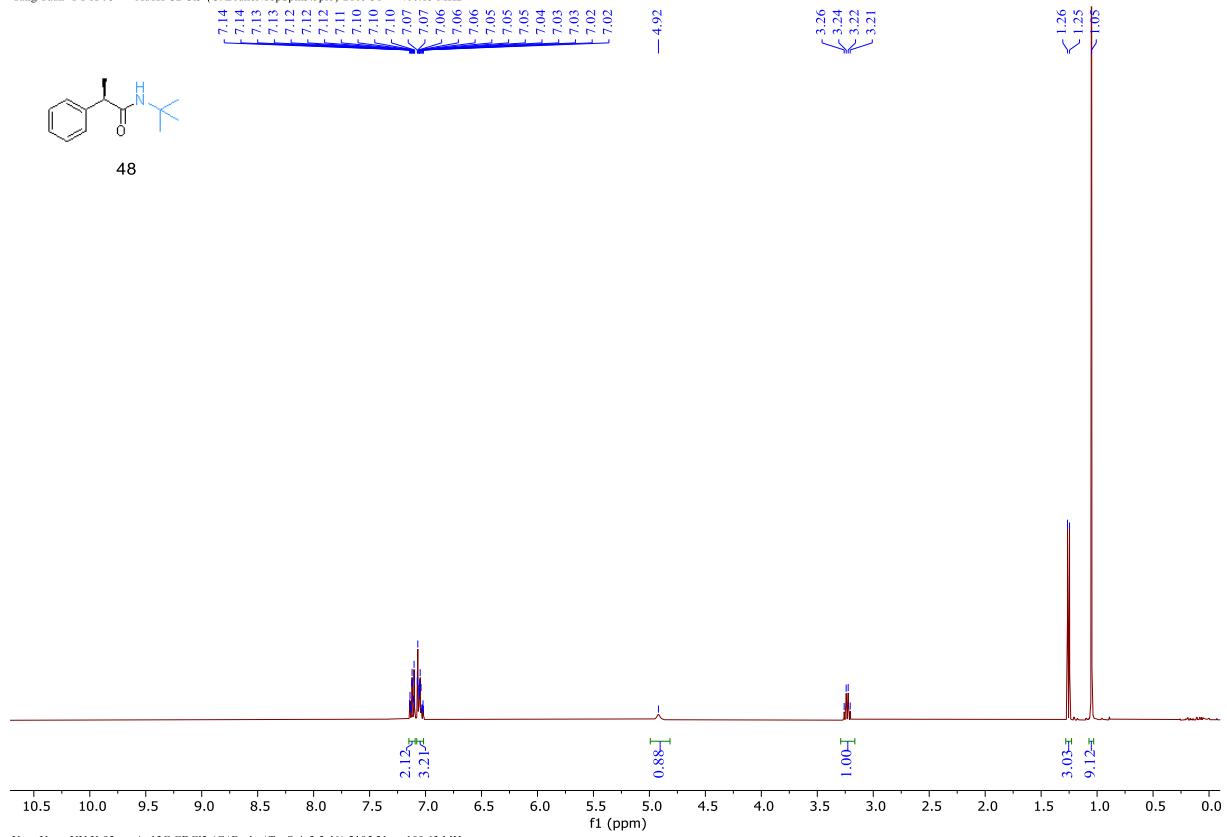
141.44
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140.79
128.88
127.52
127.20
126.74
124.73

50.56
—47.12
—40.11
39.97

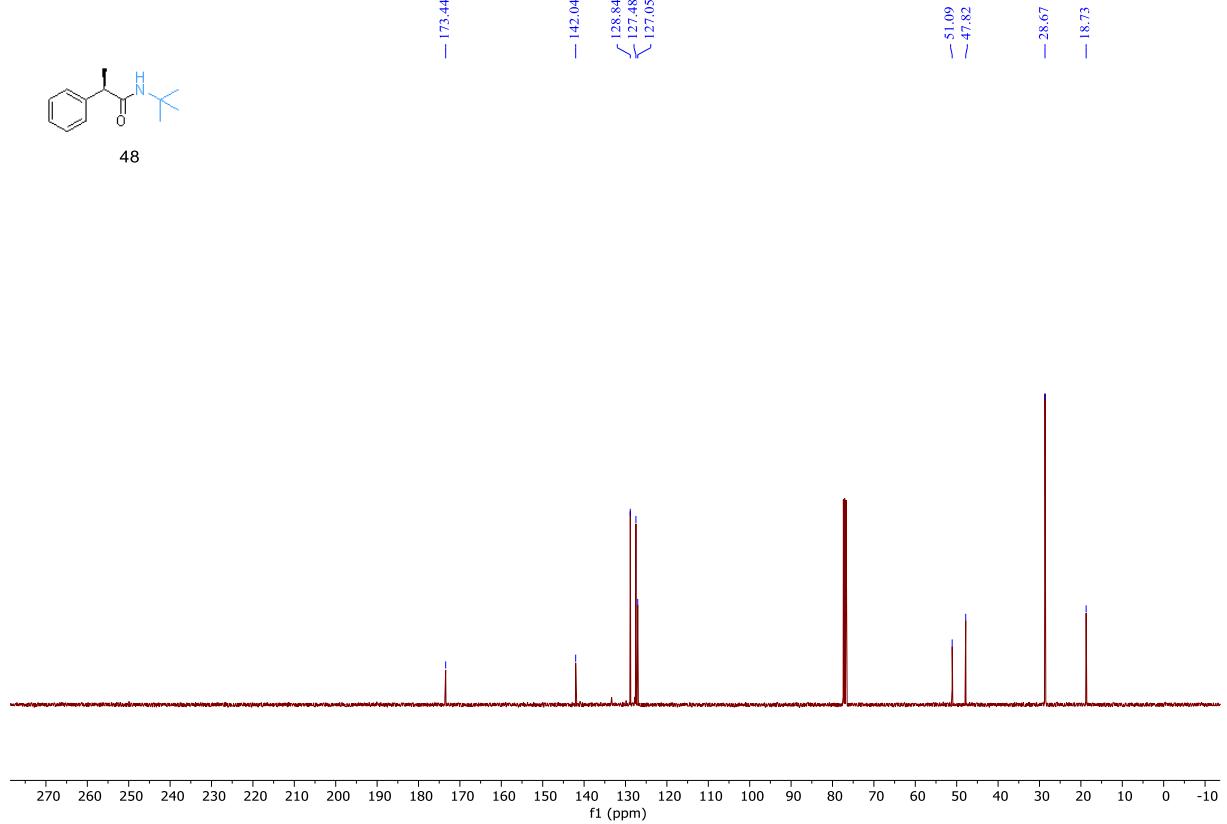
—18.75



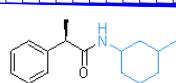
Yang Yuan YY-X-93 — Au1H CDCl₃ {C:\Bruker\TopSpin3.5pl6} 2105 31 — 400.13 MHz



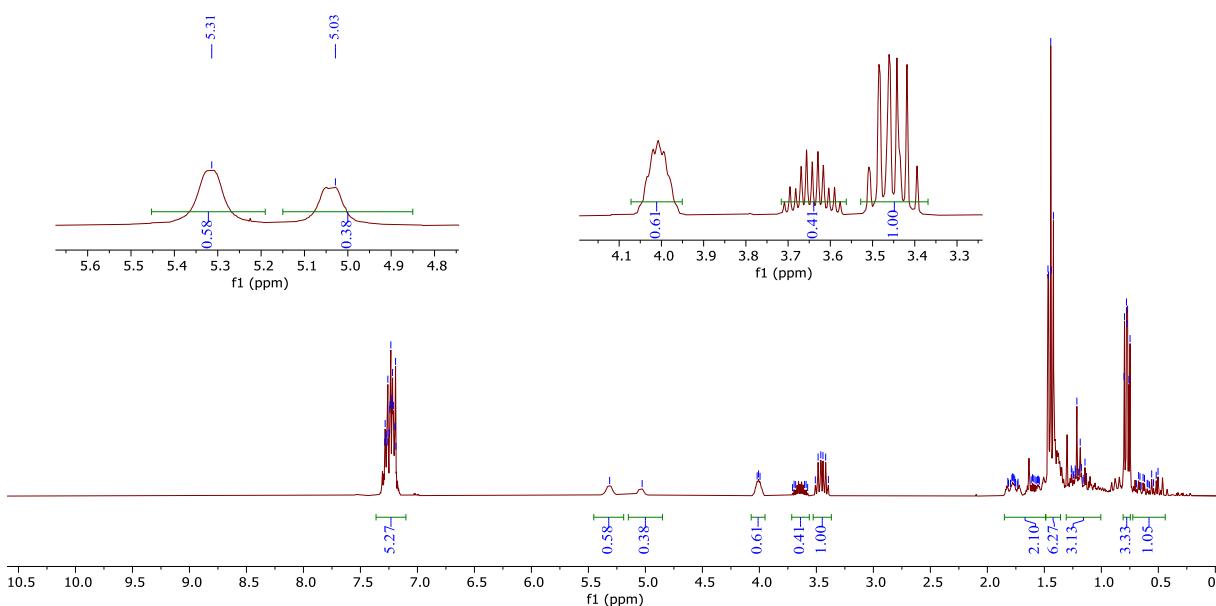
Yang Yuan YY-X-93 — Au13C CDCl₃ {C:\Bruker\TopSpin3.5pl6} 2105 31 — 100.63 MHz



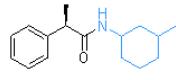
Yang Yuan, YY-X-113 — Au1H CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 27 — 300.13 MHz



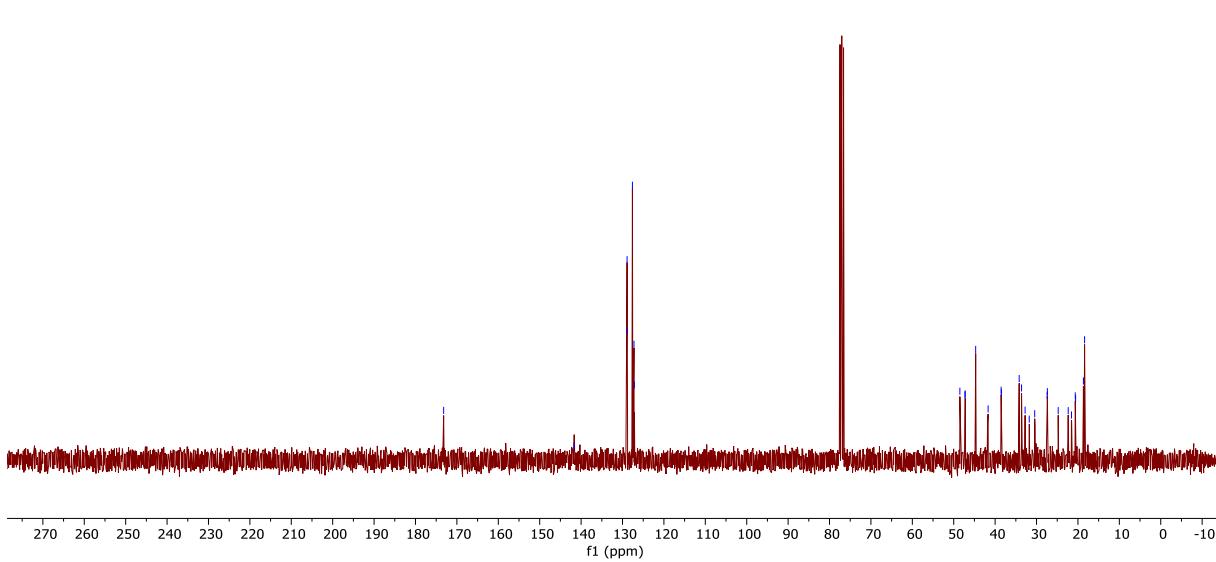
49
dr = 1.5:1

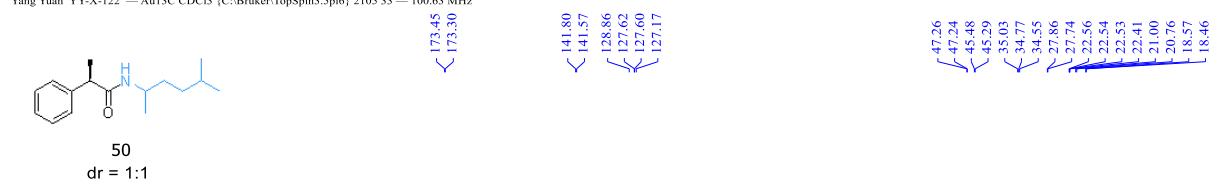
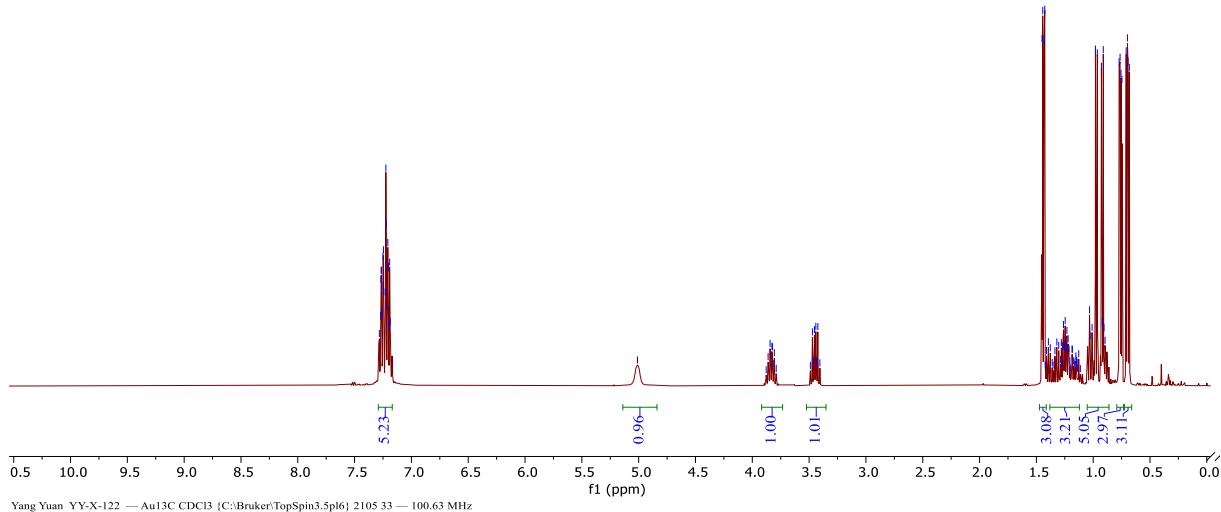
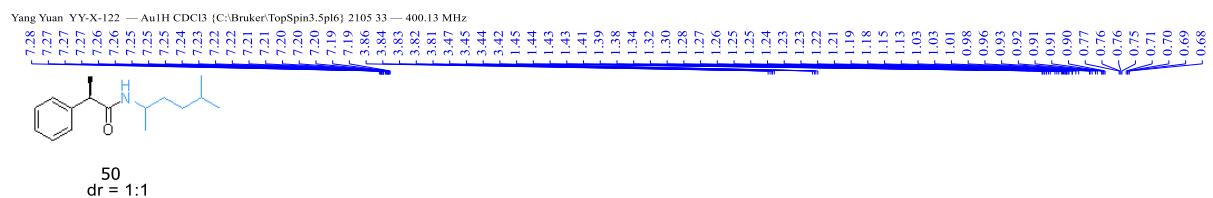


Yang Yuan, YY-X-113 — Au13C CDCl₃ {C:\Bruker\TopSpin3.6.2} 2103 27 — 75.48 MHz

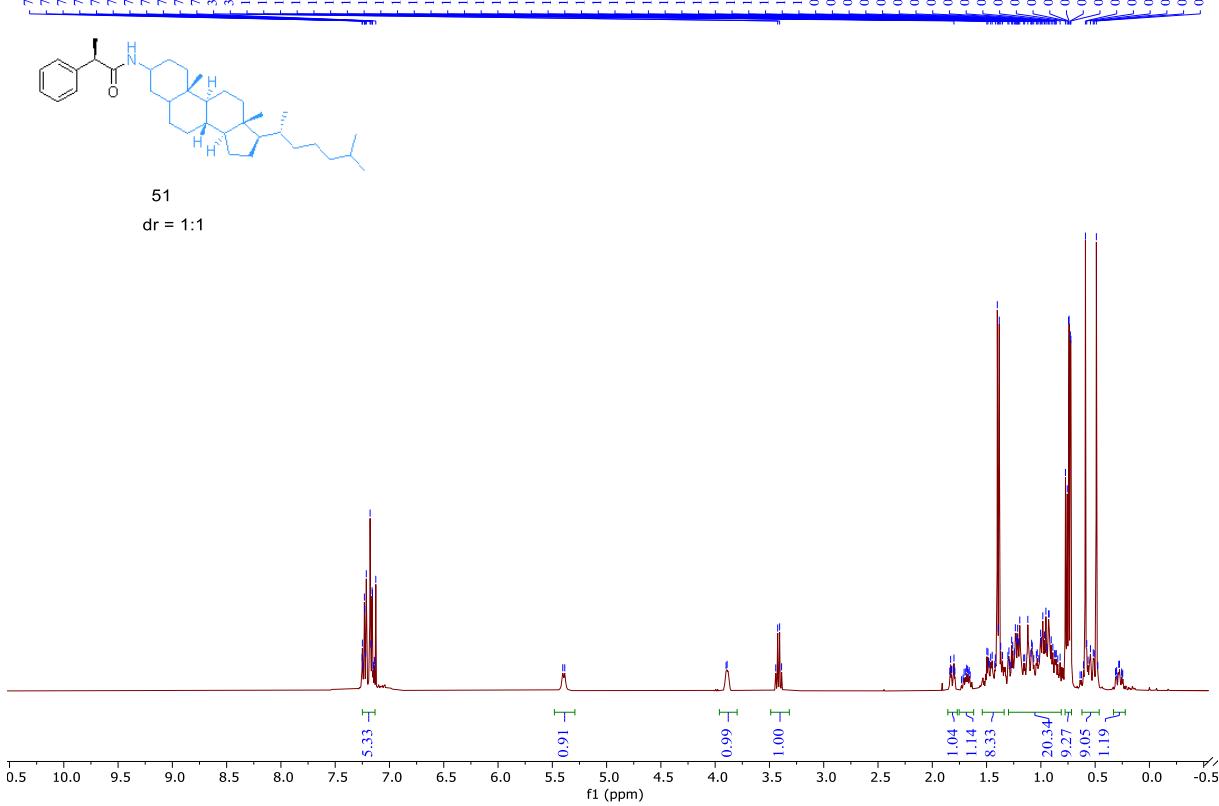


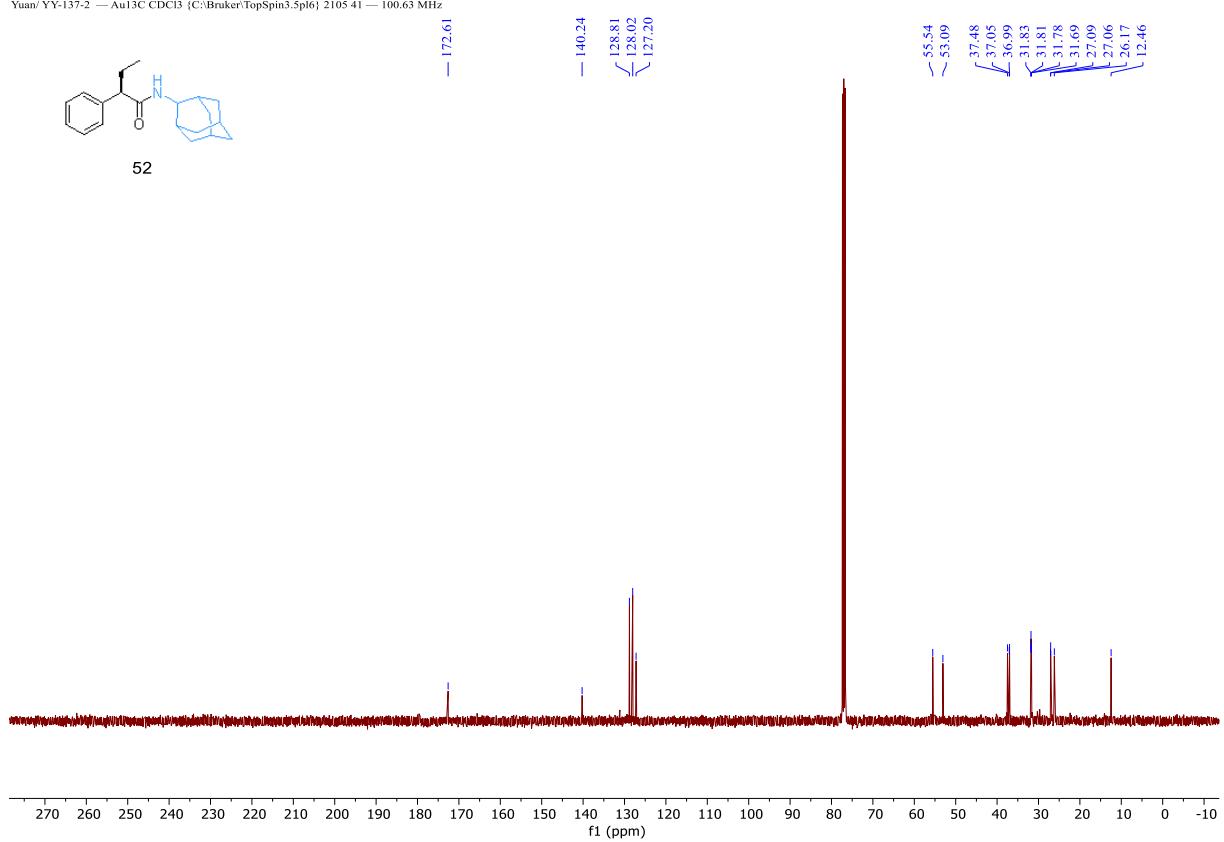
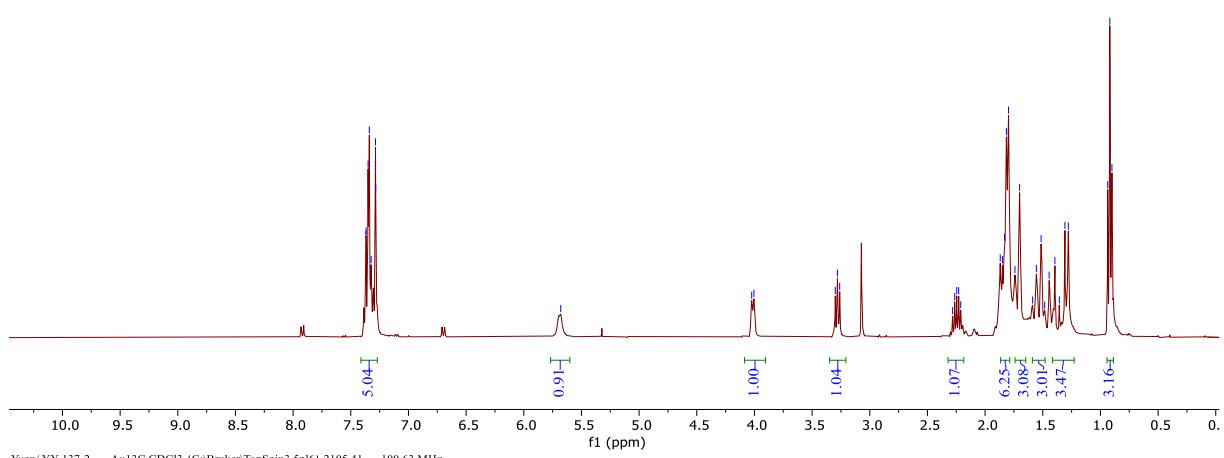
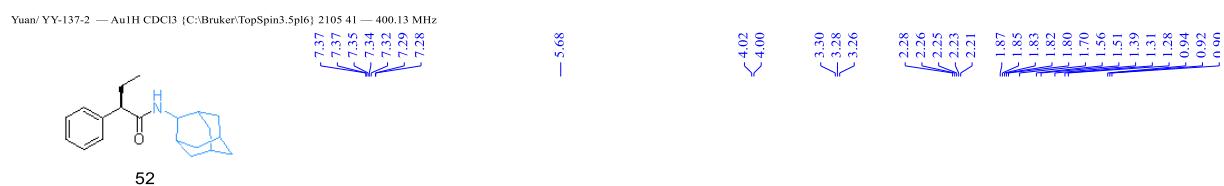
49
dr = 1.5:1

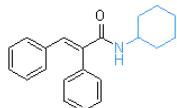
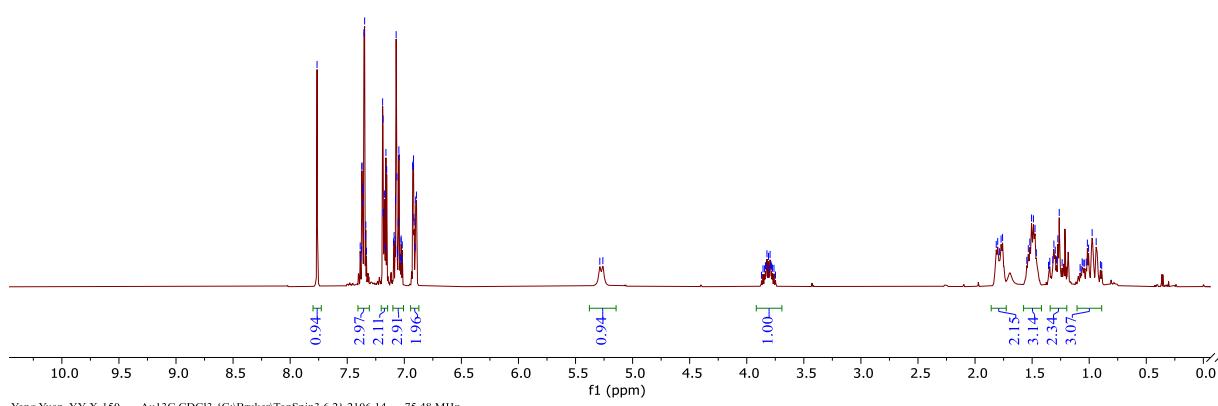
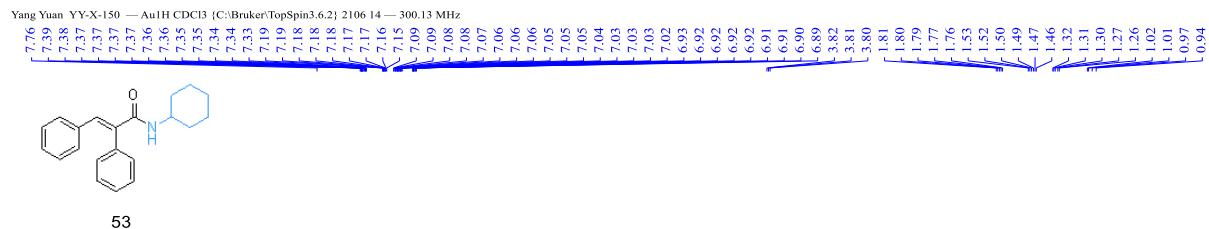




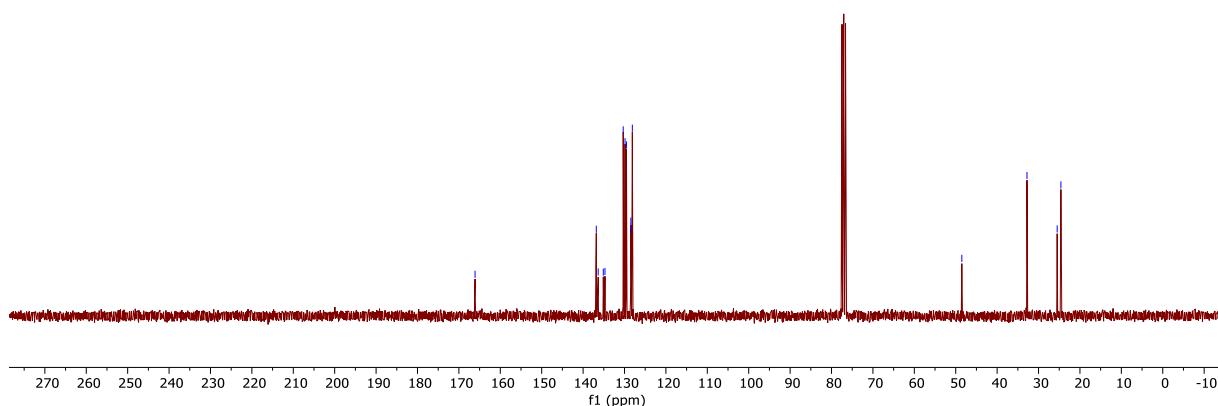
Yang Yuan YY-X-124-2 — Au1H CDCl₃ {C:\Bruker\TopSpin3.5\pl6} 2105.35 — 400.13 MHz

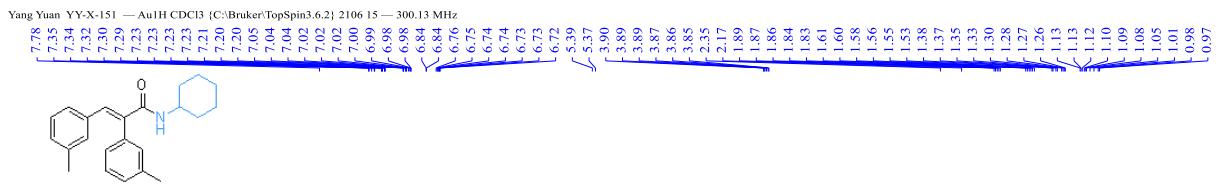




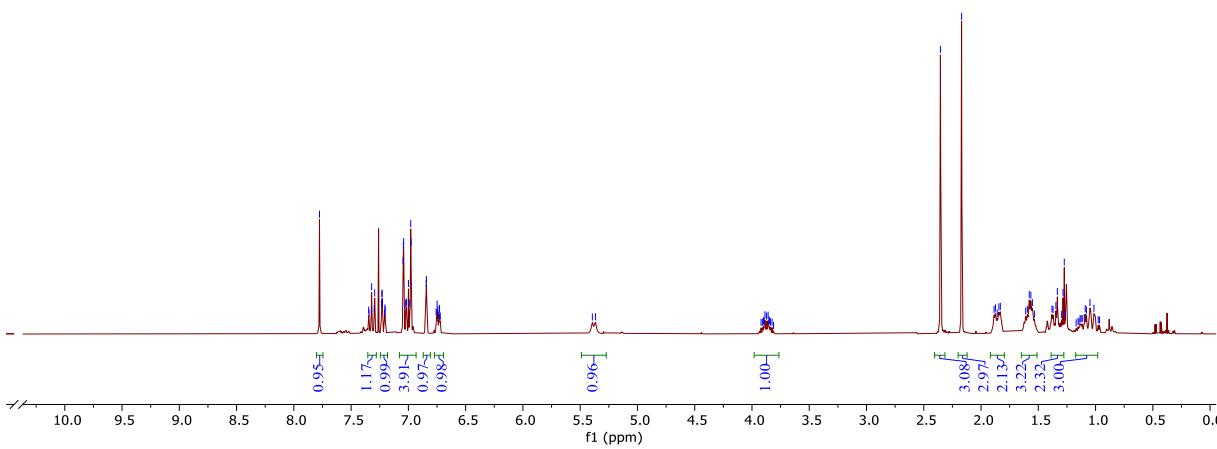


53

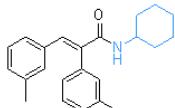




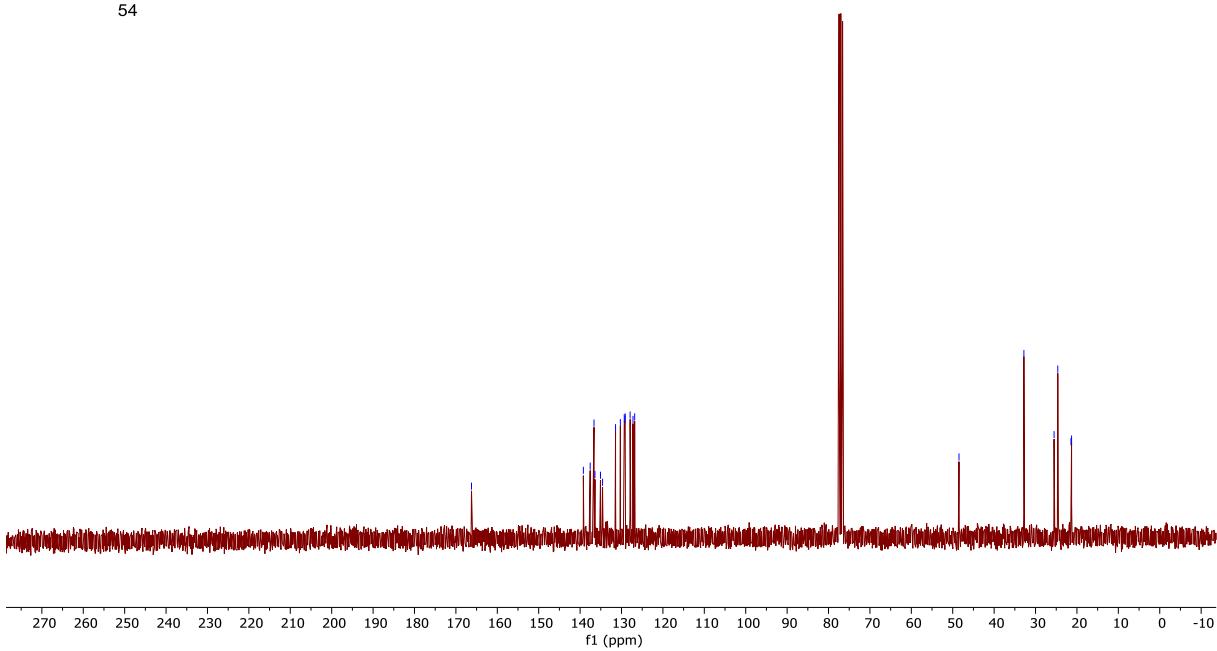
54

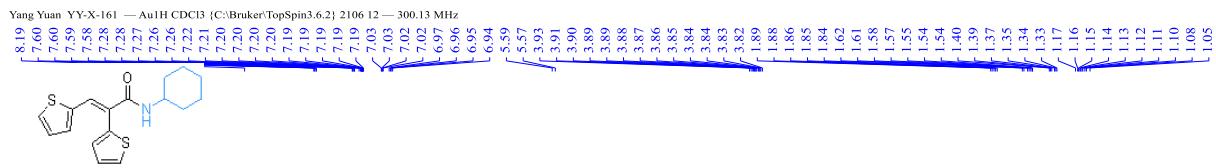


Yang Yuan YY-X-151 — Au13C CDCl₃ {C:\Bruker\TopSpin3.6.2\} 2106.15 — 75.48 MHz

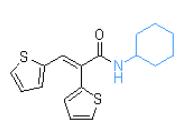
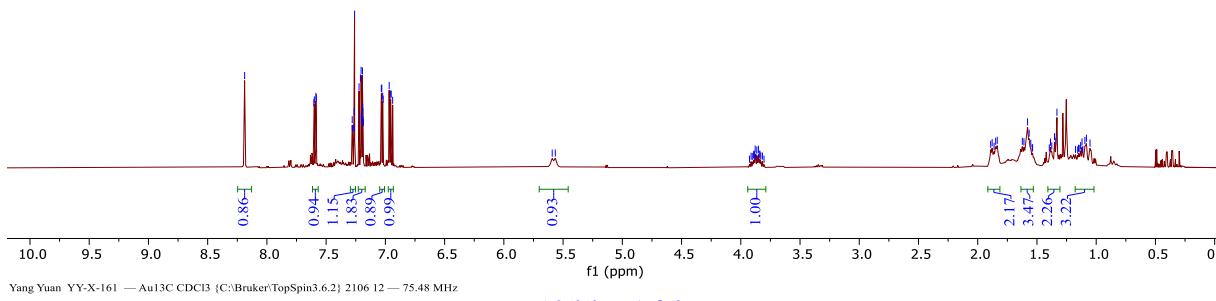


54

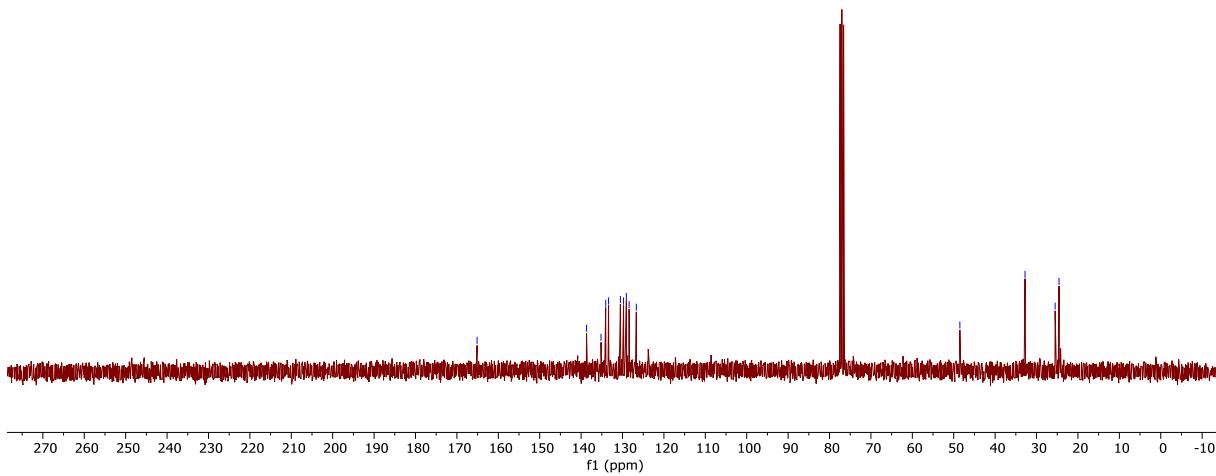




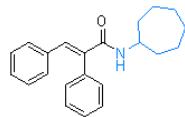
55



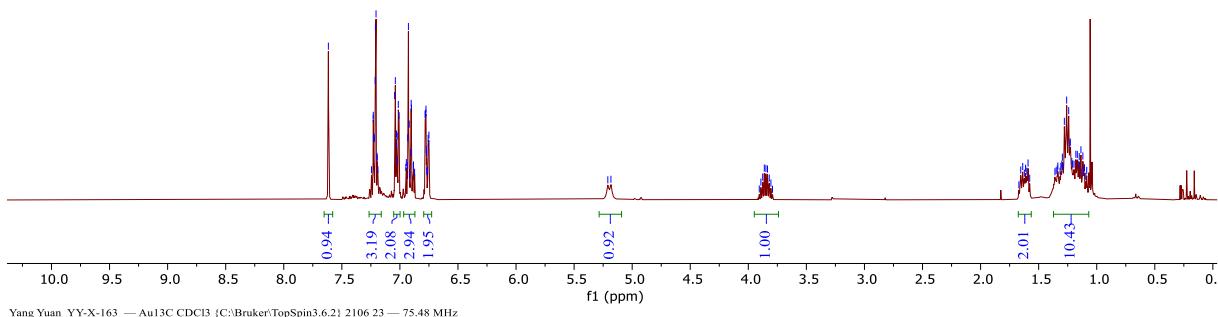
55



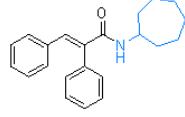
Yang Yuan YY-X-163 — Au1H CDCl₃ {C:\Bruker\TopSpin3.6.2} 2106 23 — 300.13 MHz



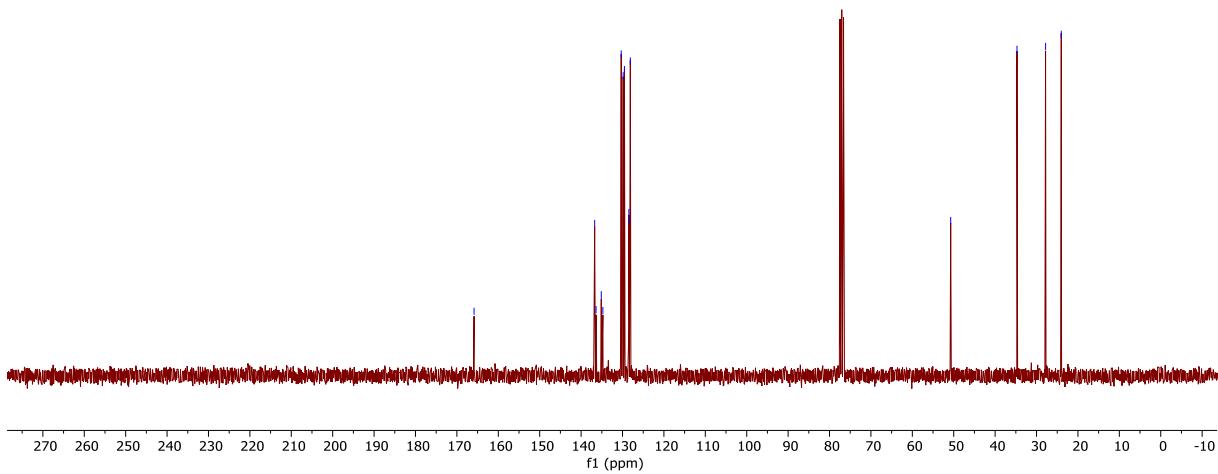
56



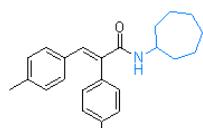
Yang Yuan YY-X-163 — Au13C CDCl₃ {C:\Bruker\TopSpin3.6.2} 2106 23 — 75.48 MHz



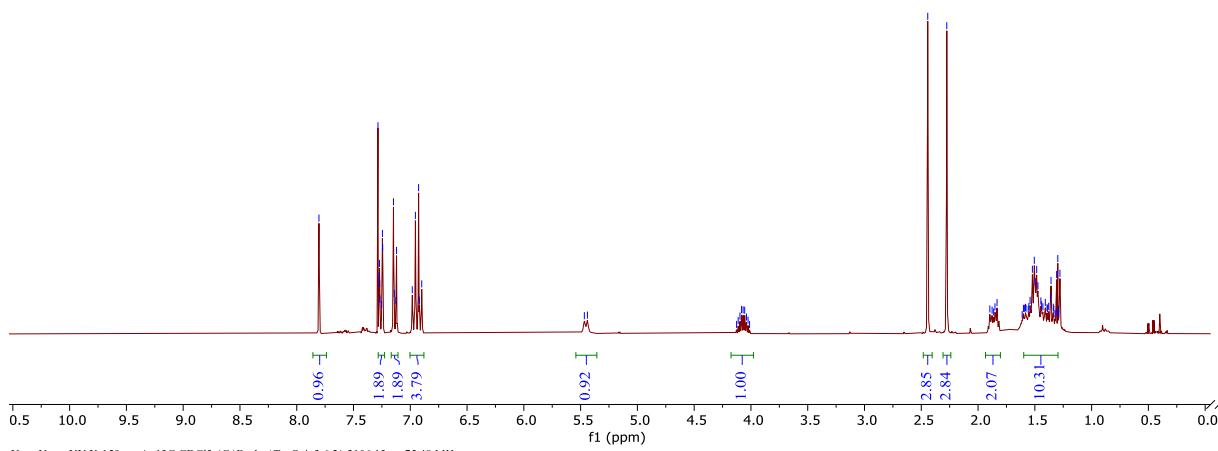
56



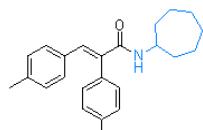
Yang Yuan YY-X-159 — Au1H CDCl₃ {C:\Bruker\TopSpin3.6.2} 2106 13 — 300.13 MHz



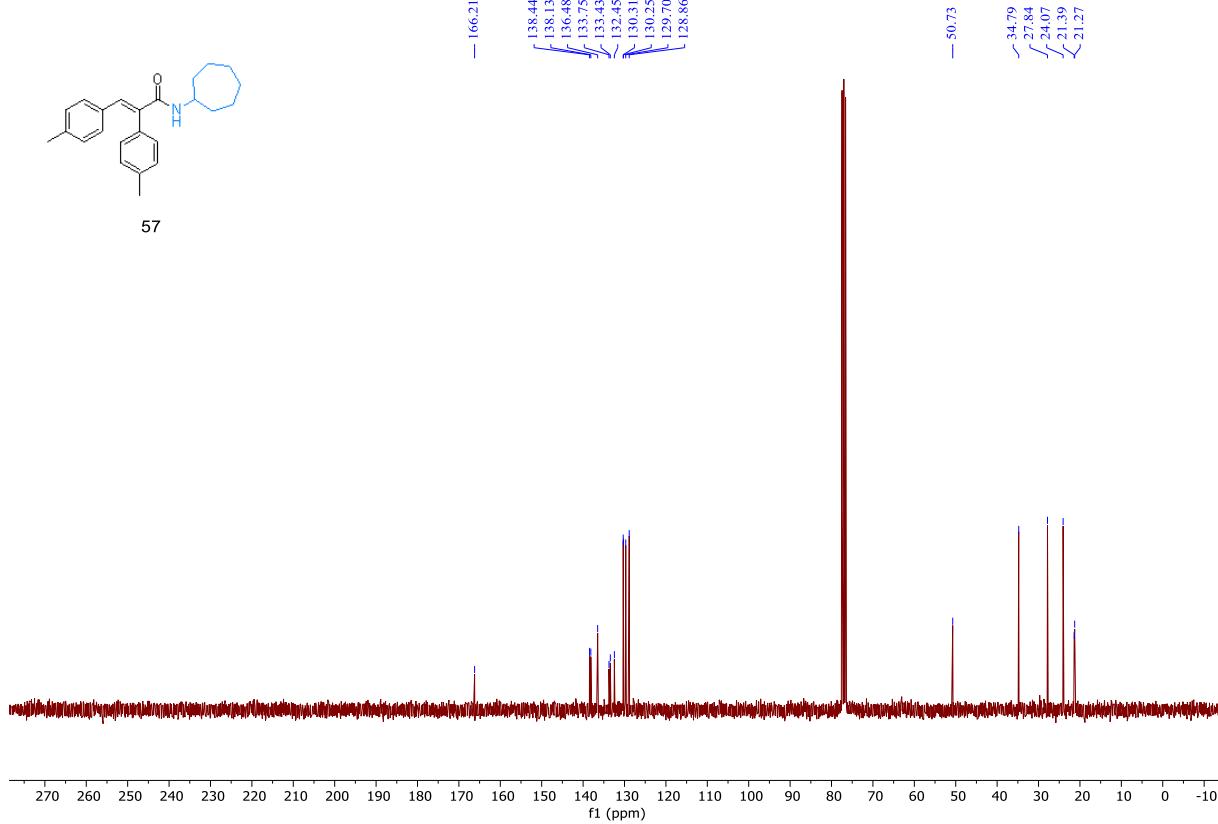
57

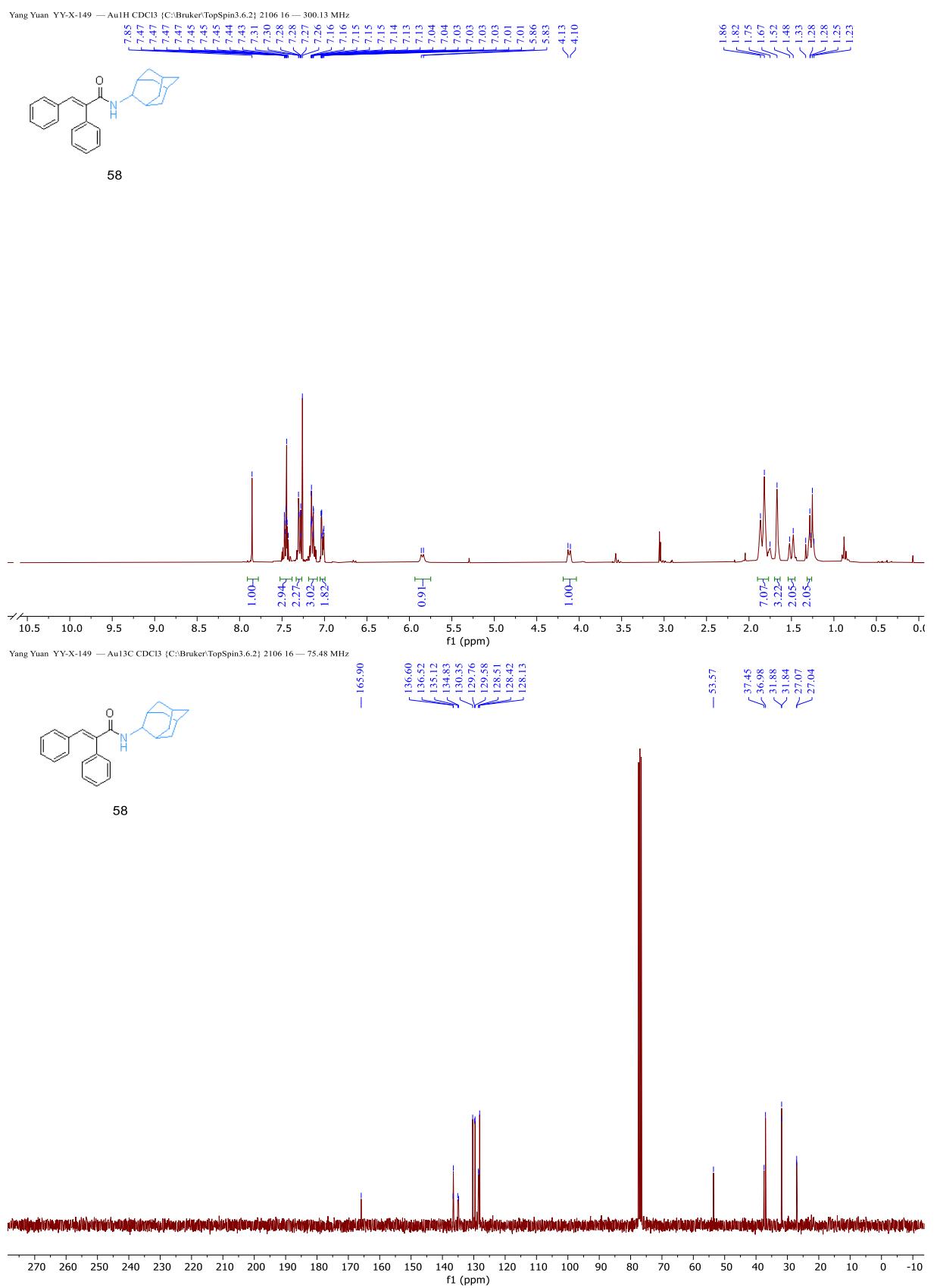


Yang Yuan YY-X-159 — Au13C CDCl₃ {C:\Bruker\TopSpin3.6.2} 2106 13 — 75.48 MHz

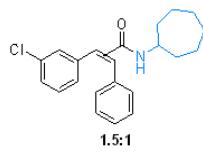


57

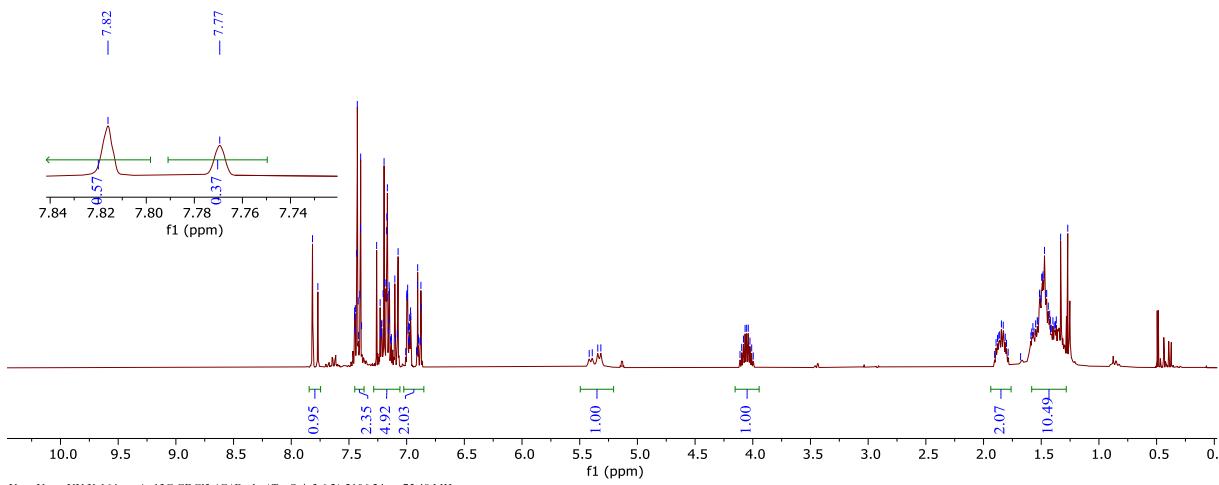




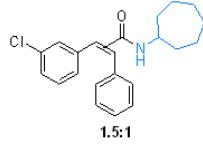
Yang Yuan YY-X-164 — Au1H CDCl₃ {C:\Bruker\TopSpin3.6.2} 2106 24 — 300.13 MHz



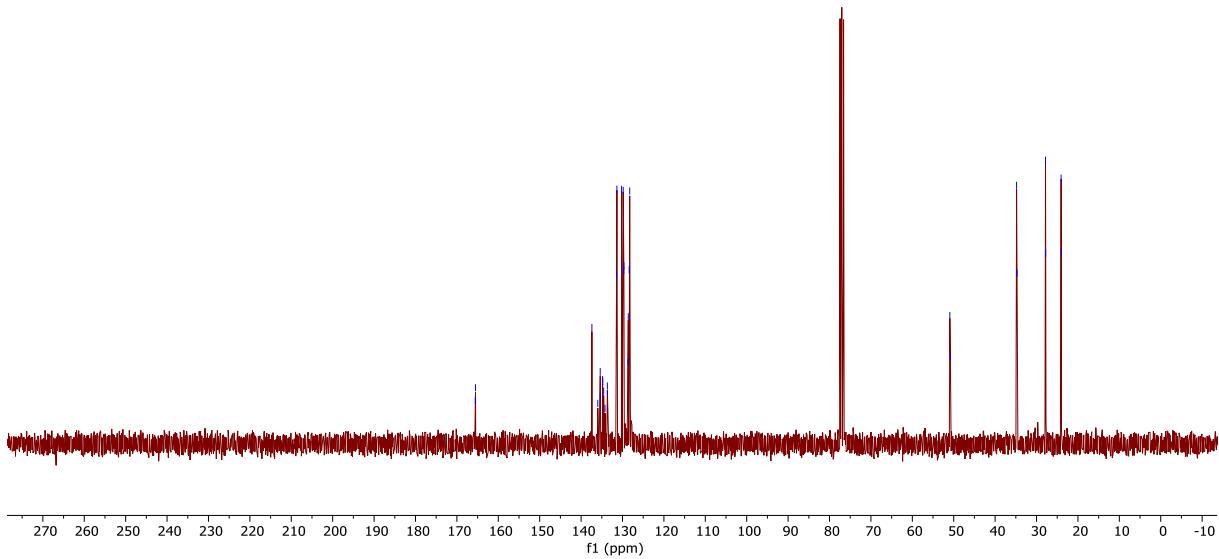
59



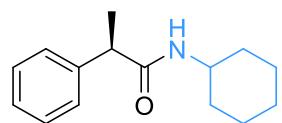
Yang Yuan YY-X-164 — Au13C CDCl₃ {C:\Bruker\TopSpin3.6.2} 2106 24 — 75.48 MHz



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7. HPLC Spectra



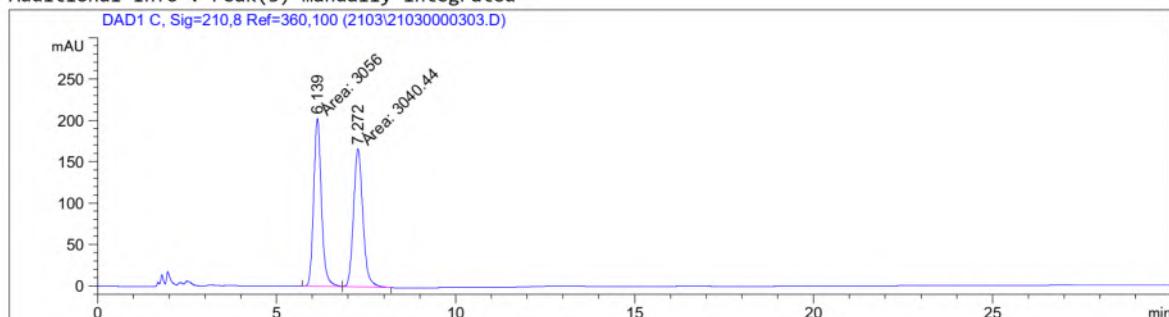
(R)-N-Cyclohexyl-2-phenylpropanamide (4)

Data File D:\CHEM32\1\DATA\2103\21030000303.D

Sample Name: YY-x-59 race

```
=====
Acq. Operator : Analytik           Seq. Line : 4
Acq. Instrument : LC5             Location : Vial 31
Injection Date : 3/3/2021 1:22:36 PM      Inj : 1
                                         Inj Volume : 0.5 µl
Different Inj Volume from Sequence !     Actual Inj Volume : 1.0 µl
Acq. Method : C:\CHEM32\1\METHODS\TEST RAC.M
Last changed : 3/3/2021 1:55:14 PM by Analytik
                         (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\TEST RAC.M
Last changed : 3/3/2021 2:51:01 PM by Analytik
                         (modified after loading)
Method Info : Cellulose4, Hept./EtOH 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

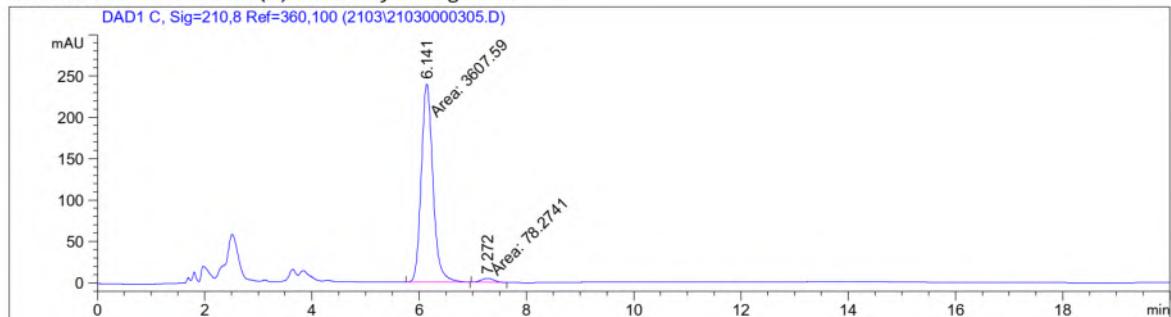
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.139	MM	0.2514	3055.99780	202.62683	50.1276
2	7.272	MM	0.3038	3040.44189	166.78253	49.8724

Data File D:\CHEM32\1\DATA\2103\21030000305.D
Sample Name: YY-x-59

```
=====
Acq. Operator   : Analytik          Seq. Line : 6
Acq. Instrument : LC5             Location  : Vial 32
Injection Date  : 3/3/2021 2:16:43 PM    Inj       : 1
                                         Inj Volume : 0.5 µl
Different Inj Volume from Sequence !    Actual Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\TEST RAC.M
Last changed    : 3/3/2021 1:55:14 PM by Analytik
                                         (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\TEST RAC.M
Last changed    : 3/3/2021 2:53:32 PM by Analytik
                                         (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated

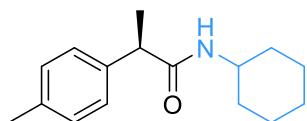


```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

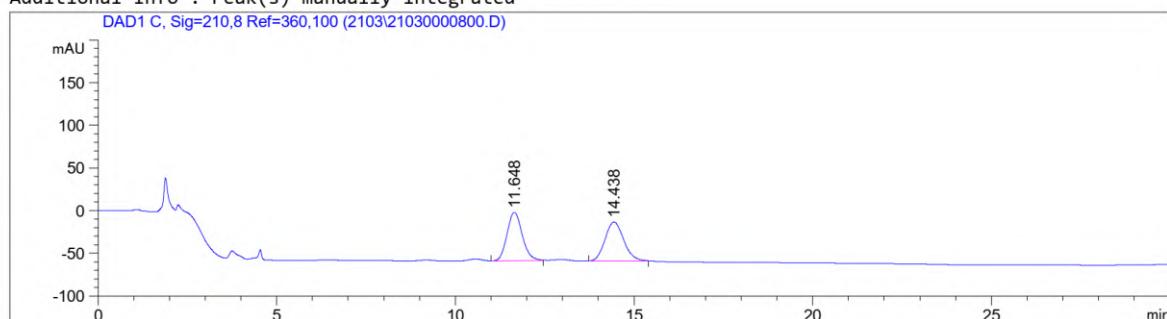
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.141	MM	0.2512	3607.59473	239.39899	97.8764
2	7.272	MM	0.2740	78.27406	4.76138	2.1236



Data File D:\CHEM32\1\DATA\2103\21030000800.D
Sample Name: YY-X-65-Race

```
=====
Acq. Operator   : Analytik          Seq. Line : 1
Acq. Instrument : LC5             Location  : Vial 31
Injection Date  : 3/8/2021 3:18:21 PM    Inj       : 1
                                         Inj Volume : 0.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/8/2021 3:18:04 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/9/2021 11:26:21 AM by Analytik
                                         (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 98:2, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

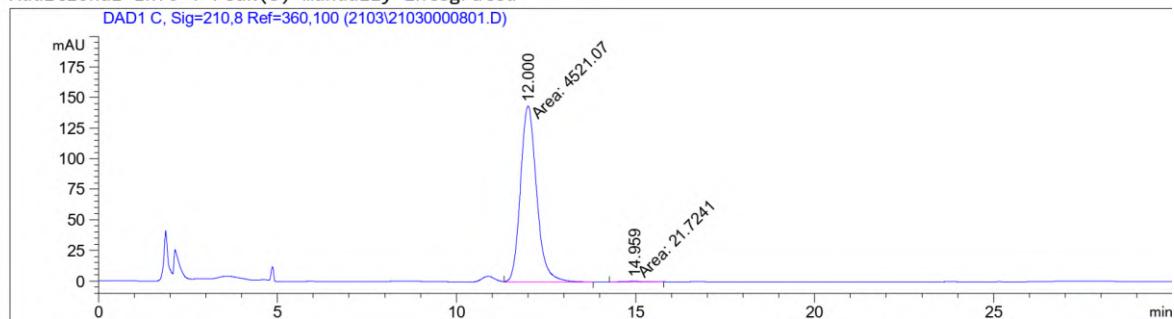
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.648	VB	0.4575	1665.84277	56.57780	49.8275
2	14.438	BB	0.5678	1677.37830	45.74106	50.1725
Totals :						3343.22107 102.31886

Data File D:\CHEM32\1\DATA\2103\21030000801.D
Sample Name: YY-X-65-C

```
=====
Acq. Operator : Analytik          Seq. Line : 2
Acq. Instrument : LC5           Location : Vial 41
Injection Date : 3/8/2021 3:59:27 PM      Inj : 1
                                         Inj Volume : 0.5 µl
Different Inj Volume from Sequence !    Actual Inj Volume : 2.0 µl
Acq. Method : C:\CHEM32\1\METHODS\YY.M
Last changed : 3/8/2021 3:18:04 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed : 3/9/2021 11:27:44 AM by Analytik
                                         (modified after loading)
Method Info : Cellulose4, Hept./EtOH 98:2, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

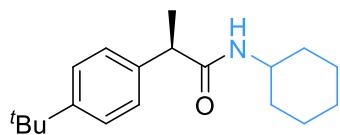
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area %
#	[min]		[min]	[mAU*s]	[mAU]	%
1	12.000	FM	0.5233	4521.06836	143.98621	99.5218
2	14.959	MM	0.6906	21.72414	5.24294e-1	0.4782

Totals : 4542.79250 144.51050

LC5 3/9/2021 11:28:45 AM Analytik

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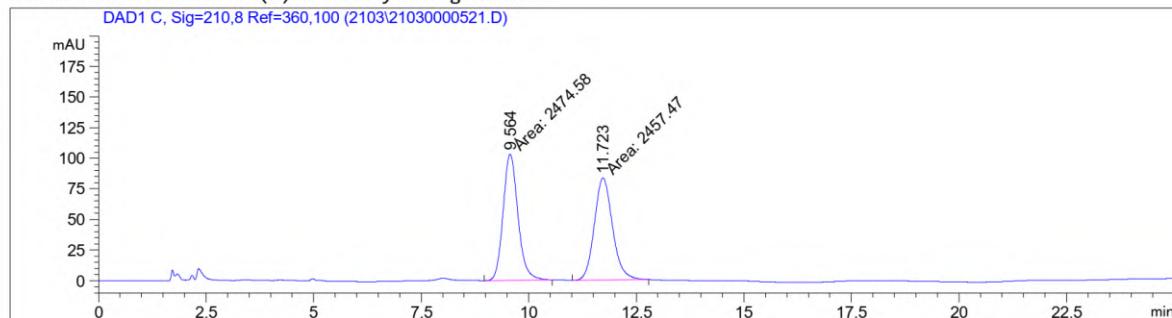
(R)-2-(4-(tert-Butyl)phenyl)-N-cyclohexylpropanamide (6)

Data File D:\CHEM32\1\DATA\2103\21030000521.D

Sample Name: YY-X-67-Race

```
=====
Acq. Operator   : Analytik           Seq. Line : 5
Acq. Instrument : LC5             Location  : Vial 33
Injection Date  : 3/5/2021 9:32:12 PM      Inj       : 1
                                                Inj Volume : 0.5 μl
Acq. Method    : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/5/2021 4:03:06 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/8/2021 12:36:39 PM by Analytik
                                                (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 98:2, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.564	MM	0.3996	2474.57935	103.22060	50.1734
2	11.723	MM	0.4906	2457.47388	83.47901	49.8266

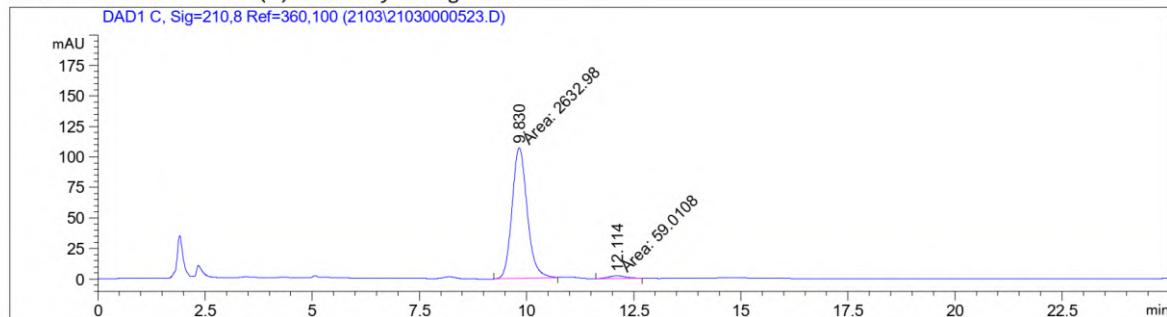
Totals : 4932.05322 186.69961

Data File D:\CHEM32\1\DATA\2103\21030000523.D

Sample Name: YY-X-67-C

```
=====
Acq. Operator   : Analytik          Seq. Line :   6
Acq. Instrument : LC5             Location  : Vial 43
Injection Date  : 3/5/2021 10:54:26 PM      Inj :   1
                                         Inj Volume : 0.5 µl
Acq. Method    : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/5/2021 4:03:06 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/8/2021 12:36:39 PM by Analytik
                                         (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 98:2, 1ml/min
```

Additional Info : Peak(s) manually integrated



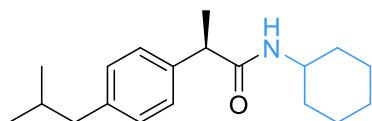
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.830	MM	0.4098	2632.97778	107.08657	97.8079
2	12.114	MM	0.4480	59.01076	2.19550	2.1921

Totals : 2691.98854 109.28207



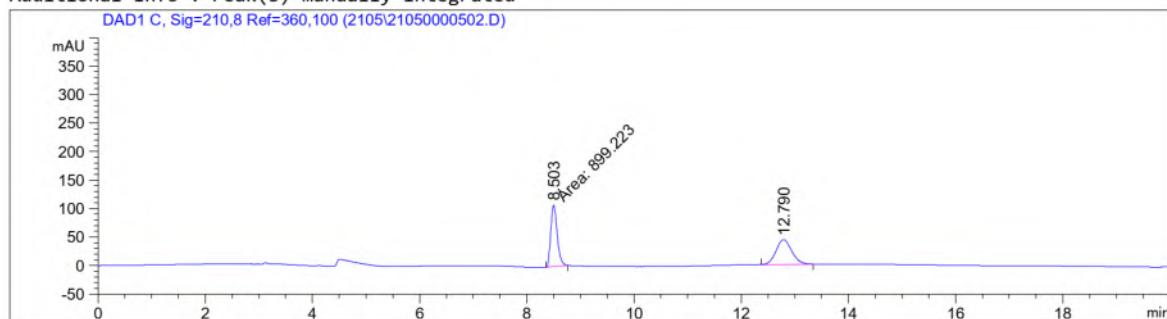
(R)-N-Cyclohexyl-2-(4-isobutylphenyl)propanamide (7)

Data File D:\CHEM32\1\DATA\2105\21050000502.D

Sample Name: YY-X-125 Race

```
=====
Acq. Operator   : Analytik           Seq. Line :  3
Acq. Instrument : LCS              Location : Vial 9
Injection Date  : 5/5/2021 11:16:14 AM    Inj :  1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/5/2021 11:15:50 AM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/5/2021 2:13:26 PM by Analytik
                  (modified after loading)
Method Info     : Chiralcel OJ, Hept./EtOH 99:1, 1.0ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.503	MM	0.1383	899.22321	108.37066	50.2277
2	12.790	BB	0.3163	891.06921	43.61647	49.7723

Totals : 1790.29242 151.98713

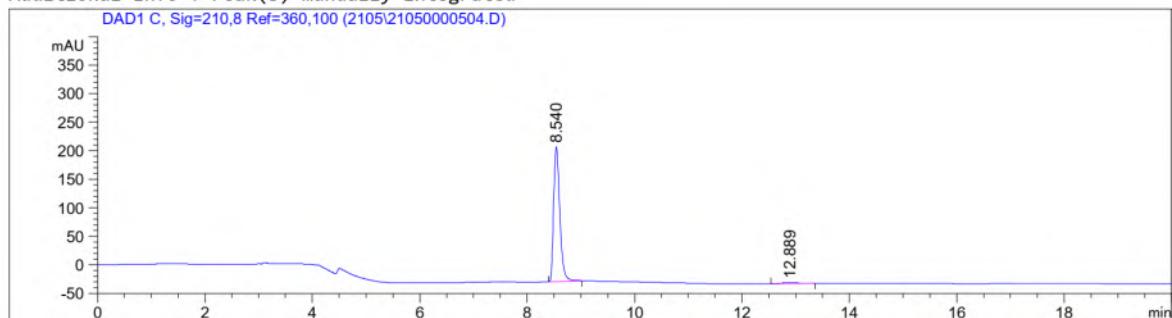
LC5 5/5/2021 2:15:06 PM Analytik

Page 1 of 2

Data File D:\CHEM32\1\DATA\2105\21050000504.D
Sample Name: YY-X-125 -C

```
=====
Acq. Operator   : Analytik          Seq. Line :  5
Acq. Instrument : LC5             Location  : Vial 8
Injection Date  : 5/5/2021 12:38:43 PM    Inj :  1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/5/2021 10:03:49 AM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/5/2021 2:13:26 PM by Analytik
                                         (modified after loading)
Method Info     : Chiralcel OJ, Hept./EtOH 99:1, 1.0ml/min
```

Additional Info : Peak(s) manually integrated



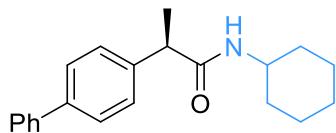
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.540	BB	0.1236	1893.06592	235.96875	98.2875
2	12.889	BB	0.3045	32.98305	1.62707	1.7125

Totals : 1926.04897 237.59582



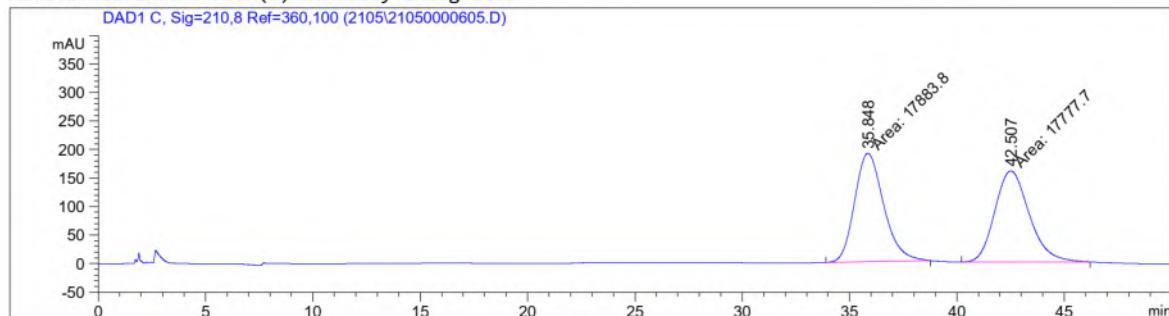
(R)-2-((1,1'-Biphenyl)-4-yl)-N-cyclohexylpropanamide (8)

Data File D:\CHEM32\1\DATA\2105\21050000605.D

Sample Name: YY-X-103 Race

```
=====
Acq. Operator   : Analytik          Seq. Line :  5
Acq. Instrument : LC5             Location  : Vial 21
Injection Date  : 5/6/2021 4:06:09 PM      Inj       : 1
                                                Inj Volume : 1.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/6/2021 1:51:06 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/7/2021 10:54:22 AM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 99:1, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	35.848	MM	1.5683	1.78838e4	190.05756	50.1488
2	42.507	MM	1.8571	1.77777e4	159.55045	49.8512

Totals : 3.56615e4 349.60800

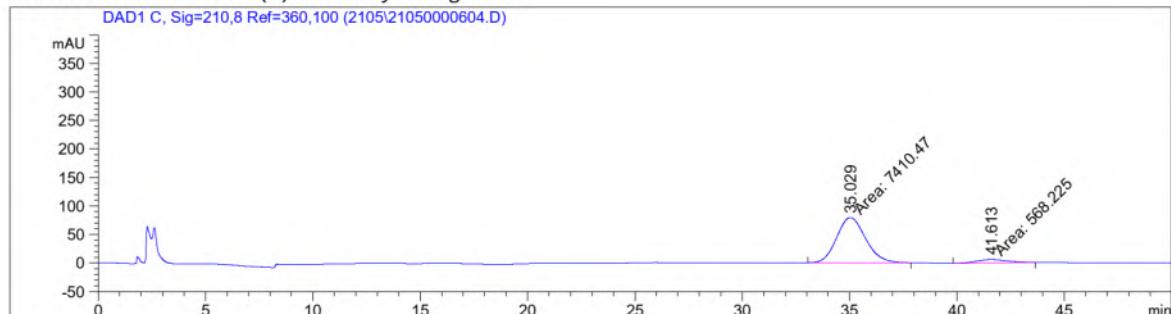
LC5 5/7/2021 11:03:46 AM Analytik

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Data File D:\CHEM32\1\DATA\2105\21050000604.D
Sample Name: YY-X-103 rt

```
=====
Acq. Operator   : Analytik          Seq. Line : 4
Acq. Instrument : LC5             Location  : Vial 22
Injection Date  : 5/6/2021 3:10:00 PM    Inj       : 1
                                         Inj Volume : 1.0 µl
Different Inj Volume from Sequence !    Actual Inj Volume : 5.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/6/2021 1:51:06 PM by Analytik
                                         (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/7/2021 10:54:22 AM by Analytik
                                         (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 99:1, 1ml/min
```

Additional Info : Peak(s) manually integrated

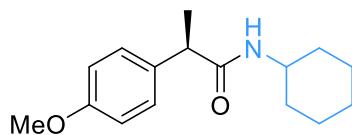


```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	35.029	MM	1.5630	7410.47363	79.01877	92.8782
2	41.613	MM	1.1961	568.22485	5.72760	7.1218



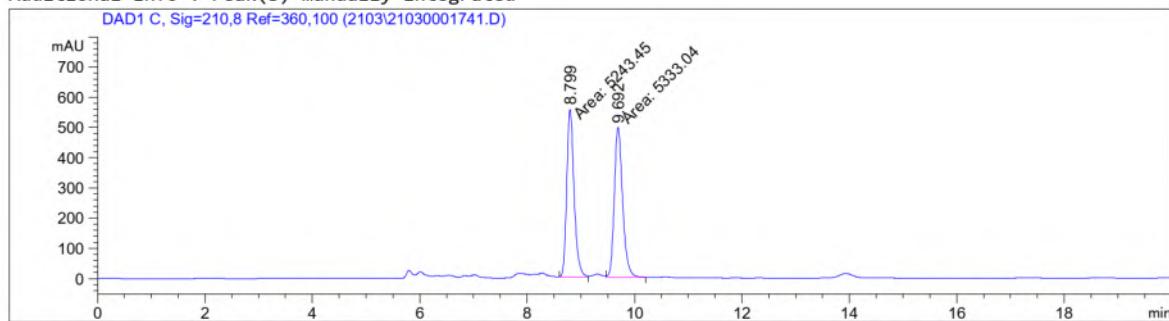
(R)-N-Cyclohexyl-2-(4-methoxyphenyl)propanamide (9)

Data File D:\CHEM32\1\DATA\2103\21030001741.D

Sample Name: YY-X-68 Race

```
=====
Acq. Operator   : Analytik           Seq. Line : 18
Acq. Instrument : LC5             Location  : Vial 34
Injection Date  : 3/18/2021 10:29:55 AM    Inj       : 1
                                                Inj Volume : 0.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY-72.M
Last changed    : 3/17/2021 2:33:11 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY-72.M
Last changed    : 3/18/2021 11:12:55 AM by Analytik
                                                (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.799	MF	0.1577	5243.45264	554.05408	49.5765
2	9.692	FM	0.1792	5333.03906	496.12857	50.4235

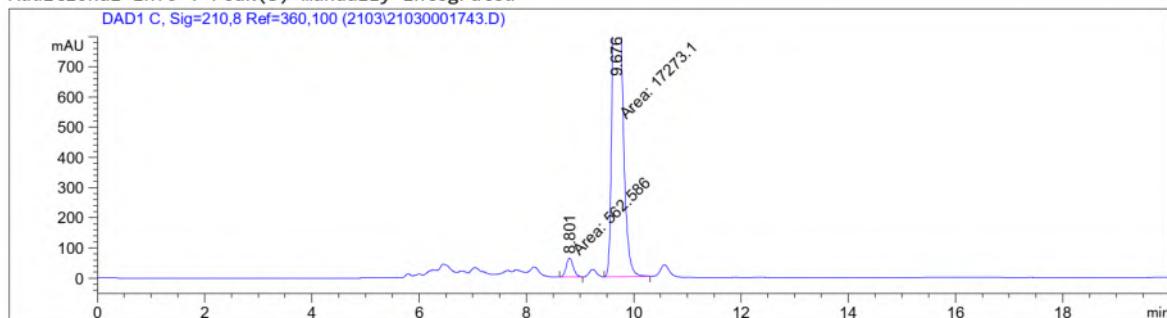
Totals : 1.05765e4 1050.18265

Data File D:\CHEM32\1\DATA\2103\21030001743.D

Sample Name: YY-X-68 C

```
=====
Acq. Operator   : Analytik          Seq. Line : 19
Acq. Instrument : LC5             Location  : Vial 44
Injection Date  : 3/18/2021 11:12:07 AM    Inj       : 2
                                                Inj Volume : 0.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY-72.M
Last changed    : 3/17/2021 2:33:11 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY-72.M
Last changed    : 3/18/2021 11:12:55 AM by Analytik
                                                (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



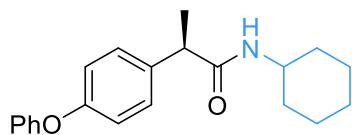
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.801	MM	0.1498	562.58582	62.57995	3.1543
2	9.676	MM	0.1842	1.72731e4	1563.02771	96.8457

Totals : 1.78357e4 1625.60766



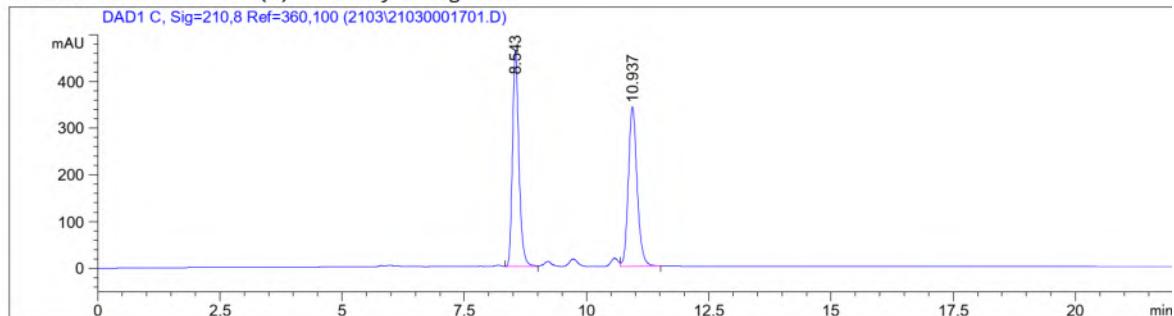
(R)-N-Cyclohexyl-2-(4-phenoxyphenyl)propanamide (10)

Data File D:\CHEM32\1\DATA\2103\21030001701.D

Sample Name: YY-X-72 Race

```
=====
Acq. Operator   : Analytik                     Seq. Line :  1
Acq. Instrument : LC5                         Location  : Vial 31
Injection Date  : 3/17/2021 11:42:23 AM          Inj       : 1
                                                Inj Volume : 0.5 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 0.1 µl
Acq. Method    : C:\CHEM32\1\METHODS\YY-72.M
Last changed   : 3/17/2021 12:16:02 PM by Analytik
                           (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY-72.M
Last changed   : 3/17/2021 12:41:09 PM by Analytik
                           (modified after loading)
Method Info    : Chiraldak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

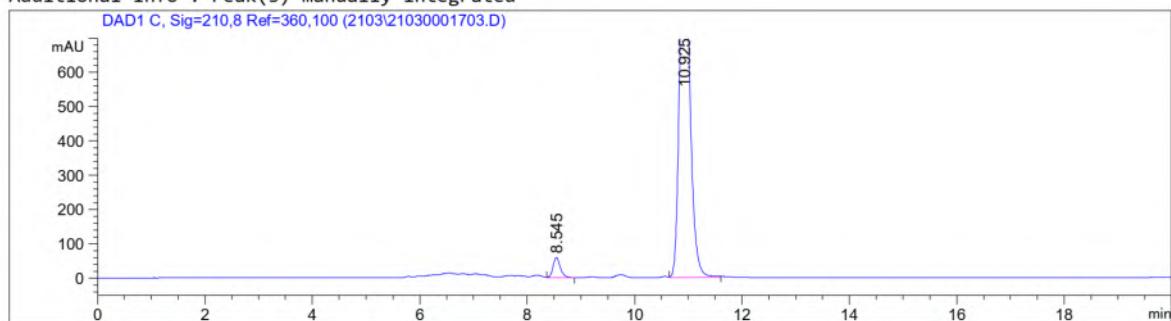
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.543	VV	0.1407	4295.44385	461.12308	49.9743
2	10.937	VB	0.1931	4299.86230	341.02795	50.0257

Data File D:\CHEM32\1\DATA\2103\21030001703.D

Sample Name: YY-X-72 C

```
=====
Acq. Operator   : Analytik          Seq. Line : 2
Acq. Instrument : LC5             Location  : Vial 41
Injection Date  : 3/17/2021 12:37:30 PM    Inj : 2
                                                Inj Volume : 0.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY-72.M
Last changed    : 3/17/2021 12:16:02 PM by Analytik
                    (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY-72.M
Last changed    : 3/17/2021 2:27:03 PM by Analytik
                    (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

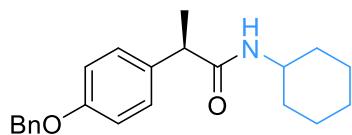
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area %
#	[min]		[min]	[mAU*s]	[mAU]	
1	8.545	VB	0.1442	573.69867	59.66669	3.8028
2	10.925	VB	0.1980	1.45124e4	1128.56689	96.1972

Totals : 1.50861e4 1188.23359

LC5 3/17/2021 2:27:13 PM Analytik

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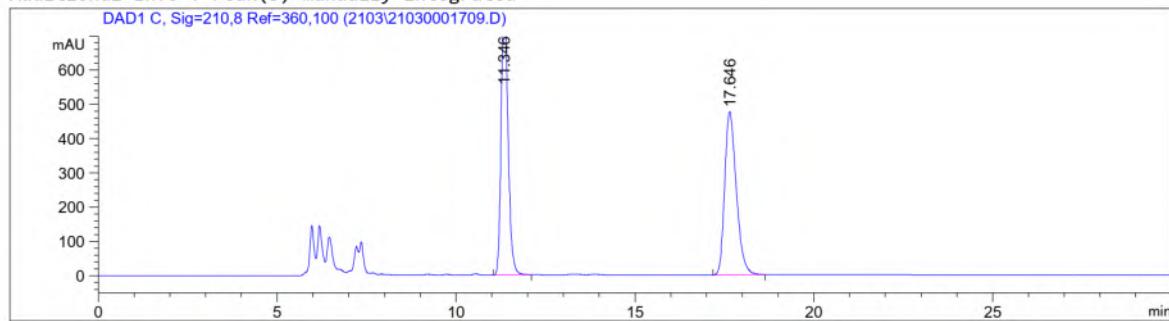
(R)-2-(4-(Benzyl)phenyl)-N-cyclohexylpropanamide (11)

Data File D:\CHEM32\1\DATA\2103\21030001709.D

Sample Name: YY-X-82-Race

```
=====
Acq. Operator   : Analytik          Seq. Line :  1
Acq. Instrument : LC5             Location  : Vial 33
Injection Date  : 3/17/2021 2:54:40 PM      Inj       : 2
                                                Inj Volume : 0.5 µl
Acq. Method    : C:\CHEM32\1\METHODS\YY-72.M
Last changed    : 3/17/2021 3:03:43 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY-72.M
Last changed    : 3/17/2021 3:48:02 PM by Analytik
                  (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.346	BB	0.2025	1.06880e4	807.14893	49.7251
2	17.646	BB	0.3458	1.08062e4	477.85327	50.2749

Totals : 2.14942e4 1285.00220

LC5 3/17/2021 3:48:50 PM Analytik

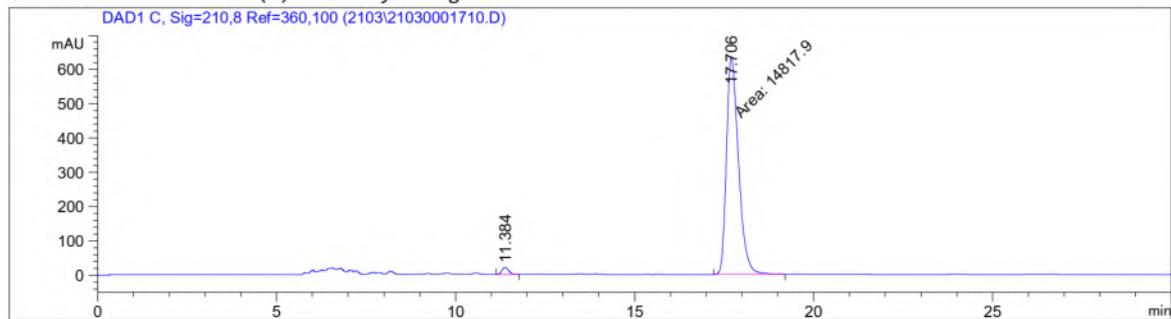
Page 1 of 2

Data File D:\CHEM32\1\DATA\2103\21030001710.D

Sample Name: YY-X-82-C

```
=====
Acq. Operator   : Analytik          Seq. Line : 2
Acq. Instrument : LC5             Location  : Vial 43
Injection Date  : 3/17/2021 3:35:48 PM    Inj : 1
                                         Inj Volume : 0.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY-72.M
Last changed    : 3/17/2021 3:03:43 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY-72.M
Last changed    : 3/17/2021 3:48:02 PM by Analytik
                  (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



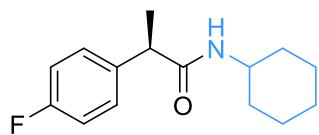
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.384	BB	0.2019	272.90213	20.68949	1.8084
2	17.706	MM	0.3878	1.48179e4	636.76093	98.1916

Totals : 1.50908e4 657.45041

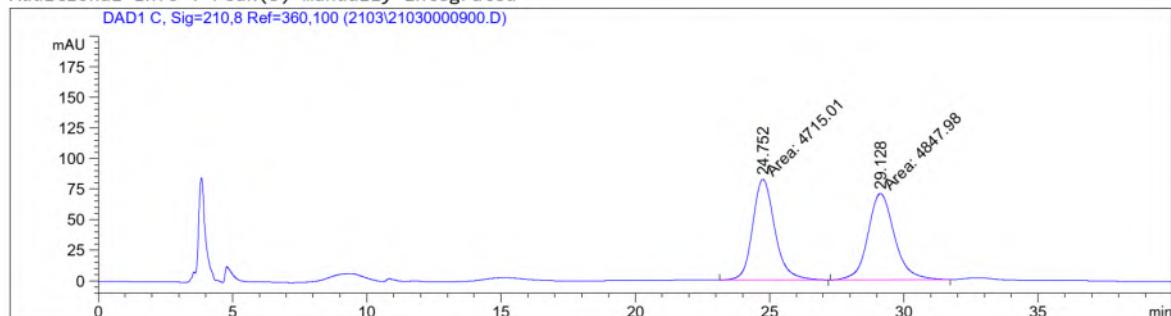


(R)-N-Cyclohexyl-2-(4-fluorophenyl)propanamide (12)

Data File D:\CHEM32\1\DATA\2103\21030000900.D
Sample Name: YY-X-66-Race

```
=====
Acq. Operator   : Analytik          Seq. Line : 1
Acq. Instrument : LC5             Location  : Vial 32
Injection Date  : 3/9/2021 11:50:50 AM    Inj       : 1
                                                Inj Volume : 0.5 μl
Acq. Method     : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/9/2021 11:50:29 AM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/9/2021 12:38:01 PM by Analytik
                                                (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 98:2, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

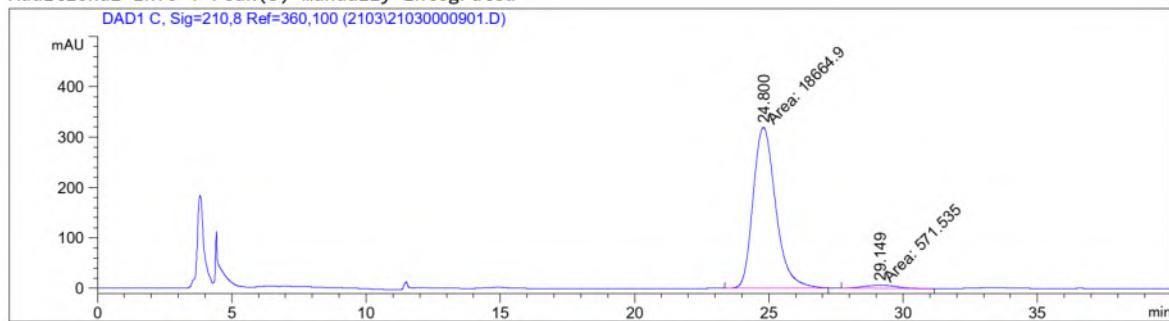
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	24.752	MM	0.9558	4715.00537	82.21693	49.3048
2	29.128	MM	1.1476	4847.97754	70.40529	50.6952

Totals : 9562.98291 152.62222

Data File D:\CHEM32\1\DATA\2103\21030000901.D
Sample Name: YY-X-66-C

```
=====
Acq. Operator   : Analytik          Seq. Line :  2
Acq. Instrument : LC5             Location  : Vial 42
Injection Date  : 3/9/2021 12:31:56 PM      Inj :  1
                                                Inj Volume : 0.5 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 2.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/9/2021 11:50:29 AM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/9/2021 2:05:42 PM by Analytik
                                                (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 98:2, 0,5ml/min
```

Additional Info : Peak(s) manually integrated



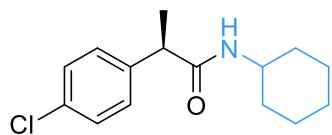
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	24.800	MM	0.9744	1.86649e4	319.26517	97.0289
2	29.149	MM	1.4684	571.53461	6.48698	2.9711

Totals : 1.92365e4 325.75215



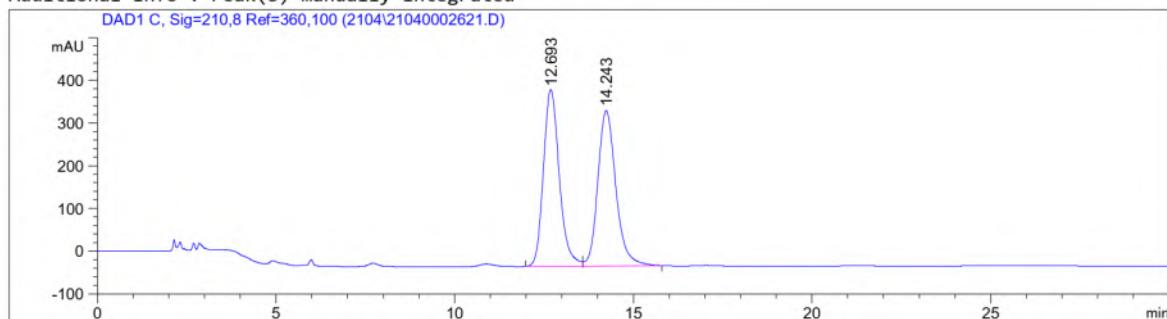
(R)-2-(4-Chlorophenyl)-N-cyclohexylpropanamide (13)

Data File D:\CHEM32\1\DATA\2104\21040002621.D

Sample Name: YY-X-81 Race

```
=====
Acq. Operator   : Analytik          Seq. Line : 13
Acq. Instrument : LC5             Location : Vial 26
Injection Date  : 4/27/2021 1:21:35 AM    Inj : 1
                                                Inj Volume : 0.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY 0,8.M
Last changed    : 4/26/2021 4:26:28 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY 0,8.M
Last changed    : 4/27/2021 11:09:04 AM by Analytik
                                                (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 98:2, 0.8ml/min
```

Additional Info : Peak(s) manually integrated



Area Percent Report

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

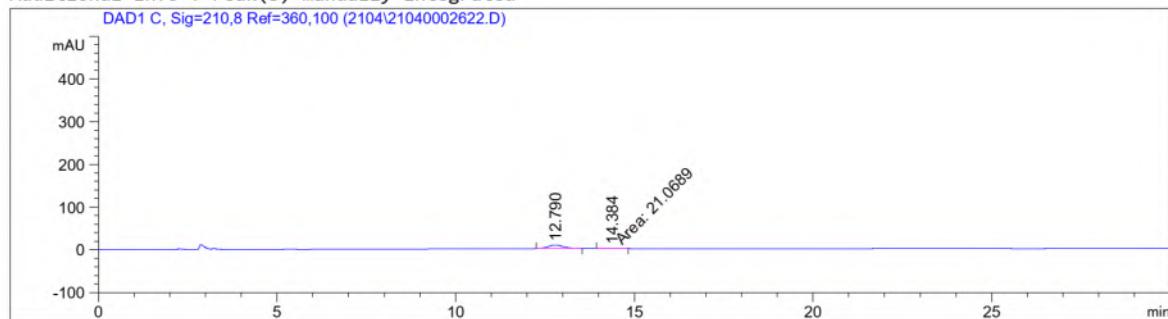
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.693	BV	0.4763	1.27643e4	413.30167	49.9141
2	14.243	VB	0.5402	1.28083e4	364.24915	50.0859

Totals : 2.55726e4 777.55081

Data File D:\CHEM32\1\DATA\2104\21040002622.D
Sample Name: YY-X-81 rt

```
=====
Acq. Operator   : Analytik          Seq. Line : 14
Acq. Instrument : LC5             Location  : Vial 36
Injection Date  : 4/27/2021 1:52:39 AM    Inj       : 1
                                         Inj Volume : 0.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY 0,8.M
Last changed    : 4/26/2021 4:26:28 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY 0,8.M
Last changed    : 4/27/2021 11:09:04 AM by Analytik
                                         (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 98:2, 0.8ml/min
```

Additional Info : Peak(s) manually integrated



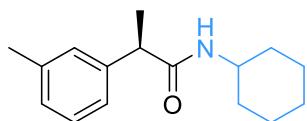
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.790	BB	0.4517	234.26813	7.95322	91.7486
2	14.384	MM	0.4592	21.06885	7.64721e-1	8.2514

Totals : 255.33698 8.71794



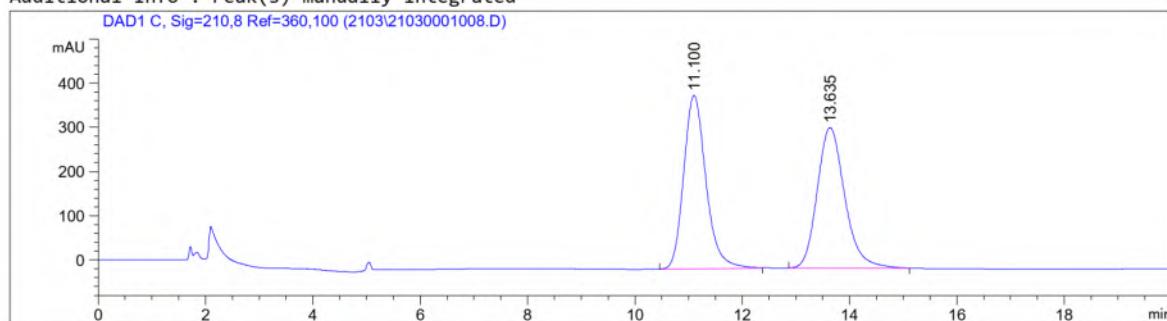
(R)-N-Cyclohexyl-2-(m-tolyl)propanamide (14)

Data File D:\CHEM32\1\DATA\2103\21030001008.D

Sample Name: YY-X-79 Race

```
=====
Acq. Operator   : Analytik           Seq. Line : 1
Acq. Instrument : LC5             Location  : Vial 56
Injection Date  : 3/10/2021 2:23:16 PM      Inj : 1
                                                Inj Volume : 3.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\YY.M
Last changed   : 3/10/2021 2:36:58 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed   : 3/10/2021 3:00:22 PM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 98:2, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.100	BB	0.4443	1.12971e4	391.91498	49.9120
2	13.635	BB	0.5503	1.13369e4	317.72333	50.0880
Totals :					2.26340e4	709.63831

LC5 3/10/2021 3:00:31 PM Analytik

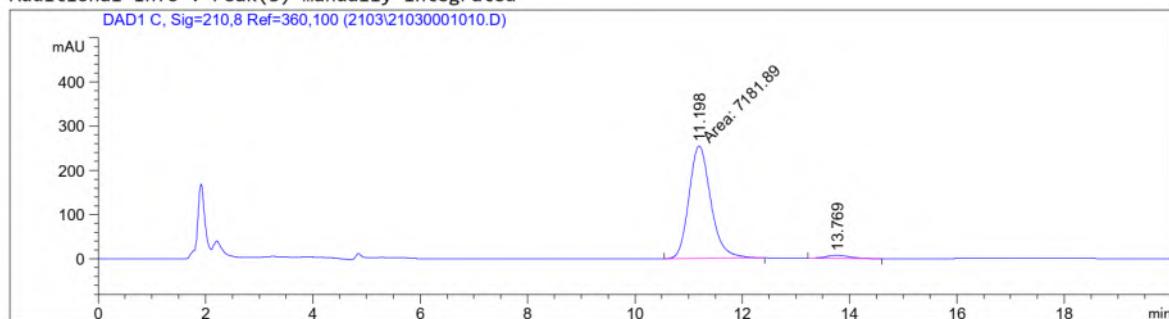
Page 1 of 2

Data File D:\CHEM32\1\DATA\2103\21030001010.D

Sample Name: YY-X-79 C

```
=====
Acq. Operator   : Analytik          Seq. Line : 3
Acq. Instrument : LC5             Location  : Vial 63
Injection Date  : 3/10/2021 3:25:33 PM    Inj : 1
                                         Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/10/2021 2:36:58 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY-81.M
Last changed    : 3/10/2021 3:58:18 PM by Analytik
Method Info     : Cellulose4, Hept./EtOH 98:2, 1ml/min
```

Additional Info : Peak(s) manually integrated



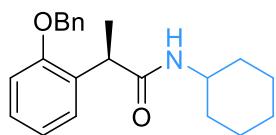
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.198	MM	0.4714	7181.89014	253.90724	97.1749
2	13.769	BB	0.4857	208.79301	6.73403	2.8251

Totals : 7390.68315 260.64127

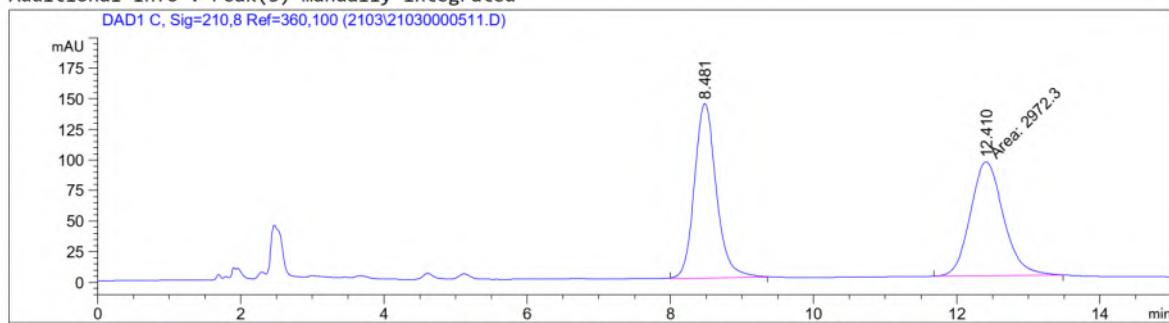


(R)-2-(2-(Benzyl)phenyl)-N-cyclohexylpropanamide (15)

Data File D:\CHEM32\1\DATA\2103\21030000511.D
Sample Name: YY-X-71-Race

```
=====
Acq. Operator   : Analytik          Seq. Line : 5
Acq. Instrument : LC5             Location  : Vial 35
Injection Date  : 3/5/2021 3:01:55 PM    Inj       : 1
                                         Inj Volume : 0.5 µl
Acq. Method    : C:\CHEM32\1\METHODS\TEST RAC.M
Last changed    : 3/5/2021 3:10:59 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\TEST RAC.M
Last changed    : 3/8/2021 12:27:53 PM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated



Area Percent Report

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area
#	[min]		[min]	[mAU*s]	[mAU]	%
1	8.481	BB	0.3262	3034.06128	142.62895	50.5141
2	12.410	MM	0.5333	2972.30225	92.89326	49.4859

Totals : 6006.36353 235.52221

LC5 3/8/2021 12:28:48 PM Analytik

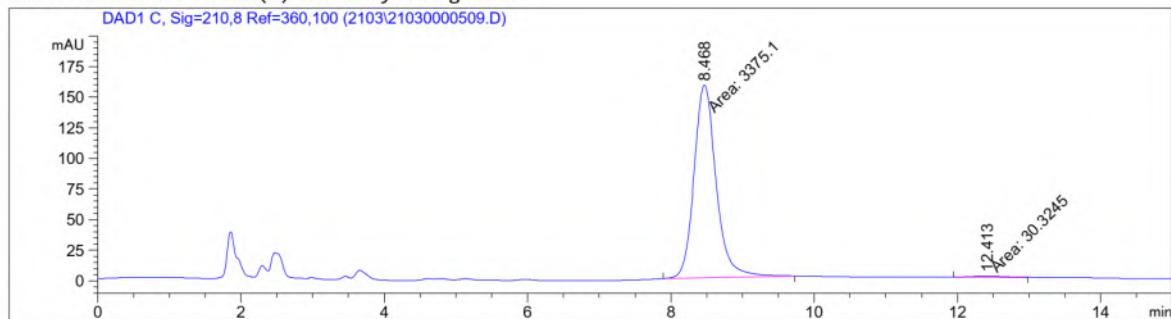
Page 1 of 2

Data File D:\CHEM32\1\DATA\2103\21030000509.D

Sample Name: YY-X-71-C

```
=====
Acq. Operator   : Analytik          Seq. Line : 3
Acq. Instrument : LC5             Location  : Vial 41
Injection Date  : 3/5/2021 2:29:46 PM    Inj : 1
                                         Inj Volume : 0.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\TEST RAC.M
Last changed    : 3/5/2021 2:13:16 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\TEST RAC.M
Last changed    : 3/8/2021 12:27:53 PM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

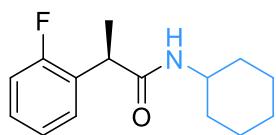
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.468	MM	0.3574	3375.10303	157.39024	99.1095
2	12.413	MM	0.5100	30.32452	9.90925e-1	0.8905

Totals : 3405.42755 158.38117

LC5 3/8/2021 12:33:09 PM Analytik

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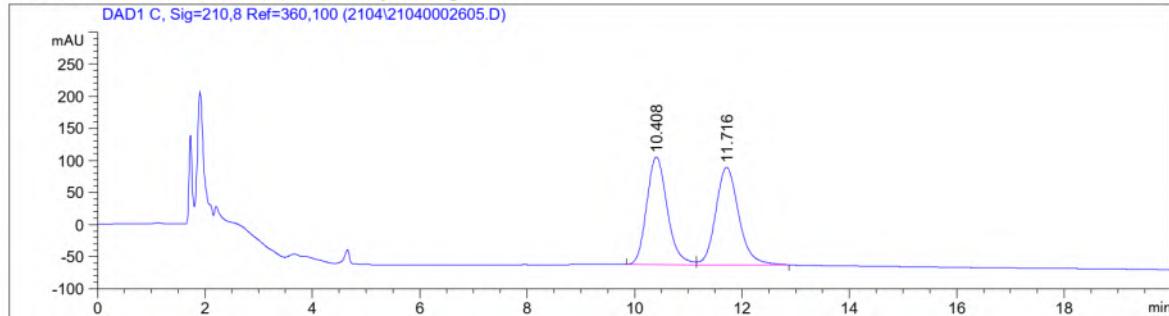
(R)-N-Cyclohexyl-2-(2-fluorophenyl)propanamide (16)

Data File D:\CHEM32\1\DATA\2104\21040002605.D

Sample Name: YY-X-100 Race

```
=====
Acq. Operator   : Analytik           Seq. Line : 1
Acq. Instrument : LC5               Location  : Vial 21
Injection Date  : 4/26/2021 2:58:27 PM    Inj       : 1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY.M
Last changed    : 4/26/2021 3:13:39 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed    : 4/27/2021 10:24:10 AM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 98:2, 1ml/min
```

Additional Info : Peak(s) manually integrated



Area Percent Report

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.408	BV	0.4134	4476.12646	167.71735	49.6773
2	11.716	VB	0.4602	4534.28613	151.91861	50.3227

Totals : 9010.41260 319.63596

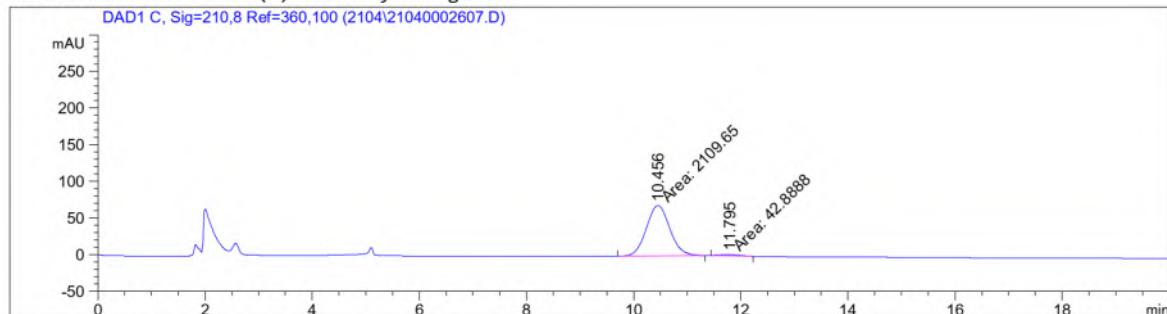
LC5 4/27/2021 10:25:30 AM Analytik

Page 1 of 2

Data File D:\CHEM32\1\DATA\2104\21040002607.D
Sample Name: YY-X-100-30

```
=====
Acq. Operator   : Analytik                     Seq. Line : 3
Acq. Instrument : LC5                         Location : Vial 31
Injection Date  : 4/26/2021 3:40:40 PM          Inj : 1
                                                Inj Volume : 1.0 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 8.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\YY.M
Last changed   : 4/26/2021 3:13:39 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed   : 4/27/2021 10:26:28 AM by Analytik
                  (modified after loading)
Method Info    : Cellulose4, Hept./EtOH 98:2, 1ml/min
```

Additional Info : Peak(s) manually integrated



=====
Area Percent Report

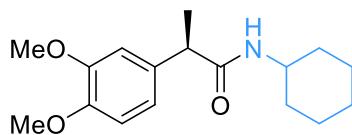
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area
#	[min]		[min]	[mAU*s]	[mAU]	%
1	10.456	MM	0.5097	2109.65234	68.97826	98.0075
2	11.795	MM	0.4103	42.88883	1.74197	1.9925

LC5 4/27/2021 10:37:07 AM Analytik

Page 1 of 2



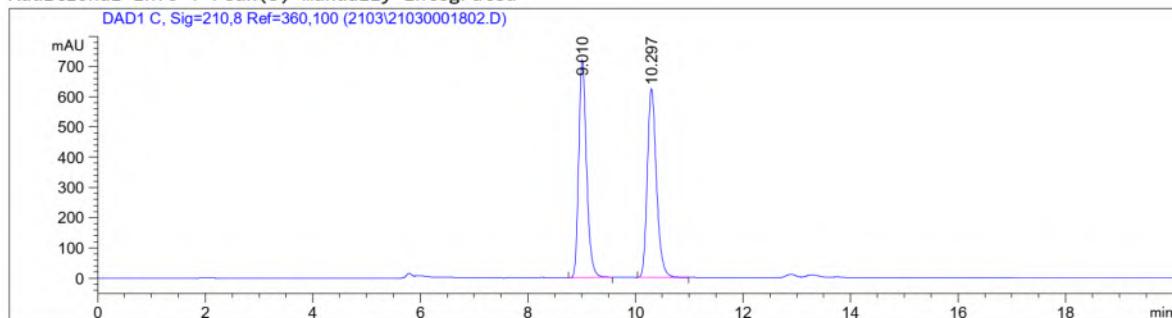
(R)-N-Cyclohexyl-2-(3,4-dimethoxyphenyl)propanamide (17)

Data File D:\CHEM32\1\DATA\2103\21030001802.D

Sample Name: YY-X-70 Race

```
=====
Acq. Operator   : Analytik           Seq. Line : 3
Acq. Instrument : LC5             Location  : Vial 35
Injection Date  : 3/18/2021 1:14:07 PM      Inj       : 1
                                                Inj Volume : 0.5 µl
Acq. Method    : C:\CHEM32\1\METHODS\YY-72.M
Last changed    : 3/17/2021 2:33:11 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY-72.M
Last changed    : 3/18/2021 3:12:26 PM by Analytik
                                                (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.010	VB	0.1542	7298.84277	720.39355	49.6158
2	10.297	BB	0.1826	7411.88623	624.03900	50.3842

Totals : 1.47107e4 1344.43256

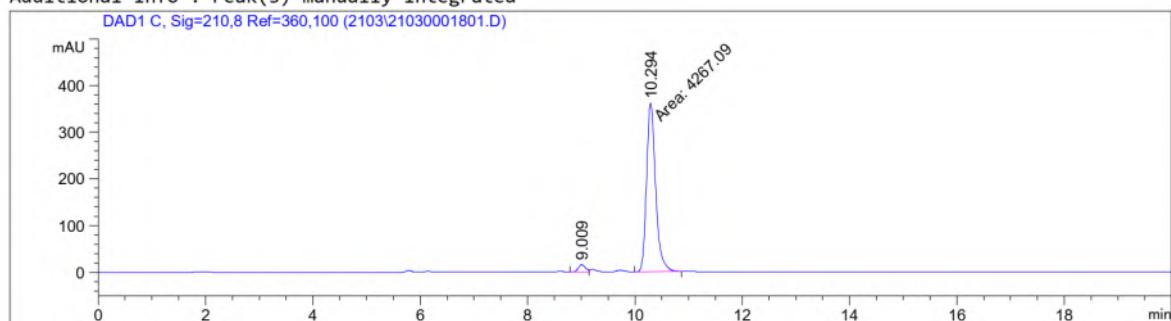
LC5 3/18/2021 3:12:29 PM Analytik

Page 1 of 2

Data File D:\CHEM32\1\DATA\2103\21030001801.D
Sample Name: YY-X-70 C

```
=====
Acq. Operator   : Analytik          Seq. Line :  2
Acq. Instrument : LC5             Location  : Vial 45
Injection Date  : 3/18/2021 12:52:58 PM    Inj       : 1
                                                Inj Volume : 0.5 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 0.1 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY-72.M
Last changed    : 3/17/2021 2:33:11 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY-72.M
Last changed    : 3/18/2021 3:07:31 PM by Analytik
                                                (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



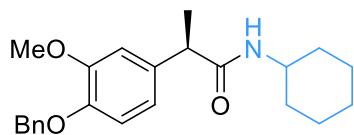
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.009	VV	0.1447	147.89467	15.58652	3.3498
2	10.294	MM	0.1970	4267.08740	360.97736	96.6502

Totals : 4414.98207 376.56387



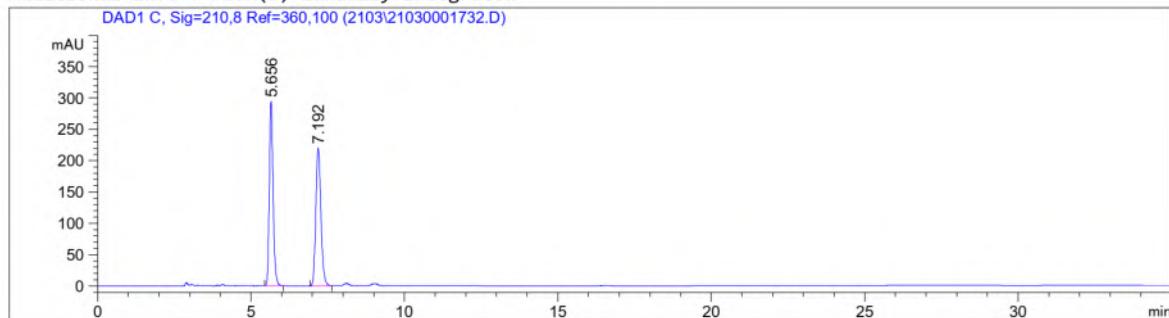
(R)-2-(4-(Benzylxy)-3-methoxyphenyl)-N-cyclohexylpropanamide (18)

Data File D:\CHEM32\1\DATA\2103\21030001732.D

Sample Name: YY-X-90 Race

```
=====
Acq. Operator   : Analytik           Seq. Line : 13
Acq. Instrument : LC5               Location  : Vial 38
Injection Date : 3/18/2021 5:20:04 AM    Inj       : 1
                                                Inj Volume : 0.5 µl
Acq. Method    : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 3/17/2021 4:50:04 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 3/18/2021 9:27:13 AM by Analytik
                                                (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 80:20, 1ml/min
```

Additional Info : Peak(s) manually integrated



Area Percent Report

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	5.656	BB	0.1307	2532.93750	293.43527	49.9898
2	7.192	BB	0.1787	2533.97144	219.47058	50.0102

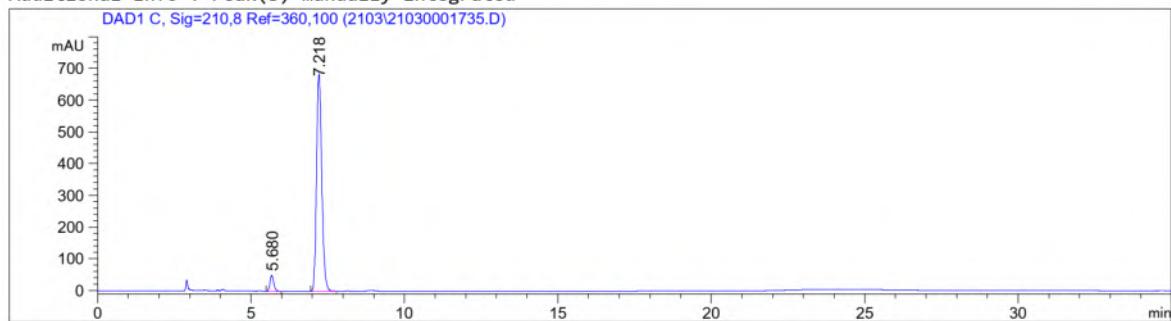
Totals : 5066.90894 512.90585

Data File D:\CHEM32\1\DATA\2103\21030001735.D

Sample Name: YY-X-90 C

```
=====
Acq. Operator   : Analytik          Seq. Line : 14
Acq. Instrument : LC5             Location  : Vial 48
Injection Date  : 3/18/2021 7:23:18 AM    Inj       : 2
                                         Inj Volume : 0.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 3/17/2021 4:50:04 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 3/18/2021 9:28:48 AM by Analytik
                                         (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 80:20, 1ml/min
```

Additional Info : Peak(s) manually integrated



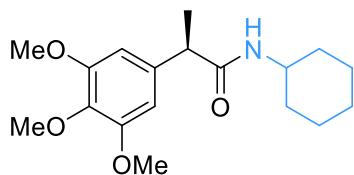
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	5.680	BB	0.1330	434.48895	50.17144	5.0875
2	7.218	BB	0.1821	8105.78857	685.00366	94.9125

Totals : 8540.27753 735.17511



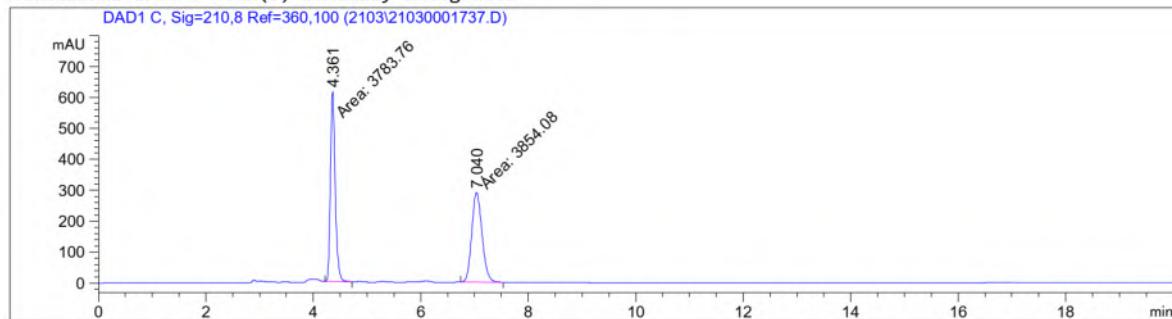
(R)-N-Cyclohexyl-2-(3,4,5-trimethoxyphenyl)propanamide (19)

Data File D:\CHEM32\1\DATA\2103\21030001737.D

Sample Name: YY-X-91 Race

```
=====
Acq. Operator   : Analytik          Seq. Line : 15
Acq. Instrument : LC5             Location  : Vial 39
Injection Date  : 3/18/2021 8:45:29 AM    Inj       : 2
                                                Inj Volume : 0.5 μl
Acq. Method     : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 3/17/2021 4:50:04 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 3/18/2021 9:53:49 AM by Analytik
                                                (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 80:20, 1ml/min
```

Additional Info : Peak(s) manually integrated



Area Percent Report

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	4.361	MM	0.1028	3783.76318	613.43744	49.5397
2	7.040	MM	0.2219	3854.08301	289.43216	50.4603

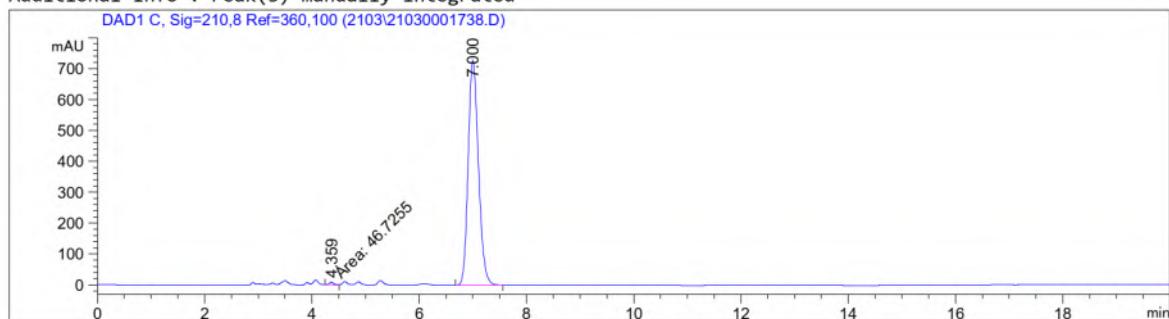
Totals : 7637.84619 902.86960

Data File D:\CHEM32\1\DATA\2103\21030001738.D

Sample Name: YY-X-91 C

```
=====
Acq. Operator   : Analytik          Seq. Line : 16
Acq. Instrument : LC5             Location  : Vial 49
Injection Date  : 3/18/2021 9:26:33 AM    Inj : 1
                                         Inj Volume : 0.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 3/18/2021 9:37:46 AM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 3/18/2021 9:53:49 AM by Analytik
                  (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 80:20, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

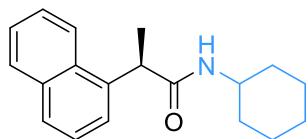
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	4.359	MM	0.0971	46.72552	8.01900	0.4806
2	7.000	BB	0.2074	9676.38086	726.53180	99.5194

Totals : 9723.10638 734.55080

LC5 3/18/2021 9:53:51 AM Analytik

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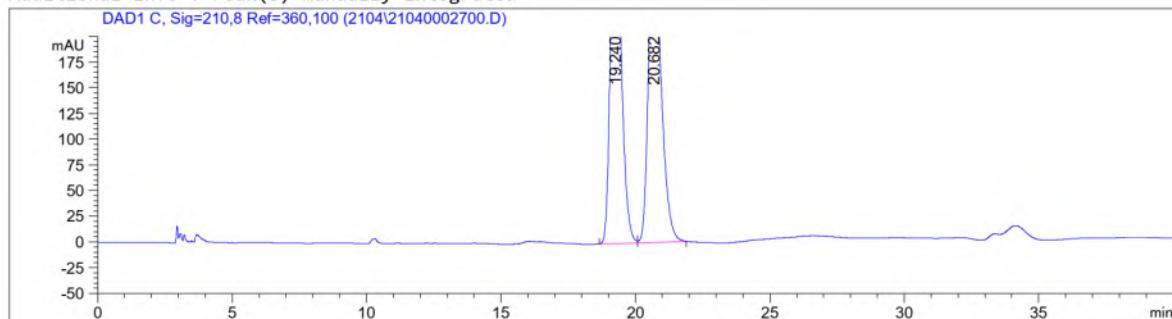
(R)-N-Cyclohexyl-2-(naphthalen-1-yl)propanamide (20)

Data File D:\CHEM32\1\DATA\2104\21040002700.D

Sample Name: YY-X-76 Race

```
=====
Acq. Operator   : Analytik          Seq. Line :  1
Acq. Instrument : LC5             Location  : Vial 1
Injection Date  : 4/27/2021 12:20:29 PM      Inj       : 1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY.M
Last changed    : 4/26/2021 2:58:10 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed    : 4/27/2021 2:51:41 PM by Analytik
                                (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 98:2, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area
#	[min]		[min]	[mAU*s]	[mAU]	%
1	19.240	BV	0.4226	9755.81445	357.19949	49.9461
2	20.682	BV	0.5035	9776.88086	299.01672	50.0539

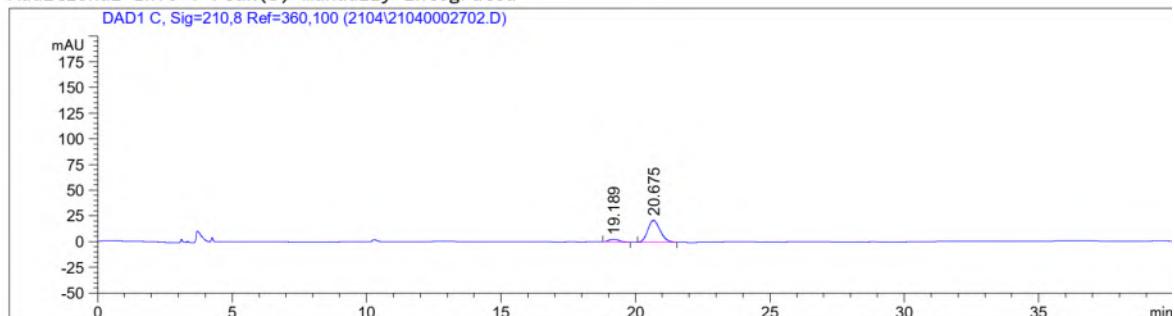
Totals : 1.95327e4 656.21622

Data File D:\CHEM32\1\DATA\2104\21040002702.D

Sample Name: YY-X-76 rt

```
=====
Acq. Operator   : Analytik          Seq. Line : 3
Acq. Instrument : LC5             Location  : Vial 2
Injection Date  : 4/27/2021 1:52:40 PM    Inj : 1
                                         Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY.M
Last changed    : 4/26/2021 2:58:10 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed    : 4/27/2021 2:51:41 PM by Analytik
                                         (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 98:2, 1ml/min
```

Additional Info : Peak(s) manually integrated



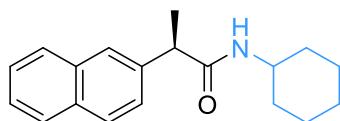
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	19.189	BB	0.3731	74.15709	2.87334	9.7107
2	20.675	BB	0.5057	689.50793	21.40855	90.2893

Totals : 763.66502 24.28189



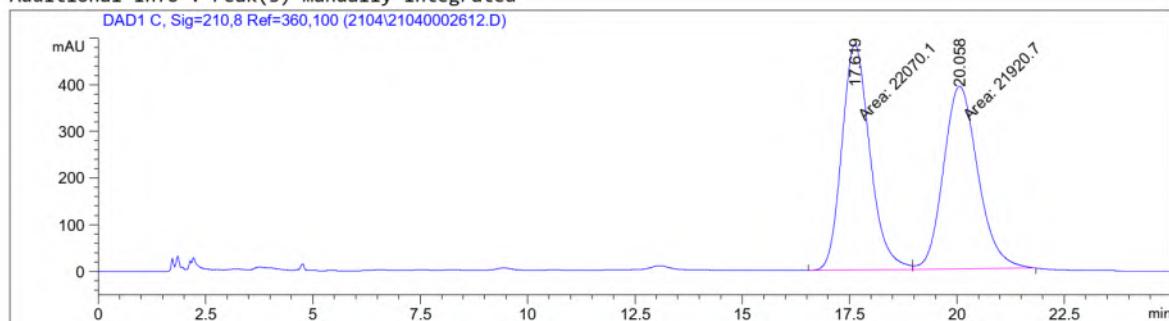
(R)-N-Cyclohexyl-2-(naphthalen-2-yl)propanamide (21)

Data File D:\CHEM32\1\DATA\2104\21040002612.D

Sample Name: YY-X-69 Race

```
=====
Acq. Operator   : Analytik          Seq. Line : 7
Acq. Instrument : LC5             Location  : Vial 23
Injection Date  : 4/26/2021 6:26:31 PM    Inj       : 1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY.M
Last changed    : 4/26/2021 4:24:04 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed    : 4/27/2021 10:44:39 AM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 98:2, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	17.619	MF	0.7615	2.2070e4	483.05246	50.1698
2	20.058	FM	0.9332	2.1920e4	391.48169	49.8302
Totals :				4.39908e4	874.53415	

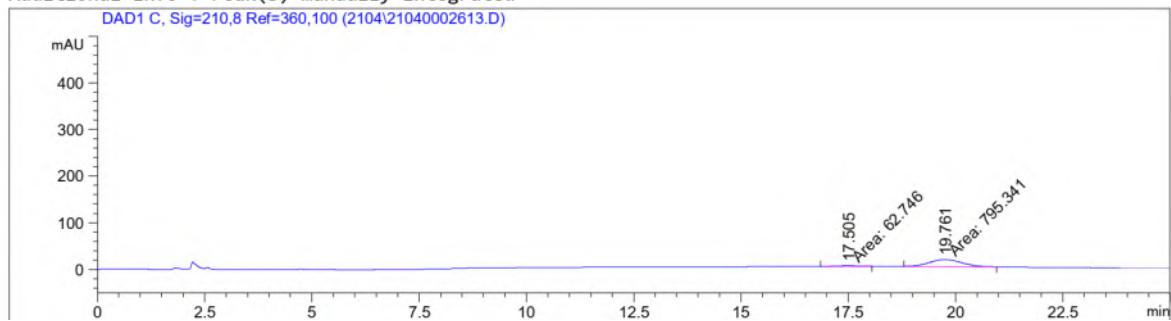
LC5 4/27/2021 10:45:05 AM Analytik

Page 1 of 2

Data File D:\CHEM32\1\DATA\2104\21040002613.D
Sample Name: YY-X-69 rt

```
=====
Acq. Operator   : Analytik          Seq. Line :  8
Acq. Instrument : LC5             Location  : Vial 33
Injection Date  : 4/26/2021 7:07:36 PM      Inj :  1
                                                Inj Volume : 1.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\YY.M
Last changed   : 4/26/2021 4:24:04 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed   : 4/27/2021 10:44:39 AM by Analytik
                  (modified after loading)
Method Info    : Cellulose4, Hept./EtOH 98:2, 1ml/min
```

Additional Info : Peak(s) manually integrated



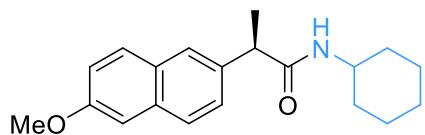
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	17.505	MM	0.5005	62.74595	1.60821	7.3123
2	19.761	MM	0.8804	795.34088	15.05666	92.6877

Totals : 858.08683 16.66487



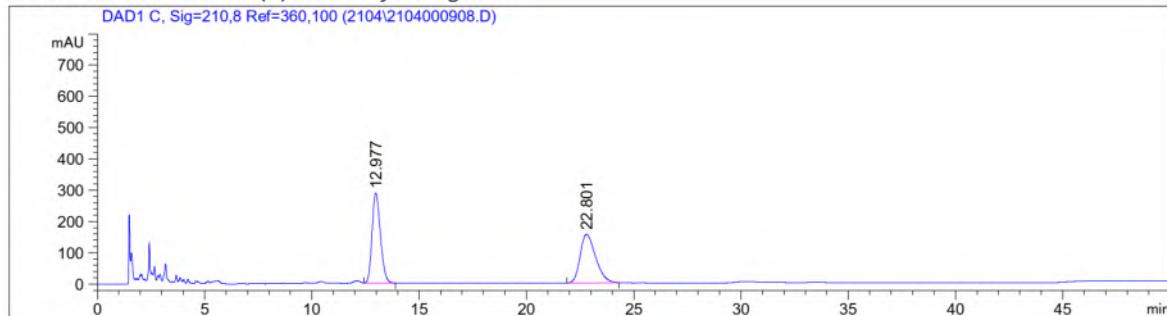
(R)-N-Cyclohexyl-2-(6-methoxynaphthalen-2-yl)propanamide (22)

Data File D:\CHEM32\1\DATA\2104\2104000908.D

Sample Name: YY-X-102 Race

```
=====
Acq. Operator   : Analytik                     Seq. Line :  8
Acq. Instrument : LC5                         Location  : Vial 11
Injection Date  : 4/9/2021 8:49:38 PM          Inj       : 1
                                                Inj Volume : 1.0 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 5.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\CHIRAL OJ-H-2.0.M
Last changed    : 4/9/2021 4:22:35 PM by Analytik
                    (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\CHIRAL OJ-H-2.0.M
Last changed    : 4/12/2021 12:41:25 PM by Analytik
                    (modified after loading)
Method Info     : Chiralcel OJ-H , Hept./EtOH 95:5 , 2.0 ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

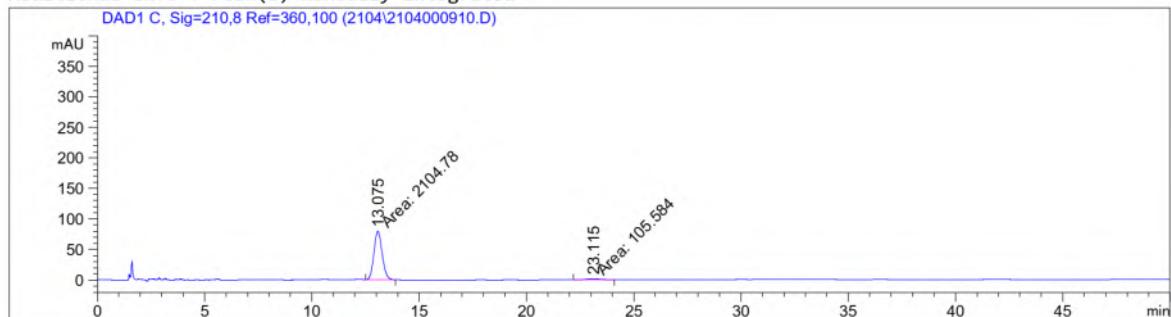
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.977	VB	0.4207	7830.07422	288.43195	50.0303
2	22.801	BB	0.7677	7820.59229	155.76659	49.9697

Data File D:\CHEM32\1\DATA\2104\2104000910.D

Sample Name: YY-X-102 C

```
=====
Acq. Operator   : Analytik          Seq. Line :  9
Acq. Instrument : LC5             Location  : Vial 31
Injection Date  : 4/9/2021 10:31:53 PM    Inj :  2
                                                Inj Volume : 1.0 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 2.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\CHIRAL OJ-H-2.0.M
Last changed    : 4/9/2021 4:22:35 PM by Analytik
                           (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\CHIRAL OJ-H-2.0.M
Last changed    : 4/12/2021 12:42:19 PM by Analytik
                           (modified after loading)
Method Info     : Chiralcel OJ-H , Hept./EtOH 95:5 , 2.0 ml/min
```

Additional Info : Peak(s) manually integrated

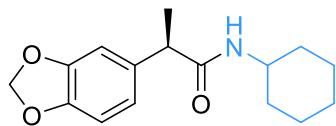


```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	13.075	MM	0.4434	2104.77808	79.11703	95.2232
2	23.115	MM	0.8175	105.58363	2.15261	4.7768



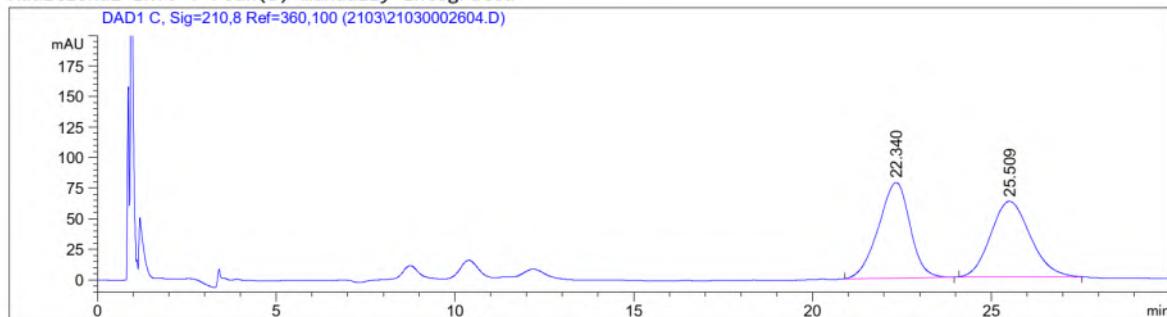
(R)-2-(Benzo[d][1,3]dioxol-5-yl)-N-cyclohexylpropanamide (23)

Data File D:\CHEM32\1\DATA\2103\21030002604.D

Sample Name: YY-X-92-Race

```
=====
Acq. Operator   : Analytik          Seq. Line : 4
Acq. Instrument : LCS             Location : Vial 11
Injection Date  : 3/26/2021 2:11:10 PM    Inj : 1
                                                Inj Volume : 0.5 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 2.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/26/2021 1:34:42 PM by Analytik
                           (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY-0,5.M
Last changed    : 3/26/2021 1:53:02 PM by Analytik
                           (modified after loading)
Method Info     : Cellulose 4, Hept./EtOH 98:2, 2.0ml/min
```

Additional Info : Peak(s) manually integrated



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

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

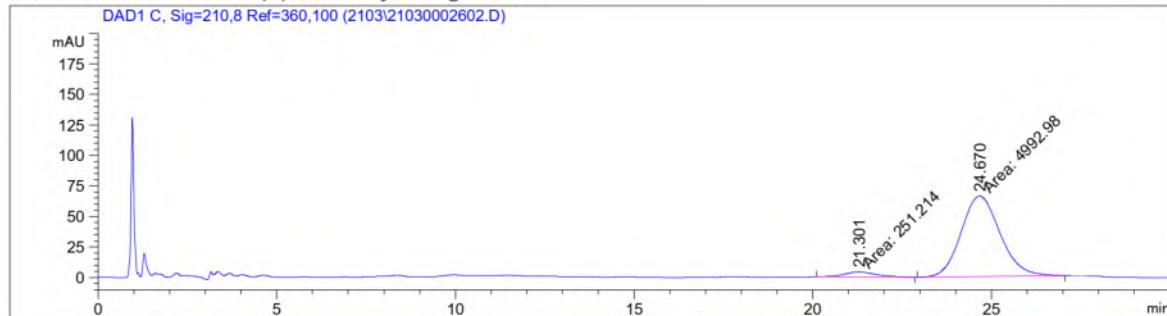
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	22.340	BB	0.9401	4817.88672	78.27844	50.1406
2	25.509	BB	1.1761	4790.86963	61.95966	49.8594

Data File D:\CHEM32\1\DATA\2103\21030002602.D
Sample Name: YY-X-92-C

```
=====
Acq. Operator   : Analytik          Seq. Line :  2
Acq. Instrument : LC5             Location  : Vial 21
Injection Date  : 3/26/2021 12:55:57 PM    Inj       : 1
                                                Inj Volume : 0.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/26/2021 1:34:42 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY-0,5.M
Last changed    : 3/26/2021 1:53:02 PM by Analytik
                  (modified after loading)
Method Info     : Cellulose 4, Hept./EtOH 98:2, 2.0ml/min
```

Additional Info : Peak(s) manually integrated



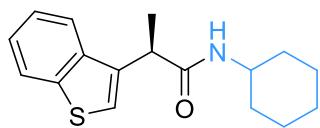
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width MM	Area [mAU*s]	Height [mAU]	Area %
1	21.301	MM	1.0427	251.21382	4.01560	4.7903
2	24.670	MM	1.2628	4992.97900	65.89839	95.2097

Totals : 5244.19283 69.91399



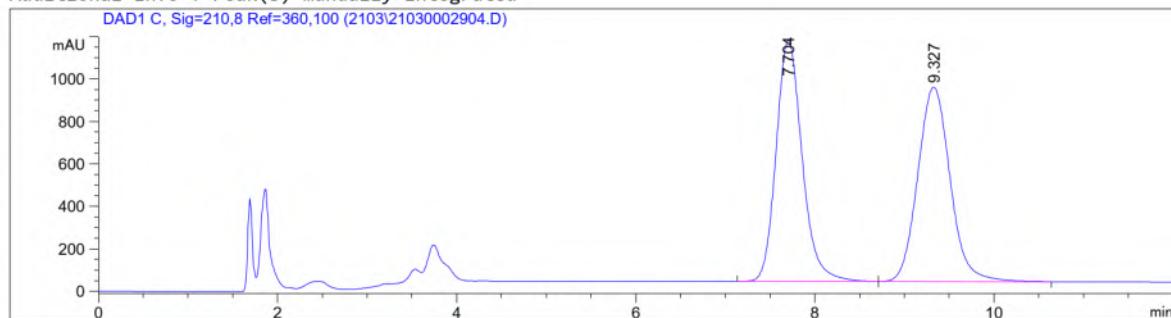
(R)-2-(Benzo[b]thiophen-3-yl)-N-cyclohexylpropanamide (24)

Data File D:\CHEM32\1\DATA\2103\21030002904.D

Sample Name: YY-X-97-Race

```
=====
Acq. Operator   : Analytik           Seq. Line : 5
Acq. Instrument : LC5               Location  : Vial 12
Injection Date  : 3/29/2021 12:07:30 PM    Inj       : 1
                                                Inj Volume : 3.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/29/2021 12:19:00 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/29/2021 12:39:26 PM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated



Area Percent Report

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area %
#	[min]		[min]	[mAU*s]	[mAU]	%
1	7.704	BV	0.3120	2.29718e4	1135.46838	49.9078
2	9.327	VB	0.3935	2.30566e4	916.47522	50.0922
Totals :						
4.60284e4 2051.94360						

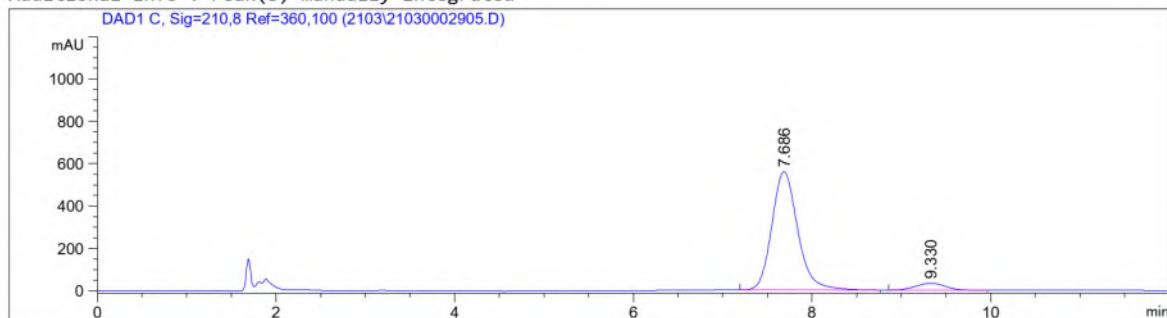
LC5 3/29/2021 12:39:31 PM Analytik

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Data File D:\CHEM32\1\DATA\2103\21030002905.D
Sample Name: YY-X-97-C-35

```
=====
Acq. Operator   : Analytik          Seq. Line :   6
Acq. Instrument : LC5             Location  : Vial 22
Injection Date  : 3/29/2021 12:23:35 PM    Inj :   1
                                                Inj Volume : 3.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/29/2021 12:19:00 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/29/2021 12:39:26 PM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated



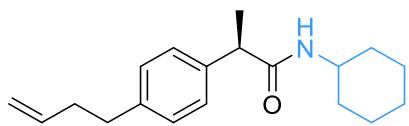
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area %
#	[min]		[min]	[mAU*s]	[mAU]	
1	7.686	BB	0.3116	1.12752e4	558.27411	93.6905
2	9.330	BB	0.3654	759.31665	32.64115	6.3095

Totals : 1.20345e4 590.91526



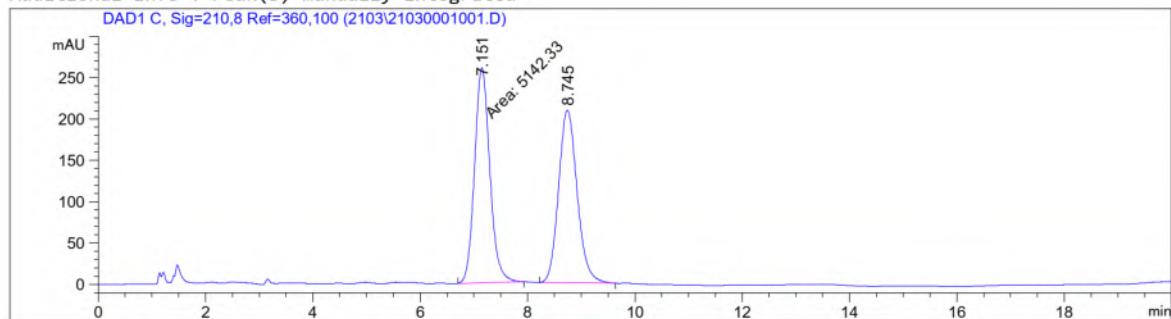
(R)-2-(4-(but-3-en-1-yl)phenyl)-N-cyclohexylpropanamide (25)

Data File D:\CHEM32\1\DATA\2103\21030001001.D

Sample Name: YY-X-74 Race

```
=====
Acq. Operator   : Analytik          Seq. Line :  2
Acq. Instrument : LC5             Location  : Vial 53
Injection Date  : 3/10/2021 11:24:23 AM    Inj :  1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/10/2021 11:23:59 AM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/10/2021 2:13:47 PM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 98:2, 1.5ml/min
```

Additional Info : Peak(s) manually integrated



Area Percent Report

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.151	MM	0.3299	5142.32764	259.77167	50.6281
2	8.745	BB	0.3703	5014.73730	208.76010	49.3719

Totals : 1.01571e4 468.53177

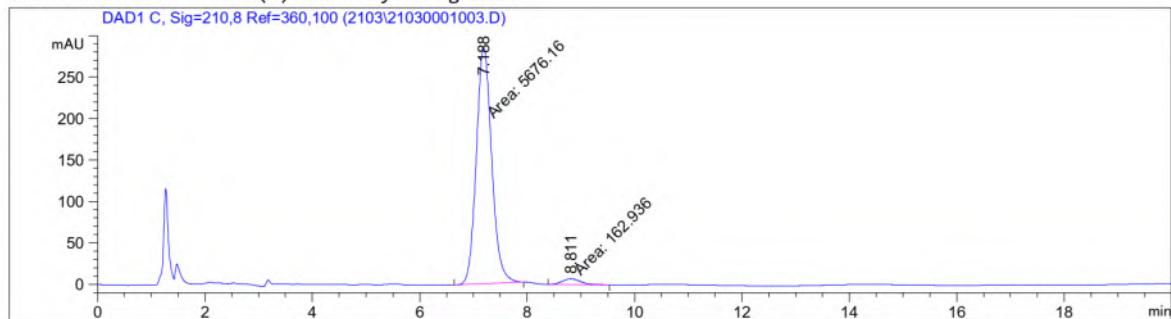
LC5 3/10/2021 2:15:22 PM Analytik

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Data File D:\CHEM32\1\DATA\2103\21030001003.D
Sample Name: YY-X-74 C

```
=====
Acq. Operator   : Analytik          Seq. Line : 4
Acq. Instrument : LC5             Location  : Vial 61
Injection Date  : 3/10/2021 12:06:41 PM    Inj       : 1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/10/2021 11:23:59 AM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed    : 3/10/2021 2:13:47 PM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 98:2, 1.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

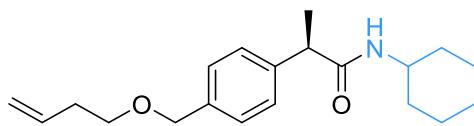
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.188	MM	0.3316	5676.15723	285.30203	97.2096
2	8.811	MM	0.3925	162.93628	6.91835	2.7904

Totals : 5839.09351 292.22038

LC5 3/10/2021 2:17:29 PM Analytik

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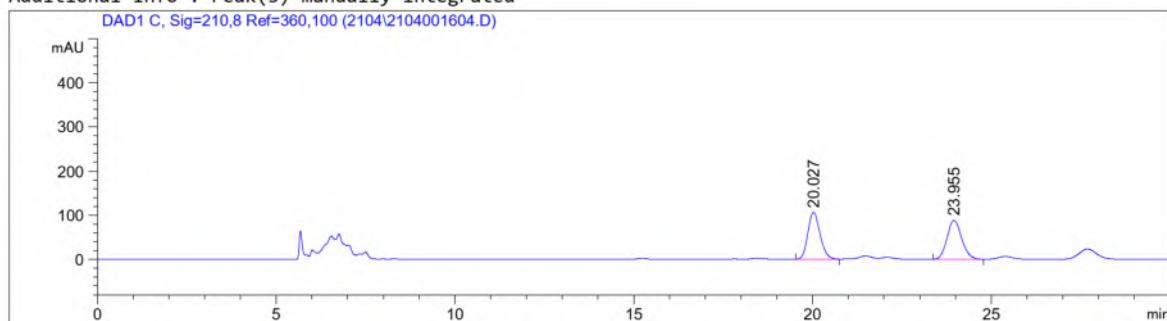
(R)-2-((4-((But-3-en-1-yloxy)methyl)phenyl)-N-cyclohexylpropanamide (26)

Data File D:\CHEM32\1\DATA\2104\2104001604.D

Sample Name: YY-X-88 Race

```
=====
Acq. Operator   : Analytik           Seq. Line :  4
Acq. Instrument : LC5             Location : Vial 1
Injection Date  : 4/16/2021 2:46:18 PM    Inj :  1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY 0.5.M
Last changed    : 3/29/2021 10:16:57 AM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY-0,5.M
Last changed    : 4/16/2021 2:49:17 PM by Analytik
                                (modified after loading)
Method Info     : AD-H, Hept./Isoprop 95:5 0,5ml/min
```

Additional Info : Peak(s) manually integrated



Area Percent Report

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	20.027	BB	0.3694	2527.19849	106.28359	49.9057
2	23.955	BB	0.4434	2536.74609	88.23108	50.0943

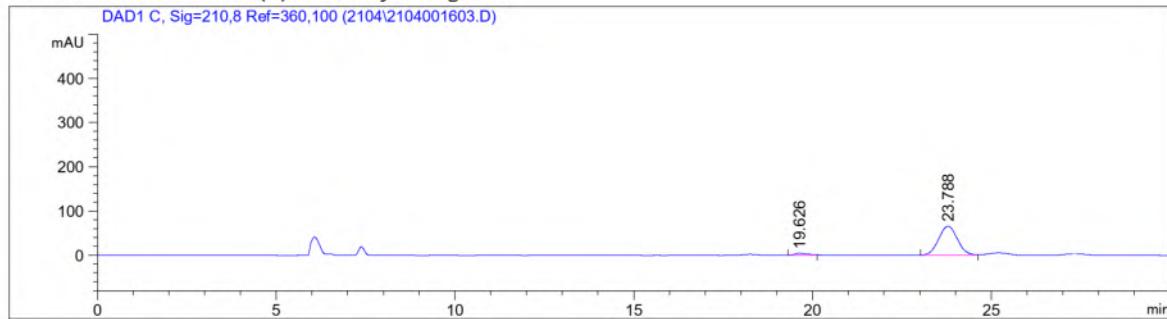
Totals : 5063.94458 194.51467

Data File D:\CHEM32\1\DATA\2104\2104001603.D

Sample Name: YY-X-88 C

```
=====
Acq. Operator   : Analytik          Seq. Line :  3
Acq. Instrument : LC5             Location  : Vial 2
Injection Date  : 4/16/2021 2:00:10 PM      Inj : 1
                                         Inj Volume : 1.0 µl
Different Inj Volume from Sequence !    Actual Inj Volume : 5.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY 0.5.M
Last changed    : 3/29/2021 10:16:57 AM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY-0,5.M
Last changed    : 4/16/2021 2:49:17 PM by Analytik
                                         (modified after loading)
Method Info     : AD-H, Hept./Isoprop 95:5 0,5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

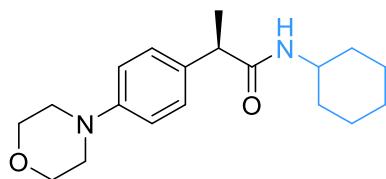
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	19.626	BB	0.3003	99.83392	4.69176	3.9910
2	23.788	BV	0.5776	2401.64136	65.53313	96.0090

Totals : 2501.47528 70.22488

LC5 4/19/2021 11:10:35 AM Analytik

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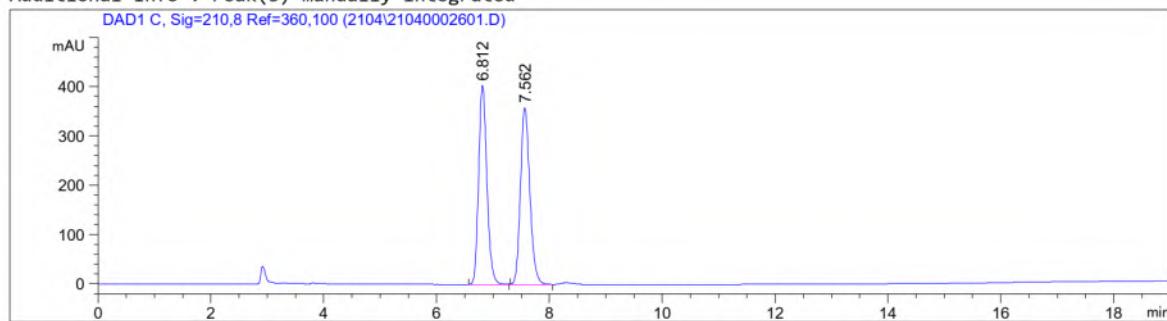
(R)-N-Cyclohexyl-2-(4-morpholinophenyl)propanamide (27)

Data File D:\CHEM32\1\DATA\2104\21040002601.D

Sample Name: YY-X-104 Race

```
=====
Acq. Operator   : Analytik           Seq. Line : 1
Acq. Instrument : LC5               Location : Vial 27
Injection Date  : 4/26/2021 12:47:30 PM    Inj : 1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 4/26/2021 12:58:28 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 4/26/2021 2:07:05 PM by Analytik
                  (modified after loading)
Method Info     : Chiralpak AD-H, Hept./EtOH 80:20, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area %
#	[min]		[min]	[mAU*s]	[mAU]	%
1	6.812	BB	0.1571	4121.67480	403.83118	49.8126
2	7.562	BB	0.1791	4152.68066	358.53961	50.1874

Totals : 8274.35547 762.37079

LC5 4/26/2021 2:07:43 PM Analytik

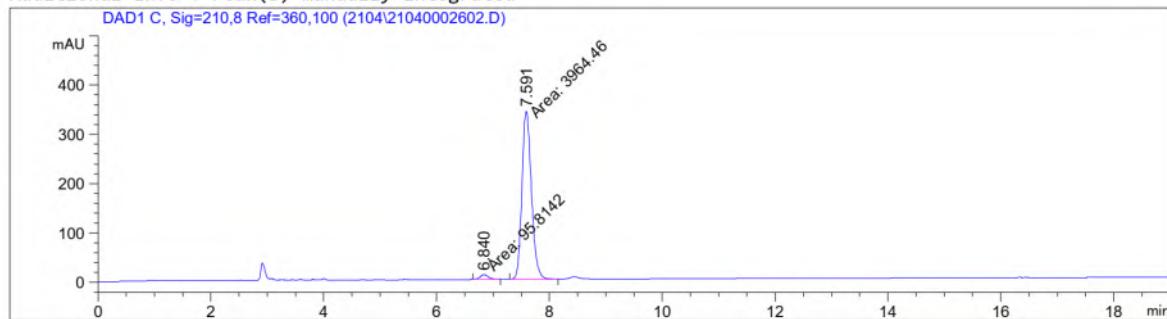
Page 1 of 2

Data File D:\CHEM32\1\DATA\2104\21040002602.D

Sample Name: YY-X-104 -C

```
=====
Acq. Operator   : Analytik          Seq. Line : 2
Acq. Instrument : LC5             Location : Vial 11
Injection Date  : 4/26/2021 1:08:35 PM      Inj : 1
                                                Inj Volume : 1.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 4/26/2021 12:58:28 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 4/26/2021 2:07:05 PM by Analytik
                  (modified after loading)
Method Info     : Chiralpak AD-H, Hept./EtOH 80:20, 1ml/min
```

Additional Info : Peak(s) manually integrated



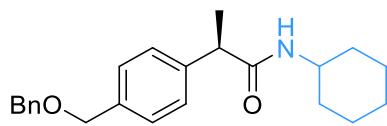
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.840	MM	0.1691	95.81420	9.44570	2.3598
2	7.591	MM	0.1939	3964.45776	340.69324	97.6402

Totals : 4060.27197 350.13894



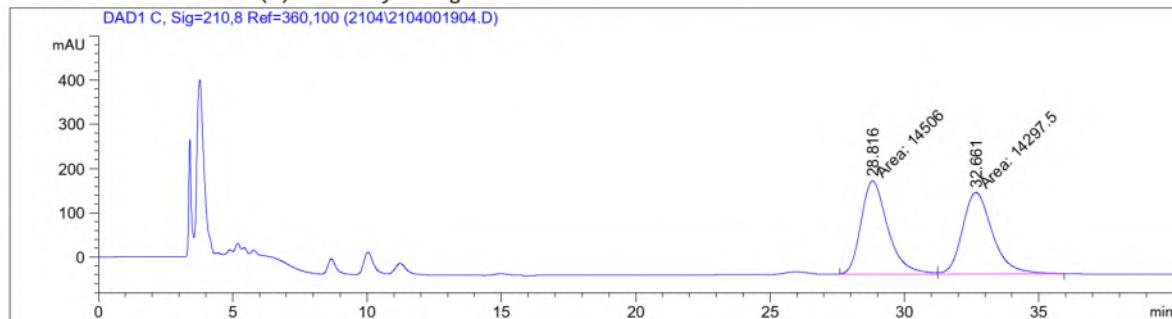
(R)-2-((Benzyl)oxy)methylphenyl-N-cyclohexylpropanamide (28)

Data File D:\CHEM32\1\DATA\2104\2104001904.D

Sample Name: YY-X-98 Race

```
=====
Acq. Operator   : Analytik          Seq. Line : 3
Acq. Instrument : LC5             Location  : Vial 11
Injection Date  : 4/19/2021 1:58:13 PM      Inj       : 1
                                                Inj Volume : 1.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\YY 0.5.M
Last changed    : 3/29/2021 10:16:57 AM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\YY-0,5.M
Last changed    : 4/19/2021 4:06:36 PM by Analytik
                                                (modified after loading)
Method Info     : Cellulose 4, Hept./Isoprop. 95:5, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

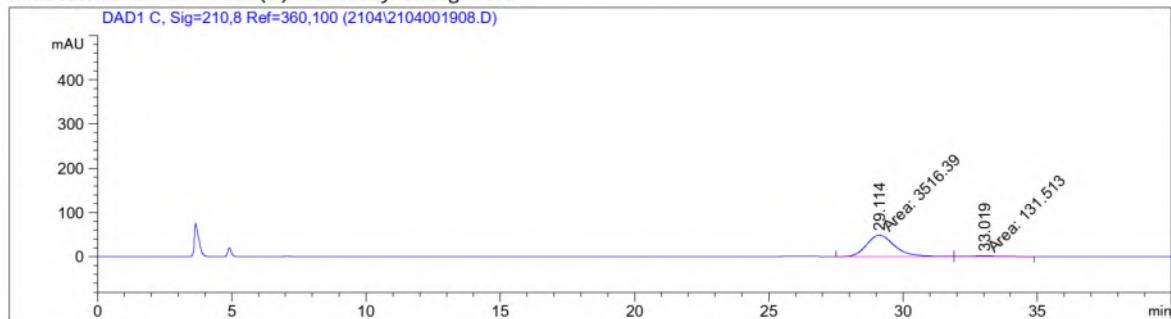
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	28.816	MF	1.1532	1.45060e4	209.65520	50.3619
2	32.661	FM	1.2953	1.42975e4	183.97166	49.6381
Totals :					2.88035e4	393.62686

Data File D:\CHEM32\1\DATA\2104\2104001908.D

Sample Name: YY-X-98 -30

```
=====
Acq. Operator   : Analytik          Seq. Line :   6
Acq. Instrument : LC5             Location  : Vial 12
Injection Date  : 4/19/2021 5:02:32 PM    Inj       : 1
                                                Inj Volume : 2.5 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 5.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY 0.5.M
Last changed    : 4/19/2021 4:10:16 PM by Analytik
                           (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY-0,5.M
Last changed    : 4/19/2021 4:06:36 PM by Analytik
                           (modified after loading)
Method Info     : Cellulose 4, Hept./Isoprop. 95:5, 0.5ml/min
```

Additional Info : Peak(s) manually integrated

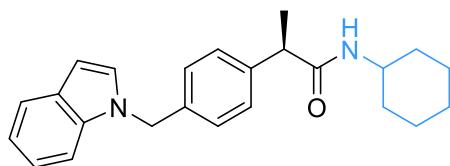


```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	29.114	MF	1.2179	3516.38525	48.12136	96.3948
2	33.019	FM	1.4214	131.51259	1.54204	3.6052



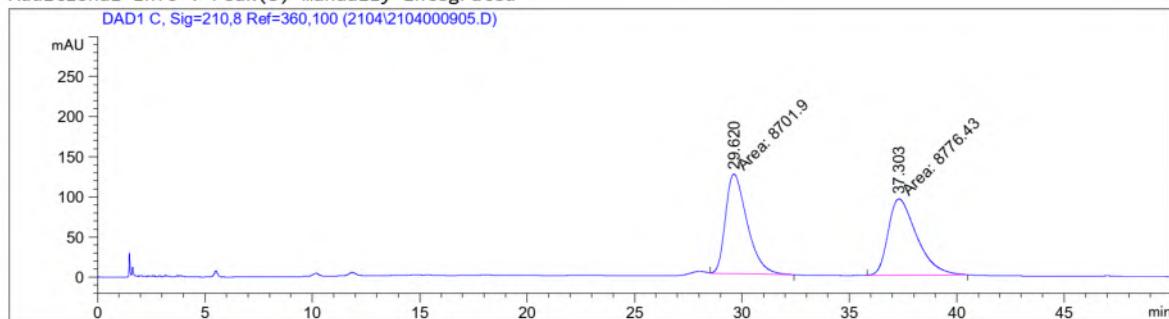
(R)-2-(4-((1*H*-Indol-1-yl)methyl)phenyl)-*N*-cyclohexylpropanamide (29)

Data File D:\CHEM32\1\DATA\2104\2104000905.D

Sample Name: YY-X-106 Race

```
=====
Acq. Operator   : Analytik           Seq. Line : 5
Acq. Instrument : LC5             Location : Vial 12
Injection Date  : 4/9/2021 6:16:24 PM    Inj : 1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\CHIRAL OJ-H-2.0.M
Last changed    : 4/9/2021 4:22:35 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\CHIRAL OJ-H-2.0.M
Last changed    : 4/12/2021 12:16:26 PM by Analytik
                  (modified after loading)
Method Info     : Chiralcel OJ-H , Hept./EtOH 95:5 , 2.0 ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	29.620	MM	1.1673	8701.90430	124.24153	49.7868
2	37.303	MM	1.5391	8776.43066	95.03600	50.2132

Totals : 1.74783e4 219.27753

LC5 4/12/2021 12:19:10 PM Analytik

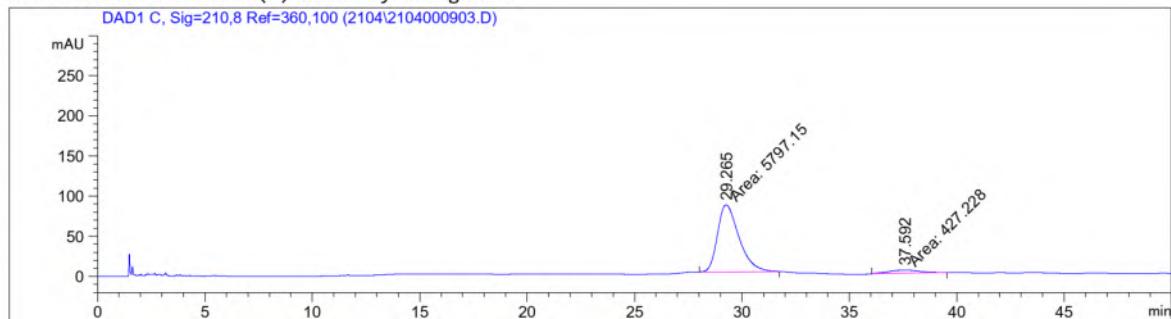
Page 1 of 2

Data File D:\CHEM32\1\DATA\2104\2104000903.D

Sample Name: YY-X-106 C

```
=====
Acq. Operator   : Analytik          Seq. Line : 4
Acq. Instrument : LC5             Location  : Vial 23
Injection Date  : 4/9/2021 4:34:12 PM    Inj : 1
                                         Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\CHIRAL OJ-H-2.0.M
Last changed    : 4/9/2021 4:22:35 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\CHIRAL OJ-H-2.0.M
Last changed    : 4/12/2021 12:16:26 PM by Analytik
                  (modified after loading)
Method Info     : Chiralcel OJ-H , Hept./EtOH 95:5 , 2.0 ml/min
```

Additional Info : Peak(s) manually integrated



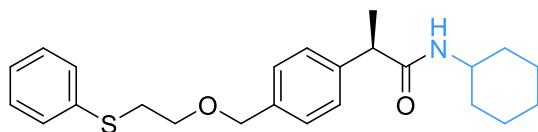
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	29.265	MM	1.1567	5797.14600	83.53152	93.1362
2	37.592	MM	1.6934	427.22800	4.20484	6.8638

Totals : 6224.37399 87.73636



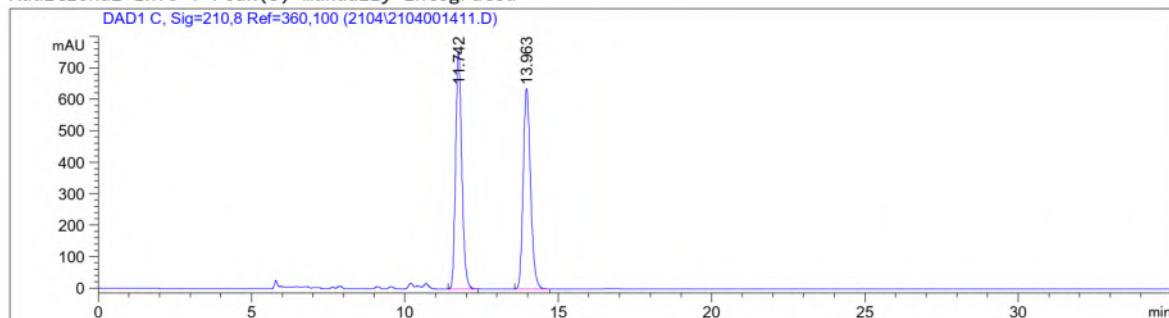
(R)-N-Cyclohexyl-2-(4-((2-(phenylthio)ethoxy)methyl)phenyl)propanamide (30)

Data File D:\CHEM32\1\DATA\2104\2104001411.D

Sample Name: YY-X-107 Race

```
=====
Acq. Operator   : Analytik          Seq. Line : 3
Acq. Instrument : LC5             Location  : Vial 11
Injection Date  : 4/14/2021 5:12:19 PM      Inj       : 2
                                                Inj Volume : 0.5 µl
Acq. Method    : C:\CHEM32\1\METHODS\HOLZ 0.5.M
Last changed   : 4/14/2021 11:19:56 AM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\HOLZ 0.5.M
Last changed   : 4/15/2021 11:14:11 AM by Analytik
                                                (modified after loading)
Method Info    : Chiraldak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

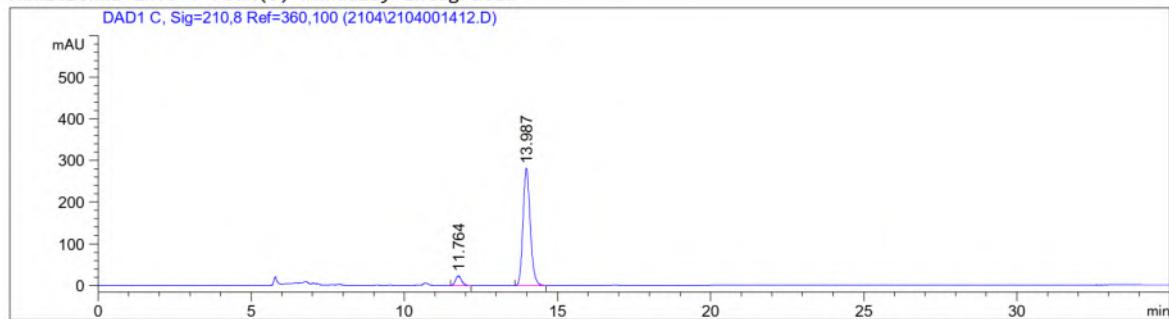
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.742	BB	0.2156	1.05899e4	755.02643	49.9776
2	13.963	BB	0.2561	1.05993e4	637.07794	50.0224

Totals : 2.11892e4 1392.10437

Data File D:\CHEM32\1\DATA\2104\2104001412.D
Sample Name: YY-X-107 C

```
=====
Acq. Operator   : Analytik          Seq. Line :  4
Acq. Instrument : LC5             Location  : Vial 21
Injection Date  : 4/14/2021 5:58:21 PM      Inj :  1
                                                Inj Volume : 0.5 µl
Acq. Method    : C:\CHEM32\1\METHODS\HOLZ 0.5.M
Last changed   : 4/14/2021 11:19:56 AM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\HOLZ 0.5.M
Last changed   : 4/15/2021 11:16:29 AM by Analytik
                                         (modified after loading)
Method Info    : Chiraldak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



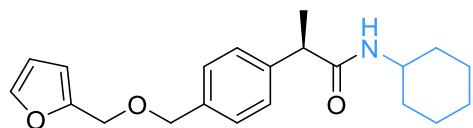
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area %
#	[min]		[min]	[mAU*s]	[mAU]	
1	11.764	BB	0.2114	325.37738	23.52243	6.5601
2	13.987	BB	0.2543	4634.57861	281.16498	93.4399

Totals : 4959.95599 304.68741



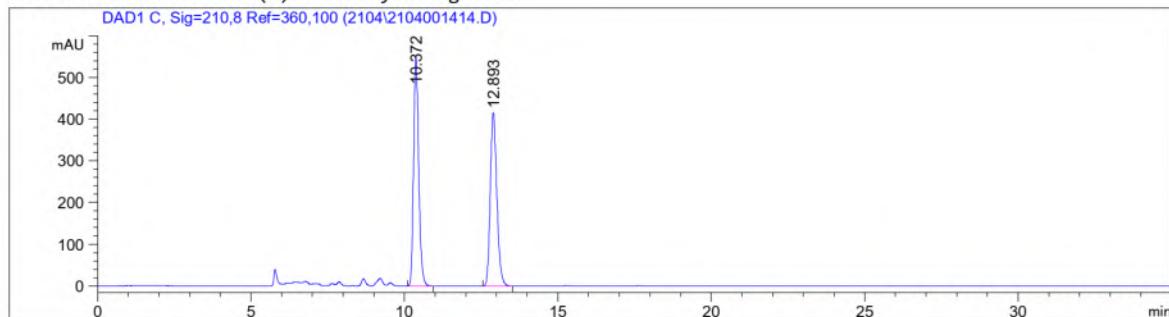
(R)-N-Cyclohexyl-2-(4-((furan-2-ylmethoxy)methyl)phenyl)propanamide (31)

Data File D:\CHEM32\1\DATA\2104\2104001414.D

Sample Name: YY-X-108 Race

```
=====
Acq. Operator   : Analytik          Seq. Line :  5
Acq. Instrument : LC5             Location  : Vial 12
Injection Date  : 4/14/2021 7:30:28 PM      Inj       : 1
                                                Inj Volume : 0.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\HOLZ 0.5.M
Last changed    : 4/14/2021 11:19:56 AM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\HOLZ 0.5.M
Last changed    : 4/15/2021 11:16:29 AM by Analytik
                                (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.372	BB	0.1767	6303.33350	546.02582	50.4632
2	12.893	BB	0.2294	6187.61523	416.30707	49.5368

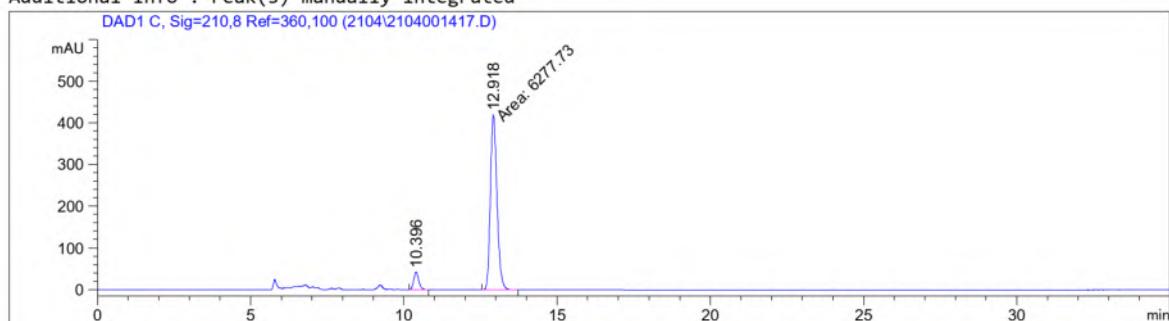
Totals : 1.24909e4 962.33289

Data File D:\CHEM32\1\DATA\2104\2104001417.D

Sample Name: YY-X-108 C

```
=====
Acq. Operator   : Analytik          Seq. Line : 6
Acq. Instrument : LC5             Location  : Vial 22
Injection Date  : 4/14/2021 9:48:40 PM      Inj : 2
                                                Inj Volume : 0.5 µl
Acq. Method    : C:\CHEM32\1\METHODS\HOLZ 0.5.M
Last changed   : 4/14/2021 11:19:56 AM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\HOLZ 0.5.M
Last changed   : 4/15/2021 11:16:29 AM by Analytik
                                                (modified after loading)
Method Info    : Chiraldak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



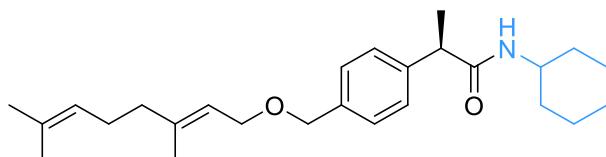
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.396	BB	0.1769	491.24588	42.47208	7.2573
2	12.918	MM	0.2493	6277.73389	419.64307	92.7427

Totals : 6768.97977 462.11514

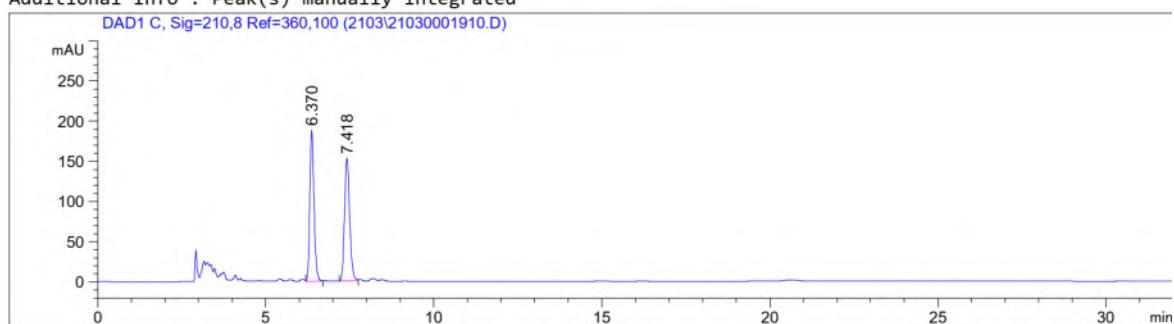


(R)-N-Cyclohexyl-2-(4-(((3,7-dimethylocta-2,6-dien-1-yl)oxy)methyl)phenyl)propanamide (32)

Data File D:\CHEM32\1\DATA\2103\21030001910.D
Sample Name: YY-X-87 Race

```
=====
Acq. Operator   : Analytik          Seq. Line :  2
Acq. Instrument : LC5             Location  : Vial 36
Injection Date  : 3/19/2021 4:10:45 PM    Inj       : 1
                                                Inj Volume : 0.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\TEST RAC.M
Last changed    : 3/19/2021 3:02:19 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\TEST RAC.M
Last changed    : 3/19/2021 3:59:49 PM by Analytik
                                                (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

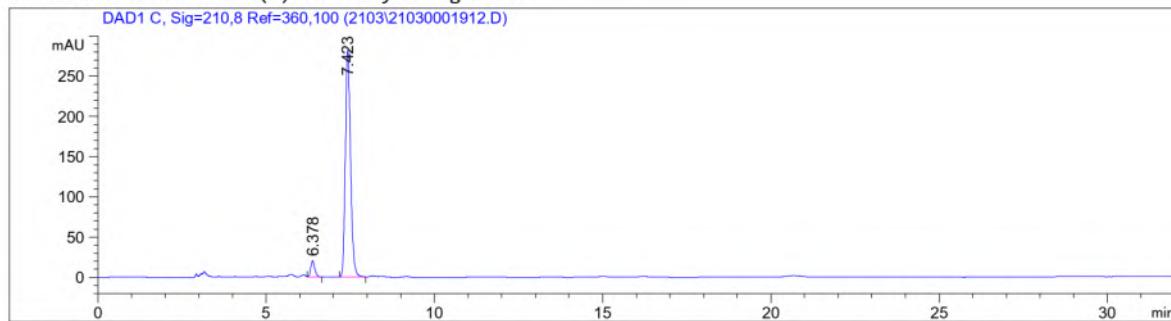
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.370	VB	0.1307	1625.06848	188.24072	50.0614
2	7.418	BB	0.1640	1621.08081	152.46329	49.9386
Totals :					3246.14929	340.70401

Data File D:\CHEM32\1\DATA\2103\21030001912.D

Sample Name: YY-X-87 C

```
=====
Acq. Operator   : Analytik          Seq. Line : 3
Acq. Instrument : LC5             Location  : Vial 46
Injection Date  : 3/19/2021 5:18:59 PM      Inj : 1
                                                Inj Volume : 0.5 µl
Acq. Method    : C:\CHEM32\1\METHODS\TEST RAC.M
Last changed    : 3/19/2021 3:02:19 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\TEST RAC.M
Last changed    : 3/19/2021 3:59:49 PM by Analytik
                                                (modified after loading)
Method Info     : Chiralpak AD-H, Hept./EtOH 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated



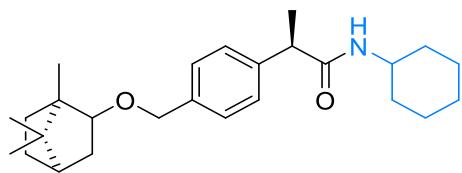
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area %
#	[min]		[min]	[mAU*s]	[mAU]	
1	6.378	VB	0.1301	172.23213	20.08257	5.3616
2	7.423	BB	0.1659	3040.10889	281.81577	94.6384

Totals : 3212.34102 301.89834



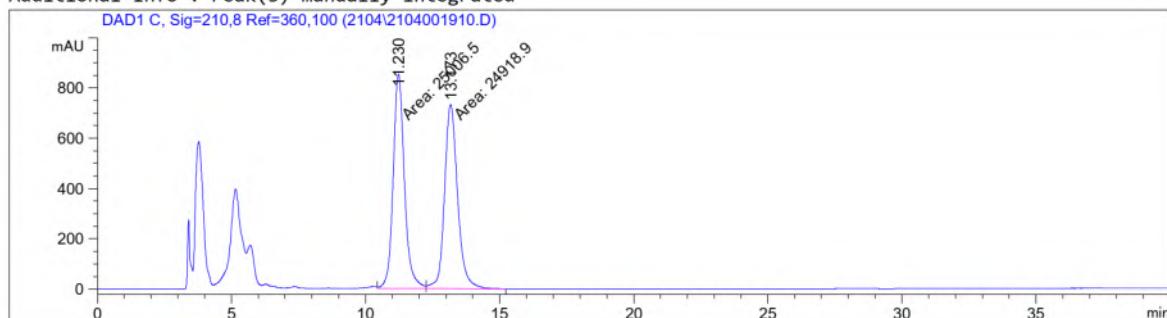
(2*R*)-*N*-cyclohexyl-2-(4-(((1*S*,4*S*)-1,7,7-trimethylbicyclo[2.2.1]heptan-2-yl)oxy)methyl)phenylpropanamide (33)

Data File D:\CHEM32\1\DATA\2104\2104001910.D

Sample Name: YY-X-99 Race

```
=====
Acq. Operator   : Analytik           Seq. Line : 7
Acq. Instrument : LC5               Location  : Vial 13
Injection Date  : 4/19/2021 6:34:49 PM    Inj       : 1
                                         Inj Volume : 2.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY 0.5.M
Last changed    : 4/19/2021 4:10:16 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY-0.5.M
Last changed    : 4/20/2021 9:16:17 AM by Analytik
                  (modified after loading)
Method Info     : Cellulose 4, Hept./Isoprop. 95:5, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.230	MF	0.4888	2.50065e4	852.73309	50.0878
2	13.173	FM	0.5672	2.49189e4	732.23187	49.9122

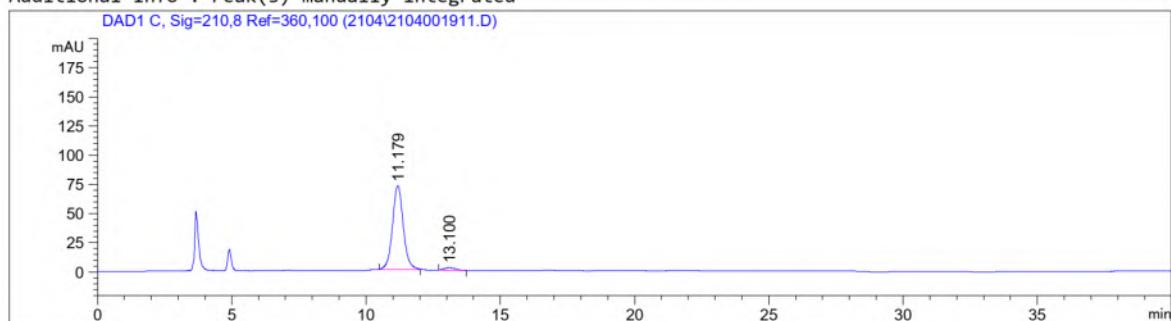
Totals : 4.99254e4 1584.96497

Data File D:\CHEM32\1\DATA\2104\2104001911.D

Sample Name: YY-X-99 -30

```
=====
Acq. Operator   : Analytik          Seq. Line :  8
Acq. Instrument : LC5             Location  : Vial 14
Injection Date  : 4/19/2021 7:20:59 PM      Inj :  1
                                                Inj Volume : 2.5 µl
Acq. Method    : C:\CHEM32\1\METHODS\YY 0.5.M
Last changed    : 4/19/2021 4:10:16 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY-0,5.M
Last changed    : 4/20/2021 9:18:11 AM by Analytik
                  (modified after loading)
Method Info     : Cellulose 4, Hept./Isoprop. 95:5, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

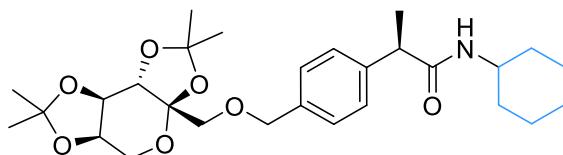
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.179	BB	0.4282	2013.56116	72.00964	97.1110
2	13.100	BB	0.3655	59.90166	2.08332	2.8890

Totals : 2073.46281 74.09295

LC5 4/20/2021 9:19:16 AM Analytik

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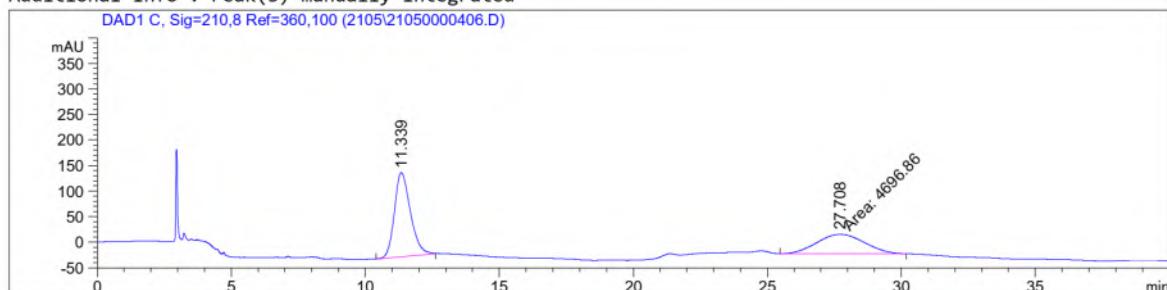
(2R)-N-Cyclohexyl-2-(4-(((5aR,8aR,8bS)-2,2,7,7-tetramethyltetrahydro-3aH-bis([1,3]dioxolo)[4,5-b:4',5'-d]pyran-3a-yl)methoxy)methyl)phenyl)propanamide (34)

Data File D:\CHEM32\1\DATA\2105\21050000406.D

Sample Name: YY-X-120 Race

```
=====
Acq. Operator   : Analytik          Seq. Line : 5
Acq. Instrument : LCS             Location : Vial 4
Injection Date  : 5/4/2021 1:23:00 PM Inj : 1
                           Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/4/2021 10:03:07 AM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/5/2021 9:00:26 AM by Analytik
                           (modified after loading)
Method Info     : Chiralcel OJ, Hept./EtOH 95:5, 1.0ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

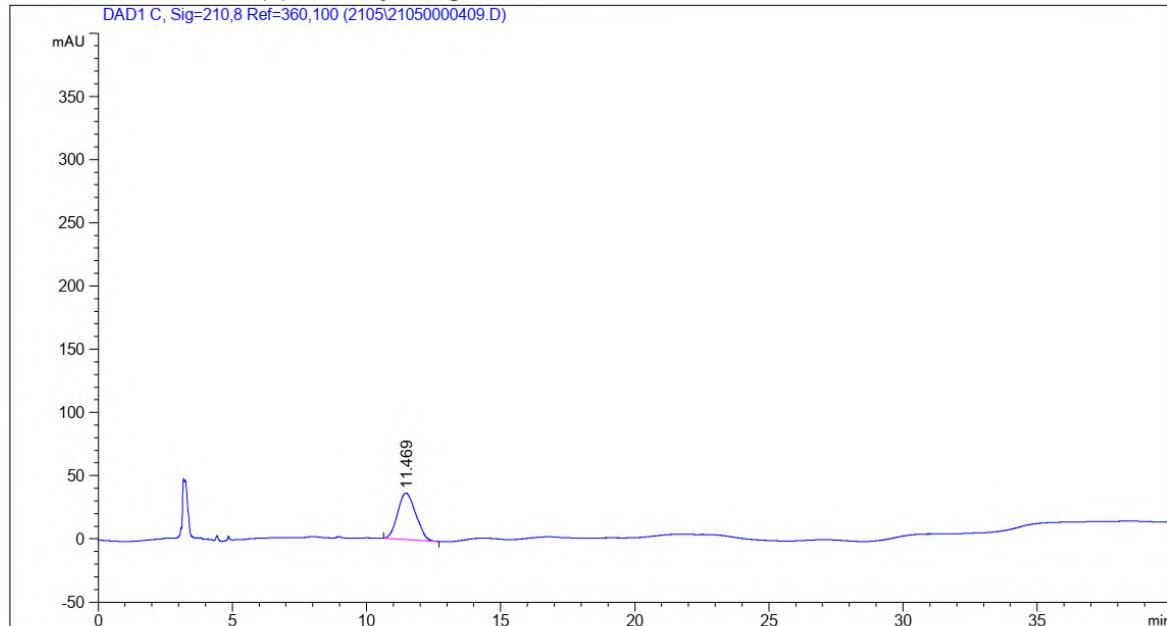
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.339	BB	0.6303	6831.67432	164.56416	59.2588
2	27.708	MM	2.0619	4696.86230	37.96592	40.7412

Totals : 1.15285e4 202.53009

Data File D:\CHEM32\1\DATA\2105\21050000409.D
Sample Name: YY-X-120-30

```
=====
Acq. Operator   : Analytik          Seq. Line :   8
Acq. Instrument : LC5             Location  : Vial 44
Injection Date  : 5/4/2021 4:22:31 PM      Inj       : 1
                                         Inj Volume : 1.0 µl
Different Inj Volume from Sequence !    Actual Inj Volume : 5.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/4/2021 4:22:07 PM by Analytik
                           (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/5/2021 9:00:26 AM by Analytik
                           (modified after loading)
Method Info     : Chiralcel OJ, Hept./EtOH 95:5, 1.0ml/min
```

Additional Info : Peak(s) manually integrated



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

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

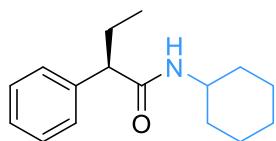
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area %
#	[min]		[min]	[mAU*s]	[mAU]	
1	11.469	BB	0.7228	1718.49609	36.79789	100.0000

Totals : 1718.49609 36.79789

LC5 5/5/2021 9:16:55 AM Analytik

Page 1 of 2



(R)-N-Cyclohexyl-2-phenylbutanamide (35)

from trans- β -methylstyrene

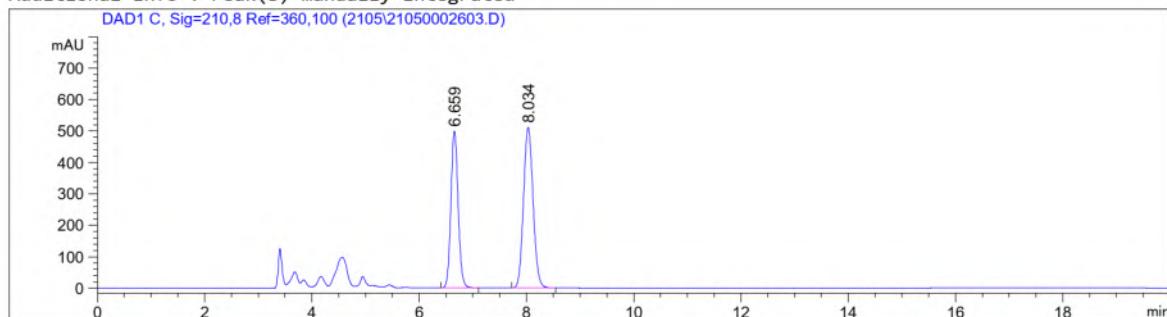
*Due to the use of ligand mixture [(R,R)-Ph-BPE & (S,S)-Ph-BPE] as the racemic ligand, the ratio of enantiomeric isomer in this case is not 50:50 er.

Data File D:\CHEM32\1\DATA\2105\21050002603.D

Sample Name: YY-X-129 Race

```
=====
Acq. Operator   : Analytik          Seq. Line :  4
Acq. Instrument : LCS             Location  : Vial 1
Injection Date  : 5/26/2021 4:47:12 PM    Inj       : 1
                                         Inj Volume : 1.0 μl
Acq. Method     : C:\CHEM32\1\METHODS\AMYLOSE2-0.5.M
Last changed    : 5/26/2021 4:26:29 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\AMYLOSE2-0.5.M
Last changed    : 5/26/2021 4:16:50 PM by Analytik
                  (modified after loading)
Method Info     : Amylose2, Hept./EtOH 90:10, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



=====

Area Percent Report

=====

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.659	BB	0.1450	4665.87354	499.49747	42.4692
2	8.034	BB	0.1925	6320.62500	510.16016	57.5308

Totals : 1.09865e4 1009.65762

LC5 5/27/2021 9:30:30 AM Analytik

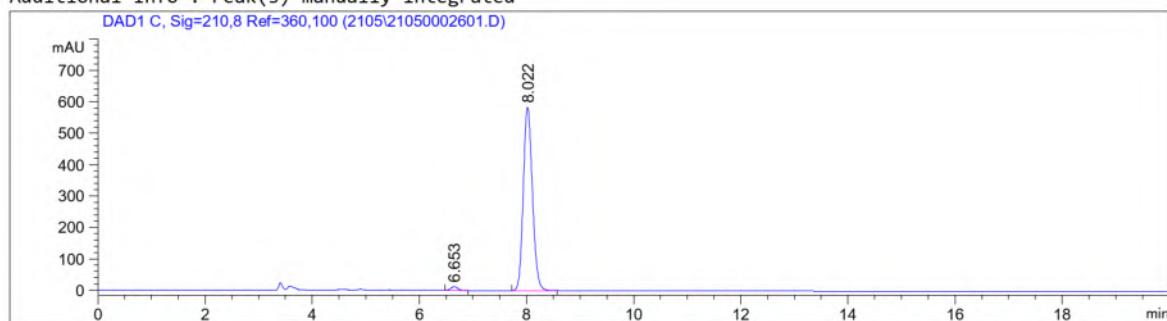
Page 1 of 2

Data File D:\CHEM32\1\DATA\2105\21050002601.D

Sample Name: YY-X-127-C

```
=====
Acq. Operator   : Analytik          Seq. Line :  2
Acq. Instrument : LC5             Location  : Vial 2
Injection Date  : 5/26/2021 3:50:05 PM    Inj :  1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\AMYLOSE2-0.5.M
Last changed    : 5/26/2021 3:49:41 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\AMYLOSE2-0.5.M
Last changed    : 5/26/2021 4:16:50 PM by Analytik
                  (modified after loading)
Method Info     : Amylose2, Hept./EtOH 90:10, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.653	BB	0.1439	114.01968	12.32931	1.6266
2	8.022	BB	0.1841	6895.87354	582.56409	98.3734

Totals : 7009.89321 594.89340

LC5 5/26/2021 4:18:11 PM Analytik

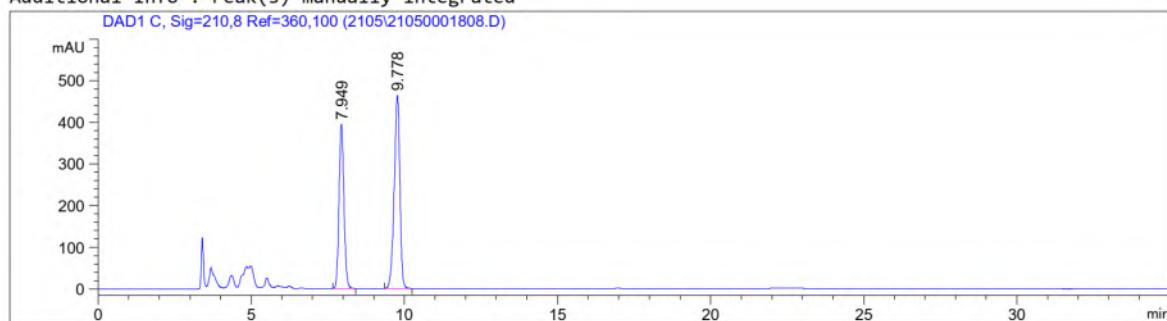
Page 1 of 2

from *cis*- β -methylstyrene

Data File D:\CHEM32\1\DATA\2105\21050001808.D
Sample Name: YY-X-129 Race

```
=====
Acq. Operator   : Analytik          Seq. Line :   6
Acq. Instrument : LC5              Location  : Vial 11
Injection Date  : 5/18/2021 1:46:04 PM      Inj :   1
                                         Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\AMYLOSE2-0.5.M
Last changed    : 5/18/2021 12:46:25 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\AMYLOSE2.M
Last changed    : 5/25/2021 3:04:44 PM by Analytik
                  (modified after loading)
Method Info     : Amylose2, Hept./EtOH 90:10, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



Area Percent Report

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area %
#	[min]		[min]	[mAU*s]	[mAU]	
1	7.949	BB	0.1721	4397.89844	394.51910	42.4912
2	9.778	BB	0.1957	5952.24609	463.77136	57.5088

Totals : 1.03501e4 858.29047

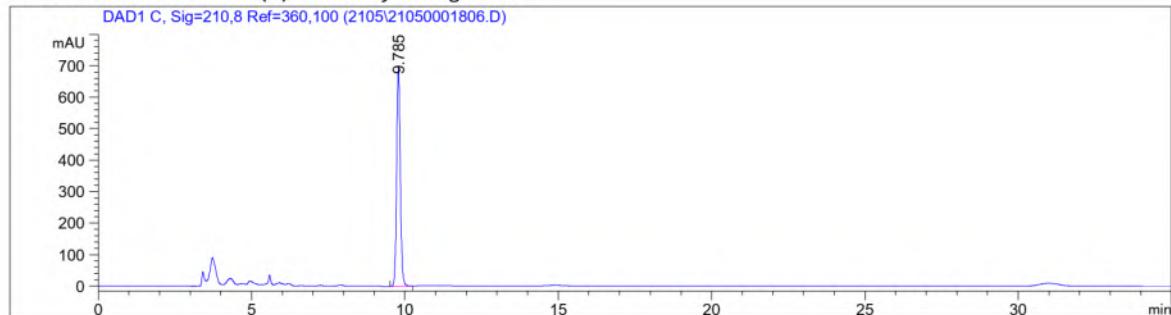
LC5 5/25/2021 3:31:18 PM Analytik

Page 1 of 2

Data File D:\CHEM32\1\DATA\2105\21050001806.D
Sample Name: YY-X-129- C

```
=====
Acq. Operator   : Analytik          Seq. Line : 4
Acq. Instrument : LC5             Location  : Vial 21
Injection Date  : 5/18/2021 12:33:54 PM    Inj       : 1
                                                Inj Volume : 1.0 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 3.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\AMYLOSE2-0.5.M
Last changed    : 5/18/2021 12:46:25 PM by Analytik
                           (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\AMYLOSE2.M
Last changed    : 5/25/2021 3:32:38 PM by Analytik
                           (modified after loading)
Method Info     : Amylose2, Hept./EtOH 90:10, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

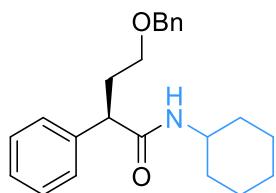
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area %
#	[min]		[min]	[mAU*s]	[mAU]	
1	9.785	BB	0.1265	5787.76855	699.40259	100.0000

Totals : 5787.76855 699.40259

LC5 5/25/2021 3:35:39 PM Analytik

Page 1 of 2



(R)-4-(Benzyl)-N-cyclohexyl-2-phenylbutanamide (36)

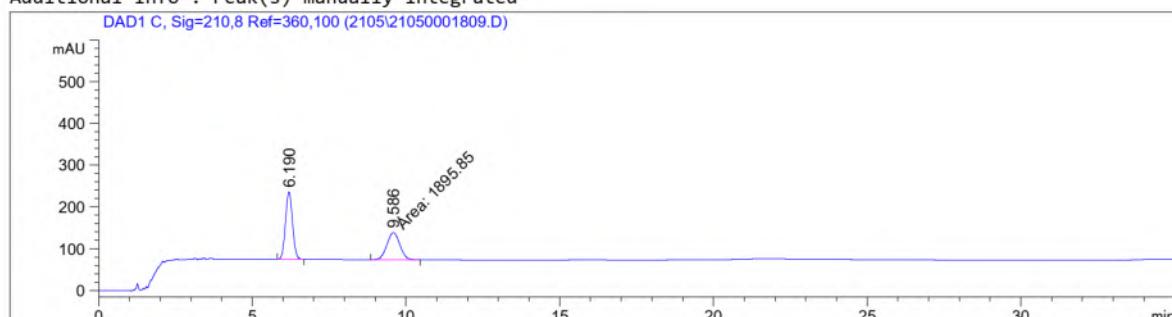
*Due to the use of ligand mixture [(R,R)-Ph-BPE & (S,S)-Ph-BPE] as the racemic ligand, the ratio of enantiomeric isomer in this case is not 50:50 er.

Data File D:\CHEM32\1\DATA\2105\21050001809.D

Sample Name: YY-X-131 Race

```
=====
Acq. Operator   : Analytik           Seq. Line : 7
Acq. Instrument : LC5               Location  : Vial 12
Injection Date  : 5/18/2021 2:22:13 PM    Inj       : 1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\AMYLOSE2.M
Last changed    : 5/17/2021 3:03:04 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\AMYLOSE2.M
Last changed    : 5/25/2021 3:04:44 PM by Analytik
                                (modified after loading)
Method Info     : Amylose2, Hept./EtOH 90:10, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

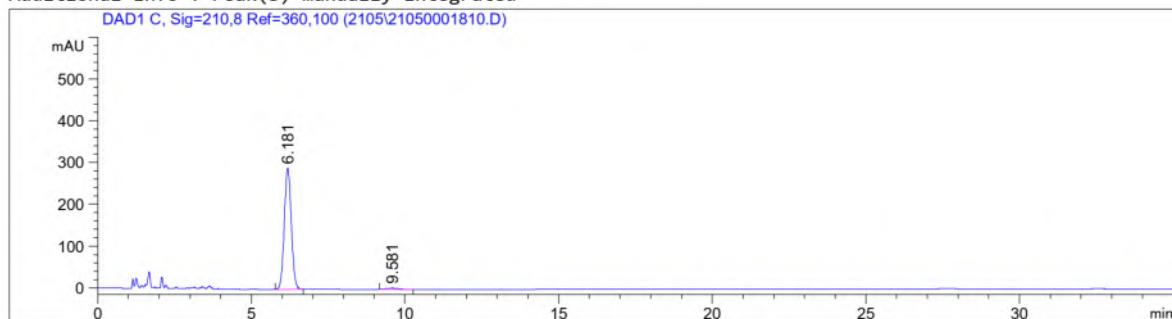
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.190	BB	0.2472	2558.36670	161.13013	57.4369
2	9.586	MM	0.4878	1895.85291	64.77711	42.5631
Totals :						4454.21960 225.90724

Data File D:\CHEM32\1\DATA\2105\21050001810.D

Sample Name: YY-X-131 C

```
=====
Acq. Operator   : Analytik          Seq. Line :  8
Acq. Instrument : LC5             Location  : Vial 22
Injection Date  : 5/18/2021 2:58:16 PM    Inj :  1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\AMYLOSE2.M
Last changed    : 5/17/2021 3:03:04 PM by Analytik
Analysis Method : C:\CHEM32\1\METHODS\AMYLOSE2.M
Last changed    : 5/25/2021 3:04:44 PM by Analytik
                                         (modified after loading)
Method Info     : Amylose2, Hept./EtOH 90:10, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



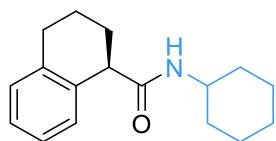
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area
#	[min]		[min]	[mAU*s]	[mAU]	%
1	6.181	BB	0.2466	4581.46045	289.61316	98.4100
2	9.581	BB	0.3748	74.02377	2.70918	1.5900

Totals : 4655.48421 292.32234



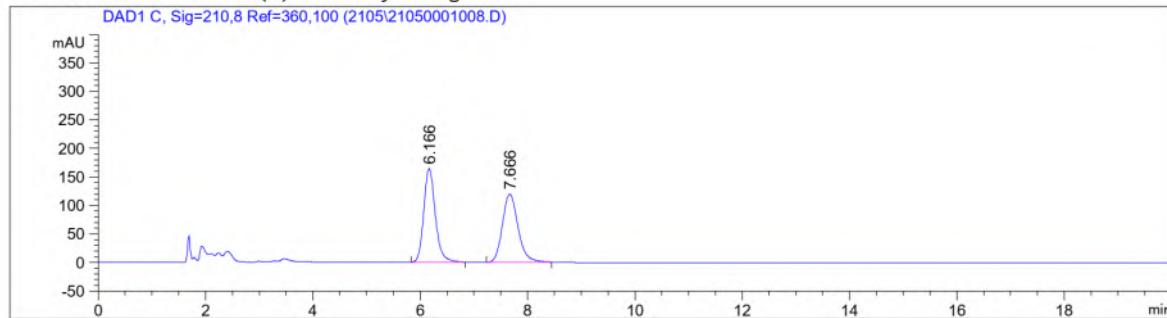
(R)-N-Cyclohexyl-1,2,3,4-tetrahydronaphthalene-1-carboxamide (37)

Data File D:\CHEM32\1\DATA\2105\21050001008.D

Sample Name: YY-X-130 Race

```
=====
Acq. Operator   : Analytik          Seq. Line :  9
Acq. Instrument : LC5             Location  : Vial 13
Injection Date  : 5/10/2021 4:33:33 PM      Inj :  1
                                         Inj Volume : 1.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/10/2021 2:26:33 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/10/2021 3:10:59 PM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area %
#	[min]		[min]	[mAU*s]	[mAU]	%
1	6.166	BB	0.2290	2468.32031	164.50887	50.4445
2	7.666	BB	0.3122	2424.81909	119.77151	49.5555

Totals : 4893.13940 284.28038

LC5 5/11/2021 12:36:45 PM Analytik

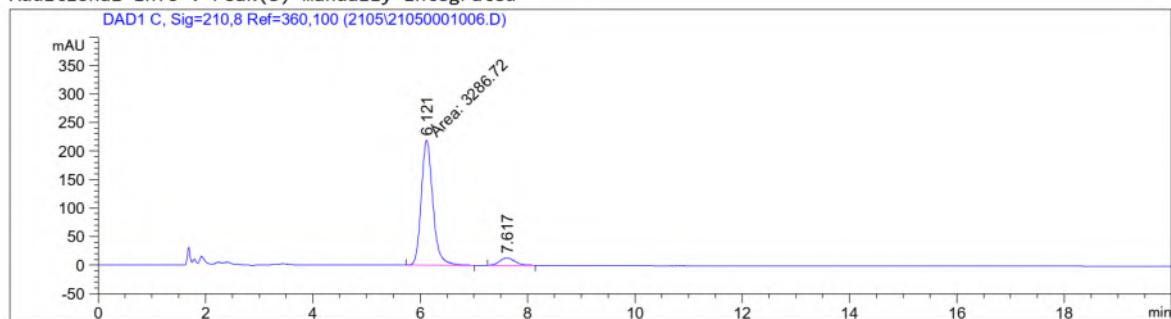
Page 1 of 2

Data File D:\CHEM32\1\DATA\2105\21050001006.D

Sample Name: YY-X-130-C

```
=====
Acq. Operator   : Analytik          Seq. Line :    7
Acq. Instrument : LC5             Location  : Vial 23
Injection Date  : 5/10/2021 3:51:23 PM      Inj :    1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/10/2021 2:26:33 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/10/2021 3:10:59 PM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

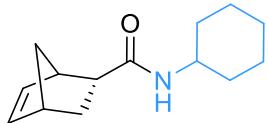
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.121	MM	0.2500	3286.71509	219.09990	92.7491
2	7.617	BB	0.3041	256.94687	13.02880	7.2509

Totals : 3543.66196 232.12870

LC5 5/11/2021 12:38:28 PM Analytik

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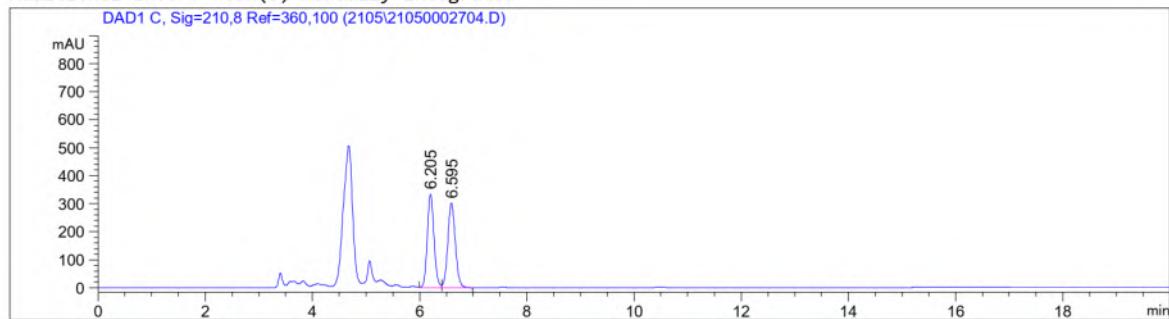
(4R)-N-Cyclohexylbicyclo[2.2.1]hept-5-ene-2-carboxamide (38)

Data File D:\CHEM32\1\DATA\2105\21050002704.D

Sample Name: YY-X-83 Race

```
=====
Acq. Operator   : Analytik           Seq. Line : 4
Acq. Instrument : LCS              Location : Vial 24
Injection Date : 5/27/2021 11:32:49 AM    Inj : 1
                                                Inj Volume : 1.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\AMYLOSE2-0.5.M
Last changed   : 5/27/2021 11:56:29 AM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\AMYLOSE2-0.5.M
Last changed   : 5/27/2021 11:11:12 AM by Analytik
                  (modified after loading)
Method Info    : Amylose2, Hept./EtOH 90:10, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



Area Percent Report

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area
#	[min]		[min]	[mAU*s]	[mAU]	%
1	6.205	VV	0.1336	2913.05566	334.43451	49.7039
2	6.595	VB	0.1517	2947.76514	302.45889	50.2961

Totals : 5860.82080 636.89340

LC5 5/27/2021 11:58:03 AM Analytik

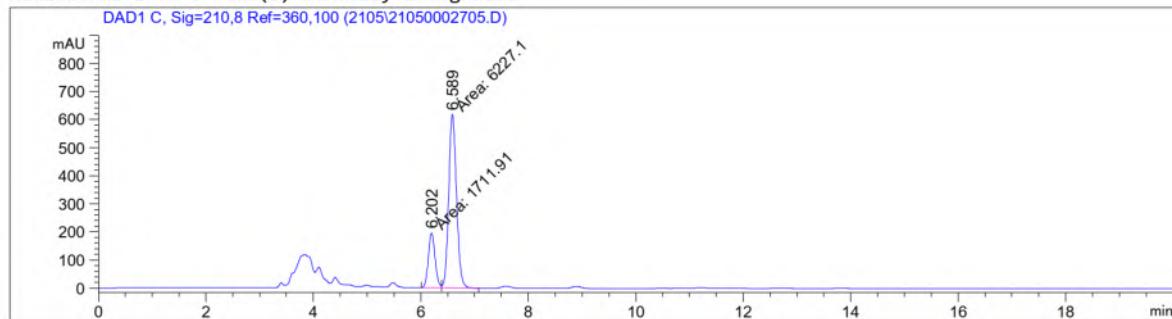
Page 1 of 2

Data File D:\CHEM32\1\DATA\2105\21050002705.D

Sample Name: YY-X-83 C

```
=====
Acq. Operator   : Analytik                     Seq. Line :  5
Acq. Instrument : LC5                         Location  : Vial 34
Injection Date  : 5/27/2021 11:56:52 AM          Inj       : 1
                                                Inj Volume : 1.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\AMYLOSE2-0.5.M
Last changed    : 5/27/2021 11:56:29 AM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\AMYLOSE2-0.5.M
Last changed    : 5/27/2021 11:11:12 AM by Analytik
                  (modified after loading)
Method Info     : Amylose2, Hept./EtOH 90:10, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

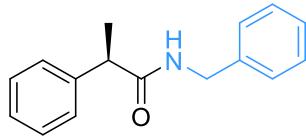
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.202	MF	0.1466	1711.90796	194.62668	21.5633
2	6.589	FM	0.1681	6227.09521	617.23059	78.4367

Totals : 7939.00317 811.85727

LC5 5/27/2021 12:36:03 PM Analytik

Page 1 of 2



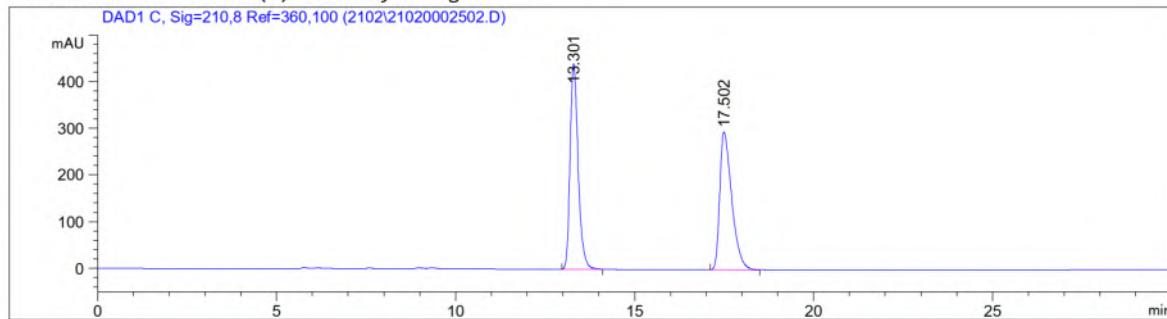
(R)-N-benzyl-2-phenylpropanamide (3)

Data File D:\CHEM32\1\DATA\2102\21020002502.D

Sample Name: YY-x-race

```
=====
Acq. Operator   : Analytik          Seq. Line :  3
Acq. Instrument : LC5             Location  : Vial 51
Injection Date  : 2/25/2021 10:56:38 AM    Inj :  1
                                         Inj Volume : 0.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 2/25/2021 11:41:45 AM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 2/25/2021 12:30:47 PM by Analytik
                  (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



Area Percent Report

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	13.301	BB	0.2385	6881.00146	439.86713	49.9153
2	17.502	BB	0.3547	6904.35840	295.35629	50.0847

Totals : 1.37854e4 735.22342

LC5 2/25/2021 12:31:07 PM Analytik

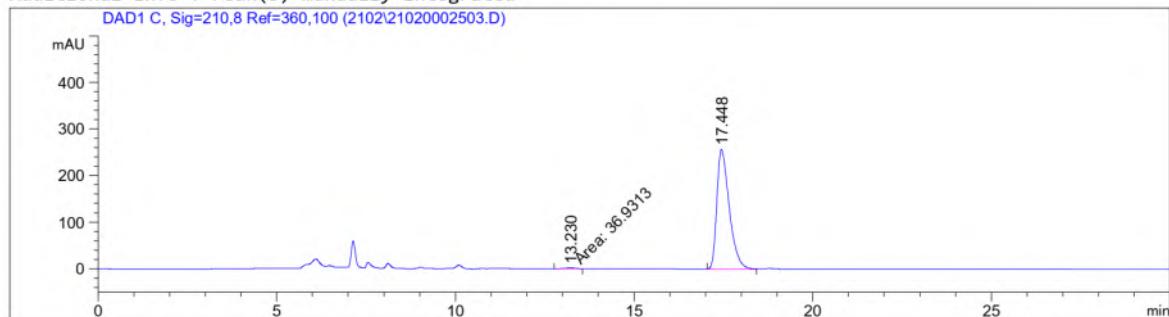
Page 1 of 2

Data File D:\CHEM32\1\DATA\2102\21020002503.D

Sample Name: YY-x-53

```
=====
Acq. Operator   : Analytik          Seq. Line :  4
Acq. Instrument : LC5             Location  : Vial 53
Injection Date  : 2/25/2021 11:42:09 AM    Inj       : 1
                                                Inj Volume : 0.5 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 3.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 2/25/2021 11:41:45 AM by Analytik
                           (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\HOLZ.M
Last changed    : 2/25/2021 12:30:47 PM by Analytik
                           (modified after loading)
Method Info     : Chiraldak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated

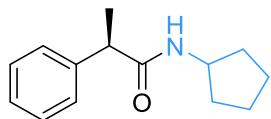


```
=====
                           Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210.8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	13.230	MM	0.3222	36.93130	1.91056	0.5913
2	17.448	BB	0.3662	6208.45605	256.65204	99.4087



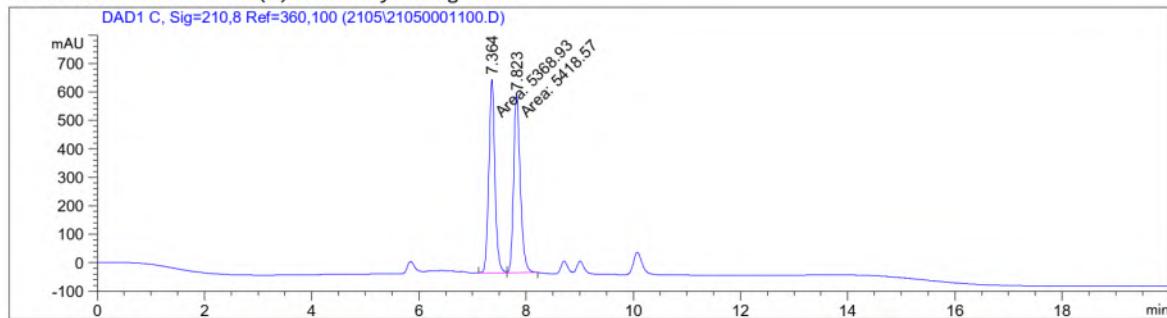
(R)-N-Cyclopentyl-2-phenylpropanamide (41)

Data File D:\CHEM32\1\DATA\2105\21050001100.D

Sample Name: YY-X-109 Race

```
=====
Acq. Operator   : Analytik                     Seq. Line :  1
Acq. Instrument : LCS                        Location : Vial 1
Injection Date  : 5/11/2021 2:42:53 PM          Inj :  1
                                                Inj Volume : 1.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\HOLZ 0.5.M
Last changed   : 5/11/2021 3:25:19 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\HOLZ 0.5.M
Last changed   : 5/12/2021 12:08:42 PM by Analytik
                  (modified after loading)
Method Info     : Chiralpak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



Area Percent Report

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.364	MF	0.1316	5368.93311	680.12775	49.7699
2	7.823	FM	0.1450	5418.57129	622.89972	50.2301

Totals : 1.07875e4 1303.02747

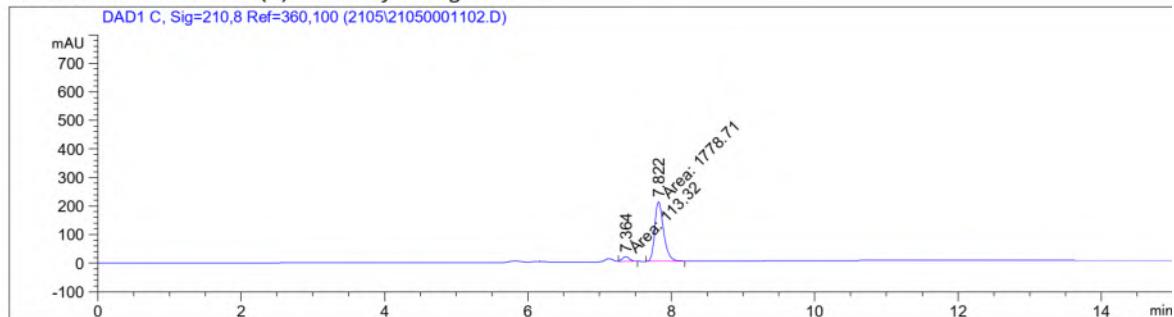
LC5 5/12/2021 12:10:18 PM Analytik

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Data File D:\CHEM32\1\DATA\2105\21050001102.D
Sample Name: YY-X-109-rt

```
=====
Acq. Operator   : Analytik           Seq. Line : 3
Acq. Instrument : LC5               Location  : Vial 2
Injection Date  : 5/11/2021 3:41:47 PM    Inj : 1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\HOLZ 0.5.M
Last changed    : 5/11/2021 3:25:19 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\HOLZ 0.5.M
Last changed    : 5/12/2021 12:11:18 PM by Analytik
                  (modified after loading)
Method Info     : Chiralpak AD-H, Hept./EtOH 80:20, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



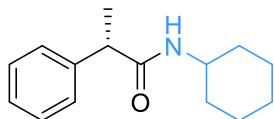
```
=====
          Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area %
#	[min]		[min]	[mAU*s]	[mAU]	
1	7.364	MM	0.1145	113.32011	16.49935	5.9893
2	7.822	MM	0.1424	1778.70959	208.19086	94.0107

Totals : 1892.02970 224.69020



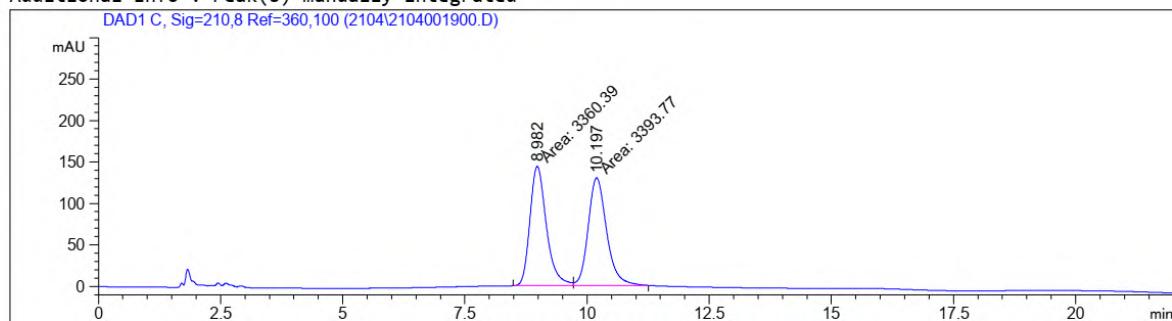
(S)-N-Cyclohexyl-2-phenylpropanamide (42)

Data File D:\CHEM32\1\DATA\2104\2104001900.D

Sample Name: YY-X-59 Race

```
=====
Acq. Operator   : Analytik           Seq. Line : 1
Acq. Instrument : LC5             Location : Vial 1
Injection Date  : 4/19/2021 12:11:17 PM      Inj : 1
                                                Inj Volume : 1.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\YY.M
Last changed   : 4/19/2021 12:39:33 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed   : 4/19/2021 3:57:24 PM by Analytik
                  (modified after loading)
Method Info     : Cellulse4, Hept./Isoprop 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated



Area Percent Report

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

Signal 1: DAD1 C, Sig=210.8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.982	MF	0.3893	3360.39331	143.85008	49.7529
2	10.197	FM	0.4359	3393.76636	129.77121	50.2471

Totals : 6754.15967 273.62129

LC5 4/19/2021 3:58:31 PM Analytik

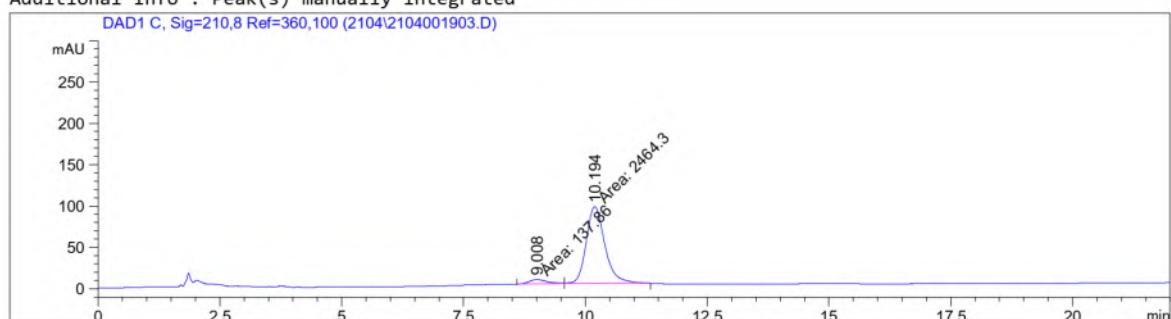
Page 1 of 2

Data File D:\CHEM32\1\DATA\2104\2104001903.D

Sample Name: YY-X-59 S

```
=====
Acq. Operator   : Analytik          Seq. Line :  2
Acq. Instrument : LC5             Location  : Vial 2
Injection Date  : 4/19/2021 1:32:05 PM      Inj :  2
                                                Inj Volume : 1.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\YY.M
Last changed   : 4/19/2021 12:39:33 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed   : 4/19/2021 3:57:24 PM by Analytik
                  (modified after loading)
Method Info    : Cellulse4, Hept./Isoprop 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

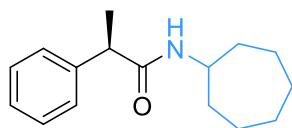
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.008	MF	0.4393	137.85976	5.22978	5.2979
2	10.194	FM	0.4421	2464.29565	92.90602	94.7021

Totals : 2602.15541 98.13580

LC5 4/19/2021 4:04:26 PM Analytik

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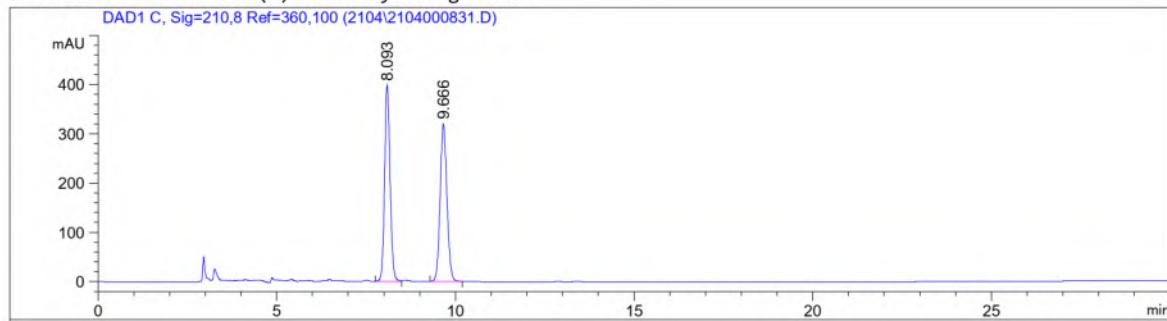
(R)-N-Cycloheptyl-2-phenylpropanamide (43)

Data File D:\CHEM32\1\DATA\2104\2104000831.D

Sample Name: YY-X-110 Race

```
=====
Acq. Operator   : Analytik                     Seq. Line : 22
Acq. Instrument : LC5                         Location : Vial 19
Injection Date  : 4/9/2021 8:58:04 AM           Inj : 1
                                                Inj Volume : 1.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\CHIRAL OJ-H.M
Last changed    : 4/9/2021 7:24:13 AM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\CHIRAL OJ-H.M
Last changed    : 4/9/2021 12:12:40 PM by Analytik
                  (modified after loading)
Method Info     : Chiralcel OJ-H , Hept./EtOH 95:5 , 1.0 ml/min
```

Additional Info : Peak(s) manually integrated



Area Percent Report

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.093	BB	0.1648	4254.18652	397.81912	50.1203
2	9.666	BB	0.2045	4233.75977	319.66937	49.8797

Totals : 8487.94629 717.48849

LC5 4/9/2021 12:13:17 PM Analytik

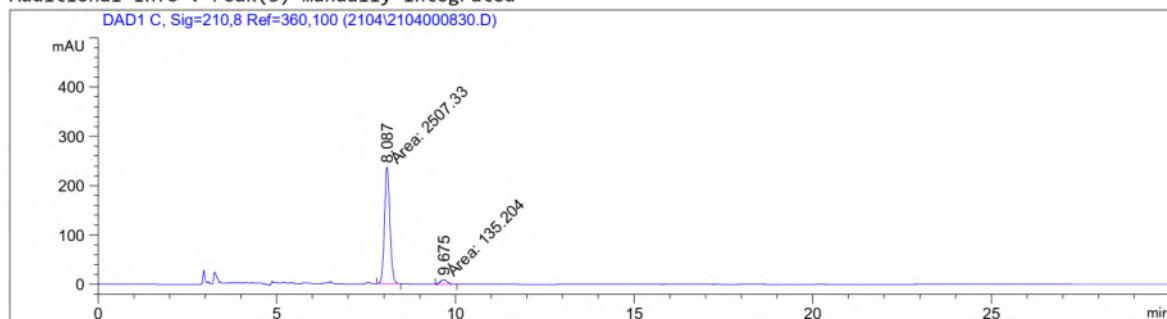
Page 1 of 2

Data File D:\CHEM32\1\DATA\2104\2104000830.D

Sample Name: YY-X-110 C

```
=====
Acq. Operator   : Analytik          Seq. Line : 21
Acq. Instrument : LC5             Location  : Vial 53
Injection Date  : 4/9/2021 8:26:54 AM    Inj       : 2
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\CHIRAL OJ-H.M
Last changed    : 4/9/2021 7:24:13 AM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\CHIRAL OJ-H.M
Last changed    : 4/9/2021 12:12:40 PM by Analytik
                  (modified after loading)
Method Info     : Chiralcel OJ-H , Hept./EtOH 95:5 , 1.0 ml/min
```

Additional Info : Peak(s) manually integrated



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

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

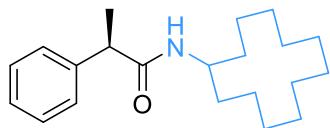
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width MM	Area [mAU*s]	Height [mAU]	Area %
1	8.087	MM	0.1766	2507.33276	236.58972	94.8835
2	9.675	MM	0.2309	135.20445	9.75895	5.1165

Totals : 2642.53722 246.34867

LC5 4/9/2021 12:16:56 PM Analytik

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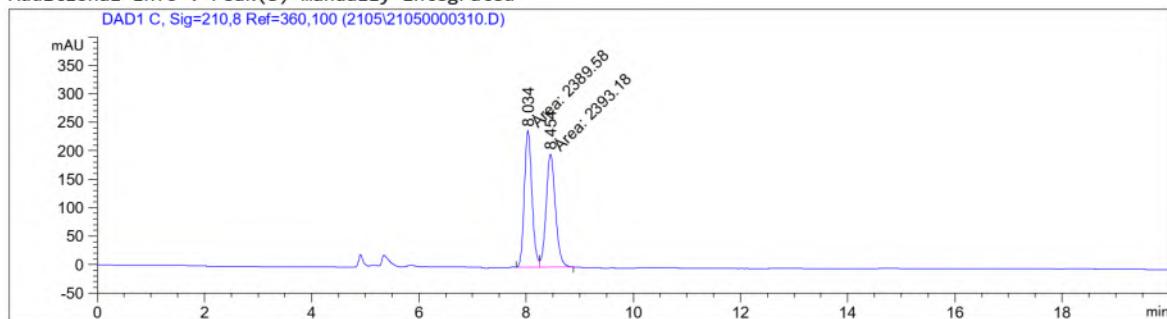
(R)-N-Cyclododecyl-2-phenylpropanamide (44)

Data File D:\CHEM32\1\DATA\2105\21050000310.D

Sample Name: YY-X-112 Race

```
=====
Acq. Operator   : Analytik          Seq. Line : 10
Acq. Instrument : LC5             Location : Vial 5
Injection Date  : 5/3/2021 3:56:06 PM    Inj : 1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\SDC 0.6.M
Last changed    : 5/3/2021 3:13:28 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\SDC 0.6.M
Last changed    : 5/4/2021 9:21:06 AM by Analytik
                  (modified after loading)
Method Info     : Chiralcel OJ, Hept./EtOH 95:5, 0.6ml/min
```

Additional Info : Peak(s) manually integrated



Area Percent Report

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.034	MF	0.1659	2389.58105	240.03227	49.9624
2	8.454	FM	0.2009	2393.17578	198.51573	50.0376

Totals : 4782.75684 438.54800

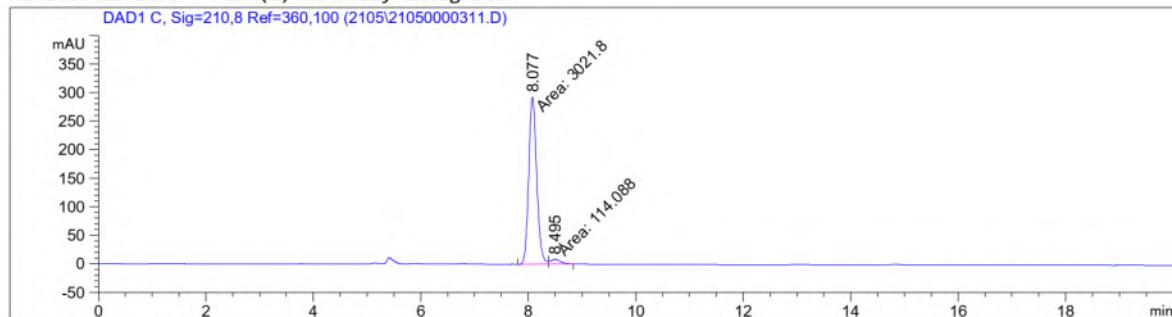
LC5 5/4/2021 9:25:06 AM Analytik

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Data File D:\CHEM32\1\DATA\2105\21050000311.D
Sample Name: YY-X-112 C

```
=====
Acq. Operator   : Analytik          Seq. Line : 11
Acq. Instrument : LC5             Location  : Vial 15
Injection Date  : 5/3/2021 4:17:15 PM    Inj       : 1
                                         Inj Volume : 1.0 µl
Different Inj Volume from Sequence !    Actual Inj Volume : 0.5 µl
Acq. Method     : C:\CHEM32\1\METHODS\SDC 0.6.M
Last changed    : 5/3/2021 3:13:28 PM by Analytik
                           (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\SDC 0.6.M
Last changed    : 5/4/2021 9:21:06 AM by Analytik
                           (modified after loading)
Method Info     : Chiralcel OJ, Hept./EtOH 95:5, 0.6ml/min
```

Additional Info : Peak(s) manually integrated

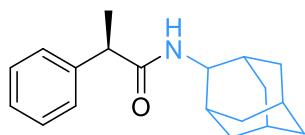


```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.077	MF	0.1719	3021.80005	292.93423	96.3619
2	8.495	FM	0.2190	114.08807	8.68151	3.6381



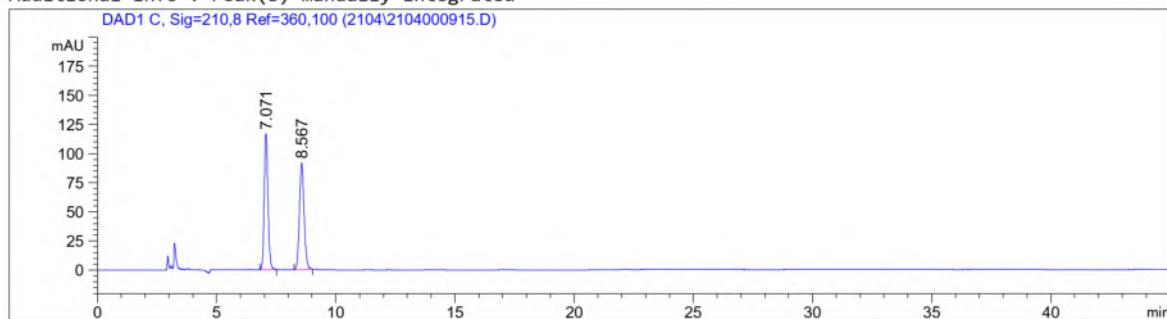
(R)-N-(Adamantan-2-yl)-2-phenylpropanamide (45)

Data File D:\CHEM32\1\DATA\2104\2104000915.D

Sample Name: YY-X-111 Race

```
=====
Acq. Operator   : Analytik          Seq. Line : 12
Acq. Instrument : LC5             Location  : Vial 17
Injection Date  : 4/10/2021 2:27:24 AM    Inj       : 1
                                                Inj Volume : 1.0 µl
Method          : C:\CHEM32\1\METHODS\CHIRAL OJ-H.M
Last changed    : 4/9/2021 2:37:18 PM by Analytik
Method Info     : Chiralcel OJ-H , Hept./EtOH 95:5 , 1.0 ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.071	BB	0.1684	1260.41406	116.35311	50.1350
2	8.567	BB	0.2123	1253.62427	91.25807	49.8650

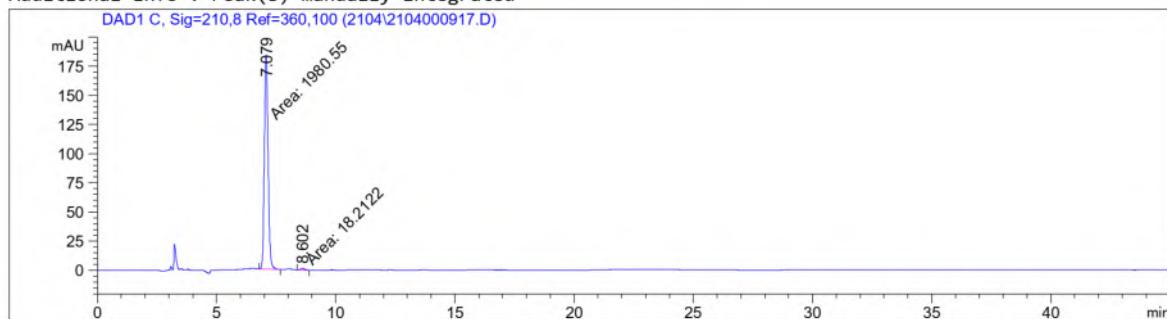
Totals : 2514.03833 207.61118

Data File D:\CHEM32\1\DATA\2104\2104000917.D

Sample Name: YY-X-111 C

```
=====
Acq. Operator   : Analytik          Seq. Line : 13
Acq. Instrument : LC5             Location  : Vial 18
Injection Date  : 4/10/2021 3:59:34 AM    Inj       : 1
                                         Inj Volume : 1.0 µl
Method          : C:\CHEM32\1\METHODS\CHIRAL OJ-H.M
Last changed    : 4/9/2021 2:37:18 PM by Analytik
Method Info     : Chiralcel OJ-H , Hept./EtOH 95:5 , 1.0 ml/min
```

Additional Info : Peak(s) manually integrated

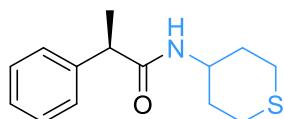


```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area %
#	[min]		[min]	[mAU*s]	[mAU]	
1	7.079	MM	0.1799	1980.55469	183.48459	99.0888
2	8.602	MM	0.2110	18.21221	1.43874	0.9112
Totals :						
				1998.76690	184.92333	



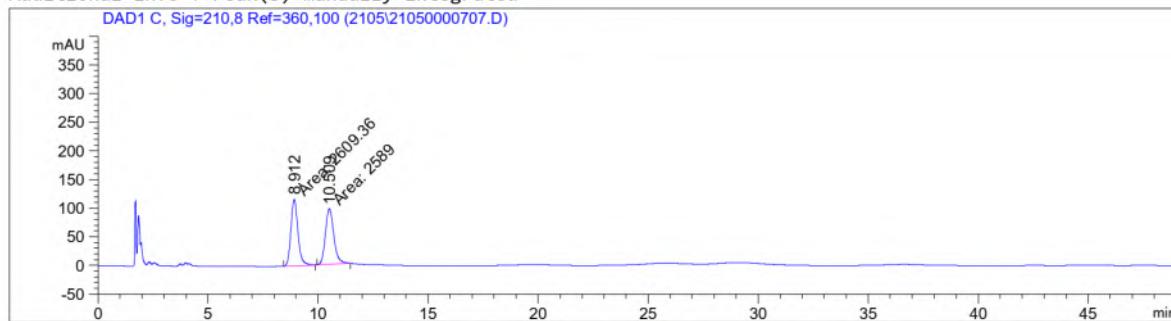
(R)-2-phenyl-N-(tetrahydro-2H-thiopyran-4-yl)propanamide (46)

Data File D:\CHEM32\1\DATA\2105\21050000707.D

Sample Name: YY-X-115-Race

```
=====
Acq. Operator   : Analytik           Seq. Line :  8
Acq. Instrument : LC5               Location  : Vial 23
Injection Date  : 5/7/2021 5:01:28 PM      Inj       : 1
                                                Inj Volume : 1.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/7/2021 3:44:32 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\SDC 1.2.M
Last changed    : 5/10/2021 11:54:58 AM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.912	MM	0.3747	2609.35938	116.07904	50.1958
2	10.509	MM	0.4453	2589.00293	96.90015	49.8042

Totals : 5198.36230 212.97919

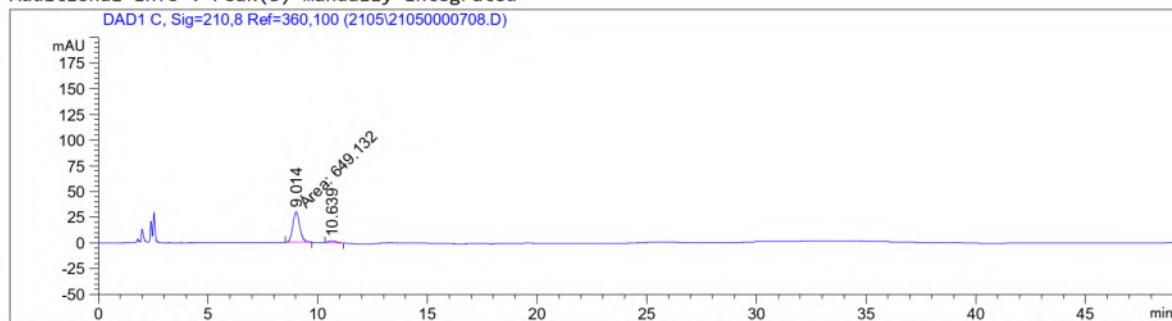
LC5 5/10/2021 12:27:16 PM Analytik

Page 1 of 2

Data File D:\CHEM32\1\DATA\2105\21050000708.D
Sample Name: YY-X-115-rt

```
=====
Acq. Operator   : Analytik           Seq. Line : 9
Acq. Instrument : LC5               Location  : Vial 24
Injection Date  : 5/7/2021 6:02:34 PM    Inj       : 1
                                         Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/7/2021 3:44:32 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\SDC 1.2.M
Last changed    : 5/10/2021 12:27:50 PM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./EtOH 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

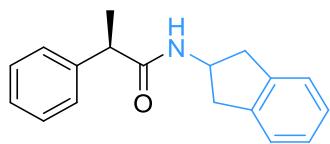
Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.014	MM	0.3655	649.13165	29.60283	93.4663
2	10.639	BB	0.3298	45.37730	1.74266	6.5337

Totals : 694.50896 31.34549

LC5 5/10/2021 12:29:10 PM Analytik

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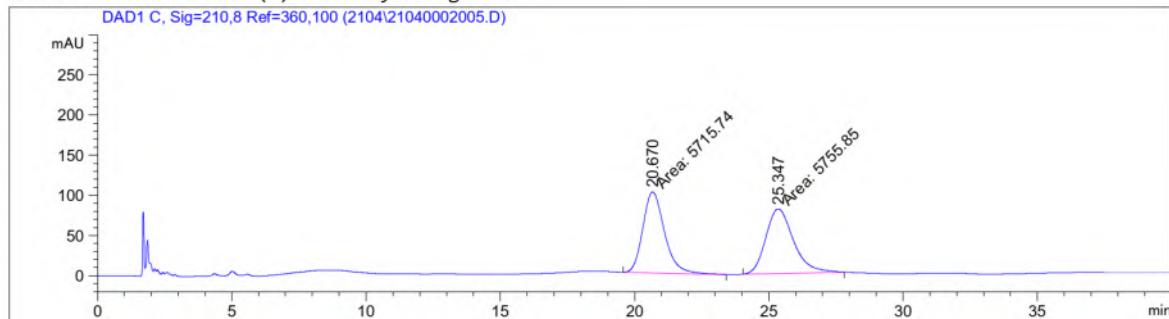
(R)-N-(2,3-Dihydro-1H-inden-2-yl)-2-phenylpropanamide (47)

Data File D:\CHEM32\1\DATA\2104\21040002005.D

Sample Name: YY-X-114 Race

```
=====
Acq. Operator   : Analytik          Seq. Line : 3
Acq. Instrument : LC5             Location  : Vial 22
Injection Date  : 4/20/2021 2:30:26 PM      Inj : 2
                                         Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\YY.M
Last changed    : 4/20/2021 2:27:56 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed    : 4/21/2021 8:56:35 AM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./Isoprop 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	20.670	MM	0.9431	5715.73584	101.01209	49.8251
2	25.347	MM	1.1921	5755.85498	80.47052	50.1749

Totals : 1.14716e4 181.48261

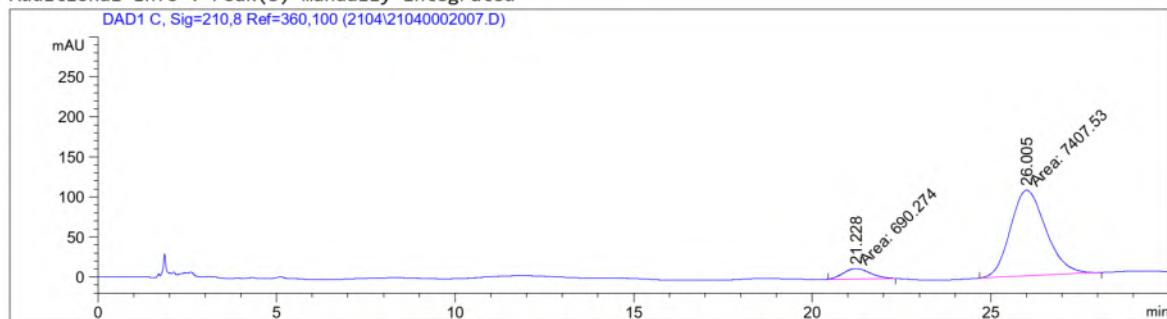
LC5 4/21/2021 8:57:46 AM Analytik

Page 1 of 2

Data File D:\CHEM32\1\DATA\2104\21040002007.D
Sample Name: YY-X-114 C

```
=====
Acq. Operator   : Analytik          Seq. Line : 4
Acq. Instrument : LC5             Location  : Vial 32
Injection Date  : 4/20/2021 3:42:37 PM      Inj : 2
                                         Inj Volume : 1.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\YY.M
Last changed    : 4/20/2021 3:52:53 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\YY.M
Last changed    : 4/21/2021 8:58:19 AM by Analytik
                  (modified after loading)
Method Info     : Cellulose4, Hept./Isoprop 95:5, 1ml/min
```

Additional Info : Peak(s) manually integrated



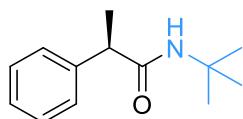
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	21.228	MM	0.8752	690.27374	13.14580	8.5242
2	26.005	MM	1.1516	7407.52930	107.20650	91.4758

Totals : 8097.80304 120.35230



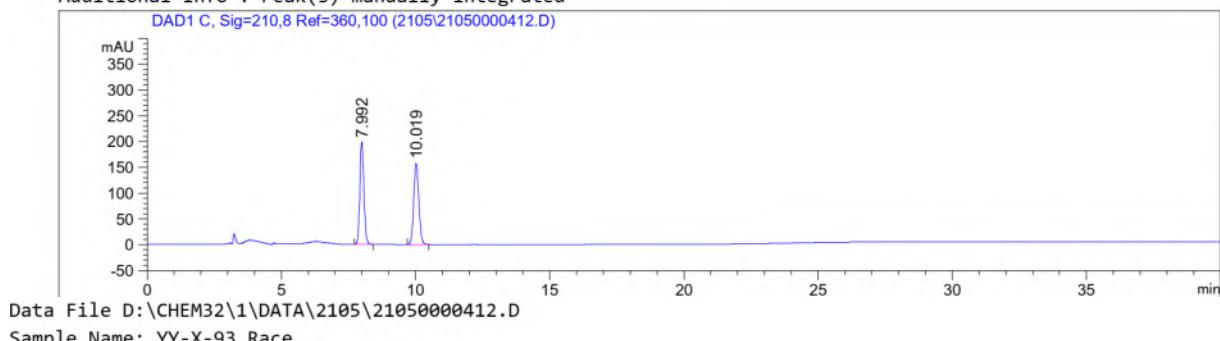
(R)-N-(tert-Butyl)-2-phenylpropanamide (48)

Data File D:\CHEM32\1\DATA\2105\21050000412.D

Sample Name: YY-X-93 Race

```
=====
Acq. Operator   : Analytik           Seq. Line : 11
Acq. Instrument : LC5             Location : Vial 8
Injection Date  : 5/4/2021 6:26:00 PM    Inj : 1
                                         Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/4/2021 4:22:07 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/5/2021 9:00:26 AM by Analytik
                  (modified after loading)
Method Info     : Chiralcel OJ, Hept./EtOH 95:5, 1.0ml/min
```

Additional Info : Peak(s) manually integrated



Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.992	BB	0.1620	2078.37427	198.67500	50.0101
2	10.019	BB	0.2037	2077.53784	157.69313	49.9899

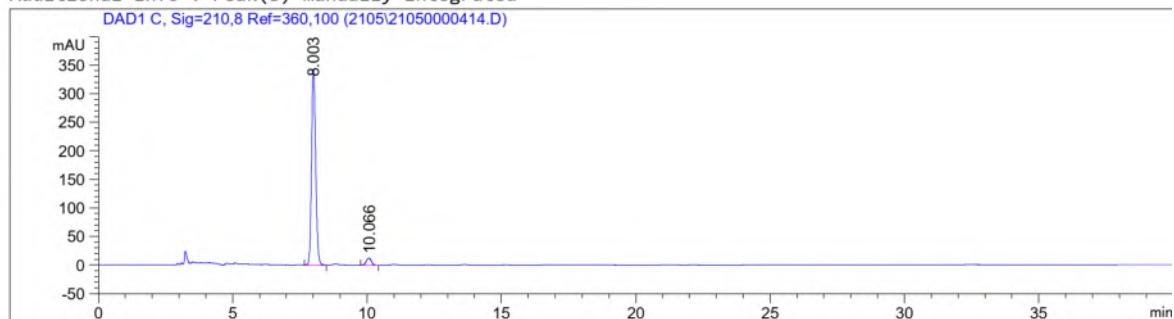
Totals : 4155.91211 356.36813

Data File D:\CHEM32\1\DATA\2105\21050000414.D

Sample Name: YY-X-93-C

```
=====
Acq. Operator   : Analytik          Seq. Line : 13
Acq. Instrument : LC5             Location  : Vial 18
Injection Date  : 5/4/2021 7:48:12 PM      Inj : 1
                                         Inj Volume : 1.0 µl
Acq. Method    : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/4/2021 4:22:07 PM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\SDC.M
Last changed    : 5/5/2021 9:00:26 AM by Analytik
                  (modified after loading)
Method Info     : Chiralcel OJ, Hept./EtOH 95:5, 1.0ml/min
```

Additional Info : Peak(s) manually integrated



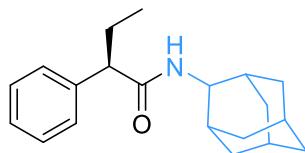
```
=====
Area Percent Report
=====
```

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.003	VB	0.1632	3628.12646	343.66391	95.3171
2	10.066	BB	0.2097	178.25041	13.02036	4.6829

Totals : 3806.37688 356.68427



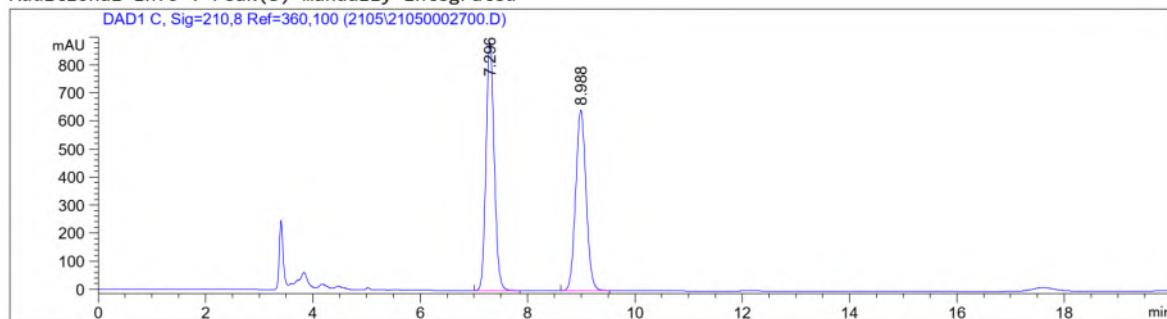
(R)-N-Adamantan-2-yl)-2-phenylbutanamide (52)

Data File D:\CHEM32\1\DATA\2105\21050002700.D

Sample Name: YY-X-137 Race

```
=====
Acq. Operator   : Analytik          Seq. Line : 1
Acq. Instrument : LC5             Location  : Vial 23
Injection Date  : 5/27/2021 9:48:32 AM      Inj       : 1
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\AMYLOSE2-0.5.M
Last changed    : 5/27/2021 10:07:58 AM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\AMYLOSE2-0.5.M
Last changed    : 5/27/2021 11:11:12 AM by Analytik
                  (modified after loading)
Method Info     : Amylose2, Hept./EtOH 90:10, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.296	BB	0.1661	9407.94336	884.61633	51.8692
2	8.988	BB	0.2102	8729.88574	643.98096	48.1308

Totals : 1.81378e4 1528.59729

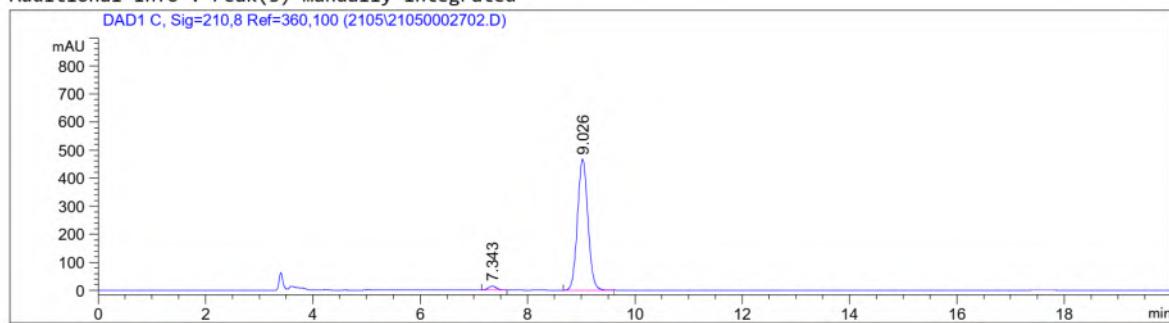
LC5 5/27/2021 11:11:31 AM Analytik

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Data File D:\CHEM32\1\DATA\2105\21050002702.D
Sample Name: YY-X-137 C

```
=====
Acq. Operator   : Analytik          Seq. Line :  2
Acq. Instrument : LC5             Location  : Vial 33
Injection Date  : 5/27/2021 10:40:42 AM    Inj :  2
                                                Inj Volume : 1.0 µl
Acq. Method     : C:\CHEM32\1\METHODS\AMYLOSE2-0.5.M
Last changed    : 5/27/2021 10:07:58 AM by Analytik
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\AMYLOSE2-0.5.M
Last changed    : 5/27/2021 11:11:12 AM by Analytik
                  (modified after loading)
Method Info     : Amylose2, Hept./EtOH 90:10, 0.5ml/min
```

Additional Info : Peak(s) manually integrated



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

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

Signal 1: DAD1 C, Sig=210,8 Ref=360,100

Peak	RetTime	Type	Width	Area	Height	Area %
#	[min]		[min]	[mAU*s]	[mAU]	%
1	7.343	BB	0.1618	141.78105	13.58352	2.1886
2	9.026	BB	0.2101	6336.25293	467.71741	97.8114

Totals : 6478.03398 481.30093

LC5 5/27/2021 11:14:24 AM Analytik

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