

Supporting Information

Synthesis of Chiral Sultams *via* Palladium-Catalyzed Intramolecular Asymmetric Reductive Amination

*Bo Song, Chang-Bin Yu, Yue Ji, Mu-Wang Chen and Yong-Gui Zhou**
Email: ygzhou@dicp.ac.cn

Table of Contents

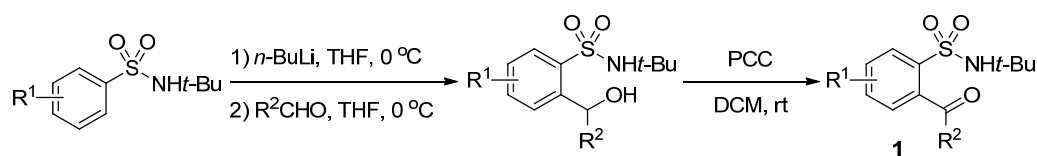
1. General.....	S1
2. The Preparation of Keto Sulfonamides.....	S1-8
3. The Procedure for Intramolecular Asymmetric Reductive Amination....	S8-19
4. Control Experiments.....	S19
5. Determination of Absolute Configuration of Products.....	S19-S20
6. References.....	S20
7. Copy of NMR and HPLC for the Compounds.....	S21-216

1. General

All reactions were carried out under an atmosphere of nitrogen using the standard Schlenk techniques, unless otherwise noted. Commercially available reagents were used without further purification. Solvents were treated prior to use according to the standard methods. ^1H NMR, ^{13}C NMR and ^{19}F NMR spectra were recorded at room temperature in CDCl_3 on 400 MHz instrument with tetramethylsilane (TMS) as internal standard. Enantiomeric excess was determined by HPLC analysis using chiral column. Optical rotations were measured by polarimeter. Flash column chromatography was performed on silica gel (200-300 mesh). All reactions were monitored by TLC analysis.

2. The Preparation of Keto Sulfonamides

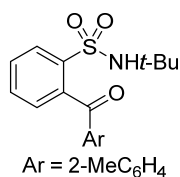
Keto sulfonamides **1** were prepared from the corresponding aldehydes and *N-tert*-butylbenzenesulfonamide according to the following procedures.



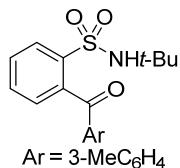
To a cooled ($0\text{ }^{\circ}\text{C}$) solution of *N-tert*-butylbenzenesulfonamide (853 mg, 4.0 mmol) in THF (20 mL) was added *n*-butyl lithium (3.5 mL, 8.8 mmol, 2.5 M in *n*-hexane). After stirring the resulting mixture for 30 min, a solution of aldehyde (4.4 mmol) in THF (5 mL) was added dropwise over a period of 20 min and stirring was continued for additional 2 hours, the mixture was poured into a saturated aqueous ammonium chloride solution (20 mL). The aqueous layer was extracted three times with ethyl acetate (20 mL $\times$ 3), and the organic extracts were dried over anhydrous sodium sulfate. After concentration in *vacuo*, the residue was finally purified by flash chromatography to afford the corresponding alcohol. To a solution of the entire amount of intermediate alcohol in dichloromethane (25 mL) was added PCC (3.10 g, 13.9 mmol). The resulting dark-brown solution was stirred for overnight at ambient temperature. After addition of diethyl ether (10 mL) and additional stirring (30 min), the mixture was filtered through a pad of silica gel. Concentration in *vacuo* afforded the analytically pure keto sulfonamides **1**. The yields given are overall yields for two steps.

2-Benzoyl-*N-tert*-butylbenzenesulfonamide (1a): 0.951 g, 75% yield (4 mmol scale), white solid, mp $82\text{--}83\text{ }^{\circ}\text{C}$. ^1H NMR (400 MHz, CDCl_3) δ 8.14–8.12 (m, 1H), 7.80–7.78 (m, 2H), 7.61–7.57 (m, 3H), 7.44 (t, $J = 7.7\text{ Hz}$, 2H), 7.36 (dd, $J = 7.1, 1.5\text{ Hz}$, 1H), 5.24 (br, 1H), 1.30 (s, 9H); ^{13}C NMR (100 MHz, CDCl_3) δ 198.1, 142.7, 137.6, 136.4, 134.1, 131.2, 130.7, 130.3, 129.3, 128.9, 128.6, 55.3, 30.3. HRMS Calculated for $\text{C}_{17}\text{H}_{20}\text{NO}_3\text{S}$ $[\text{M}+\text{H}]^+$ 340.0978, found: 340.0971.

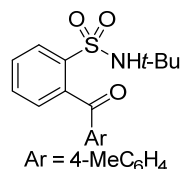
***N-tert*-Butyl-2-(2-methylbenzoyl)-benzenesulfonamide (1b):** 0.857 g, 65% yield (4 mmol scale), white solid, mp $136\text{--}137\text{ }^{\circ}\text{C}$. ^1H NMR (400 MHz, CDCl_3) δ 8.14 (d, $J = 7.7\text{ Hz}$, 1H), 7.59 (t, $J = 7.3\text{ Hz}$, 1H), 7.52 (t, $J = 7.3\text{ Hz}$, 1H), 7.41 (dd, $J = 13.2, 7.2\text{ Hz}$, 2H), 7.27 (d, $J = 7.7\text{ Hz}$, 2H), 7.20 (t, $J = 7.5\text{ Hz}$, 1H), 5.47 (br, 1H), 2.49 (s, 3H), 1.32 (s, 9H); ^{13}C NMR (100 MHz, CDCl_3) δ 200.1, 142.6, 140.0, 139.4, 136.7, 132.6, 132.4, 131.8, 131.4, 130.7, 129.9, 129.1, 125.7, 55.3, 30.4, 21.2. HRMS Calculated for $\text{C}_{18}\text{H}_{22}\text{NO}_3\text{S}$ $[\text{M}+\text{H}]^+$ 354.1134, found: 354.1136.



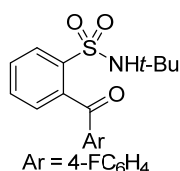
***N*-tert-Butyl-2-(3-methylbenzoyl)-benzenesulfonamide (1c):** 1.127 g, 85% yield (4 mmol scale), white solid, mp 84-85 °C. ¹H NMR (400 MHz, CDCl₃) δ 8.14-8.12 (m, 1H), 7.62-7.56 (m, 4H), 7.40 (d, *J* = 7.6 Hz, 1H), 7.37-7.33 (m, 2H), 5.25 (br, 1H), 2.37 (s, 3H), 1.30 (s, 9H); ¹³C NMR (100 MHz, CDCl₃) δ 198.3, 142.7, 138.5, 137.8, 136.5, 135.0, 131.2, 131.0, 130.3, 129.3, 128.9, 128.5, 128.2, 55.3, 30.3, 21.4. HRMS Calculated for C₁₈H₂₂NO₃S [M+H]⁺ 354.1134, found: 354.1131.



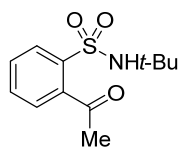
***N*-tert-Butyl-2-(4-methylbenzoyl)-benzenesulfonamide (1d):** 0.805 g, 61% yield (4 mmol scale), white solid, mp 120-121 °C. ¹H NMR (400 MHz, CDCl₃) δ 8.14-8.11 (m, 1H), 7.70 (d, *J* = 8.2 Hz, 2H), 7.62-7.57 (m, 2H), 7.37-7.35 (m, 1H), 7.25 (d, *J* = 8.5 Hz, 2H), 5.26 (br, 1H), 2.42 (s, 3H), 1.30 (s, 9H); ¹³C NMR (100 MHz, CDCl₃) δ 197.8, 145.3, 142.7, 137.9, 134.0, 131.2, 130.9, 130.2, 129.4, 129.3, 128.9, 55.3, 30.3, 21.9. HRMS Calculated for C₁₈H₂₂NO₃S [M+H]⁺ 354.1134, found: 354.1140.



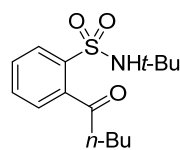
***N*-tert-Butyl-2-(4-fluorobenzoyl)-benzenesulfonamide (1e):** 1.192g, 89% yield (4 mmol scale), white solid, mp 82-83 °C. ¹H NMR (400 MHz, CDCl₃) δ 8.13 (d, *J* = 7.3 Hz, 1H), 7.83 (dd, *J* = 7.9, 5.7 Hz, 2H), 7.65-7.58 (m, 2H), 7.35 (d, *J* = 7.2 Hz, 1H), 7.12 (t, *J* = 8.1 Hz, 2H), 5.19 (br, 1H), 1.29 (s, 9H); ¹³C NMR (100 MHz, CDCl₃) δ 196.5, 166.3 (d, *J* = 256.8 Hz), 142.6, 137.4, 133.4 (d, *J* = 9.7 Hz), 132.9 (d, *J* = 2.9 Hz), 131.3, 130.4, 129.0, 116.0, 115.9 (d, *J* = 22.1 Hz), 55.3, 30.2; ¹⁹F NMR (376 MHz, CDCl₃) δ -130.17. HRMS Calculated for C₁₇H₁₈FNO₃Na [M+Na]⁺ 358.0884, found: 358.0881.



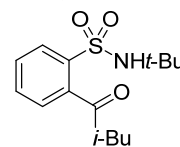
2-Acetyl-*N*-tert-butylbenzenesulfonamide (1f): 0.458 g, 45% yield (4 mmol scale), white solid, mp 84-85 °C. ¹H NMR (400 MHz, CDCl₃) δ 8.06 (dd, *J* = 7.8, 1.2 Hz, 1H), 7.62-7.54 (m, 3H), 5.42 (br, 1H), 2.63 (s, 3H), 1.24 (s, 9H); ¹³C NMR (100 MHz, CDCl₃) δ 204.5, 141.4, 139.2, 131.9, 130.8, 129.1, 127.8, 55.2, 30.5, 30.3. HRMS Calculated for C₁₂H₁₇NO₃Na [M+Na]⁺ 278.0821, found: 278.0825.



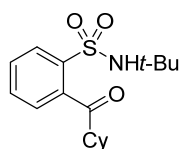
***N*-tert-Butyl-2-pentanoylbenzenesulfonamide (1g):** 0.482 g, 41% yield (4 mmol scale), colorless oil. ¹H NMR (400 MHz, CDCl₃) δ 8.04 (dd, *J* = 7.6, 1.3 Hz, 1H), 7.60-7.52 (m, 2H), 7.48 (dd, *J* = 7.3, 1.4 Hz, 1H), 5.34 (br, 1H), 2.92 (t, *J* = 7.3 Hz, 2H), 1.75-1.67 (m, 2H), 1.44-1.35 (m, 2H), 1.25 (s, 9H), 0.93 (t, *J* = 7.3 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 207.5, 141.4, 139.7, 131.9, 130.4, 129.0, 127.4, 55.2, 42.8, 30.3, 25.8, 22.3, 14.0. HRMS Calculated for C₁₅H₂₃NO₃Na [M+Na]⁺ 320.1291, found: 320.1283.



***N*-tert-Butyl-2-(3-methylbutanoyl)-benzenesulfonamide (1h):** 0.508 g, 43% yield (4 mmol scale), colorless oil. ¹H NMR (400 MHz, CDCl₃) δ 8.05 (dd, *J* = 7.7, 1.3 Hz, 1H), 7.60-7.48 (m, 3H), 5.43 (br, 1H), 2.81 (d, *J* = 6.7 Hz, 2H), 2.36-2.26 (m, 1H), 1.25 (s, 9H), 1.00 (d, *J* = 6.7 Hz, 6H); ¹³C NMR (100 MHz, CDCl₃) δ 206.6, 141.4, 139.6, 131.8, 130.5, 129.0, 127.8, 55.2, 51.7, 30.3, 24.3, 22.6. HRMS Calculated for C₁₅H₂₃NO₃Na [M+Na]⁺ 320.1291, found: 320.1303.



***N*-tert-Butyl-2-(cyclohexanecarbonyl)-benzenesulfonamide (1i):** 0.426 g, 33% yield (4 mmol scale), white solid, mp 110-111 °C. ¹H NMR (400 MHz, CDCl₃) δ 8.07 (dd, *J* = 7.6, 1.3 Hz, 1H), 7.61-7.51 (m, 3H), 5.49 (br, 1H), 3.02-2.94 (m, 1H), 1.91 (d, *J* = 12.4 Hz, 2H), 1.82-1.77 (m, 2H), 1.68-1.66 (m, 1H), 1.54-1.48 (m, 2H), 1.30-1.18 (m, 12H); ¹³C NMR (100 MHz, CDCl₃) δ 209.9, 142.1, 138.9, 131.8,



130.5, 129.2, 127.6, 55.2, 50.1, 30.3, 28.7, 25.9, 25.9. HRMS Calculated for $C_{17}H_{25}NO_3SNa$ $[M+Na]^+$ 346.1447, found: 346.1438.

2-Benzoyl-*N*-(*tert*-butyl)-4-methylbenzenesulfonamide (1j): 0.690 g, 52% yield (4 mmol scale), white solid, mp 158-159 °C. 1H NMR (400 MHz, $CDCl_3$) δ 8.00 (d, J = 8.1 Hz, 1H), 7.81-7.79 (m, 2H), 7.62-7.59 (m, 1H), 7.48-7.44 (m, 2H), 7.40 (d, J = 8.1 Hz, 1H), 7.15 (s, 1H), 5.14 (s, 1H), 2.42 (s, 3H), 1.29 (s, 9H); ^{13}C NMR (100 MHz, $CDCl_3$) δ 198.2, 142.0, 139.9, 137.7, 136.5, 133.9, 130.7, 130.7, 129.6, 129.0, 128.5, 55.1, 30.2, 21.4. HRMS Calculated for $C_{18}H_{21}NO_3SNa$ $[M+Na]^+$ 354.1134, found: 354.1137.

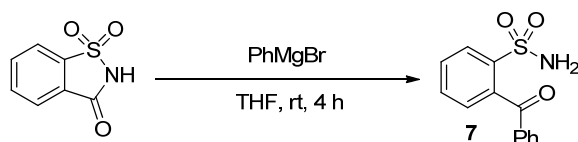
2-Benzoyl-*N*-(*tert*-butyl)-4-methoxybenzenesulfonamide (1k): 1.013 g, 73% yield (4 mmol scale), white solid, mp 142-143 °C. 1H NMR (400 MHz, $CDCl_3$) δ 8.04 (d, J = 8.8 Hz, 1H), 7.82 (d, J = 7.5 Hz, 2H), 7.60 (t, J = 7.4 Hz, 1H), 7.47-7.43 (m, 2H), 7.05 (dd, J = 8.8, 2.4 Hz, 1H), 6.83-6.82 (m, 1H), 5.05 (s, 1H), 3.84 (s, 3H), 1.29 (s, 9H); ^{13}C NMR (100 MHz, $CDCl_3$) δ 197.5, 161.4, 139.5, 136.2, 134.3, 134.1, 131.1, 130.7, 128.6, 115.3, 114.2, 55.8, 55.0, 30.2. HRMS Calculated for $C_{18}H_{21}NO_4SNa$ $[M+Na]^+$ 370.1083, found: 370.1086.

2-Benzoyl-*N*-(*tert*-butyl)-4-chlorobenzenesulfonamide (1l): 0.841 g, 60% yield (4 mmol scale), white solid, mp 103-104 °C. 1H NMR (400 MHz, $CDCl_3$) δ 8.06 (d, J = 8.5 Hz, 1H), 7.80 (d, J = 7.3 Hz, 2H), 7.65-7.57 (m, 2H), 7.50-7.46 (m, 2H), 7.34-7.33 (m, 1H), 5.11 (s, 1H), 1.31 (s, 9H); ^{13}C NMR (100 MHz, $CDCl_3$) δ 196.4, 141.2, 139.3, 137.7, 135.8, 134.4, 130.7, 130.4, 130.2, 129.0, 128.7, 55.4, 30.2. HRMS Calculated for $C_{17}H_{18}ClNO_3SNa$ $[M+Na]^+$ 374.0588, found: 374.0592.

***N*-(*tert*-Butyl)-4-methyl-2-pentanoylbenzenesulfonamide (1m):** 0.469 g, 38% yield (4 mmol scale), white solid, mp 57-58 °C. 1H NMR (400 MHz, $CDCl_3$) δ 7.92 (d, J = 8.1 Hz, 1H), 7.33 (d, J = 8.1 Hz, 1H), 7.26 (s, 1H), 5.27 (s, 1H), 2.92 (t, J = 7.3 Hz, 2H), 2.45 (s, 3H), 1.75-1.68 (m, 2H), 1.46-1.36 (m, 2H), 1.25 (s, 9H), 0.95 (t, J = 7.3 Hz, 3H); ^{13}C NMR (100 MHz, $CDCl_3$) δ 207.5, 142.5, 139.7, 138.5, 130.7, 129.0, 127.9, 55.0, 42.7, 30.2, 25.7, 22.2, 21.4, 13.9. HRMS Calculated for $C_{16}H_{25}NO_3SNa$ $[M+Na]^+$ 334.1449, found: 334.1449.

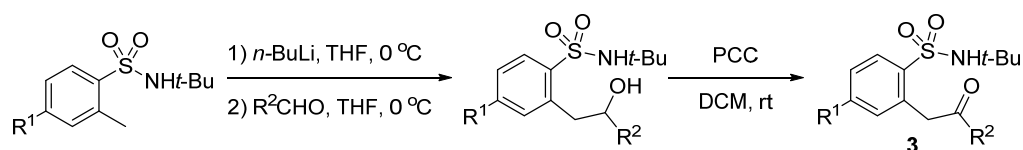
2-Benzoyl-*N*-(*tert*-butyl)-3-fluoro-6-methylbenzenesulfonamide (1n): 0.456 g, 33% yield (4 mmol scale), white solid, mp 179-180 °C. 1H NMR (400 MHz, $CDCl_3$) δ 7.80 (d, J = 7.3 Hz, 2H), 7.58-7.54 (m, 1H), 7.46-7.42 (m, 2H), 7.39-7.36 (m, 1H), 7.26-7.20 (m, 1H), 4.75 (s, 1H), 2.72 (s, 3H), 1.27 (s, 9H); ^{13}C NMR (100 MHz, $CDCl_3$) δ 192.9, 157.7 (d, J = 248.0 Hz), 140.9 (d, J = 2.3 Hz), 137.5, 134.9 (d, J = 7.5 Hz), 133.7, 133.6 (d, J = 3.8 Hz), 129.3, 128.7, 128.2 (d, J = 20.5 Hz), 119.5 (d, J = 21.7 Hz), 55.9, 30.2, 20.7; ^{19}F NMR (376 MHz, $CDCl_3$) δ -115.29. HRMS Calculated for $C_{18}H_{24}FN_2O_3S$ $[M+NH_4]^+$ 367.1486, found: 367.1490.

Keto sulfonamide **7** was prepared from saccharin and phenylmagnesium bromide according to the following procedures. The compound **7** is known.¹



To a degassed THF (20 mL) solution of saccharin (0.915 g, 5.0 mmol) at 25 °C was added fresh prepared phenylmagnesium bromide (15.0 mmol, 1.0 M in THF) dropwise. The resulting mixture was further stirred at the same temperature for 4 hours before being quenched with a saturated aqueous ammonium chloride solution. The aqueous layer was extracted further with dichloromethane (30 mL×3); then the combined organic layer was washed with brine and dried over sodium sulfate. Evaporation of the solvent in ice-water bath gave the crude product. The solid obtained was washed two times with cold *n*-hexane and hot *n*-hexane (60-70 °C), then, crystallization of the insoluble part from toluene gave 0.146 g keto sulfonamide **7** in 11% yield.

Keto sulfonamides **3** were prepared from the corresponding aldehydes and *N*-*tert*-butyl-2-methyl benzenesulfonamide according to the following procedures.

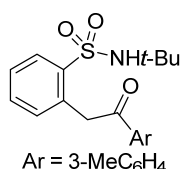


To a cooled (0 °C) solution of *N*-*tert*-butyl-2-methylbenzenesulfonamide (4.0 mmol) in THF (15 mL) was added *n*-butyl lithium (3.5 mL, 8.8 mmol, 2.5 M in *n*-hexane). After stirring the resulting mixture for 45 min, a solution of aldehyde (4.4 mmol) in THF (10 mL) was added dropwise over a period of 20 min and stirring was continued for additional 3 hours, the mixture was poured into a saturated aqueous ammonium chloride solution (20 mL). The aqueous layer was extracted three times with ethyl acetate (20 mL×3), and the organic extracts were dried over anhydrous sodium sulfate. After concentration in *vacuo*, the residue was finally purified by flash chromatography to afford the corresponding alcohol. To a solution of the entire amount of intermediate alcohol in dichloromethane (25 mL) was added PCC (1.724 g, 8.0 mmol). The resulting dark-brown solution was stirred for overnight at ambient temperature. After addition of diethyl ether (10 mL) and additional stirring (30 min), the mixture was filtered through a pad of silica gel. Concentration in *vacuo* afforded the analytically pure keto sulfonamide compounds **3**. The yields given are overall yields for two steps.

***N*-*tert*-Butyl-2-(2-oxo-2-phenylethyl)-benzenesulfonamide (3a):** 0.873 g, 53% yield (5 mmol scale), white solid, mp 88- 89 °C. ^1H NMR (400 MHz, CDCl_3) δ 8.12-8.05 (m, 3H), 7.62-7.58 (m, 1H), 7.53-7.48 (m, 3H), 7.41 (td, $J = 7.8, 1.2$ Hz, 1H), 7.30-7.27 (m, 1H), 4.83 (s, 2H), 4.55 (br, 1H), 1.17 (s, 9H); ^{13}C NMR (100 MHz, CDCl_3) δ 197.5, 141.4, 136.6, 133.8, 133.7, 133.5, 132.5, 129.7, 128.9, 128.6, 127.5, 55.5, 43.1, 30.0. HRMS Calculated for $\text{C}_{18}\text{H}_{22}\text{NO}_3\text{S}$ $[\text{M}+\text{H}]^+$ 332.1315, found: 332.1318.

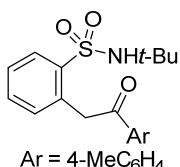
***N*-*tert*-Butyl-2-(2-oxo-2-*o*-tolylethyl)-benzenesulfonamide (3b):** 0.728 g, 53% yield (4 mmol scale), white solid, mp 81-82 °C. ^1H NMR (400 MHz, CDCl_3) δ 8.09 (dd, $J = 7.9, 1.3$ Hz, 1H), 7.90 (d, $J = 7.7$ Hz, 1H), 7.53-7.49 (m, 1H), 7.43-7.39 (m, 2H), 7.33-7.25 (m, 3H), 4.76 (br, 2H), 2.52 (s, 3H), 1.22 (s, 9H); ^{13}C NMR (100 MHz, CDCl_3) δ 201.0, 141.6, 139.1, 137.0, 133.8, 133.5, 132.5, 132.4, 132.0, 129.6, 129.3, 127.6, 126.1, 55.5, 46.1, 30.2, 21.7. HRMS Calculated for $\text{C}_{19}\text{H}_{24}\text{NO}_3\text{S}$ $[\text{M}+\text{H}]^+$ 346.1471, found: 346.1472.

***N*-*tert*-Butyl-2-(2-oxo-2-*m*-tolylethyl)-benzenesulfonamide (3c):** 0.504 g, 37% yield (4 mmol scale), white solid, mp 97-98 °C. ^1H NMR (400 MHz, CDCl_3) δ 8.10 (dd, $J = 7.9, 1.2$ Hz, 1H), 7.87-7.85 (m, 2H), 7.51 (td, $J = 7.5, 1.3$ Hz, 1H), 7.42-7.35 (m, 3H), 7.28-7.26 (m, 1H), 4.81 (s,



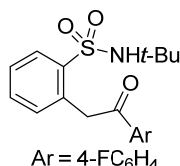
2H), 4.57 (br, 1H), 2.42 (s, 3H), 1.17 (s, 9H); ^{13}C NMR (100 MHz, CDCl_3) δ 197.7, 141.5, 138.8, 136.6, 134.5, 133.9, 133.5, 132.5, 129.6, 129.1, 128.8, 127.5, 125.9, 55.5, 43.2, 30.0, 21.5. HRMS Calculated for $\text{C}_{19}\text{H}_{24}\text{NO}_3\text{S}$ $[\text{M}+\text{H}]^+$ 346.1471, found: 346.1475.

***N*-tert-Butyl-2-(2-oxo-2-*p*-tolylethyl)-benzenesulfonamide (3d):** 0.613 g, 44% yield (4 mmol scale), white solid, mp 84-85 °C. ^1H NMR (400 MHz, CDCl_3) δ 8.10 (dd, J = 8.0, 1.3 Hz, 1H),



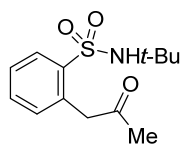
7.96 (d, J = 8.2 Hz, 2H), 7.50 (td, J = 7.5, 1.4 Hz, 1H), 7.39 (td, J = 7.8, 1.3 Hz, 1H), 7.29-7.26 (m, 3H), 4.79 (s, 2H), 4.59 (br, 1H), 2.42 (s, 3H), 1.16 (s, 9H); ^{13}C NMR (100 MHz, CDCl_3) δ 197.1, 144.6, 141.4, 134.1, 134.0, 133.5, 132.5, 129.6, 129.6, 128.8, 127.4, 55.4, 42.9, 30.0, 21.8. HRMS Calculated for $\text{C}_{19}\text{H}_{24}\text{NO}_3\text{S}$ $[\text{M}+\text{H}]^+$ 346.1471, found: 346.1470.

***N*-tert-Butyl-2-(2-(4-fluorophenyl)-2-oxo-ethyl)-benzenesulfonamide (3e):** 0.728 g, 52% yield (4 mmol scale), white solid, mp 115-116 °C. ^1H NMR (400 MHz, CDCl_3) δ 8.04-8.00 (m, 3H),



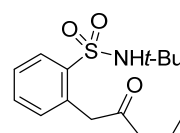
7.45 (td, J = 7.5, 1.2 Hz, 1H), 7.35 (td, J = 7.8, 1.1 Hz, 1H), 7.21-7.19 (m, 1H), 7.11-7.07 (m, 2H), 4.71 (s, 2H), 4.43 (br, 1H), 1.10 (s, 9H); ^{19}F NMR (376 MHz, CDCl_3) δ -104.31; ^{13}C NMR (100 MHz, CDCl_3) δ 195.9, 166.1 (d, J = 255.6 Hz), 141.3, 133.5, 133.4, 133.0 (d, J = 3.1 Hz), 132.6, 131.3 (d, J = 9.4 Hz), 129.7, 127.7, 116.1 (d, J = 21.9 Hz), 55.5, 43.2, 30.0. HRMS Calculated for $\text{C}_{18}\text{H}_{20}\text{FNO}_3\text{S}$ $[\text{M}+\text{H}]^+$ 350.1221, found: 350.1226.

***N*-tert-Butyl-2-(2-oxo-propyl)-benzenesulfonamide (3f):** 0.461 g, 57% yield (3 mmol scale), white solid, mp 87-88 °C. ^1H NMR (400 MHz, CDCl_3) δ 8.05 (dd, J = 7.9, 1.2 Hz, 1H), 7.51 (td, J = 7.5, 1.3 Hz, 1H), 7.40 (td, J = 7.8, 1.2 Hz, 1H), 7.27-7.24 (m, 1H), 4.95 (br, 1H), 4.22 (s, 2H), 2.26 (s, 3H), 1.23 (s, 9H); ^{13}C NMR (100 MHz, CDCl_3) δ



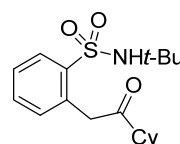
206.1, 141.5, 133.4, 133.2, 132.5, 129.5, 127.6, 55.4, 48.4, 30.2, 30.1. HRMS Calculated for $\text{C}_{13}\text{H}_{20}\text{NO}_3\text{S}$ $[\text{M}+\text{H}]^+$ 270.1158, found: 270.1158.

***N*-tert-Butyl-2-(2-oxo-pentyl)-benzenesulfonamide (3g):** 0.319 g, 36% yield (3 mmol scale), white solid, mp 86-87 °C. ^1H NMR (400 MHz, CDCl_3) δ 8.05 (dd, J = 7.9, 1.1 Hz, 1H), 7.50 (td, J = 7.5, 1.2 Hz, 1H), 7.39 (td, J = 7.8, 1.1 Hz, 1H), 7.24 (d, J = 7.5 Hz, 1H), 4.94



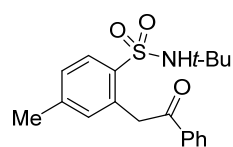
(br, 1H), 4.20 (s, 2H), 2.54 (t, J = 7.3 Hz, 2H), 1.66-1.61 (m, 2H), 1.24 (s, 9H), 0.93 (t, J = 7.4 Hz, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 208.4, 141.5, 133.3, 133.3, 132.4, 129.5, 127.5, 55.4, 47.7, 44.7, 30.2, 17.2, 13.8. HRMS Calculated for $\text{C}_{15}\text{H}_{24}\text{NO}_3\text{S}$ $[\text{M}+\text{H}]^+$ 298.1471, found: 298.1472.

***N*-tert-Butyl-2-(2-cyclohexyl-2-oxo-ethyl)-benzenesulfonamide (3h):** 0.516 g, 51% yield (3 mmol scale), white solid, mp 79-80 °C. ^1H NMR (400 MHz, CDCl_3) δ 8.05 (dd, J = 7.9, 1.1 Hz,



1H), 7.48 (td, J = 7.5, 1.3 Hz, 1H), 7.38 (td, J = 7.8, 1.2 Hz, 1H), 7.19 (d, J = 7.5 Hz, 1H), 4.81 (s, 1H), 4.29 (s, 2H), 2.58-2.51 (m, 1H), 1.96-1.93 (m, 2H), 1.82-1.79 (m, 2H), 1.70-1.64 (m, 1H), 1.42-1.23 (m, 14H); ^{13}C NMR (100 MHz, CDCl_3) δ 211.6, 141.6, 133.6, 133.3, 132.4, 129.6, 127.5, 55.4, 50.8, 46.0, 30.2,

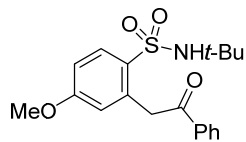
28.7, 25.9, 25.7. HRMS Calculated for $\text{C}_{18}\text{H}_{28}\text{NO}_3\text{S}$ $[\text{M}+\text{H}]^+$ 338.1784, found: 338.1783.



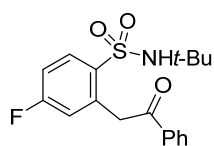
***N*-(tert-Butyl)-4-methyl-2-(2-oxo-phenylethyl)-benzenesulfonamide (3i):** 0.976 g, 71% yield (4 mmol scale), white solid, mp 119-120 °C. ^1H NMR (400 MHz, CDCl_3) δ 8.07 (d, J = 7.8 Hz, 2H), 7.97 (d, J = 8.1 Hz, 1H), 7.62-7.58 (m, 1H), 7.51-7.48 (m, 2H), 7.20 (d, J = 8.1 Hz, 1H), 7.08 (s,

1H), 4.78 (s, 2H), 4.46 (s, 1H), 2.38 (s, 3H), 1.16 (s, 9H); ¹³C NMR (100 MHz, CDCl₃) δ 197.5, 143.1, 138.4, 136.5, 134.0, 133.5, 133.5, 129.7, 128.8, 128.5, 128.1, 55.3, 43.0, 29.9, 21.3. HRMS Calculated for C₁₉H₂₃NO₃SNa [M+Na]⁺ 368.1291, found: 368.1294.

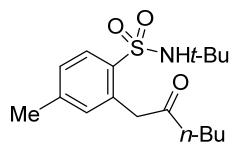
***N*-(*tert*-Butyl)-4-methoxy-2-(2-oxo-2-phenylethyl)-benzenesulfonamide (3j):** 0.345 g, 60% yield (1.6 mmol scale), white solid, mp 134-135 °C. ¹H NMR (400 MHz, CDCl₃) δ 8.07-8.02 (m, 3H), 7.62-7.58 (m, 1H), 7.51-7.49 (m, 2H), 6.87 (dd, *J* = 8.9, 2.4 Hz, 1H), 6.78-6.77 (m, 1H), 4.78 (s, 2H), 4.40 (s, 1H), 3.83 (s, 3H), 1.16 (s, 9H); ¹³C NMR (100 MHz, CDCl₃) δ 197.2, 162.4, 136.5, 135.8, 133.6, 133.1, 131.9, 128.8, 128.5, 118.9, 111.9, 55.5, 55.2, 43.2, 29.9. HRMS Calculated for C₁₉H₂₄NO₄S [M+H]⁺ 362.1421, found: 362.1424.



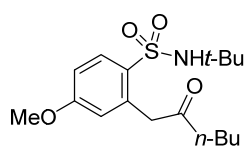
***N*-(*tert*-Butyl)-4-fluoro-2-(2-oxo-2-phenylethyl)-benzenesulfonamide (3k):** 0.404 g, 46% yield (2.5 mmol scale), colorless oil. ¹H NMR (400 MHz, CDCl₃) δ 8.11 (dd, *J* = 8.9, 5.7 Hz, 1H), 8.07-8.05 (m, 2H), 7.65-7.61 (m, 1H), 7.53-7.51 (m, 2H), 7.12-7.07 (m, 1H), 7.01 (dd, *J* = 9.1, 2.6 Hz, 1H), 4.81 (s, 2H), 4.39 (br, 1H), 1.16 (s, 9H); ¹³C NMR (100 MHz, CDCl₃) δ 196.6, 164.5 (d, *J* = 254.7 Hz), 137.3 (d, *J* = 3.4 Hz), 137.0 (d, *J* = 8.9 Hz), 136.2, 133.8, 132.3 (d, *J* = 9.3 Hz), 129.0, 128.5, 120.5 (d, *J* = 22.6 Hz), 114.4 (d, *J* = 21.6 Hz), 55.5, 43.0, 29.8; ¹⁹F NMR (376 MHz, CDCl₃) δ -106.54. HRMS Calculated for C₁₈H₂₄FN₂O₃S [M+NH₄]⁺ 367.1486, found: 367.1490.



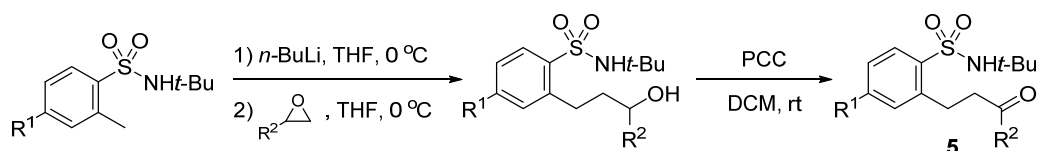
***N*-(*tert*-Butyl)-4-methyl-2-(2-oxohexyl)-benzenesulfonamide (3l):** 0.765 g, 59% yield (4 mmol scale), white solid, mp 99-100 °C. ¹H NMR (400 MHz, CDCl₃) δ 7.92 (d, *J* = 8.1 Hz, 1H), 7.18 (d, *J* = 8.1 Hz, 1H), 7.03 (s, 1H), 4.84 (s, 1H), 4.16 (br, 2H), 2.56 (t, *J* = 7.4 Hz, 2H), 2.39 (s, 3H), 1.62-1.55 (m, 2H), 1.36-1.29 (m, 2H), 1.23 (s, 9H), 0.90 (t, *J* = 7.3 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 208.6, 143.0, 138.5, 133.9, 133.1, 129.6, 128.1, 55.2, 47.5, 42.5, 30.1, 25.7, 22.3, 21.3, 13.9. HRMS Calculated for C₁₇H₂₈NO₃S [M+H]⁺ 326.1784, found: 326.1787.



***N*-(*tert*-Butyl)-4-methoxy-2-(2-oxohexyl)-benzenesulfonamide (3m):** 0.435 g, 67% yield (1.9 mmol scale), white solid, mp 72-73 °C. ¹H NMR (400 MHz, CDCl₃) δ 7.98 (d, *J* = 8.9 Hz, 1H), 6.85 (dd, *J* = 8.8, 2.4 Hz, 1H), 6.72 (d, *J* = 2.4 Hz, 1H), 4.77 (s, 1H), 4.16 (s, 2H), 3.85 (s, 3H), 2.56 (t, *J* = 7.4 Hz, 2H), 1.62-1.55 (m, 2H), 1.35-1.28 (m, 2H), 1.23 (s, 9H), 0.90 (t, *J* = 7.3 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 208.2, 162.3, 135.4, 133.2, 131.8, 118.7, 111.9, 55.5, 55.1, 47.7, 42.5, 30.1, 25.7, 22.3, 13.9. HRMS Calculated for C₁₇H₂₈NO₄S [M+H]⁺ 342.1734, found: 342.1734.



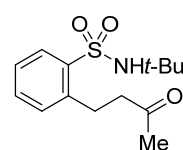
Keto sulfonamides **5** were prepared from the corresponding substituted oxiranes and *N*-*tert*-butyl-2-methylbenzenesulfonamide according to the following procedures.



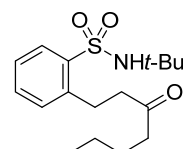
To a cooled (0 °C) solution of *N*-*tert*-butyl-2-methylbenzenesulfonamide (4.0 mmol) in THF (15 mL) was added *n*-butyl lithium (3.5 mL, 8.8 mmol, 2.5 M in *n*-hexane). After stirring the resulting mixture for 30 min, a solution of substituted oxiranes (4.4 mmol) in THF (10 mL) was

added dropwise over a period of 20 min and stirring was continued for additional 3 hours, the mixture was poured into a saturated aqueous ammonium chloride solution (20 mL). The aqueous layer was extracted three times with ethyl acetate (20 mL×3), and the organic extracts were dried over anhydrous sodium sulfate. After concentration in *vacuo*, the residue was finally purified by flash chromatography to afford the corresponding alcohol. To a solution of the entire amount of intermediate alcohol in dichloromethane (25 mL) was added PCC (1.724 g, 8.0 mmol). The resulting dark- brown solution was stirred for overnight at ambient temperature. After addition of diethyl ether (10 mL) and additional stirring (30 min), the mixture was filtered through a pad of silica gel. Concentration in *vacuo* afforded the analytically pure keto sulfonamide compounds **5**. The yields given are overall yields for two steps.

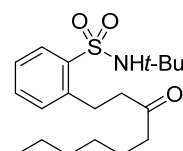
***N*-tert-Butyl-2-(3-oxo-butyl)-benzenesulfonamide (5a):** 1.040 g, 61% yield (6 mmol scale), white solid, mp 106-107 °C. ¹H NMR (400 MHz, CDCl₃) δ 8.03-8.00 (m, 1H), 7.45 (t, *J* = 7.5 Hz, 1H), 7.32-7.27 (m, 2H), 4.90 (br, 1H), 3.24 (t, *J* = 7.6 Hz, 2H), 2.87 (t, *J* = 7.6 Hz, 2H), 2.15 (s, 3H), 1.22 (s, 9H); ¹³C NMR (100 MHz, CDCl₃) δ 208.2, 141.2, 140.2, 132.6, 131.5, 129.6, 126.6, 55.1, 45.0, 30.3, 30.1, 27.1. HRMS Calculated for C₁₄H₂₁NO₃S [M+H]⁺ 284.1315, found: 284.1320.



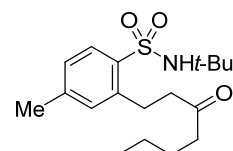
***N*-tert-Butyl-2-(3-oxo-heptyl)-benzenesulfonamide (5b):** 0.287 g, 35% yield (2.5 mmol scale), colorless oil. ¹H NMR (400 MHz, CDCl₃) δ 8.01 (d, *J* = 7.8 Hz, 1H), 7.44 (t, *J* = 7.2 Hz, 1H), 7.30-7.28 (m, 2H), 4.86 (br, 1H), 3.24 (t, *J* = 7.4 Hz, 2H), 2.84 (t, *J* = 7.5 Hz, 2H), 2.40 (t, *J* = 7.4 Hz, 2H), 1.55-1.52 (m, 2H), 1.29-1.25 (m, 2H), 1.22 (s, 9H), 0.87 (t, *J* = 7.3 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 210.6, 141.2, 140.4, 132.6, 131.5, 129.6, 126.5, 55.0, 44.0, 42.7, 30.3, 27.1, 26.0, 22.4, 13.9. HRMS Calculated for C₁₇H₂₈NO₃S [M+H]⁺ 326.1784, found: 326.1786.



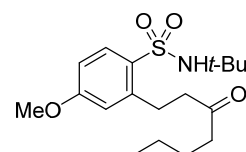
***N*-tert-Butyl-2-(3-oxo-nonyl)-benzenesulfonamide (5c):** 0.829 g, 59% yield (4 mmol scale), colorless oil. ¹H NMR (400 MHz, CDCl₃) δ 8.03 (d, *J* = 7.7 Hz, 1H), 7.47-7.43 (m, 1H), 7.32-7.28 (m, 2H), 5.25-5.19 (br, 1H), 3.28 (t, *J* = 7.5 Hz, 2H), 2.86 (t, *J* = 7.4 Hz, 2H), 2.41 (t, *J* = 7.3 Hz, 2H), 1.60-1.53 (m, 2H), 1.26-1.24 (m, 16H), 0.87-0.85 (m, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 210.6, 141.2, 140.3, 132.4, 131.4, 129.3, 126.3, 54.8, 43.9, 42.9, 31.6, 30.1, 28.8, 26.9, 23.7, 22.4, 14.0. HRMS Calculated for C₁₉H₃₂NO₃S [M+H]⁺ 354.2097, found: 354.2101.



***N*-tert-Butyl-4-methyl-2-(3-oxo-heptyl)-benzenesulfonamide (5d):** 0.875 g, 65% yield (4 mmol scale), colorless oil. ¹H NMR (400 MHz, CDCl₃) δ 7.90 (d, *J* = 8.5 Hz, 1H), 7.10-7.09 (m, 2H), 4.97 (br, 1H), 3.21 (t, *J* = 7.6 Hz, 2H), 2.85 (t, *J* = 7.5 Hz, 2H), 2.41 (t, *J* = 7.4 Hz, 2H), 2.36 (s, 3H), 1.57-1.53 (m, 2H), 1.32-1.27 (m, 2H), 1.23 (s, 9H), 0.89 (t, *J* = 7.3 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 210.6, 143.0, 140.1, 138.3, 132.1, 129.6, 127.0, 54.8, 44.0, 42.6, 30.1, 27.0, 25.9, 22.3, 21.3, 13.9. HRMS Calculated for C₁₈H₃₀NO₃S [M+H]⁺ 340.1941, found: 340.1940.

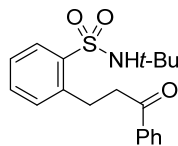


***N*-tert-Butyl-4-methoxy-2-(3-oxo-heptyl)-benzenesulfonamide (5e):** 1.071 g, 75% yield (4 mmol scale), colorless oil. ¹H NMR (400 MHz, CDCl₃) δ 7.96 (d, *J* = 8.5 Hz, 1H), 6.80-6.76 (m, 2H), 4.76 (br, 1H), 3.84 (s, 3H), 3.21 (t, *J* = 7.6 Hz, 2H), 2.85 (t, *J* = 7.6 Hz, 2H), 2.41 (t, *J* = 7.5 Hz, 2H), 1.57-1.53 (m, 2H), 1.31-1.26 (m, 2H), 1.22 (s, 9H), 0.89 (t, *J* = 7.3 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 210.5, 162.6, 142.6, 133.0, 132.0, 117.0, 111.0, 55.6, 54.8, 44.0, 42.7, 30.2, 27.3,



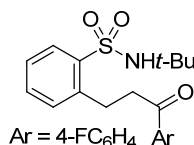
26.0, 22.4, 13.9. HRMS Calculated for $C_{18}H_{30}NO_4S$ $[M+H]^+$ 356.1890, found: 356.1898.

***N*-tert-Butyl-2-(3-oxo-3-phenylpropyl)-benzenesulfonamide (5f):** 1.284 g, 62% yield (6 mmol scale), colorless oil. 1H NMR (400 MHz, $CDCl_3$) δ 8.06 (d, J = 7.9 Hz, 1H), 7.96 (d, J = 7.5 Hz,



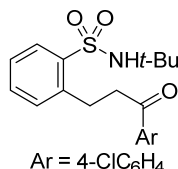
2H), 7.56-7.53 (m, 1H), 7.48-7.30 (m, 5H), 4.92 (br, 1H), 3.43 (s, 4H), 1.24 (s, 9H); ^{13}C NMR (100 MHz, $CDCl_3$) δ 199.4, 141.2, 140.3, 136.7, 133.2, 132.6, 131.6, 129.5, 128.6, 128.2, 126.5, 55.0, 40.2, 30.2, 27.5. HRMS Calculated for $C_{19}H_{24}NO_3S$ $[M+H]^+$ 346.1471, found: 346.1479.

***N*-tert-Butyl-2-(3-(4-fluorophenyl)-3-oxo-propyl)-4-methoxybenzenesulfonamide (5g):** 0.317 g, 49% yield (1.8 mmol scale), colorless oil. 1H NMR (400 MHz, $CDCl_3$) δ 8.11-7.97 (m, 3H),



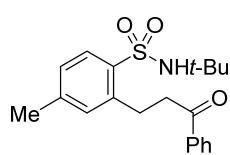
7.48-7.45 (m, 1H), 7.38-7.28 (m, 2H), 7.09 (t, J = 8.5 Hz, 2H), 5.01 (br, 1H), 3.41 (dd, J = 12.0, 5.4 Hz, 4H), 1.24 (s, 9H); ^{13}C NMR (100 MHz, $CDCl_3$) δ 197.9, 165.8 (d, J = 254.7 Hz), 141.2, 140.2, 133.1 (d, J = 3.0 Hz), 132.6, 131.7, 130.8 (d, J = 9.3 Hz), 129.5, 126.6, 115.7 (d, J = 21.9 Hz), 55.0, 40.2, 30.2, 27.6; ^{19}F NMR (376 MHz, $CDCl_3$) δ -105.11. HRMS Calculated for $C_{19}H_{23}FNO_3S$ $[M+H]^+$ 364.1377, found: 364.1375.

***N*-tert-Butyl-2-(3-(4-chlorophenyl)-3-oxopropyl)-benzenesulfonamide (5h):** 0.489 g, 54% yield (2.4 mmol scale), colorless oil. 1H NMR (400 MHz, $CDCl_3$) δ 8.05 (d, J = 7.9 Hz, 1H), 7.89 (d, J = 8.5 Hz, 2H), 7.48-7.30 (m, 5H), 4.98 (br, 1H), 3.45-3.36 (m, 4H), 1.24 (s, 9H);



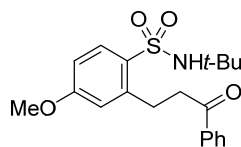
^{13}C NMR (100 MHz, $CDCl_3$) δ 198.2, 141.2, 140.1, 139.6, 135.0, 132.6, 131.7, 129.6, 129.5, 128.9, 126.6, 55.0, 40.2, 30.2, 27.6. HRMS Calculated for $C_{19}H_{23}ClNO_3S$ $[M+H]^+$ 380.1082, found: 380.1084.

***N*-tert-Butyl-4-methyl-2-(3-oxo-3-phenylpropyl)-benzenesulfonamide (5i):** 0.714 g, 66% yield (3 mmol scale), colorless oil. 1H NMR (400 MHz, $CDCl_3$) δ 7.97-7.91 (m, 3H), 7.55-7.51 (m, 1H),



7.45-7.41 (m, 2H), 7.17-7.15 (m, 1H), 7.11 (d, J = 8.1 Hz, 1H), 4.87 (br, 1H), 3.44-3.35 (m, 4H), 2.36 (s, 3H), 1.23 (s, 9H); ^{13}C NMR (100 MHz, $CDCl_3$) δ 199.6, 143.2, 140.3, 138.4, 136.8, 133.2, 132.4, 129.8, 128.7, 128.2, 127.2, 54.9, 40.4, 30.3, 27.6, 21.4. HRMS Calculated for $C_{20}H_{26}NO_3S$ $[M+H]^+$ 360.1628, found: 360.1626.

***N*-tert-Butyl-4-methoxy-2-(3-oxo-3-phenylpropyl)-benzenesulfonamide (5j):** 0.735 g, 65% yield (3 mmol scale), colorless oil. 1H NMR (400 MHz, $CDCl_3$) δ 8.01-7.96 (m, 3H), 7.56 (t, J =



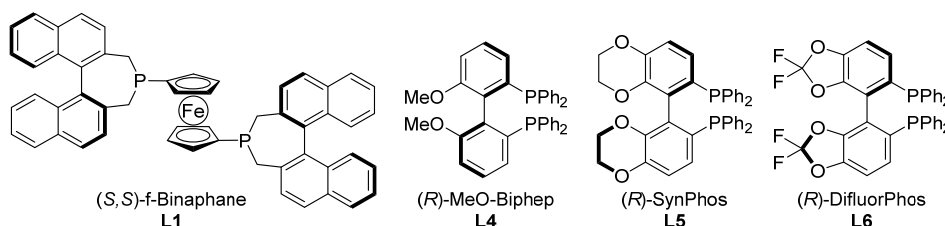
7.4 Hz, 1H), 7.45 (t, J = 7.6 Hz, 2H), 6.87 (d, J = 2.6 Hz, 1H), 6.80 (dd, J = 8.9, 2.6 Hz, 1H), 4.65 (br, 1H), 3.84 (s, 3H), 3.46-3.36 (m, 4H), 1.23 (s, 9H); ^{13}C NMR (100 MHz, $CDCl_3$) δ 199.5, 162.7, 142.7, 136.8, 133.3, 133.1, 132.2, 128.8, 128.3, 117.2, 111.1, 55.6, 54.9, 40.3, 30.3, 27.9. HRMS Calculated for $C_{20}H_{26}NO_4S$ $[M+H]^+$ 376.1577, found: 376.1583.

3. The Procedure for Intramolecular Asymmetric Reductive Amination

Table S1. Optimization of the reaction conditions for the synthesis of γ -sultams **2a**^a

entry	solvent	L	acid	yield (%) ^b	ee (%) ^c
1	TFE	L4	--	--	--
2	TFE	L4	PhCOOH	--	--
3	TFE	L4	TFA	--	--
4	TFE	L4	TsOH·H ₂ O	90	88
5	TFE	L4	L-CSA	94	88
6	TFE	L4	D-CSA	96	89
7	DCM	L4	D-CSA	51	86
8	Toluene	L4	D-CSA	82	80
9	TFE	L5	D-CSA	94	88
10	TFE	L6	D-CSA	92	93
11	TFE	L1	D-CSA	96	97

^a Conditions: **1a** (0.2 mmol), Pd(OCOCF₃)₂ (3.0 mol %), ligand (3.3 mol %), acid (100 mol %), H₂ (600 psi), solvent (3.0 mL), 50 °C, 24 h. ^b Isolated yield. ^c Determined by HPLC.



The optimal reaction conditions were established: Pd(OCOCF₃)₂/(*S,S*)-f-Binaphane as the catalyst, D-CSA as the additive, TFE as the solvent and reaction temperature 50 °C.

Intramolecular asymmetric reductive amination of keto sulfonamides **1** were performed according to the following procedures.

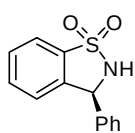


Bisphosphine ligand (*S,S*)-f-Binaphane (5.3 mg, 0.0066 mmol) and Pd(OCOCF₃)₂ (2.0 mg, 0.006 mmol) were placed in a dried Schlenk tube under nitrogen atmosphere, and degassed anhydrous acetone was added. The mixture was stirred at room temperature for 1 h. The solvent was removed under vacuum to give the catalyst. This catalyst was taken into a glove box filled with nitrogen and dissolved in 2,2,2-trifluoroethanol (TFE, 1.0 mL). To the mixture of keto sulfonamides **1** (0.2 mmol) and D-(-)-camphorsulfonic acid (46 mg, 0.2 mmol) in 2,2,2-trifluoroethanol was added this catalyst solution, and then the mixture was transferred to an autoclave, which was charged hydrogen gas (600 psi). The autoclave was stirred under directed condition (oil bath temperature was showed if it was heated). After release of the hydrogen, the autoclave was

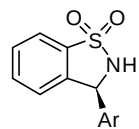
opened and the reaction mixture was evaporated. Purification was performed on silica gel using *n*-hexane/ ethyl acetate as the eluent to give the chiral products **2**. The enantiomeric excesses were determined by chiral HPLC. The absolute configurations of **2** were determined by comparison of rotation sign with literature data.

Racemates of **2** were prepared by the reduction of the corresponding imines using sodium borohydride in methanol.

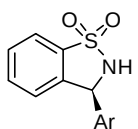
(+)-3-Phenyl-1,2-benzisothiazoline 1,1-dioxide (2a): (Known compound),^{1,2} 47 mg, 96% yield (0.2 mmol scale), 98% ee (S), white solid, $[\alpha]_D^{20} = +88.33$ (c 0.54, CHCl₃) [lit.¹ $[\alpha]_D^{20} = +101.50$ (c 1.50, CHCl₃), 97% ee (S)]. ¹H NMR (400 MHz, CDCl₃) δ 7.82-7.79 (m, 1H), 7.56-7.50 (m, 2H), 7.40-7.33 (m, 5H), 7.15-7.12 (m, 1H), 5.71 (d, *J* = 4.0 Hz, 1H), 5.10 (br, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 139.9, 138.8, 134.9, 133.4, 129.6, 129.4, 129.2, 127.7, 125.5, 121.2, 61.5. HPLC: Chiracel OJ-H column, 254 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 21.9 min (maj) and 24.3 min.



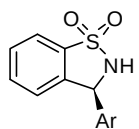
(+)-3-(*o*-Tolyl)-1,2-benzisothiazoline 1,1-dioxide (2b): (Known compound),^{1,2} 51 mg, 98% yield (0.2 mmol scale), 94% ee (S), white solid, $[\alpha]_D^{20} = +10.80$ (c 1.10, CHCl₃) [lit.¹ $[\alpha]_D^{20} = +12.50$ (c 1.00, CHCl₃), 99% ee (S)]. ¹H NMR (400 MHz, CDCl₃) δ 7.83-7.81 (m, 1H), 7.59-7.53 (m, 2H), 7.28-7.23 (m, 2H), 7.19-7.15 (m, 1H), 7.13-7.10 (m, 2H), 6.00 (d, *J* = 4.4 Hz, 1H), 5.00 (d, *J* = 4.2 Hz, 1H), 2.44 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 140.2, 136.8, 136.3, 135.6, 133.4, 131.3, 129.5, 129.1, 128.2, 127.1, 125.3, 121.3, 58.4, 19.5. HPLC: Chiracel OD-H column, 254 nm, 30 °C, *n*-hexane /*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 13.9 min (maj) and 16.8 min.



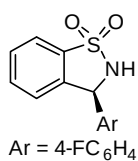
(+)-3-(*m*-Tolyl)-1,2-benzisothiazoline 1,1-dioxide (2c): (Known compound),^{1,2} 47 mg, 90% yield (0.2 mmol scale), 97% ee (S), white solid, $[\alpha]_D^{20} = +100.27$ (c 1.10, CHCl₃) [lit.¹ $[\alpha]_D^{20} = +93.49$ (c 1.00, CHCl₃), 96% ee (S)]. ¹H NMR (400 MHz, CDCl₃) δ 7.79-7.77 (m, 1H), 7.55-7.48 (m, 2H), 7.27-7.23 (m, 1H), 7.16-7.12 (m, 4H), 5.67 (s, 1H), 5.30 (s, 1H), 2.32 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 140.0, 139.1, 138.8, 134.7, 133.3, 129.8, 129.4, 129.1, 128.1, 125.4, 124.7, 121.1, 61.3, 21.4. HPLC: Chiracel OJ-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 18.2 min (maj) and 20.5 min.



(+)-3-(*p*-Tolyl)-1,2-benzisothiazoline 1,1-dioxide (2d): (Known compound),^{1,2} 48 mg, 92% yield (0.2 mmol scale), 83% ee (S), white solid, $[\alpha]_D^{20} = +75.54$ (c 1.10, CHCl₃) [lit.¹ $[\alpha]_D^{20} = +90.89$ (c 1.00, CHCl₃), 98% ee (S)]. ¹H NMR (400 MHz, CDCl₃) δ 7.82-7.78 (m, 1H), 7.56-7.50 (m, 2H), 7.27-7.23 (m, 2H), 7.18 (d, *J* = 7.8 Hz, 2H), 7.14-7.12 (d, *J* = 7.9 Hz, 1H), 5.68 (d, *J* = 4.0 Hz, 1H), 5.11 (d, *J* = 3.3 Hz, 1H), 2.34 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 140.2, 139.1, 135.8, 134.9, 133.4, 130.0, 129.5, 127.6, 125.5, 121.2, 61.3, 21.3. HPLC: Chiracel OD-H column, 254 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 22.1 min and 26.1 min (maj).

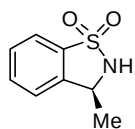


(+)-3-(4-Fluoro-phenyl)-1,2-benzisothiazoline 1,1-dioxide (2e): (Known compound),^{1,2} 49 mg, 92% yield (0.2 mmol scale), 96% ee (S), white solid, $[\alpha]_D^{20} = +94.49$ (c 1.20, CHCl₃) [lit.¹ $[\alpha]_D^{20} = +92.10$ (c 1.00, CHCl₃), 97% ee (S)]. ¹H NMR (400 MHz, CDCl₃) δ 7.85-7.80 (m, 1H), 7.60-7.53 (m, 2H), 7.38-7.33 (m, 2H), 7.15-7.12 (m, 1H), 7.10-7.04 (m, 2H), 5.73 (d, *J* = 4.2 Hz, 1H), 5.15 (d, *J* = 3.5 Hz, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 163.1 (d, *J* = 248.5 Hz), 139.7 (d, *J*_{C-F} = 0.9 Hz), 134.9, 134.7 (d, *J* = 3.2 Hz),

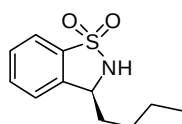


133.6, 129.8, 129.6 (d, J_{C-F} = 8.4 Hz), 125.4, 121.3, 116.4 (d, J_{C-F} = 21.8 Hz), 60.7 (d, J_{C-F} = 0.7 Hz); ^{19}F NMR (376 MHz, CDCl_3) δ -112.2. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 13.1 min (maj) and 20.4 min.

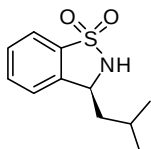
(-)-3-Methyl-1,2-benzisothiazoline 1,1-dioxide (2f): (Known compound),^{1,2} 37 mg, 95% yield (0.2 mmol scale), 95% ee (S), white solid, $[\alpha]_D^{20}$ = -18.96 (c 1.45, CHCl_3) [lit.^{2a} $[\alpha]_D^{20}$ = -20.70 (c 1.00, CHCl_3), 93% ee (S)]. ^1H NMR (400 MHz, CDCl_3) δ 7.74 (d, J = 7.8 Hz, 1H), 7.63-7.59 (m, 1H), 7.52-7.48 (m, 1H), 7.38 (d, J = 7.7 Hz, 1H), 5.09 (br, 1H), 4.81-4.75 (m, 1H), 1.59 (d, J = 6.7 Hz, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 141.9, 135.5, 133.3, 129.2, 124.0, 121.2, 53.5, 21.5. HPLC: Chiracel OD-H column, 254 nm, 30 °C, *n*-hexane/*i*-propanol = 80/20, flow = 0.8 mL/min, retention time 12.9 min (maj) and 16.7 min.



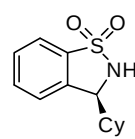
(-)-3-*n*-Butyl-1,2-benzisothiazoline 1,1-dioxide (2g): (Known compound),^{1,2} 43 mg, 96% yield (0.2 mmol scale), 94% ee (S), colorless oil, $[\alpha]_D^{20}$ = -46.92 (c 0.78, CHCl_3) [lit.¹ $[\alpha]_D^{20}$ = -48.70 (c 1.00, CHCl_3), 94% ee (S)]. ^1H NMR (400 MHz, CDCl_3) δ 7.75 (d, J = 7.8 Hz, 1H), 7.62-7.59 (m, 1H), 7.52-7.50 (m, 1H), 7.38 (d, J = 7.8 Hz, 1H), 5.06 (br, 1H), 4.69 (dt, J = 8.7, 4.3 Hz, 1H), 1.99-1.93 (m, 1H), 1.76-1.72 (m, 1H), 1.47-1.32 (m, 4H), 0.91 (t, J = 6.9 Hz, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 140.8, 135.7, 133.2, 129.3, 124.2, 121.4, 58.0, 35.6, 27.9, 22.5, 14.0. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 80/20, flow = 0.7 mL/min, retention time 11.3 min (maj) and 21.2 min.



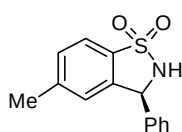
(-)-3-*i*-Butyl-1,2-benzisothiazoline 1,1-dioxide (2h): (Known compound),^{1,2} 44 mg, 98% yield (0.2 mmol scale), 96% ee (S), white solid, $[\alpha]_D^{20}$ = -52.38 (c 0.42, CHCl_3) [lit.¹ $[\alpha]_D^{20}$ = -68.80 (c 0.60, CHCl_3), 97% ee (S)]. ^1H NMR (400 MHz, CDCl_3) δ 7.72 (d, J = 7.7 Hz, 1H), 7.61-7.57 (m, 1H), 7.50-7.46 (m, 1H), 7.35 (d, J = 7.7 Hz, 1H), 5.18 (d, J = 4.8 Hz, 1H), 4.72-4.67 (m, 1H), 1.93-1.82 (m, 1H), 1.74-1.63 (m, 2H), 0.99 (dd, J = 18.4, 6.6 Hz, 6H); ^{13}C NMR (100 MHz, CDCl_3) δ 141.5, 135.6, 133.1, 129.2, 124.2, 121.3, 56.2, 45.2, 25.5, 23.5, 21.4. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 80/20, flow = 0.7 mL/min, retention time 12.2 min (maj) and 25.9 min.



(-)-3-Cyclohexyl-1,2-benzisothiazoline 1,1-dioxide (2i): (Known compound),^{1,2} 49 mg, 96% yield (0.2 mmol scale), 90% ee (S), white solid, $[\alpha]_D^{20}$ = -46.99 (c 1.23, CHCl_3) [lit.¹ $[\alpha]_D^{20}$ = -34.00 (c 0.70, CHCl_3), 97% ee (S)]. ^1H NMR (400 MHz, CDCl_3) δ 7.75 (d, J = 7.8 Hz, 1H), 7.62-7.57 (m, 1H), 7.52-7.48 (t, J = 7.5 Hz, 1H), 7.37 (d, J = 7.8 Hz, 1H), 5.11 (d, J = 4.8 Hz, 1H), 4.64-4.62 (m, 1H), 1.86-1.79 (m, 3H), 1.67-1.63 (m, 2H), 1.31-1.12 (m, 6H); ^{13}C NMR (100 MHz, CDCl_3) δ 139.1, 135.8, 133.1, 129.3, 124.5, 121.4, 63.0, 42.8, 30.8, 26.4, 26.0, 25.9, 25.8. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 75/25, flow = 0.7 mL/min, retention time 9.6 min (maj) and 24.9 min.

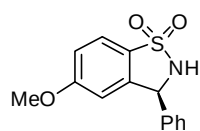


(+)-5-Methyl-3-phenyl-1,2-benzisothiazoline 1,1-dioxide (2j): 49 mg, 94% yield (0.2 mmol), 97% ee (S), white solid, 171-172 °C, $[\alpha]_D^{20}$ = +64.54 (c 1.23, CHCl_3). ^1H NMR (400 MHz, CDCl_3) δ 7.66 (d, J = 8.0 Hz, 1H), 7.39-7.34 (m, 5H), 7.30 (d, J = 8.0 Hz, 1H), 6.89 (s, 1H), 5.66 (d, J = 4.1 Hz, 1H), 5.23 (d, J = 3.7 Hz, 1H), 2.34 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 144.4, 140.2, 139.1, 132.0, 130.5, 129.2, 129.0, 127.6, 125.5, 120.9, 61.2, 21.7. HPLC: Chiracel OD-H column, 220 nm, 30 °C,

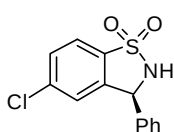


n-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 16.5 min and 17.8 min (maj). HRMS Calculated for C₁₄H₁₄NO₂S [M+H]⁺ 260.0740, found: 260.0742.

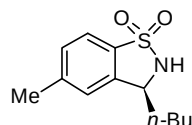
(+)-5-Methoxy-3-phenyl-1,2-benzisothiazoline 1,1-dioxide (2k): 53 mg, 96% yield (0.2 mmol scale), 97% ee (S), white solid, 120-121 °C, [α]_D²⁰ = +21.93 (c 1.40, CHCl₃). ¹H NMR (400 MHz, CDCl₃) δ 7.69 (d, *J* = 8.7 Hz, 1H), 7.36-7.33 (m, 5H), 7.01 (dd, *J* = 8.6, 1.9 Hz, 1H), 6.51 (d, *J* = 1.4 Hz, 1H), 5.64 (s, 1H), 5.17 (s, 1H), 3.75 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 163.8, 142.5, 138.9, 129.3, 129.0, 127.6, 126.9, 122.6, 116.5, 109.2, 61.2, 55.8. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 20.3 min and 22.5 min (maj). HRMS Calculated for C₁₄H₁₄NO₃S [M+H]⁺ 276.0689, found: 276.0691.



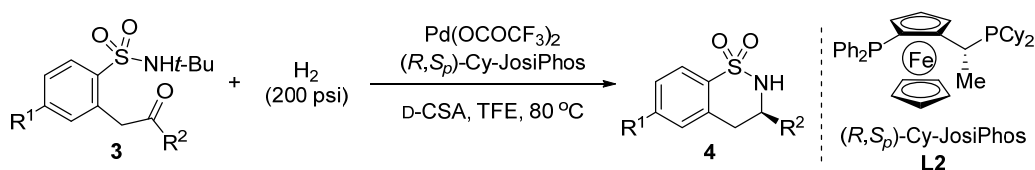
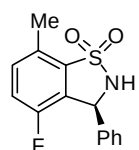
(+)-5-Chloro-3-phenyl-1,2-benzisothiazoline 1,1-dioxide (2l): 53 mg, 95% yield (0.2 mmol scale), 67% ee (S), white solid, 143-144 °C, [α]_D²⁰ = +16.52 (c 1.58, CHCl₃). ¹H NMR (400 MHz, CDCl₃) δ 7.73 (d, *J* = 8.3 Hz, 1H), 7.48 (dd, *J* = 8.3, 1.4 Hz, 1H), 7.42-7.34 (m, 5H), 7.09 (s, 1H), 5.67 (s, 1H), 5.28 (s, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 141.9, 139.8, 138.0, 133.3, 130.1, 129.4, 129.4, 127.6, 125.6, 122.5, 60.9. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 18.4 min and 22.0 min (maj). HRMS Calculated for C₁₃H₁₁Cl NO₂S [M+H]⁺ 280.0194, found: 280.0194.



(-)-5-Methyl-3-butyl-1,2-benzisothiazoline 1,1-dioxide (2m): 46 mg, 96% yield (0.2 mmol scale), 95% ee (S), white solid, 67-68 °C, [α]_D²⁰ = -66.52 (c 1.15, CHCl₃). ¹H NMR (400 MHz, CDCl₃) δ 7.63 (d, *J* = 8.0 Hz, 1H), 7.30 (d, *J* = 8.0 Hz, 1H), 7.16 (s, 1H), 5.00 (d, *J* = 4.8 Hz, 1H), 4.66-4.62 (m, 1H), 2.45 (s, 3H), 2.00-1.92 (m, 1H), 1.74-1.72 (m, 1H), 1.44-1.35 (m, 4H), 0.93-0.90 (m, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 144.1, 141.1, 132.9, 130.2, 124.4, 121.0, 57.8, 35.6, 27.9, 22.4, 21.8, 13.9. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 7.9 min (maj) and 9.4 min. HRMS Calculated for C₁₂H₁₈NO₂S [M+H]⁺ 240.1053, found: 240.1055.



(+)-4-Fluoro-7-methyl-3-phenyl-1,2-benzisothiazoline 1,1-dioxide (2n): 54 mg, 96% yield (0.2 mmol scale), 89% ee (S), white solid, 162-164 °C, [α]_D²⁰ = +113.33 (c 1.08, CHCl₃). ¹H NMR (400 MHz, CDCl₃) δ 7.38-7.33 (m, 5H), 7.29-7.26 (m, 1H), 7.12-7.08 (m, 1H), 5.73 (d, *J* = 3.5 Hz, 1H), 5.14-5.11 (br, 1H), 2.63 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 156.1 (d, *J* = 252.6 Hz), 137.9, 135.2 (d, *J* = 3.2 Hz), 133.3 (d, *J* = 6.5 Hz), 129.9 (d, *J* = 4.4 Hz), 129.1, 129.1, 127.5 (d, *J* = 0.7 Hz), 126.6 (d, *J* = 18.7 Hz), 120.2 (d, *J* = 19.6 Hz), 58.1, 16.3; ¹⁹F NMR (376 MHz, CDCl₃) δ -119.92. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 11.5 min and 16.3 min (maj). HRMS Calculated for C₁₄H₁₃FNO₂S [M+H]⁺ 278.0646, found: 278.0647.



Bisphosphine ligand (*R,S*)-Cy-JosiPhos (3.9 mg, 0.0066 mmol) and Pd(OCOCF₃)₂ (2.0 mg, 0.006 mmol) were placed in a dried Schlenk tube under nitrogen atmosphere, and degassed

anhydrous acetone was added. The mixture was stirred at room temperature for 1 h. The solvent was removed under vacuum to give the catalyst. This catalyst was taken into a glove box filled with nitrogen and dissolved in 2,2,2-trifluoroethanol (TFE, 1.0 mL). To the mixture of keto sulfonamides **3** (0.2 mmol) and D-(-)-camphorsulfonic acid (46 mg, 0.2 mmol) in 2,2,2-trifluoroethanol was added this catalyst solution, and then the mixture was transferred to an autoclave, which was charged hydrogen gas (200 psi). The autoclave was stirred under directed condition (oil bath temperature was showed if it was heated). After release of the hydrogen, the reaction mixture was evaporated. Purification was performed on silica gel using *n*-hexane/ethyl acetate as the eluent to give the chiral products **4**. The enantiomeric excesses were determined by chiral HPLC. The absolute configurations of aryl substituents of **4** were determined by comparison of rotation sign with literature data. The absolute configurations of alkyl substituents of **4** were based on single-crystal X-ray diffraction analysis.

Racemates of **4** were prepared by the reduction amination of the corresponding keto sulfonamides catalyzed by racemic catalyst.

(+)-3,4-Dihydro-3-phenyl-2H-1λ⁶-benzo[e][1,2]thiazine 1,1-dioxide (4a): (Known compound),³ 50 mg, 96% yield (0.2 mmol scale), 97% ee (*R*), white solid, $[\alpha]_{\text{D}}^{20} = +37.28$ (*c* 1.25, CHCl₃), [lit.³ $[\alpha]_{\text{D}}^{25} = +51.70$ (*c* 0.93, CHCl₃), 98% ee (*R*)]. ¹H NMR (400 MHz, CDCl₃) δ 7.77-7.75 (m, 1H), 7.44-7.39 (m, 5H), 7.36-7.34 (m, 2H), 7.25-7.21 (m, 1H), 5.03 (d, *J* = 10.5 Hz, 1H), 4.93 (td, *J* = 10.3, 5.7 Hz, 1H), 3.22-3.19 (m, 2H); ¹³C NMR (100 MHz, CDCl₃) δ 139.6, 137.8, 135.1, 132.3, 129.5, 129.1, 128.5, 127.9, 126.3, 123.7, 56.9, 35.7. HPLC: Chiracel OD-H column, 230 nm, 30 °C, *n*-hexane/ *i*-propanol = 70/30, flow = 0.7 mL/min, retention time 14.4 min and 18.1 min (maj).

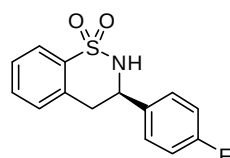
(+)-3,4-Dihydro-3-(*o*-tolyl)-2H-1λ⁶-benzo[e][1,2]thiazine 1,1-dioxide (4b): (Known compound),³ 49 mg, 89% yield (0.2 mmol scale), 96% ee (*R*), white solid, $[\alpha]_{\text{D}}^{20} = +8.86$ (*c* 1.23, CHCl₃), [lit.³ $[\alpha]_{\text{D}}^{24} = +11.40$ (*c* 0.77, CHCl₃), 98% ee (*R*)]. ¹H NMR (400 MHz, CDCl₃) δ 7.73 (d, *J* = 7.8 Hz, 1H), 7.43-7.40 (m, 2H), 7.34-7.20 (m, 5H), 5.20-5.00 (m, 2H), 3.34-3.27 (m, 1H), 3.06-3.01 (m, 1H), 2.41 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 137.5, 137.2, 136.2, 135.3, 132.3, 131.1, 129.6, 128.4, 127.9, 126.7, 125.2, 123.9, 53.4, 34.3, 19.2. HPLC: Chiracel OD-H column, 230 nm, 30 °C, *n*-hexane/ *i*-propanol = 70/30, flow = 0.7 mL/min, retention time 13.9 min and 18.1 min (maj).

(+)-3,4-Dihydro-3-(*m*-tolyl)-2H-1λ⁶-benzo[e][1,2]thiazine 1,1-dioxide (4c): (Known compound),³ 52 mg, 95% yield (0.2 mmol scale), 95% ee (*R*), white solid, $[\alpha]_{\text{D}}^{20} = +39.15$ (*c* 0.94, CHCl₃), [lit.³ $[\alpha]_{\text{D}}^{24} = +46.20$ (*c* 0.63, CHCl₃), 97% ee (*R*)]. ¹H NMR (400 MHz, CDCl₃) δ 7.75 (d, *J* = 7.8 Hz, 1H), 7.43 (td, *J* = 7.6, 1.2 Hz, 1H), 7.35-7.28 (m, 2H), 7.27-7.16 (m, 4H), 5.09-5.06 (br, 1H), 4.94-4.87 (m, 1H), 3.25-3.13 (m, 2H), 2.39 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 139.6, 138.9, 137.8, 135.3, 132.4, 129.6, 129.3, 129.1, 127.9, 127.2, 123.8, 123.3, 56.9, 35.7, 21.7. HPLC: Chiracel OD-H column, 230 nm, 30 °C, *n*-hexane/ *i*-propanol = 70/30, flow = 0.7 mL/min, retention time 12.4 min and 15.5 min (maj).

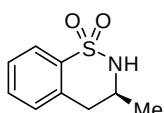
(+)-3,4-Dihydro-3-(*p*-tolyl)-2H-1λ⁶-benzo[e][1,2]thiazine 1,1-dioxide (4d): (Known compound),³ 52 mg, 95% yield (0.2 mmol scale), 79% ee (*R*), white solid, $[\alpha]_{\text{D}}^{20} = +31.44$ (*c* 1.04, CHCl₃), [lit.³ $[\alpha]_{\text{D}}^{24} = +18.70$ (*c* 0.47, CHCl₃), 97% ee (*R*)]. ¹H NMR (400 MHz, CDCl₃) δ 7.73 (d, *J* = 7.8

Hz, 1H), 7.42 (t, $J = 7.5$ Hz, 1H), 7.34-7.28 (m, 3H), 7.25-7.18 (m, 3H), 5.10-5.07 (br, 1H), 4.91-4.84 (m, 1H), 3.23-3.10 (m, 2H), 2.35 (s, 3H); ^{13}C NMR (10 MHz, CDCl_3) δ 138.2, 137.7, 136.7, 135.2, 132.2, 129.7, 129.6, 127.8, 126.2, 123.8, 56.5, 35.6, 21.2. HPLC: Chiracel OD-H column, 230 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 12.9 min and 18.1 min (maj).

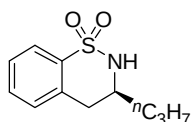
(+)-3,4-Dihydro-3-(4-fluorophenyl)-2H-1 λ ⁶-benzo[*e*][1,2]thiazine 1,1-dioxide (4e): (Known compound),³ 55 mg, 98% yield (0.2 mmol scale), 98% ee (*R*), white solid, $[\alpha]_{\text{D}}^{20} = +34.74$ (c 1.35, CHCl_3), [lit.³ $[\alpha]_{\text{D}}^{24} = +59.20$ (c 0.50, CHCl_3), 98% ee (*R*)]. ^1H NMR (400 MHz, CDCl_3) δ 7.73-7.71 (m, 1H), 7.45-7.40 (m, 3H), 7.34-7.31 (m, 1H), 7.22 (d, $J = 7.7$ Hz, 1H), 7.12-7.06 (m, 2H), 5.19-5.16 (br, 1H), 4.92-4.86 (m, 1H), 3.24-3.14 (m, 2H); ^{13}C NMR (100 MHz, CDCl_3) δ 162.5 (d, $J = 247.2$ Hz), 137.5, 135.5 (d, $J = 3.1$ Hz), 134.9, 132.3, 129.4, 128.1 (d, $J = 8.3$ Hz), 127.8, 123.5, 115.9 (d, $J = 21.6$ Hz), 56.2, 35.5; ^{19}F NMR (376 MHz, CDCl_3) δ -113.56. HPLC: Chiracel OD-H column, 230 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 12.3 min and 17.4 min (maj).



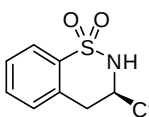
(+)-3-Methyl-3,4-dihydro-2H-benzo[*e*][1,2]thiazine 1,1-dioxide (4f): (Known compound),³ 35 mg, 90% yield (0.2 mmol scale), 94% ee, white solid, $[\alpha]_{\text{D}}^{20} = +80.88$ (c 0.68, CHCl_3), [lit.³ $[\alpha]_{\text{D}}^{24} = -69.40$ (c 0.53, CHCl_3), 98% ee]. ^1H NMR (400 MHz, CDCl_3) δ 7.70 (d, $J = 7.7$ Hz, 1H), 7.37 (t, $J = 7.4$ Hz, 1H), 7.29 (t, $J = 7.3$ Hz, 1H), 7.12 (d, $J = 7.6$ Hz, 1H), 4.74-4.71 (br, 1H), 4.02-3.93 (m, 1H), 2.90-2.85 (m, 1H), 2.75-2.68 (m, 1H), 1.37 (d, $J = 6.5$ Hz, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 137.0, 135.3, 132.0, 129.4, 127.6, 123.8, 49.4, 36.0, 21.7. HPLC: Chiracel OD-H column, 230 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 9.3 min and 17.7 min (maj).



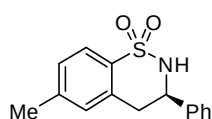
(+)-3-*n*-Propyl-3,4-dihydro-2H-benzo[*e*][1,2]thiazine 1,1-dioxide (4g): (Known compound),³ 42 mg, 93% yield (0.2 mmol scale), 95% ee, white solid, $[\alpha]_{\text{D}}^{20} = +57.89$ (c 0.76, CHCl_3), [lit.³ $[\alpha]_{\text{D}}^{24} = -60.40$ (c 0.87, CHCl_3), 96% ee]. ^1H NMR (400 MHz, CDCl_3) δ 7.69-7.67 (m, 1H), 7.35 (t, $J = 7.3$ Hz, 1H), 7.29-7.27 (m, 1H), 7.11 (d, $J = 7.3$ Hz, 1H), 4.65 (t, $J = 11.7$ Hz, 1H), 3.83 (br, 1H), 2.88-2.83 (m, 1H), 2.74-2.67 (m, 1H), 1.61-1.54 (m, 4H), 0.96 (t, $J = 6.8$ Hz, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 137.4, 135.4, 131.9, 129.5, 127.5, 123.8, 53.2, 37.8, 34.5, 18.6, 13.7. HPLC: Chiracel OD-H column, 230 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 8.3 min & 17.8 min (maj).



(+)-3-Cyclohexyl-3,4-dihydro-2H-benzo[*e*][1,2]thiazine 1,1-dioxide (4h): (Known compound),⁴ 49 mg, 92% yield (0.2 mmol scale), 96% ee, white solid, $[\alpha]_{\text{D}}^{20} = +61.00$ (c 0.90, CHCl_3). ^1H NMR (400 MHz, CDCl_3) δ 7.65 (d, $J = 7.7$ Hz, 1H), 7.34 (t, $J = 7.4$ Hz, 1H), 7.25 (t, $J = 7.7$ Hz, 1H), 7.10 (d, $J = 7.7$ Hz, 1H), 4.73-4.70 (m, 1H), 3.64-3.55 (br, 1H), 2.90-2.73 (m, 2H), 2.03-1.98 (m, 1H), 1.81-1.68 (m, 4H), 1.55-1.48 (m, 1H), 1.37-1.02 (m, 5H); ^{13}C NMR (100 MHz, CDCl_3) δ 137.7, 135.8, 132.0, 129.8, 127.6, 123.9, 58.3, 42.5, 31.9, 29.2, 28.9, 26.4, 26.1, 26.0. HPLC: Chiracel OD-H column, 230 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 7.9 min and 14.7 min (maj).

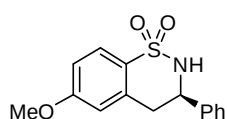


(+)-6-Methyl-3-phenyl-3,4-dihydro-2H-benzo[*e*][1,2]thiazine 1,1-dioxide (4i): 50 mg, 91% yield (0.2 mmol scale), 88% ee (*R*), white solid, mp 210-211 °C, $[\alpha]_{\text{D}}^{20} = +42.24$ (c 1.25, CHCl_3). ^1H NMR (400 MHz, CDCl_3) δ 7.73 (d, $J = 8.0$ Hz, 1H), 7.47-7.36 (m, 5H), 7.21 (d, $J = 8.0$ Hz,



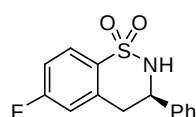
1H), 7.10 (s, 1H), 5.01-4.83 (m, 1H), 4.86-4.83 (br, 1H), 3.29-3.17 (m, 2H), 2.41 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 142.9, 139.6, 135.0, 134.9, 129.8, 129.0, 128.6, 128.4, 126.2, 123.7, 56.8, 35.7, 21.5. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 13.0 min and 15.5 min (maj). HRMS Calculated for C₁₅H₁₆NO₂S [M+H]⁺ 274.0896, found: 274.0896.

(+)-6-Methoxy-3-phenyl-3,4-dihydro-2H-benzo[e][1,2]thiazine 1,1-dioxide (4j): 54 mg, 93% yield (0.2 mmol scale), 89% ee (*R*), white solid, mp 186-187 °C, [α]_D²⁰ = +42.81 (*c* 1.35, CHCl₃).



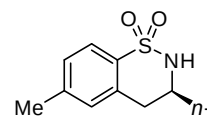
¹H NMR (400 MHz, CDCl₃) δ 7.70 (d, *J* = 8.7 Hz, 1H), 7.43- 7.32 (m, 5H), 6.83 (dd, *J* = 8.7, 2.4 Hz, 1H), 6.68 (d, *J* = 2.2 Hz, 1H), 5.05-5.02 (br, 1H), 4.93 (td, *J* = 10.9, 4.7 Hz, 1H), 3.82 (s, 3H), 3.23-3.10 (m, 2H); ¹³C NMR (100 MHz, CDCl₃) δ 162.1, 139.5, 137.2, 129.9, 129.0, 128.3, 126.2, 125.7, 114.0, 113.6, 56.7, 55.5, 36.1. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 17.5 min and 21.9 min (maj). HRMS Calculated for C₁₅H₁₆NO₃S [M+H]⁺ 290.0845, found: 290.0845.

(+)-6-Fluoro-3-phenyl-3,4-dihydro-2H-benzo[e][1,2]thiazine 1,1-dioxide (4k): 52 mg, 98% yield (0.2 mmol scale), 96% ee (*R*), white solid, mp 172-174 °C, [α]_D²⁰ = +55.18 (*c* 1.08, CHCl₃).



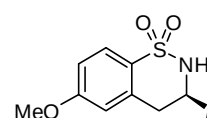
¹H NMR (400 MHz, CDCl₃) δ 7.77 (dd, *J* = 8.6, 5.4 Hz, 1H), 7.42- 7.34 (m, 5H), 7.07-7.02 (m, 1H), 6.97 (d, *J* = 8.9 Hz, 1H), 5.08-5.06 (br, 1H), 4.95 (td, *J* = 10.8, 4.7 Hz, 1H), 3.30-3.16 (m, 2H); ¹³C NMR (100 MHz, CDCl₃) δ 164.3 (d, *J* = 254.4 Hz), 139.1, 138.3 (d, *J* = 8.7 Hz), 133.9 (d, *J* = 3.3 Hz), 129.1, 128.6, 126.4 (d, *J* = 9.4 Hz), 126.2, 116.2 (d, *J* = 22.4 Hz), 115.4 (d, *J* = 22.6 Hz), 56.6, 35.8; ¹⁹F NMR (376 MHz, CDCl₃) δ -105.43. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 13.7 min and 15.6 min (maj). HRMS Calculated for C₁₄H₁₃FNO₂S [M+H]⁺ 278.0646, found: 278.0646.

(+)-3-Butyl-6-methyl-3,4-dihydro-2H-benzo[e][1,2]thiazine 1,1-dioxide (4l): 49 mg, 96% yield (0.2 mmol scale), 96% ee (*S*), colorless oil, [α]_D²⁰ = +57.26 (*c* 1.50, CHCl₃). ¹H NMR (400 MHz,



CDCl₃) δ 7.61 (d, *J* = 8.0 Hz, 1H), 7.11 (d, *J* = 8.0 Hz, 1H), 6.96 (s, 1H), 4.54 (d, *J* = 11.6 Hz, 1H), 3.85- 3.74 (m, 1H), 2.86-2.81 (m, 1H), 2.74-2.67 (m, 1H), 2.34 (s, 3H), 1.66-1.59 (m, 2H), 1.57-1.29 (m, 4H), 0.92 (t, *J* = 7.1 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 142.4, 135.4, 134.7, 129.9, 128.3, 123.9, 53.6, 35.5, 34.7, 27.5, 22.3, 21.5, 14.0. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 6.8 min and 9.9 min (maj). HRMS Calculated for C₁₃H₂₀NO₂S [M+H]⁺ 254.1209, found: 254.1211.

(+)-3-Butyl-6-methoxy-3,4-dihydro-2H-benzo[e][1,2]thiazine 1,1-dioxide (4m): 50 mg, 93% yield (0.2 mmol scale), 96% ee (*S*), colorless oil, [α]_D²⁰ = +67.81 (*c* 1.28, CHCl₃). ¹H NMR (400

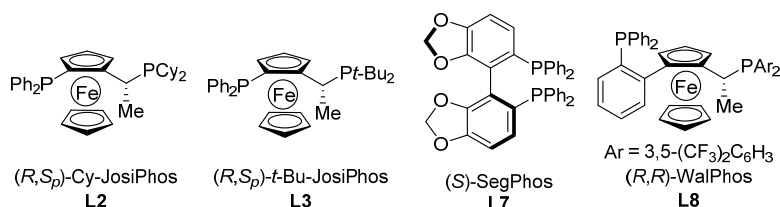


MHz, CDCl₃) δ 7.64 (d, *J* = 8.7 Hz, 1H), 6.80 (dd, *J* = 8.7, 2.5 Hz, 1H), 6.59 (d, *J* = 2.3 Hz, 1H), 4.62-4.59 (m, 1H), 3.84-3.74 (m, 4H), 2.85-2.80 (m, 1H), 2.74-2.67 (m, 1H), 1.68-1.57 (m, 2H), 1.53-1.30 (m, 4H), 0.93 (t, *J* = 7.2 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 161.8, 137.6, 129.8, 125.8, 113.7, 113.6, 55.4, 53.5, 35.5, 35.1, 27.5, 22.3, 14.0. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 70/30, flow = 0.7 mL/min, retention time 8.3 min and 11.3 min (maj). HRMS Calculated for C₁₃H₂₀NO₃S [M+H]⁺ 270.1158, found: 270.1161.

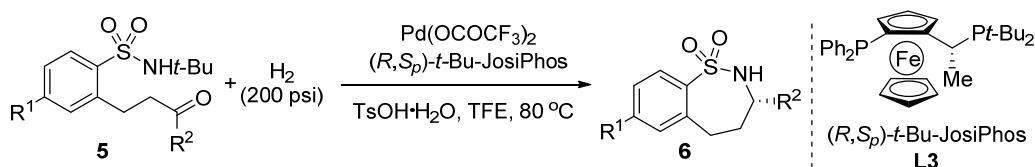
Table S2. Ligand screening for the synthesis of δ -sultams **6a^a**

entry	ligand	conv (%) ^b	ee (%) ^c
1	L2 [(<i>R,S</i> _p)-Cy-JosiPhos]	>95	80 (+)
2	L3 [(<i>R,S</i> _p)- <i>t</i> -Bu-JosiPhos]	>95	98 (+)
3	L4 [(<i>R</i>)-MeO-Biphep]	>95	36 (+)
4	L5 [(<i>R</i>)-SynPhos]	>95	34 (+)
5	L6 [(<i>R</i>)-DifluorPhos]	>95	29 (+)
6	L7 [(<i>S</i>)-SegPhos]	>95	36 (-)
7	L8 [(<i>R,R</i>)-WalPhos]	>95	95 (+)

^a Conditions: **5a** (0.1 mmol), Pd(OCOCF₃)₂ (3.0 mol %), ligand (3.3 mol %), TsOH·H₂O (100 mol %), H₂ (200 psi), TFE (2.0 mL), 80 °C, 24 h. ^b Determined by ¹H NMR. ^c Determined by HPLC.



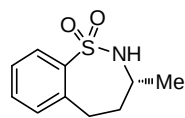
The optimal reaction conditions were established: Pd(OCOCF₃)₂/*(R,S)*_p-*t*-Bu-JosiPhos as the catalyst, TsOH·H₂O as the additive, TFE as the solvent and reaction temperature 80 °C.



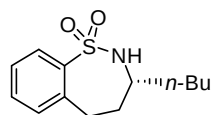
Bisphosphine ligand (*(R,S)*_p)-*t*-Bu-JosiPhos (3.6 mg, 0.0066 mmol) and Pd(OCOCF₃)₂ (2.0 mg, 0.006 mmol) were placed in a dried Schlenk tube under nitrogen atmosphere, and degassed anhydrous acetone was added. The mixture was stirred at room temperature for 1 h. The solvent was removed under vacuum to give the catalyst. This catalyst was taken into a glove box filled with nitrogen and dissolved in 2,2,2-trifluoroethanol (TFE, 1.0 mL). To the mixture of keto sulfonamides **5** (0.2 mmol) and *para*-toluenesulfonic acid monohydrate (38 mg, 0.2 mmol) in 2,2,2-trifluoroethanol was added this catalyst solution, and then the mixture was transferred to an autoclave, which was charged hydrogen gas (200 psi). The autoclave was stirred under directed condition. After release of the hydrogen, the reaction mixture was evaporated. Purification was performed on silica gel using *n*-hexane/ethyl acetate as the eluent to give the chiral products **6**. The enantiomeric excesses were determined by chiral HPLC. The absolute configurations of **6** were based on single-crystal X-ray diffraction analysis.

Racemates of **6** were prepared by the reduction amination of the corresponding keto sulfonamides catalyzed by racemic catalyst.

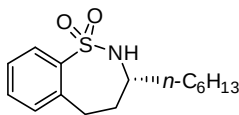
(+)-3-Methyl-2,3,4,5-tetrahydrobenzo[f][1,2]thiazepine 1,1-dioxide (6a): 20 mg, 95% yield (0.1 mmol scale), 98% ee (*R*), colorless oil, $[\alpha]_{\text{D}}^{20} = +28.30$ (*c* 1.00, CHCl₃). ¹H NMR (400 MHz, CDCl₃) δ 7.90-7.81 (m, 1H), 7.43-7.39 (m, 1H), 7.31-7.23 (m, 2H), 4.27 (br, 1H), 4.04-3.99 (m, 1H), 3.66-3.59 (m, 1H), 2.93-2.88 (m, 1H), 2.08-2.03 (m, 1H), 1.49-1.40 (m, 1H), 1.26 (d, *J* = 6.7 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 142.4, 139.4, 132.7, 131.3, 127.0, 126.5, 53.1, 35.6, 34.3, 22.1. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 80/20, flow = 0.7 mL/min, retention time 10.2 min and 13.1 min (maj). HRMS Calculated for C₁₀H₁₄NO₂S [M+H]⁺ 212.0740, found: 121.0740.



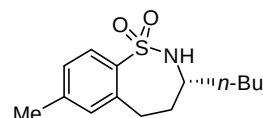
(+)-3-*n*-Butyl-2,3,4,5-tetrahydrobenzo[f][1,2]thiazepine 1,1-dioxide (6b): 50 mg, 98% yield (0.2 mmol scale), 99% ee (*R*), colorless oil, $[\alpha]_{\text{D}}^{20} = +20.46$ (*c* 1.93, CHCl₃). ¹H NMR (400 MHz, CDCl₃) δ 7.71 (d, *J* = 7.8 Hz, 1H), 7.32 (t, *J* = 7.4 Hz, 1H), 7.19-7.15 (m, 2H), 4.26-4.24 (m, 1H), 3.74 (br, 1H), 3.55-3.48 (m, 1H), 2.87-2.82 (m, 1H), 2.0-1.95 (m, 1H), 1.41-1.18 (m, 7H), 0.83 (t, *J* = 6.8 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 142.5, 139.4, 132.5, 131.2, 127.0, 126.4, 57.2, 35.7, 34.3, 34.0, 28.0, 22.4, 14.0. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 90/10, flow = 0.7 mL/min, retention time 9.7 min and 14.2 min (maj). HRMS Calculated for C₁₃H₂₀NO₂S [M+H]⁺ 254.1209, found: 254.1212.



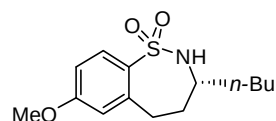
(+)-3-*n*-Hexyl-2,3,4,5-tetrahydrobenzo[f][1,2]thiazepine 1,1-dioxide (6c): 54 mg, 96% yield (0.2 mmol scale), 98% ee (*R*), colorless oil, $[\alpha]_{\text{D}}^{20} = +19.25$ (*c* 2.15, CHCl₃). ¹H NMR (400 MHz, CDCl₃) δ 7.73 (d, *J* = 7.5 Hz, 1H), 7.33 (t, *J* = 7.2 Hz, 1H), 7.19-7.15 (m, 2H), 4.19-4.17 (m, 1H), 3.75 (br, 1H), 3.56-3.50 (m, 1H), 2.87-2.82 (m, 1H), 2.00-1.95 (m, 1H), 1.41-1.21 (m, 11H), 0.80 (d, *J* = 6.6 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 142.5, 139.4, 132.5, 131.2, 127.0, 126.4, 57.2, 36.0, 34.3, 34.0, 31.7, 29.0, 25.8, 22.6, 14.1. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 90/10, flow = 0.7 mL/min, retention time 9.2 min and 15.0 min (maj). HRMS Calculated for C₁₅H₂₄NO₂S [M+H]⁺ 282.1522, found: 282.1525.



(+)-3-*n*-Butyl-7-methyl-2,3,4,5-tetrahydrobenzo[f][1,2]thiazepine 1,1-dioxide (6d): 49 mg, 92% yield (0.2 mmol scale), 98% ee (*R*), white solid, mp 86-88 °C, $[\alpha]_{\text{D}}^{20} = +25.48$ (*c* 0.93, CHCl₃). ¹H NMR (400 MHz, CDCl₃) δ 7.69-7.66 (m, 1H), 7.04-7.01 (m, 2H), 4.28-4.20 (m, 1H), 3.81 (br, 1H), 3.59-3.52 (m, 1H), 2.89-2.83 (m, 1H), 2.35 (s, 3H), 2.05-2.01 (m, 1H), 1.48-1.35 (m, 7H), 0.91 (t, *J* = 7.0 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 143.1, 139.8, 139.3, 132.1, 127.2, 126.9, 57.2, 35.8, 34.3, 34.2, 28.0, 22.4, 21.4, 14.0. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 85/15, flow = 0.7 mL/min, retention time 8.1 min and 9.7 min (maj). HRMS Calculated for C₁₄H₂₂NO₂S [M+H]⁺ 268.1366, found: 268.1367.



(+)-3-*n*-Butyl-7-methoxy-2,3,4,5-tetrahydrobenzo[f][1,2]thiazepine 1,1-dioxide (6e): 56 mg, 98% yield (0.2 mmol scale), >99% ee (*R*), colorless oil, $[\alpha]_{\text{D}}^{20} = +27.33$ (*c* 1.65, CHCl₃). ¹H NMR (400 MHz, CDCl₃) δ 7.72 (d, *J* = 8.6 Hz, 1H), 6.73-6.72 (m, 1H), 6.67 (dd, *J* = 8.7, 2.5 Hz, 1H), 4.22-4.19 (m, 1H), 3.82 (s, 3H), 3.79-3.74 (br, 1H), 3.58-3.52 (m, 1H), 2.86-2.81 (m, 1H), 2.07-2.01 (m, 1H), 1.49-1.24 (m, 7H), 0.90 (t, *J* = 7.1 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 162.5, 141.5, 134.7, 129.3, 117.3, 110.3, 57.1, 55.6, 35.8, 34.5, 34.3, 28.1, 22.4, 14.0. HPLC: Chiracel OD-H column,



220 nm, 30 °C, *n*-hexane/*i*-propanol = 85/15, flow = 0.7 mL/min, retention time 12.1 min and 13.6 min (maj). HRMS Calculated for C₁₄H₂₂NO₃S [M+H]⁺ 284.1315, found: 284.1315.

(+)-3-Phenyl-2,3,4,5-tetrahydrobenzo[f][1,2]thiazepine 1,1-dioxide (6f): 24 mg, 89% yield (0.1 mmol scale), 94% ee (*S*), colorless oil, [α]_D²⁰ = +25.92 (*c* 0.98, CHCl₃). ¹H NMR (400 MHz, CDCl₃) δ 7.92-7.90 (m, 1H), 7.48-7.42 (m, 1H), 7.36-7.25 (m, 7H), 5.03-4.97 (m, 1H), 4.53-4.51 (br, 1H), 3.82-3.75 (m, 1H), 3.06-3.00 (m, 1H), 2.31-2.26 (m, 1H), 2.02-1.90 (m, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 142.1, 140.5, 138.9, 132.8, 131.4, 129.0, 128.3, 126.9, 126.7, 126.6, 60.7, 34.3, 34.3. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 85/15, flow = 0.7 mL/min, retention time 17.2 min and 17.9 min (maj). HRMS Calculated for C₁₅H₁₆NO₂S [M+H]⁺ 274.0896, found: 274.0898.

(+)-3-(4-Fluorophenyl)-2,3,4,5-tetrahydrobenzo[f][1,2]thiazepine 1,1-dioxide (6g): 56 mg, 95% yield (0.2 mmol scale), 93% ee (*S*), colorless oil, [α]_D²⁰ = +13.69 (*c* 1.68, CHCl₃). ¹H NMR (400 MHz, CDCl₃) δ 7.86-7.84 (m, 1H), 7.47-7.44 (m, 1H), 7.35-7.25 (m, 4H), 7.03-6.99 (m, 2H), 4.99-4.94 (m, 1H), 4.69-4.62 (m, 1H), 3.78-3.71 (m, 1H), 3.05-2.99 (m, 1H), 2.29-2.24 (m, 1H), 1.97-1.88 (m, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 162.5 (*d*, *J* = 247.3 Hz), 142.1, 138.9, 136.5 (*d*, *J* = 3.3 Hz), 132.9, 131.5, 128.5 (*d*, *J* = 8.2 Hz), 127.0, 126.8, 115.9 (*d*, *J* = 21.5 Hz), 60.0, 34.4, 34.3; ¹⁹F NMR (376 MHz, CDCl₃) δ -113.46. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 80/20, flow = 0.7 mL/min, retention time 12.9 min and 17.5 min (maj). HRMS Calculated for C₁₅H₁₅FNO₂S [M+H]⁺ 292.0802, found: 292.0803.

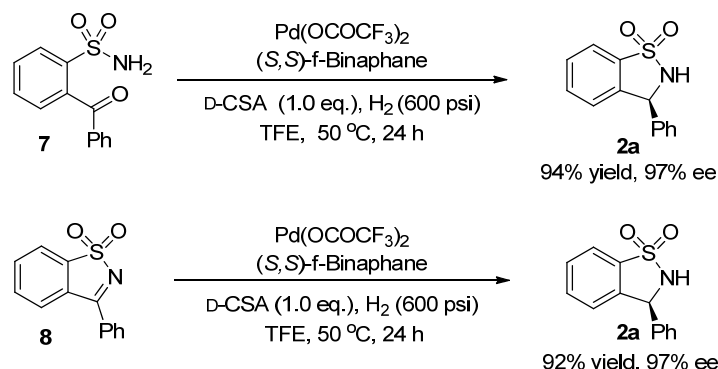
(S)-(+)-3-(4-Chlorophenyl)-2,3,4,5-tetrahydrobenzo[f][1,2]thiazepine 1,1-dioxide (6h): 54 mg, 87% yield (0.2 mmol scale), 97% ee (*S*), white solid, mp 79-80 °C, [α]_D²⁰ = +26.53 (*c* 0.98, CHCl₃). ¹H NMR (400 MHz, CDCl₃) δ 7.92-7.90 (m, 1H), 7.49-7.45 (m, 1H), 7.36-7.26 (m, 6H), 5.02-4.96 (m, 1H), 4.49-4.47 (m, 1H), 3.80-3.74 (m, 1H), 3.07-3.01 (m, 1H), 2.31-2.25 (m, 1H), 1.97-1.87 (m, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 141.9, 138.9, 138.7, 134.1, 132.9, 131.4, 129.1, 128.1, 127.0, 126.8, 59.9, 34.2, 34.1. HPLC: Chiracel OD-H column, 210 nm, 30 °C, *n*-hexane/*i*-propanol = 85/15, flow = 0.7 mL/min, retention time 16.4 min and 22.1 min (maj). HRMS Calculated for C₁₅H₁₅ClNO₂S [M+H]⁺ 308.0507, found: 308.0506.

(+)-3-Phenyl-7-methyl-2,3,4,5-tetrahydrobenzo[f][1,2]thiazepine 1,1-dioxide (6i): 50 mg, 88% yield (0.2 mmol scale), 94% ee (*S*), colorless oil, [α]_D²⁰ = +7.81 (*c* 1.55, CHCl₃). ¹H NMR (400 MHz, CDCl₃) δ 7.80-7.78 (m, 1H), 7.35-7.25 (m, 5H), 7.11-7.10 (m, 2H), 5.00-4.94 (m, 1H), 4.51-4.49 (m, 1H), 3.77-3.70 (m, 1H), 2.99-2.94 (m, 1H), 2.38 (s, 3H), 2.28-2.24 (m, 1H), 1.97-1.87; ¹³C NMR (100 MHz, CDCl₃) δ 143.4, 140.7, 139.4, 138.8, 132.2, 129.0, 128.3, 127.1, 126.7, 60.7, 34.5, 34.3, 21.5. HPLC: Chiracel OJ-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 85/15, flow = 0.7 mL/min, retention time 30.9 min (maj) and 35.8 min. HRMS Calculated for C₁₆H₁₈NO₂S [M+H]⁺ 288.1053, found: 288.1053.

(+)-3-Phenyl-7-methoxy-2,3,4,5-tetrahydrobenzo[f][1,2]thiazepine 1,1-dioxide (6j): 57 mg, 93% yield (0.2 mmol scale), 95% ee (*S*), colorless oil, [α]_D²⁰ = +7.16 (*c* 1.48, CHCl₃). ¹H NMR (400 MHz, CDCl₃) δ 7.90-7.88 (m, 1H), 7.37-7.28 (m, 5H), 6.81-6.78 (m, 2H), 5.01-4.95 (m, 1H), 4.38-4.36 (br, 1H), 3.86 (s, 3H), 3.80-3.73 (m, 1H), 2.99-2.93 (m, 1H), 2.32-2.27 (m, 1H),

1.98-1.89 (m, 1H); ^{13}C NMR (100 MHz, CDCl_3) δ 162.7, 141.0, 140.6, 134.2, 129.2, 129.0, 128.3, 126.6, 117.5, 110.4, 60.5, 55.6, 34.5, 34.5. HPLC: Chiracel OD-H column, 220 nm, 30 °C, *n*-hexane/*i*-propanol = 85/15, flow = 0.7 mL/min, retention time 22.1 min (maj) and 28.2 min. HRMS Calculated for $\text{C}_{16}\text{H}_{18}\text{NO}_3\text{S}$ $[\text{M}+\text{H}]^+$ 304.1002, found: 304.1001.

4. Control Experiments



The *in situ* prepared bisphosphine ligand and $\text{Pd}(\text{OCOCF}_3)_2$ (2.0 mg, 0.006 mmol) complex was taken into a glove box and added to the mixture of keto sulfonamide **7** or imine **8** (0.2 mmol) and acid in dichloromethane or 2,2,2-trifluoroethanol. Then, the mixture was transferred to an autoclave, which was charged hydrogen gas (600 psi). The autoclave was stirred for 24 h. After release of the hydrogen, the autoclave was opened and the reaction mixture was evaporated. The reaction mixture was evaporated. Purification was performed on silica gel using *n*-hexane/ethyl acetate as the eluent to give the chiral product **2a**. The enantiomeric excesses were determined by chiral HPLC.

5. Determination of Absolute Configurations of Products

The single crystal of (+)-3-cyclohexyl-3,4-dihydro-2*H*-benzo[*e*][1,2]thiazine 1,1-dioxide **4h** was grown from the solution of *n*-hexane and ethyl acetate, which is suitable for X-ray diffraction analysis. The structure in **Figure S1** showed that the absolute configuration of (+)-**4h** is (*R*). [CCDC 1507240] contains the structure and supplementary crystallographic data for (+)-**4h**. These data can be obtained free of charge from the Cambridge Crystallographic Data Centre via www.ccdc.cam.ac.uk.

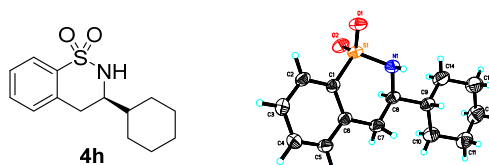


Figure S1. X-ray crystallographic analysis of (*R*)-(+)-**4h**

To determine the absolute configuration of the product (+)-3-(4-chlorophenyl)-2,3,4,5-tetrahydrobenzo[*f*][1,2]thiazepine 1,1-dioxide (+)-**6h**, *n*-hexane (3.0 mL) was slowly added into the solution of (+)-**6h** in ethyl acetate (1.0 mL) at 50 °C, then the solution was cooled down to room temperature. The crystal of (+)-**6h** was grown from the solution, which is suitable for X-ray diffraction analysis. The structure in **Figure S2** showed that the absolute configuration of (+)-**6h** is

S. [CCDC 1414841] contains the structure and supplementary crystallographic data for (+)-**6h**. These data can be obtained free of charge from the Cambridge Crystallographic Data Centre via www.ccdc.cam.ac.uk.

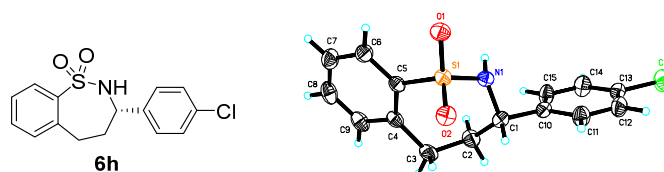
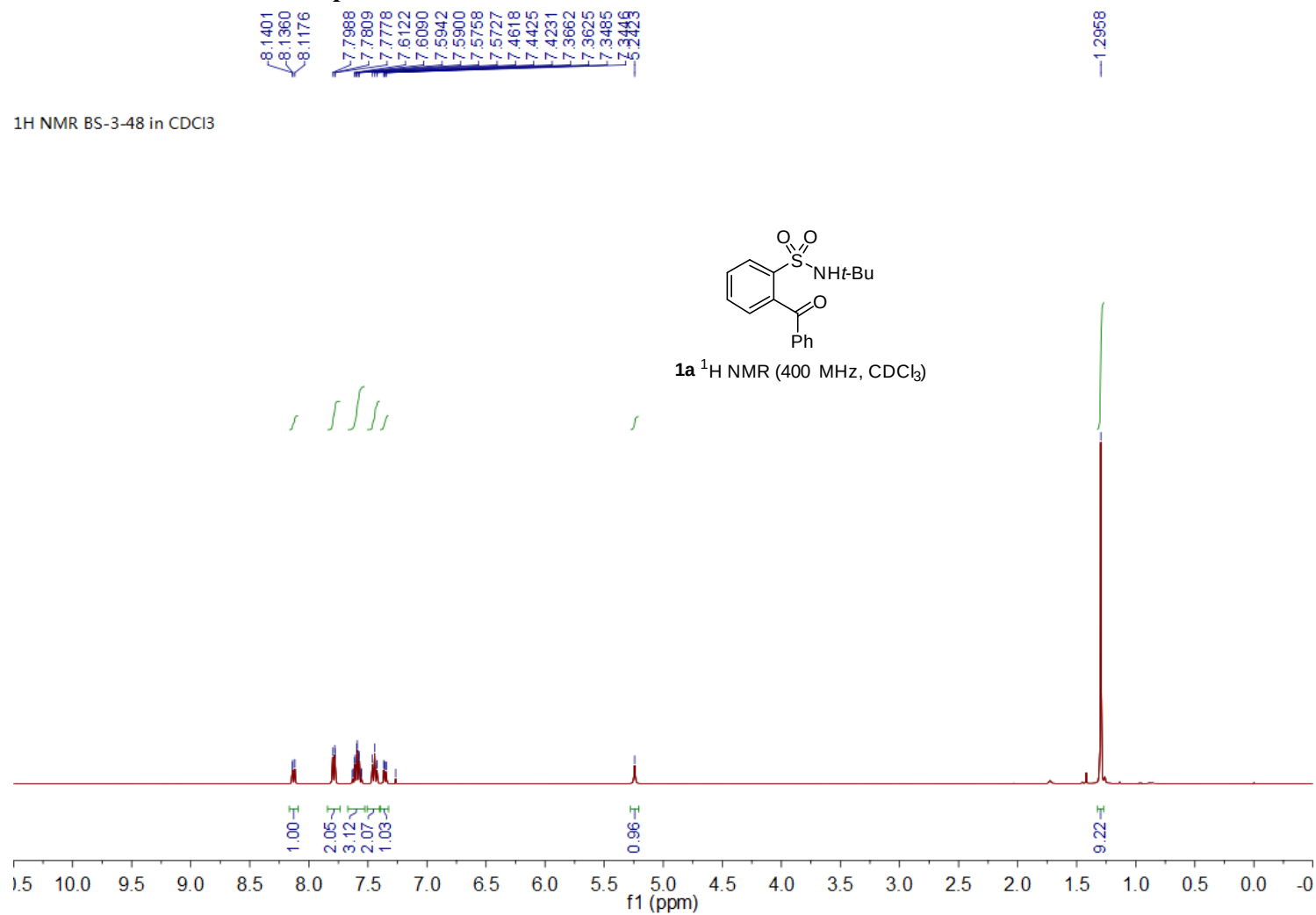


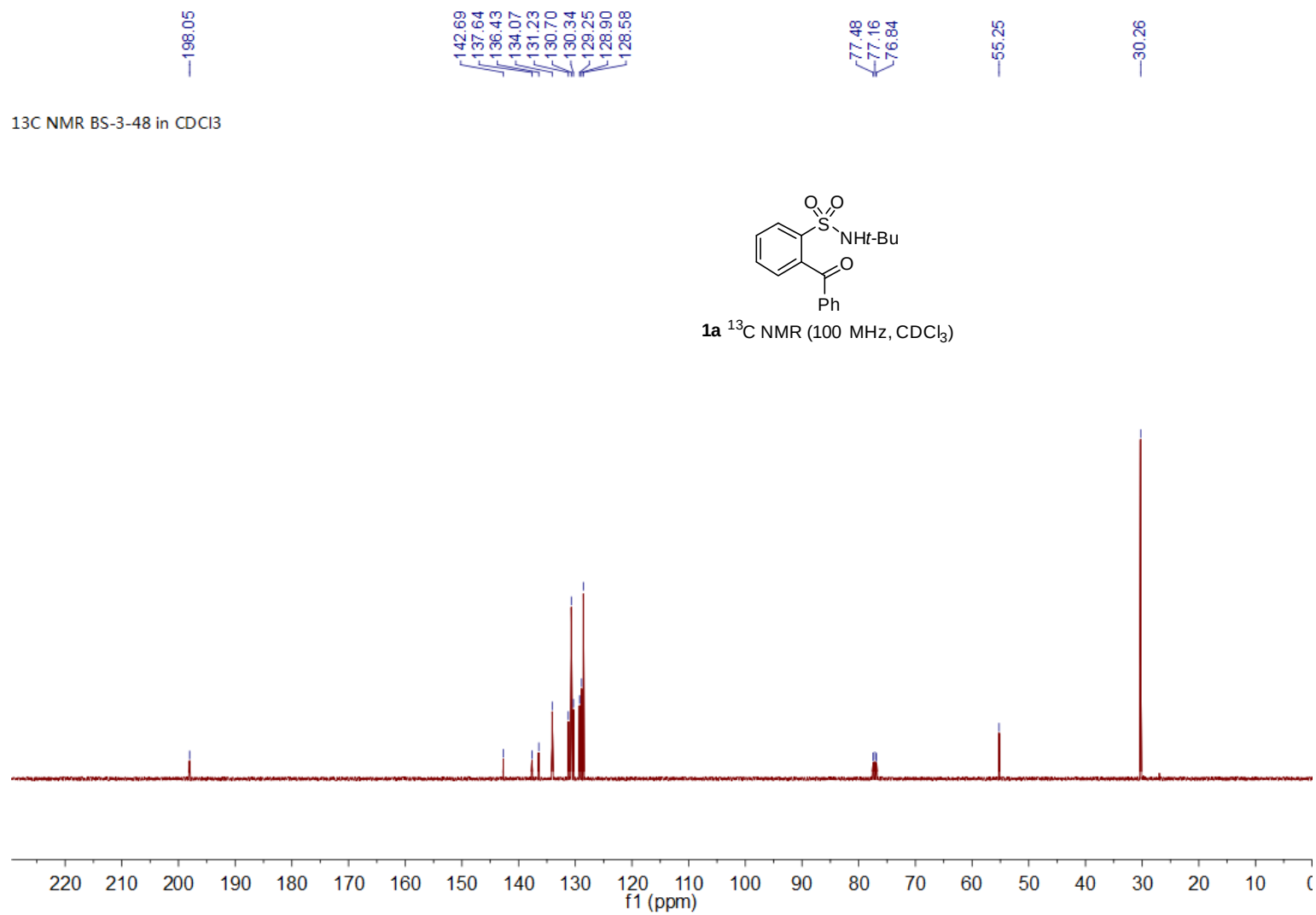
Figure S2. X-ray crystallographic analysis of (S)-(+)-**6h**

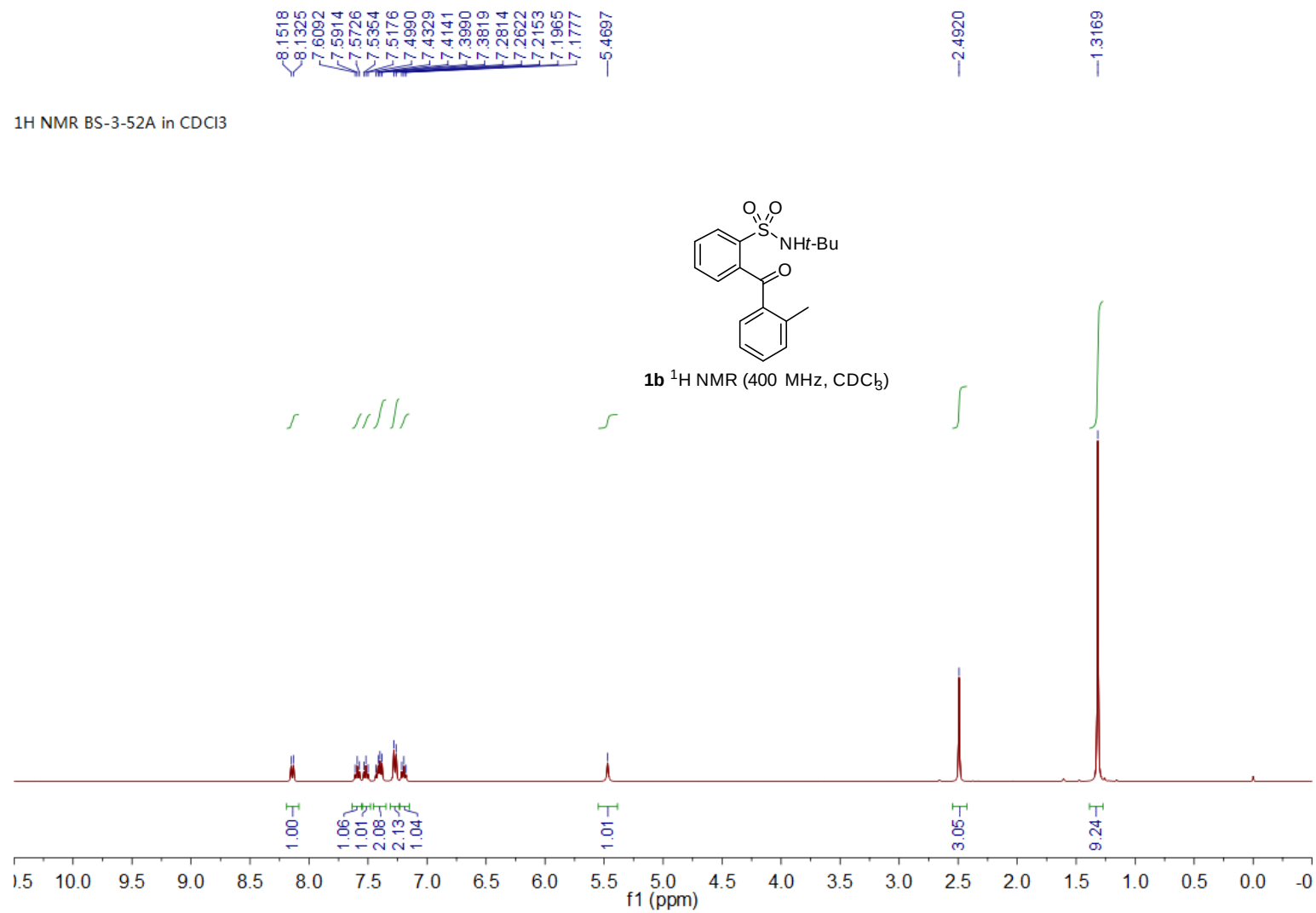
6. References

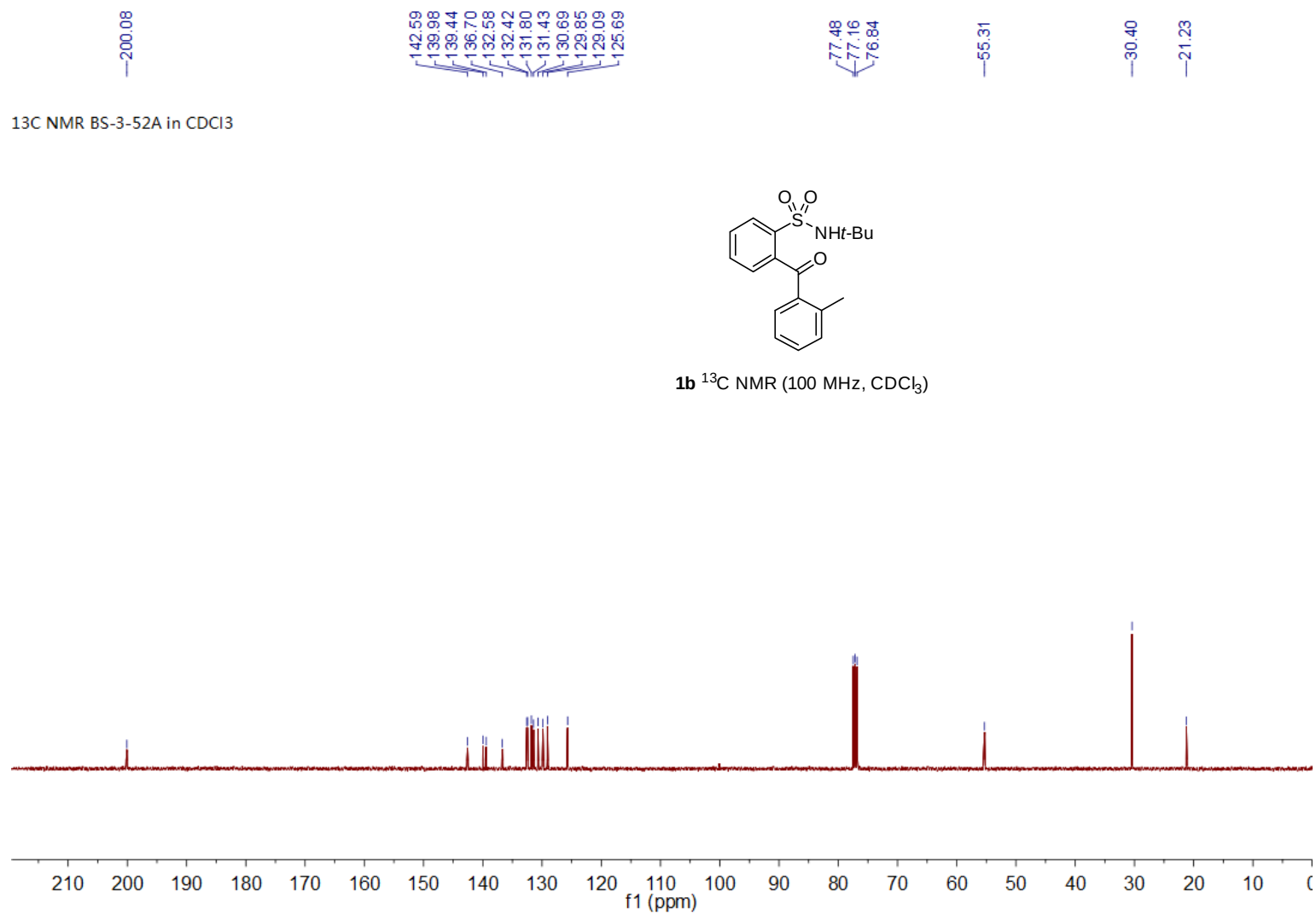
- 1) B. Song, C.-B. Yu, W.-X. Huang, M.-W. Chen and Y.-G. Zhou, *Org. Lett.*, 2015, **17**, 190.
- 2) (a) Y.-Q. Wang, S.-M. Lu and Y.-G. Zhou, *J. Org. Chem.*, 2007, **72**, 3729; (b) C.-B. Yu, D.-W. Wang and Y.-G. Zhou, *J. Org. Chem.*, 2009, **74**, 5633.
- 3) C.-B. Yu, K. Gao, D.-S. Wang, L. Shi and Y.-G. Zhou, *Chem. Commun.*, 2011, **47**, 5052.
- 4) F. Fang, R. Wang, Z.-P. Liu and A.-G. Ji, *Heterocycles*, 2007, **71**, 2377.

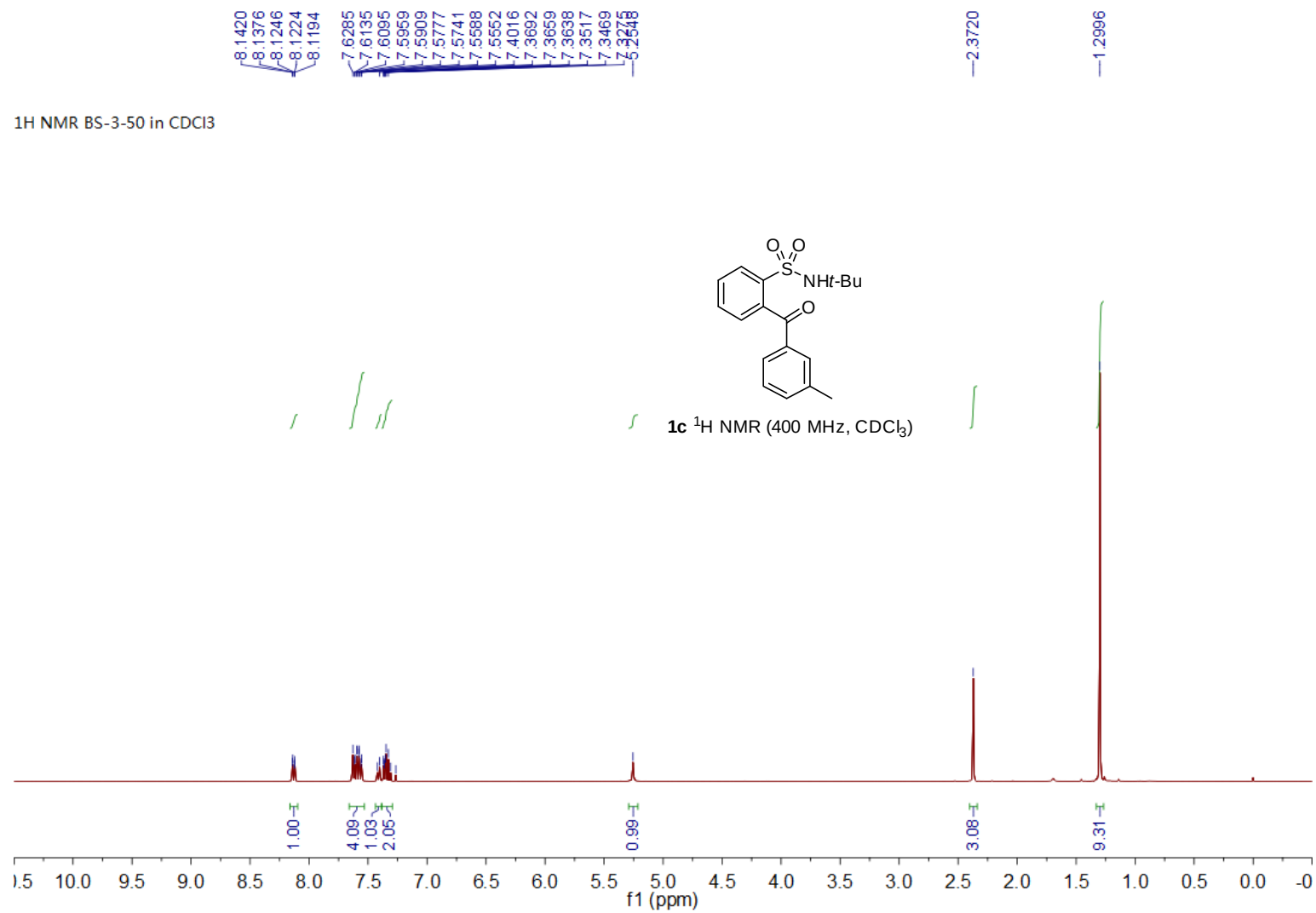
7. Copy of NMR and HPLC for the Compounds

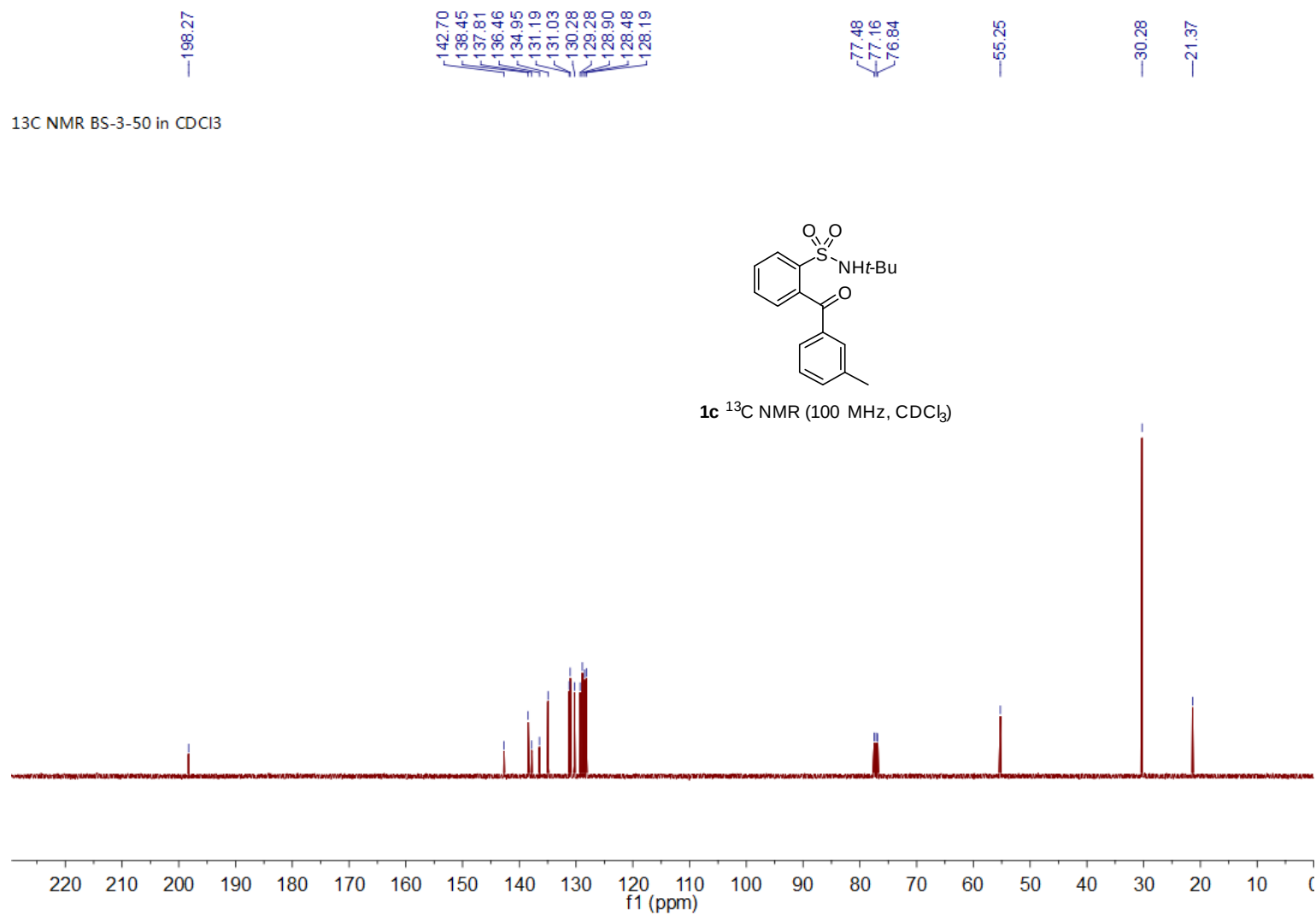




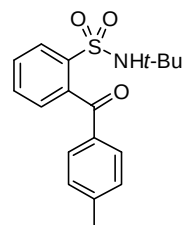
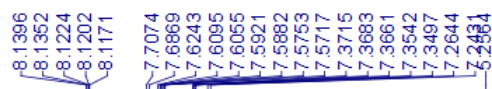




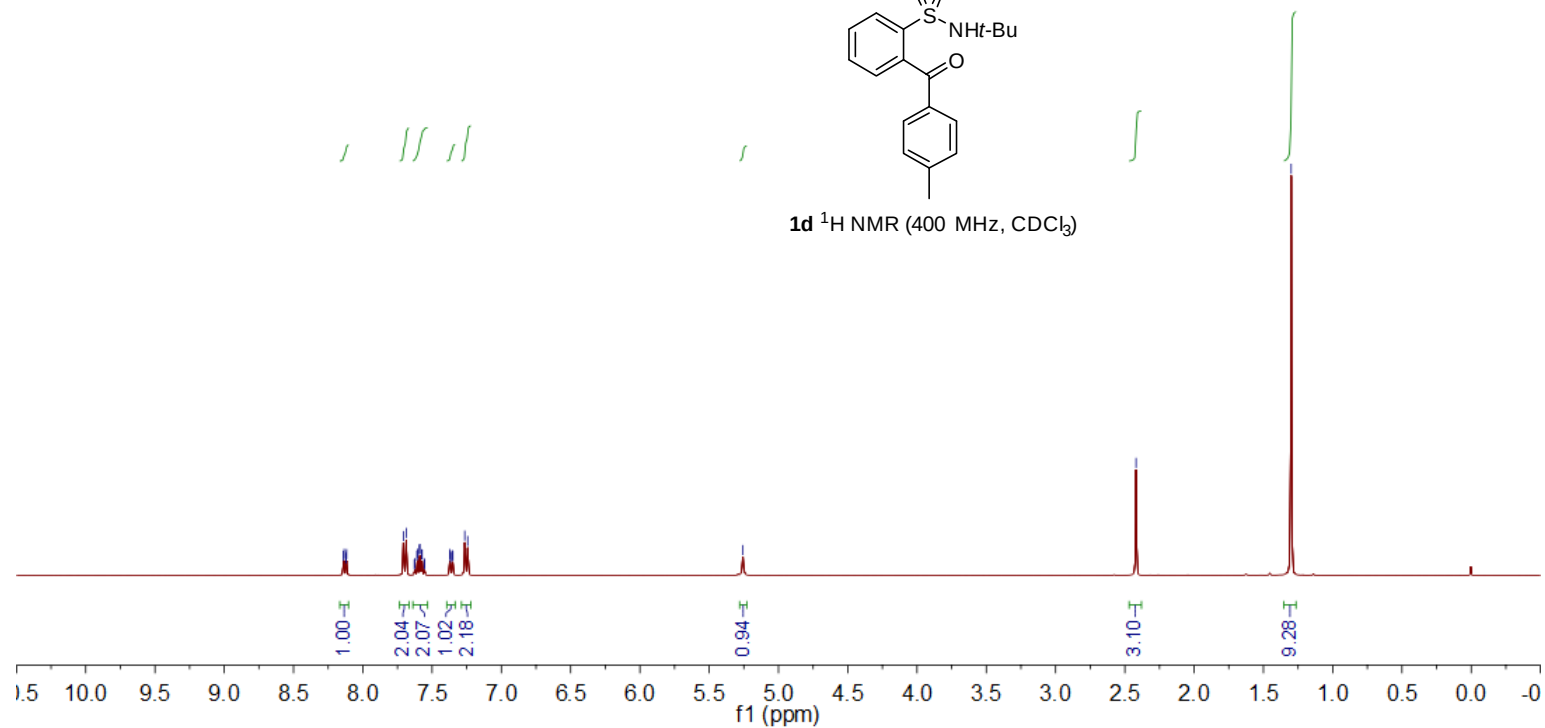


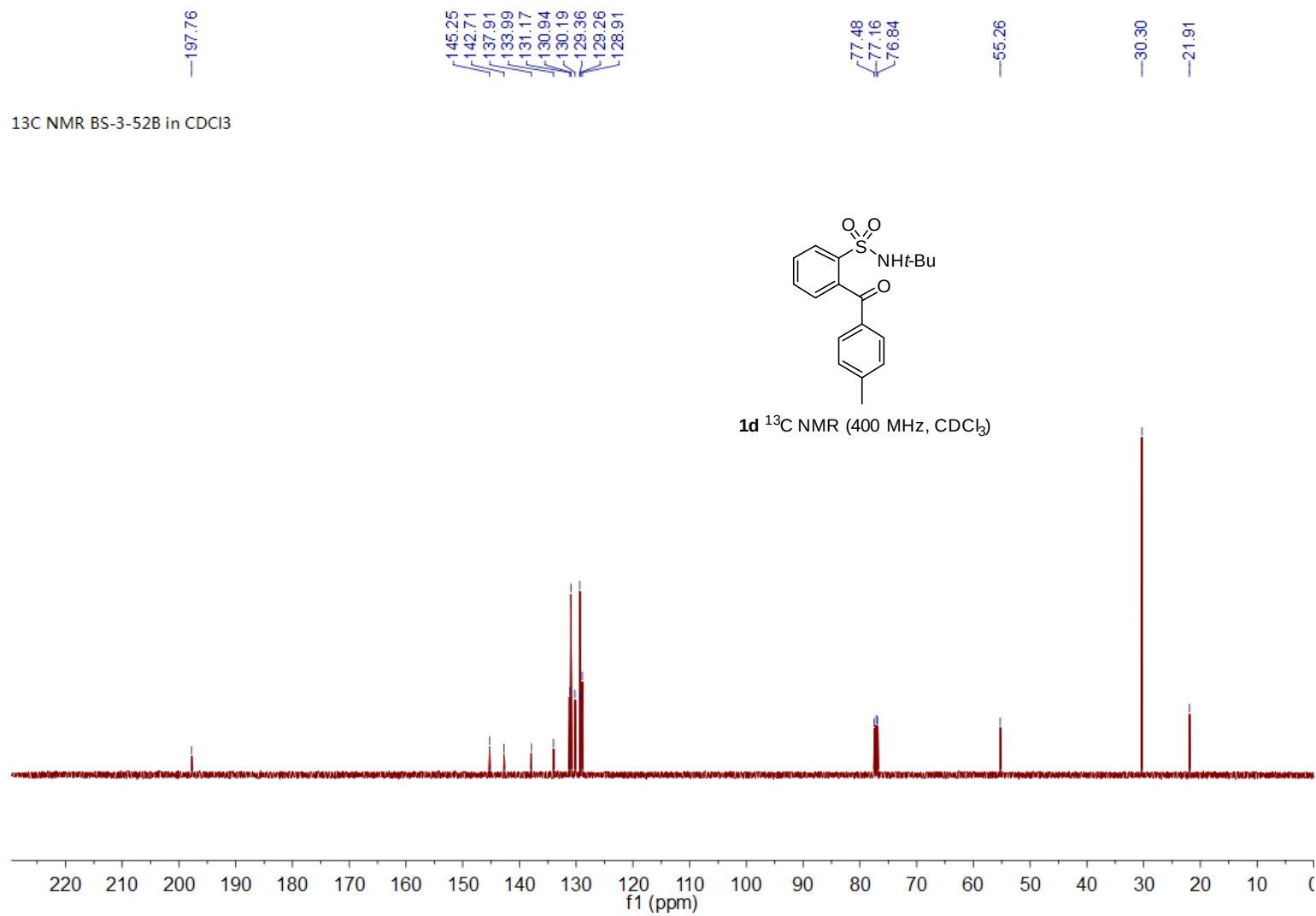


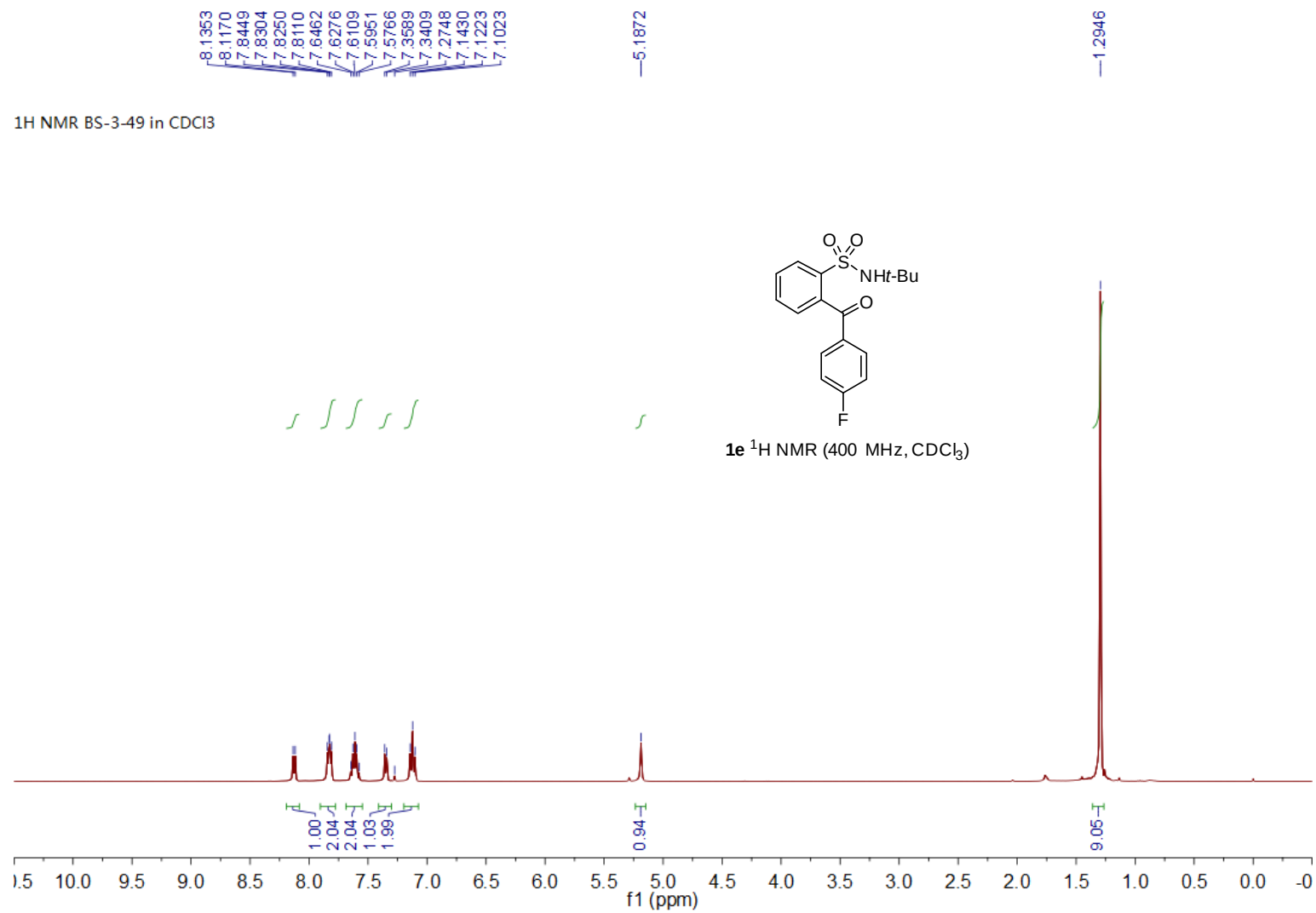
¹H NMR BS-3-52B in CDCl₃



1d ¹H NMR (400 MHz, CDCl₃)

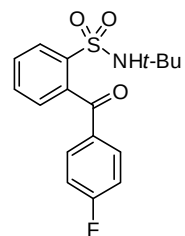




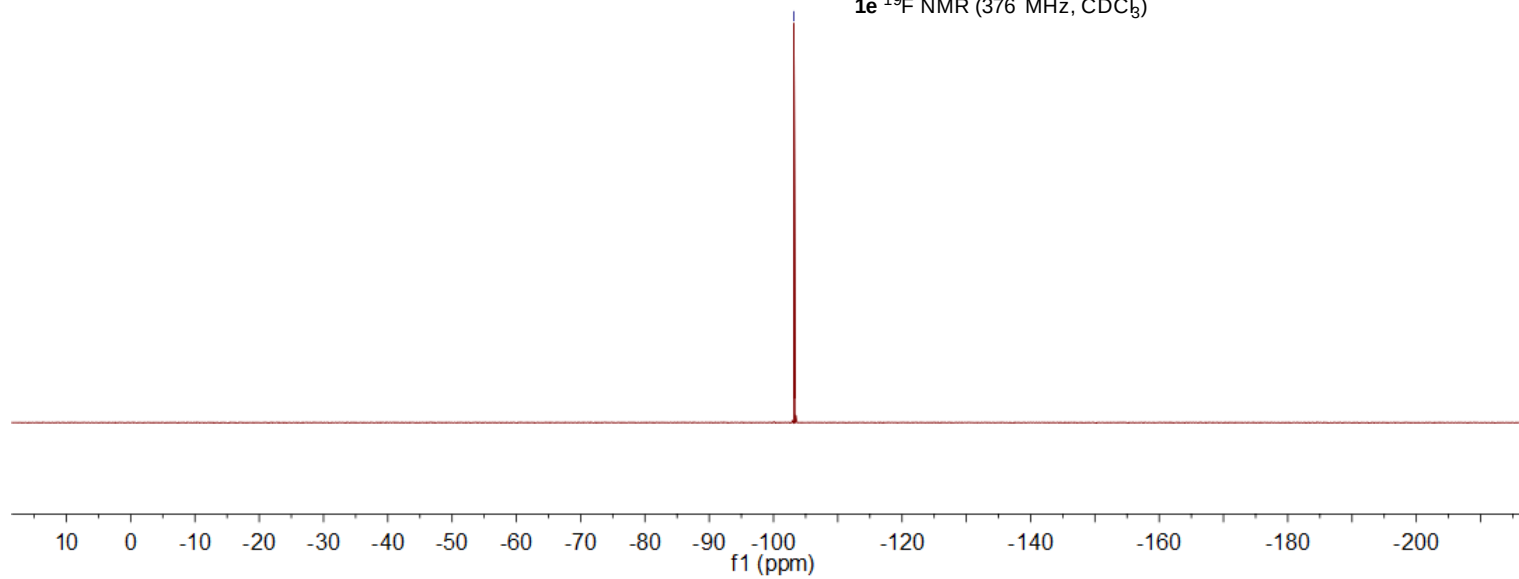


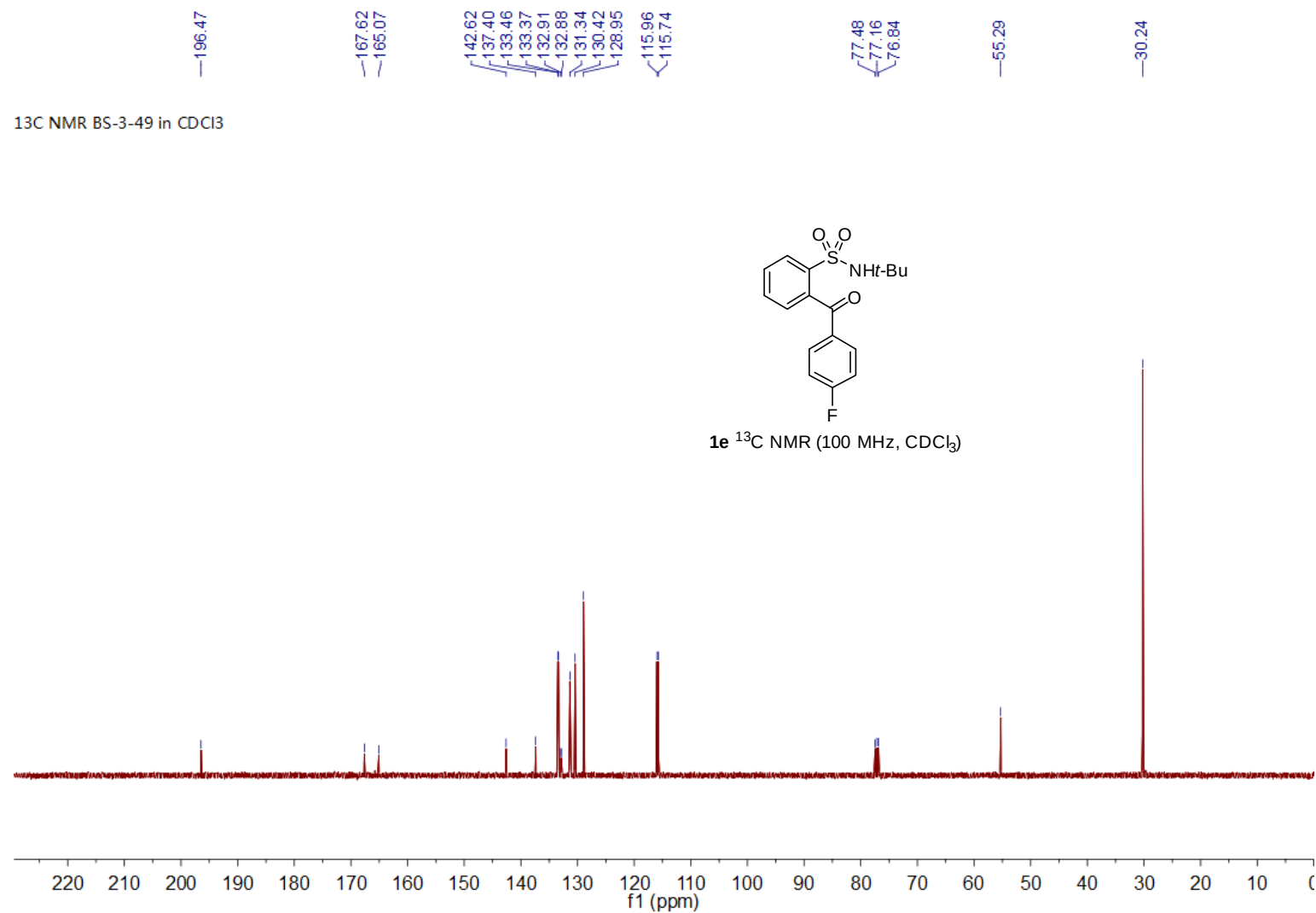
¹⁹F NMR BS-3-49 in CDCl₃

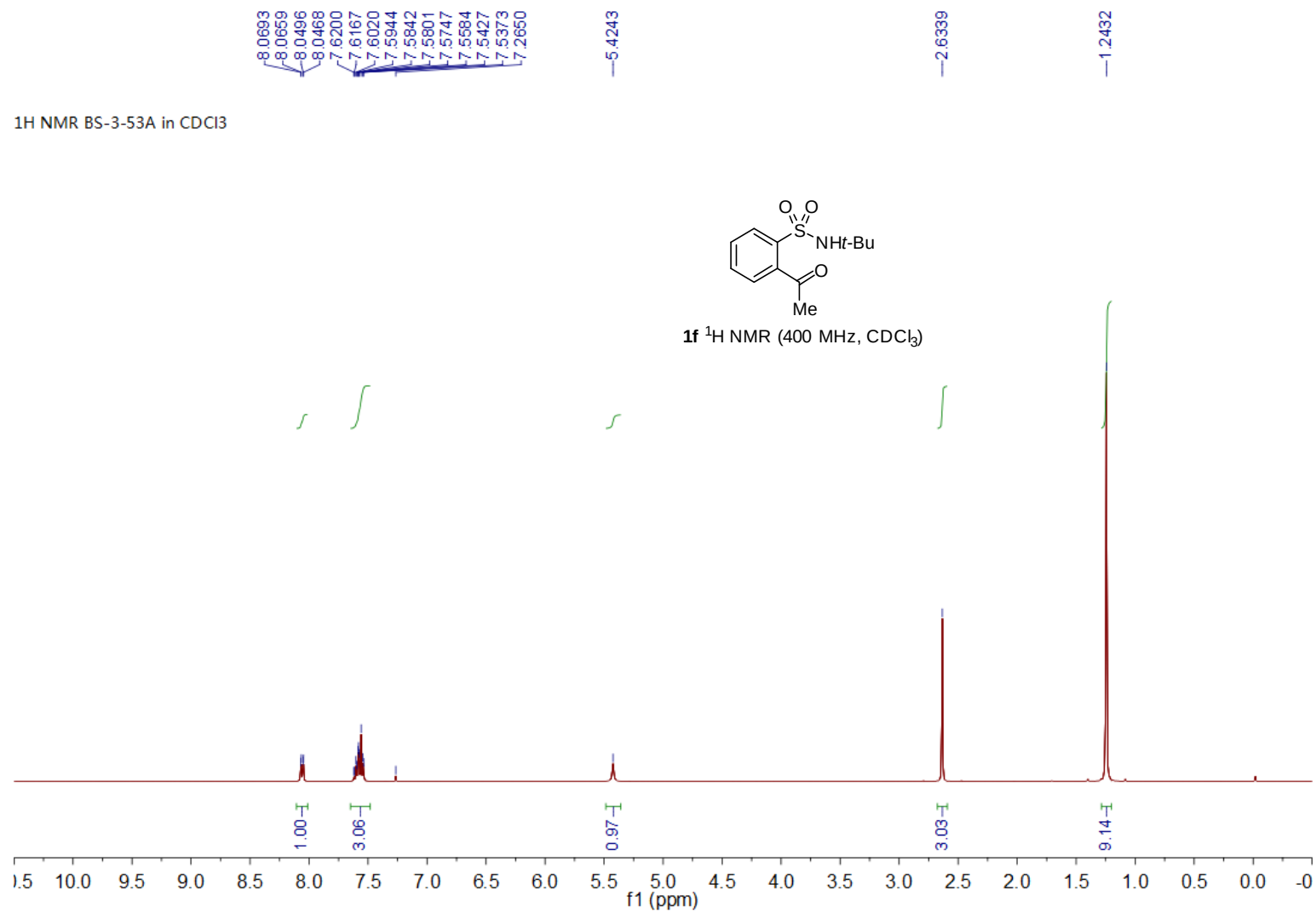
103.1704

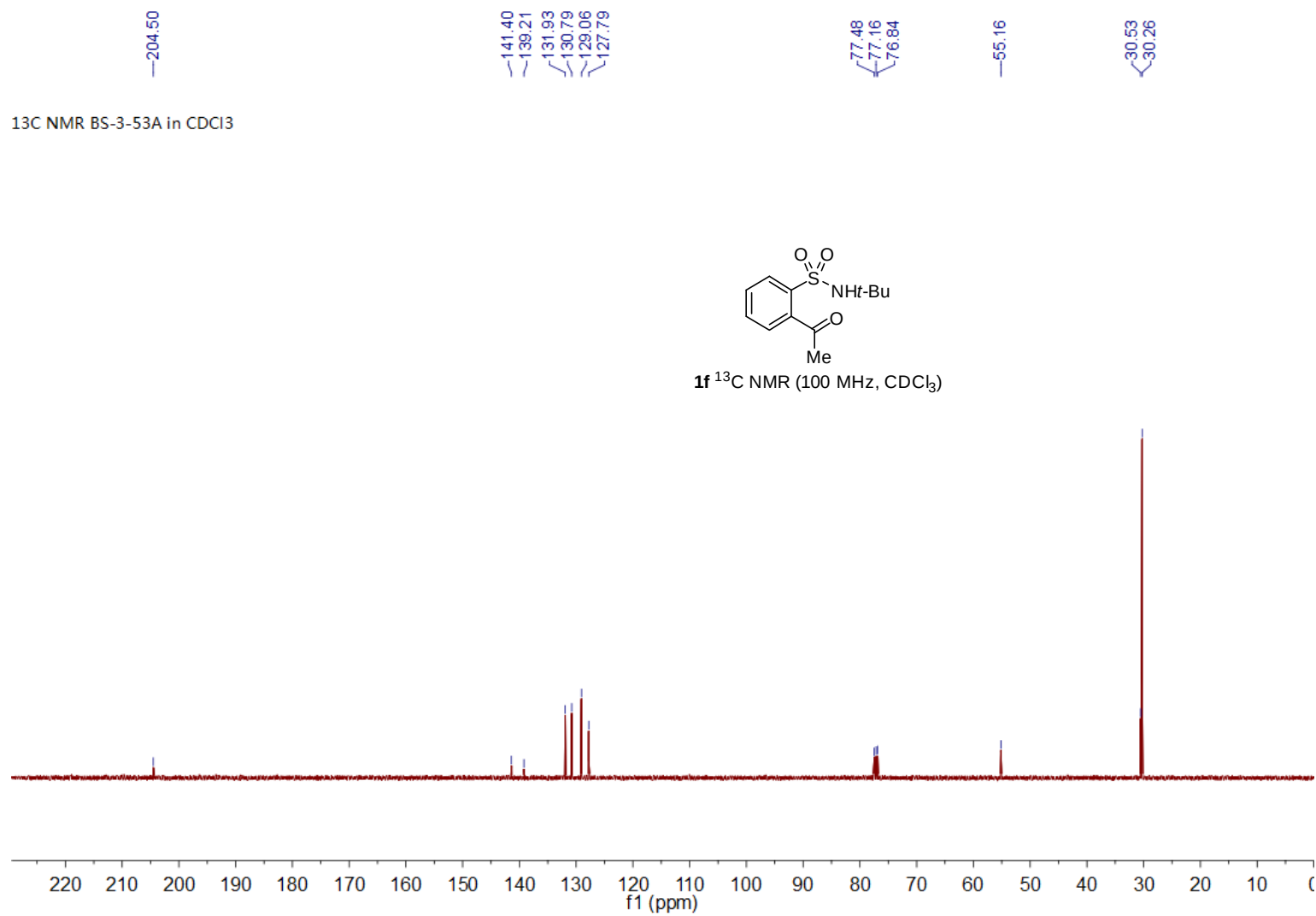


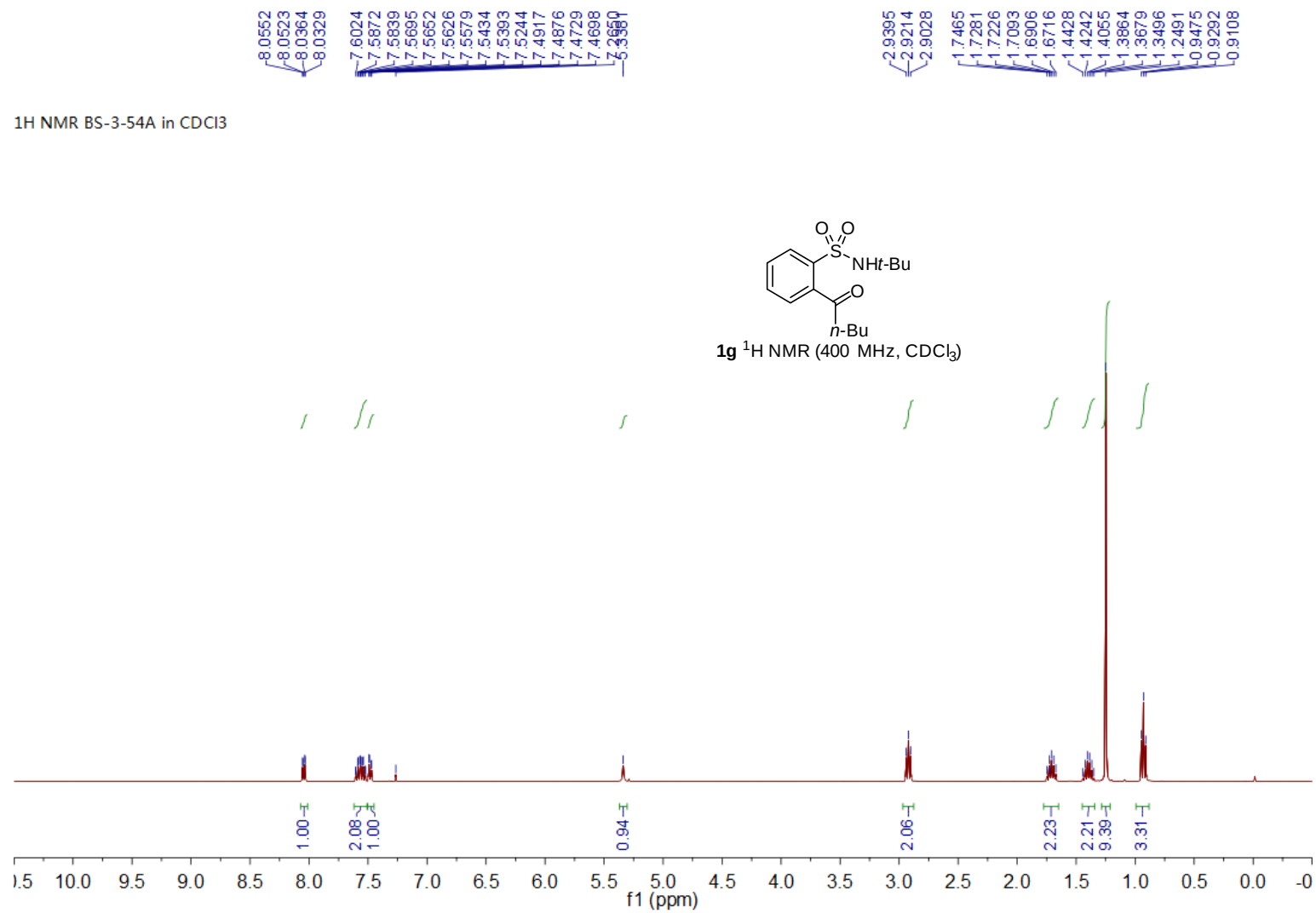
1e ¹⁹F NMR (376 MHz, CDCl₃)

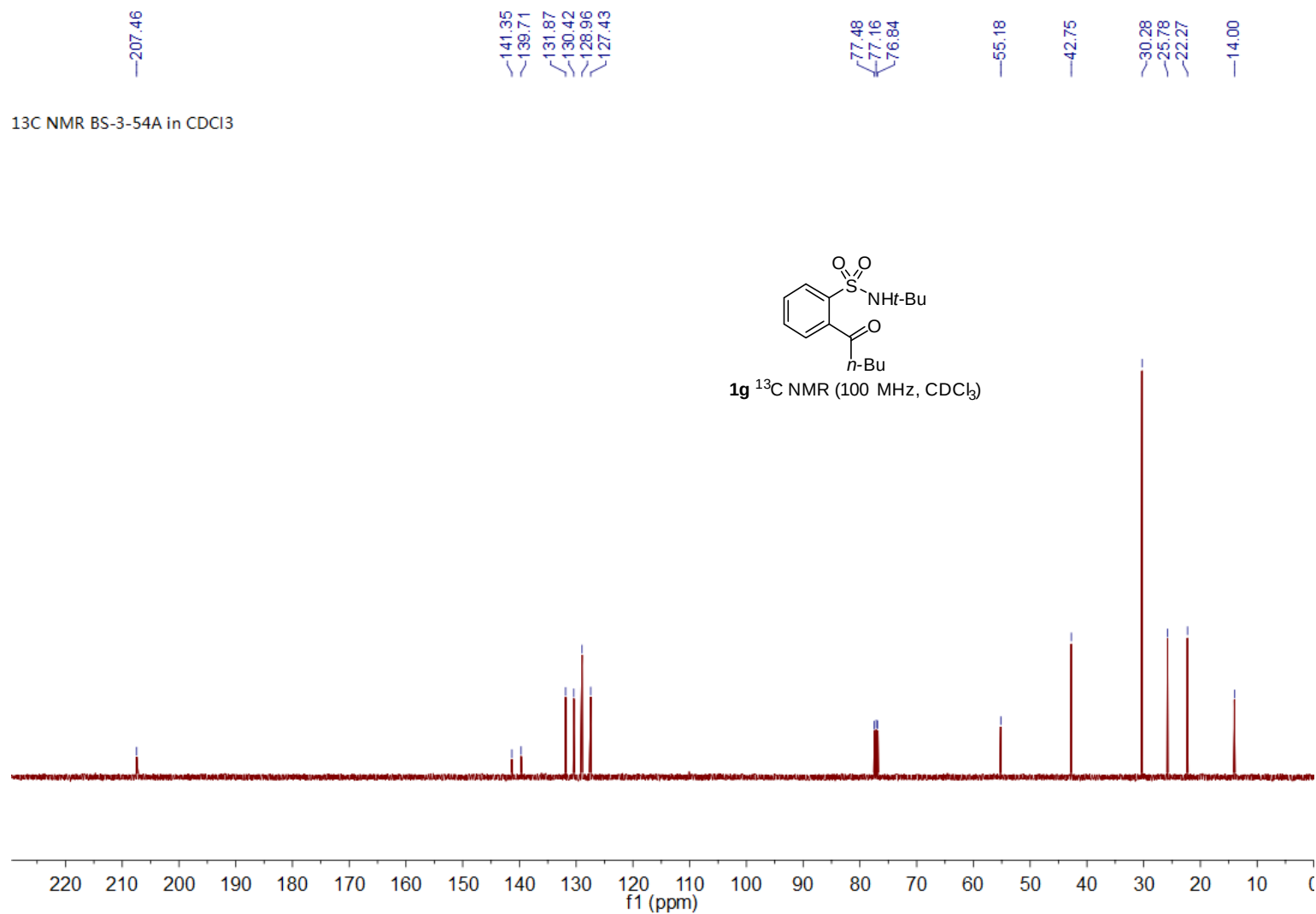


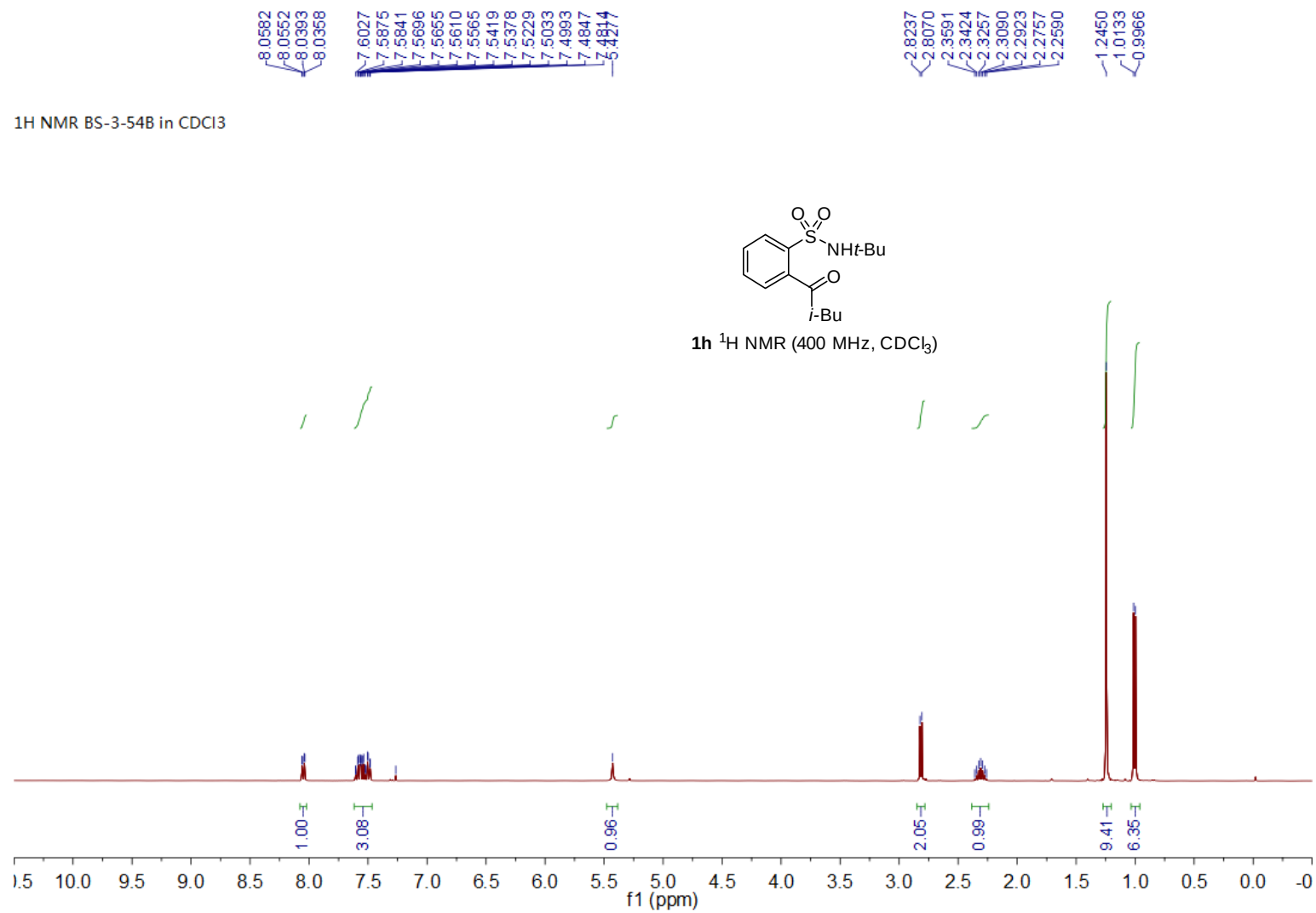


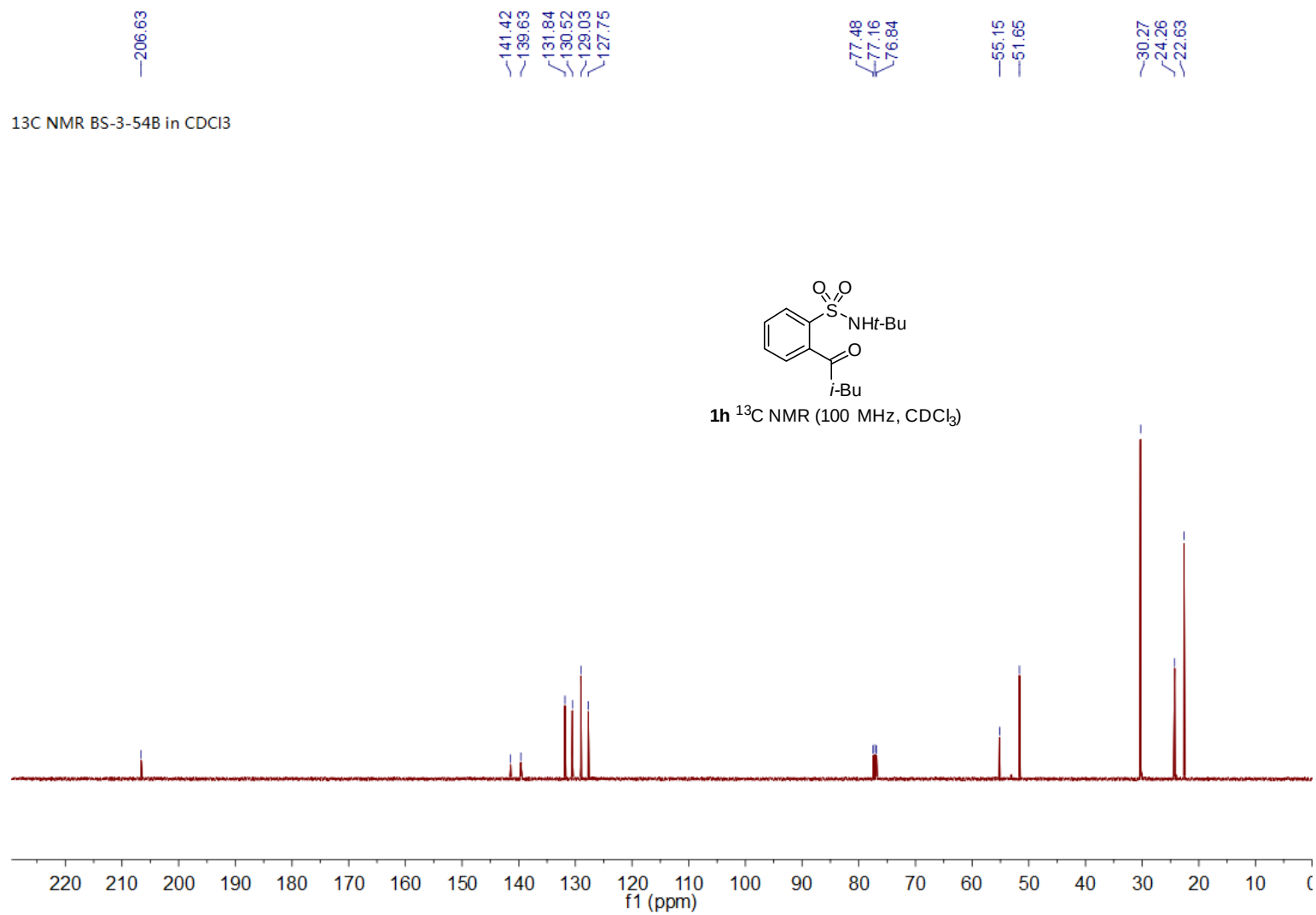


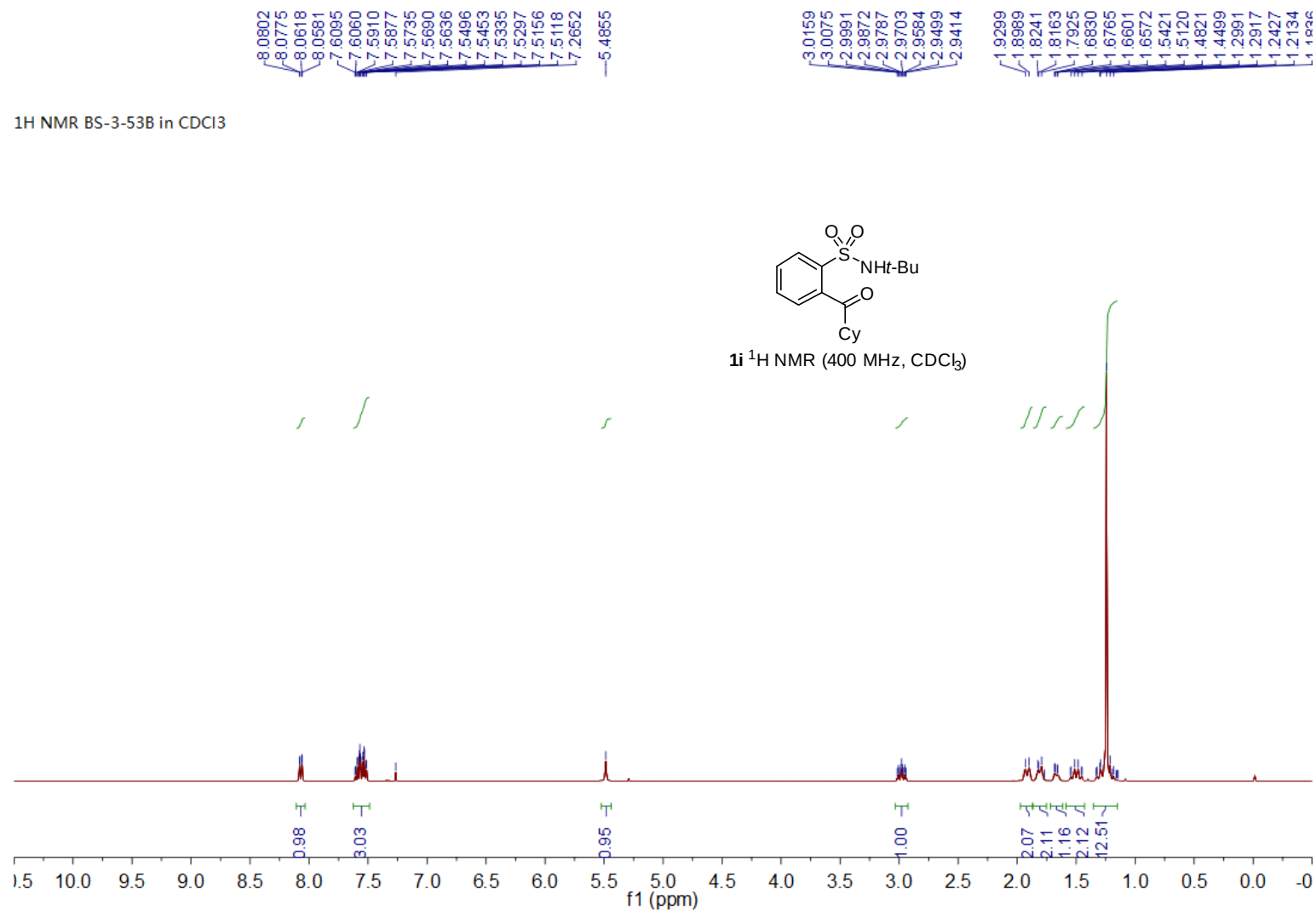


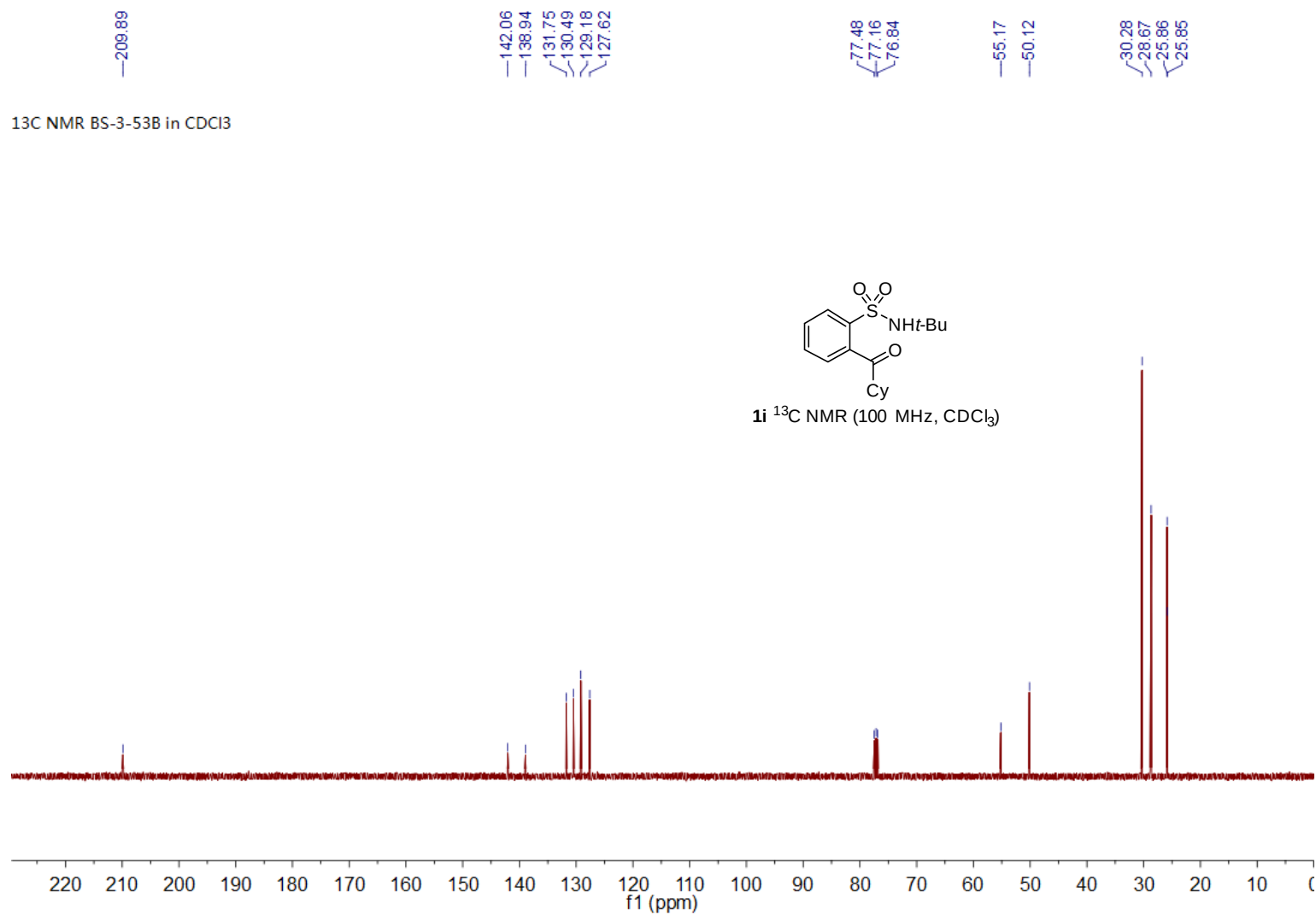


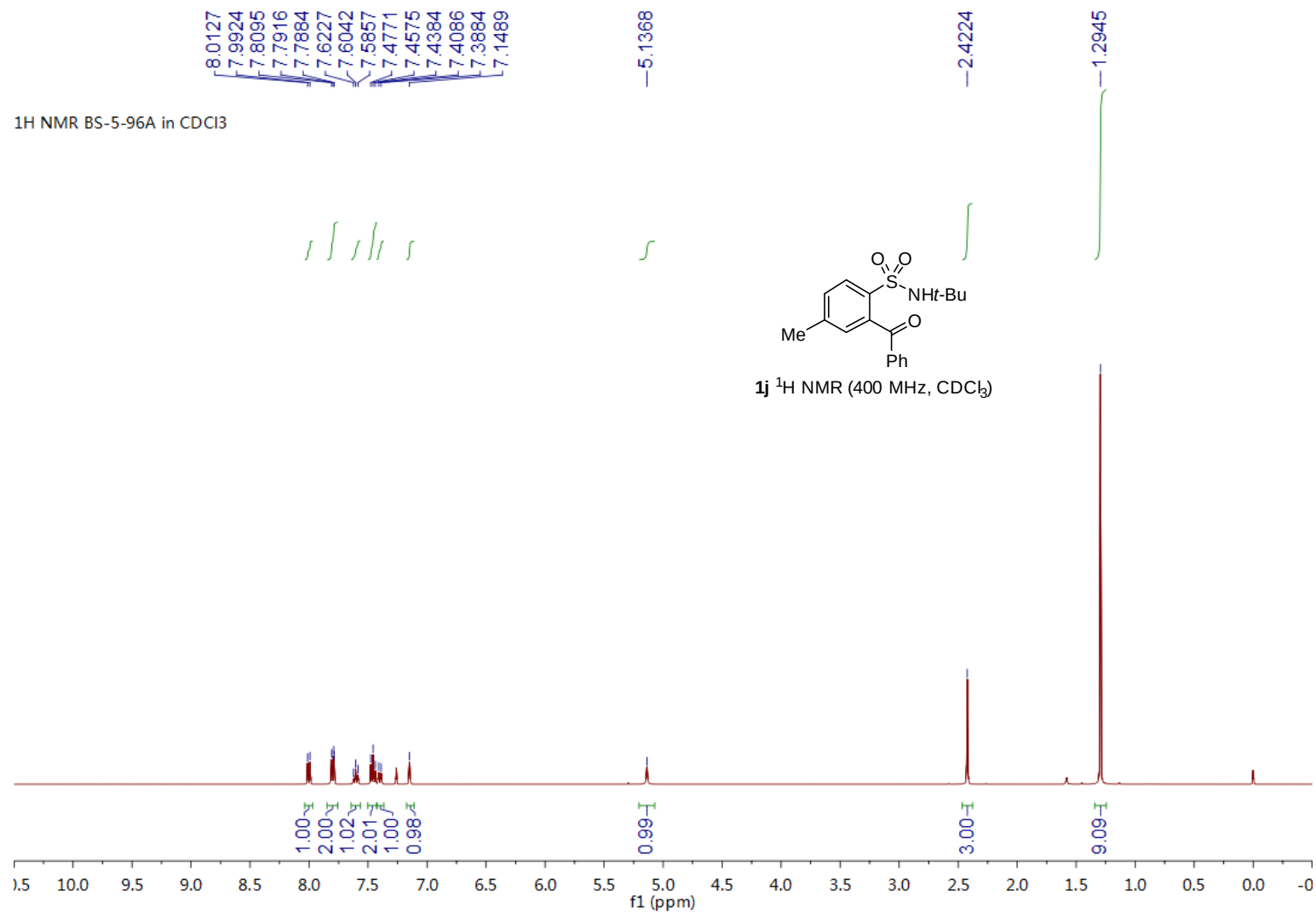


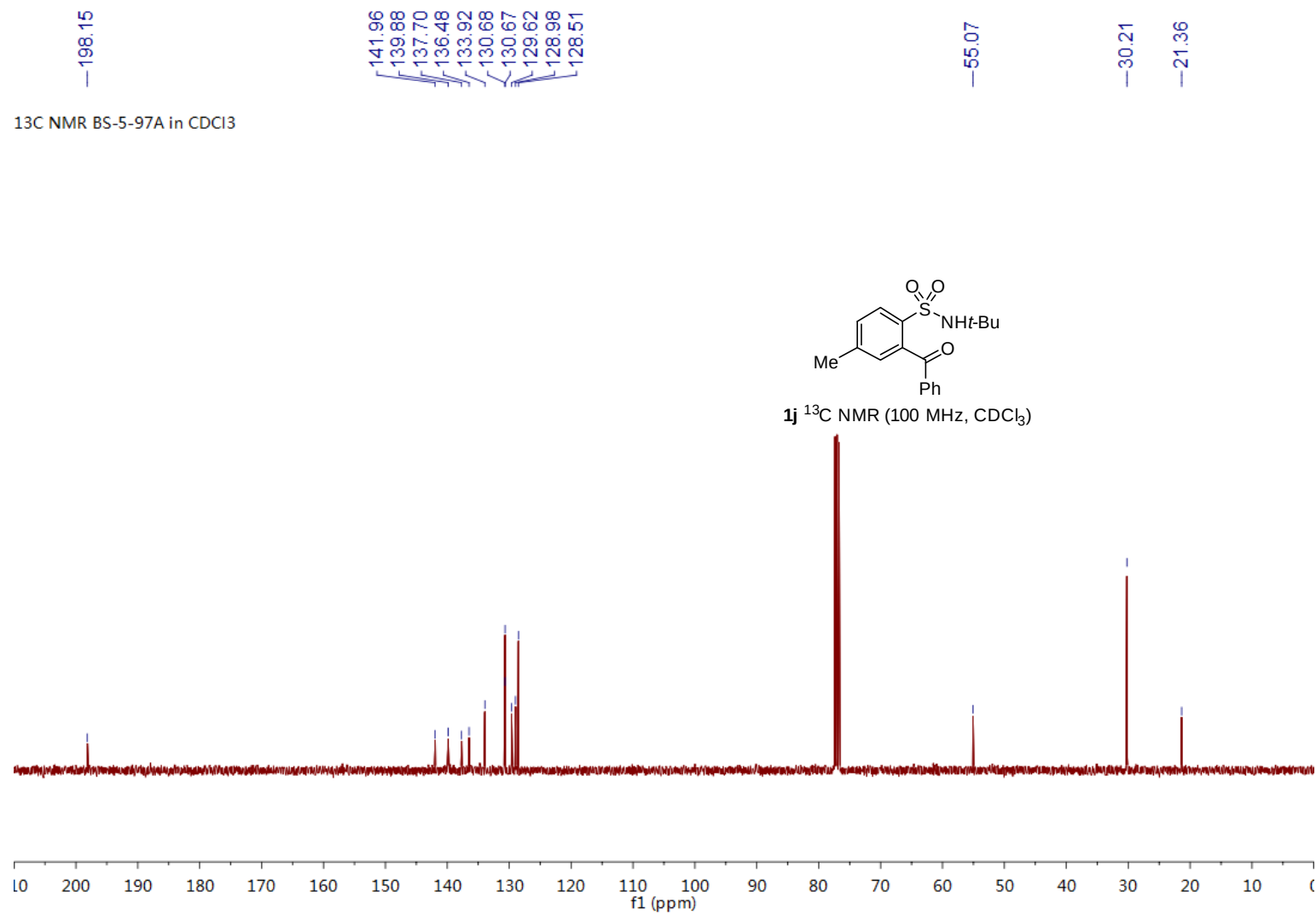


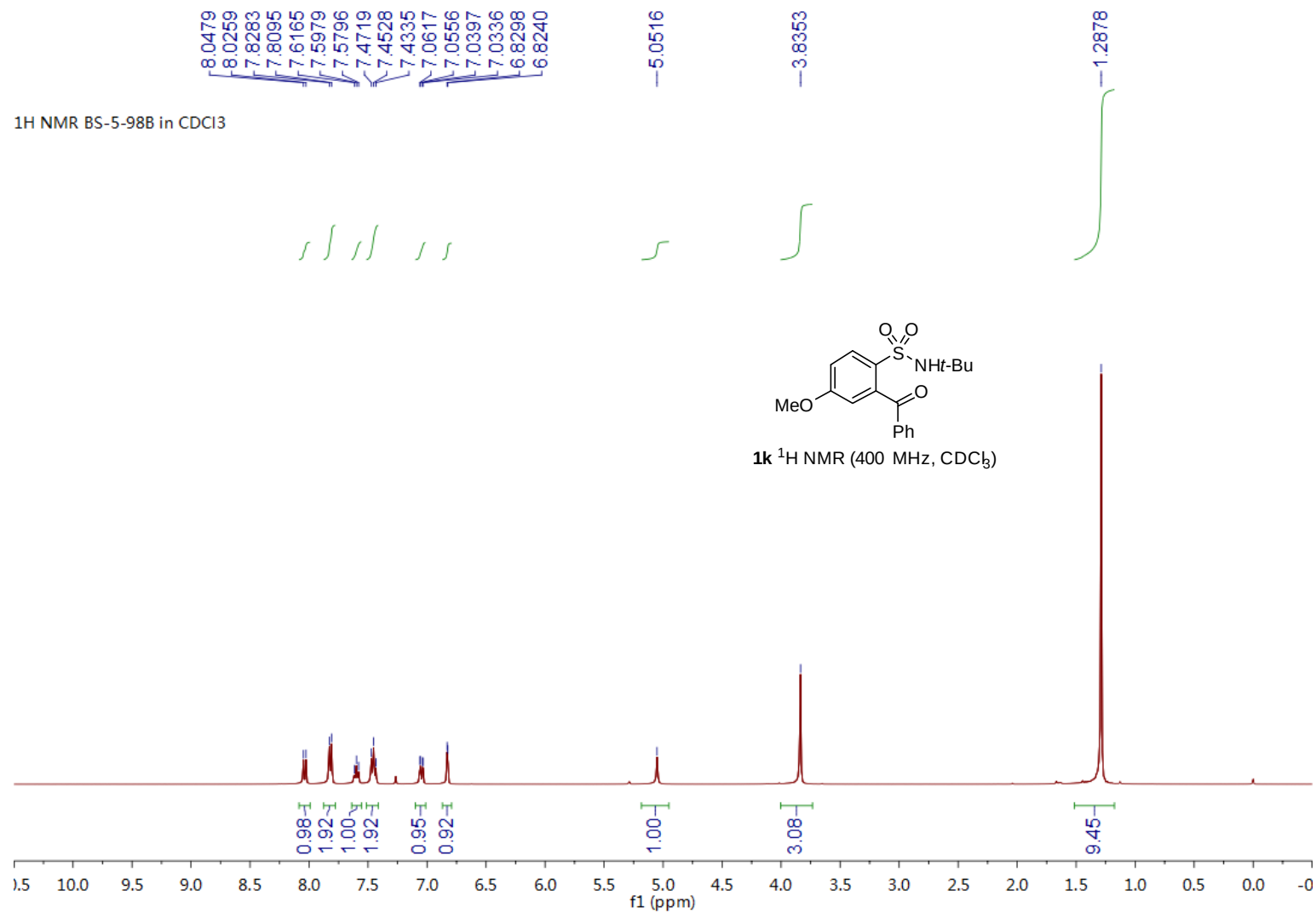


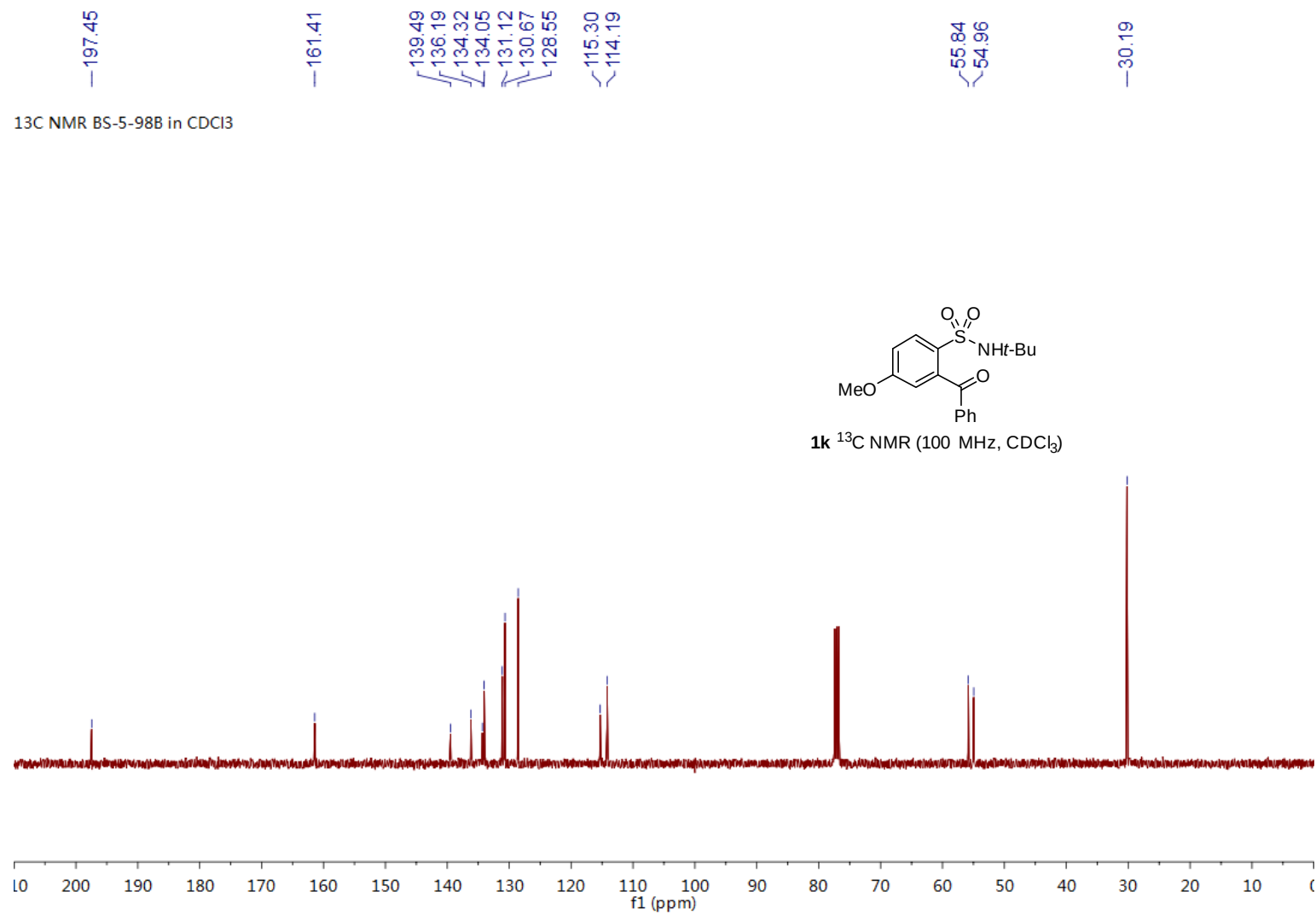


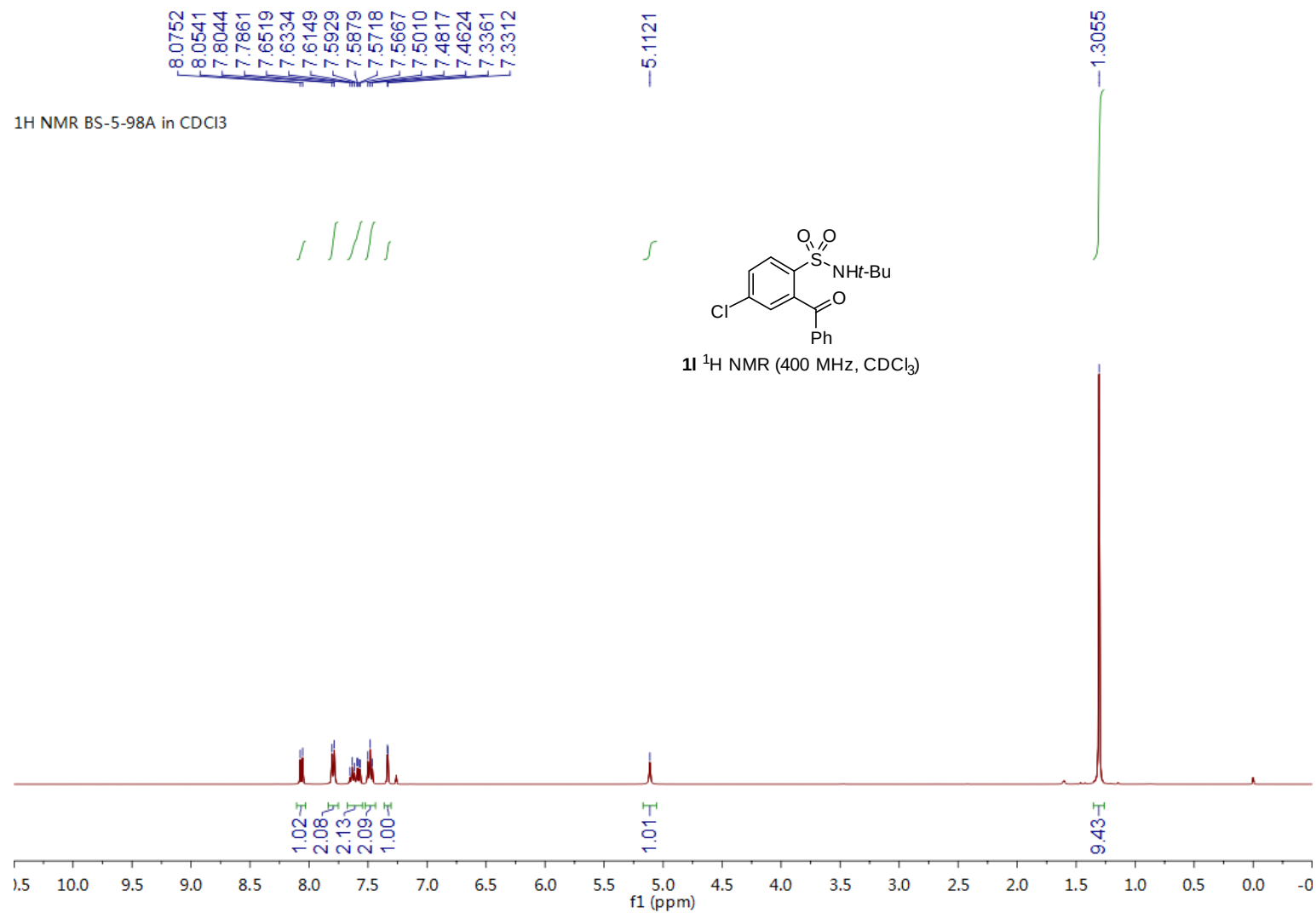


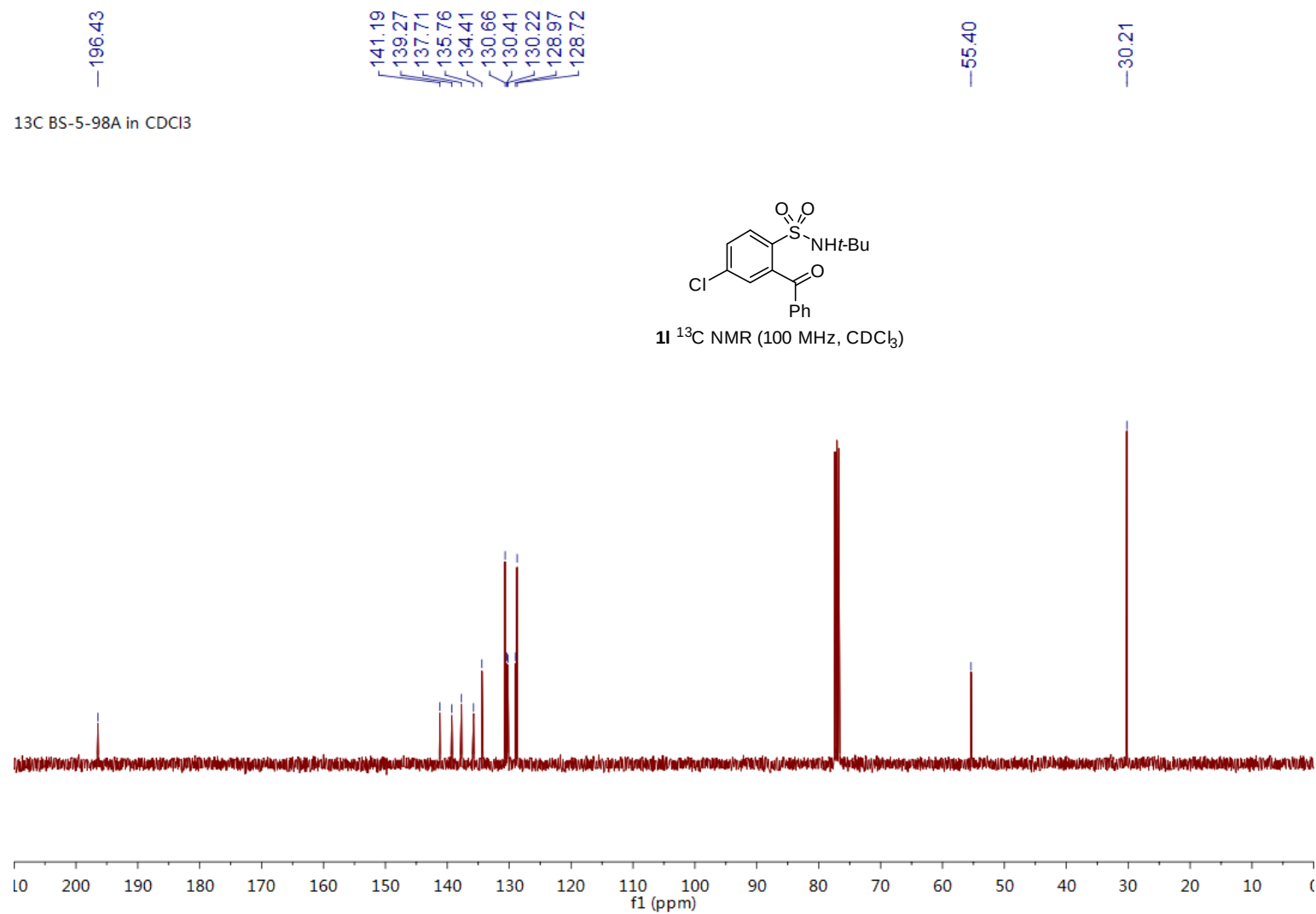


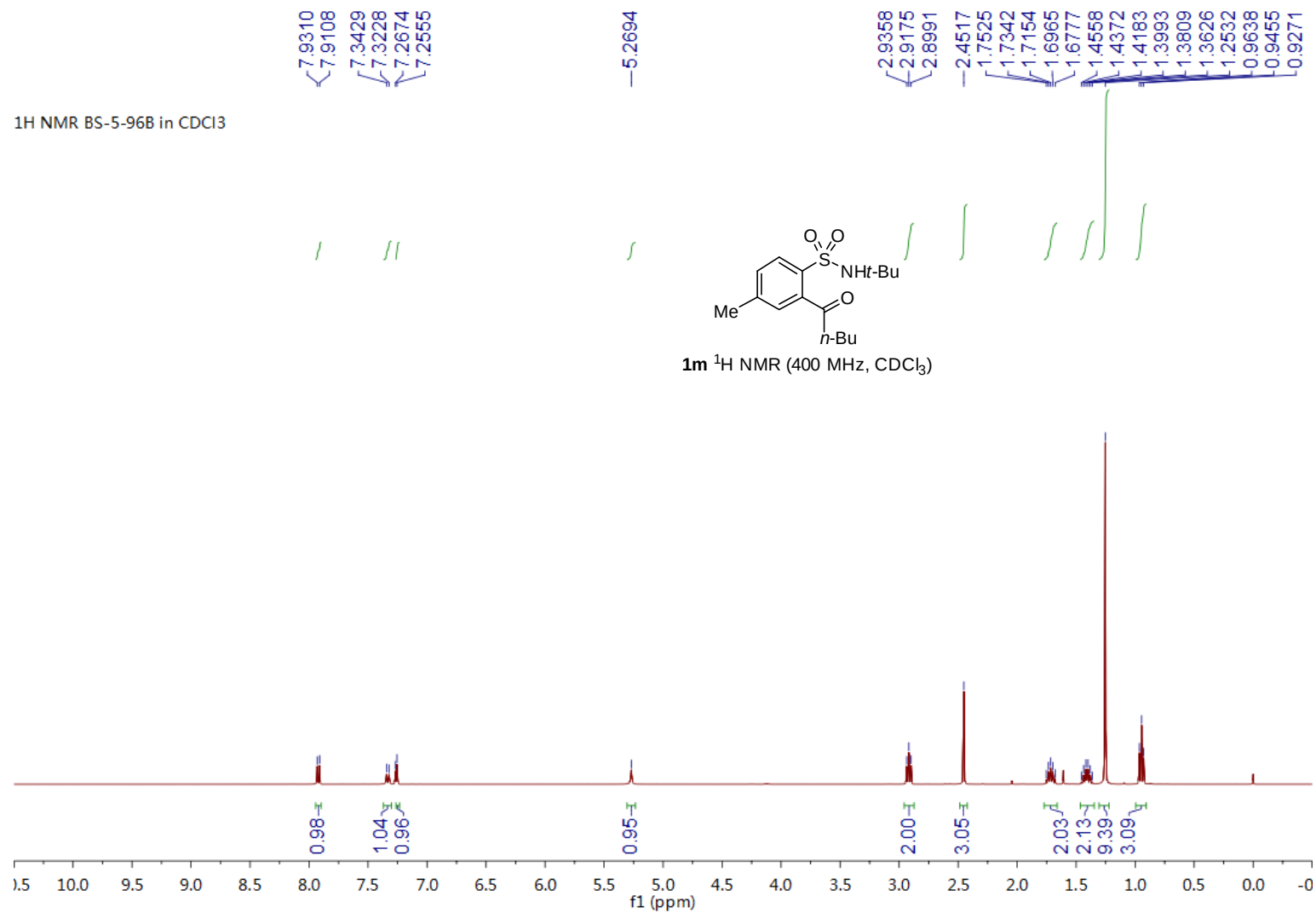


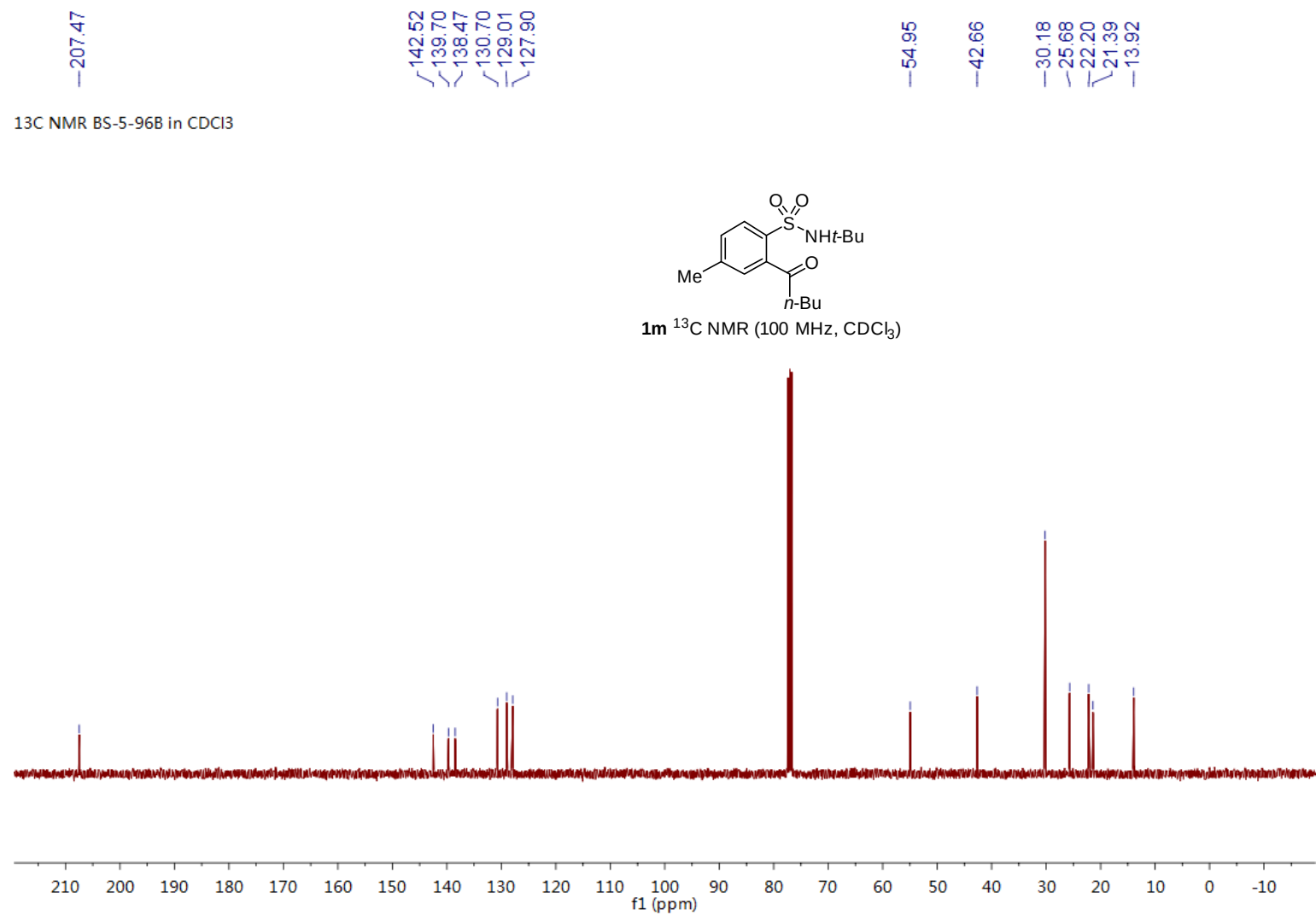


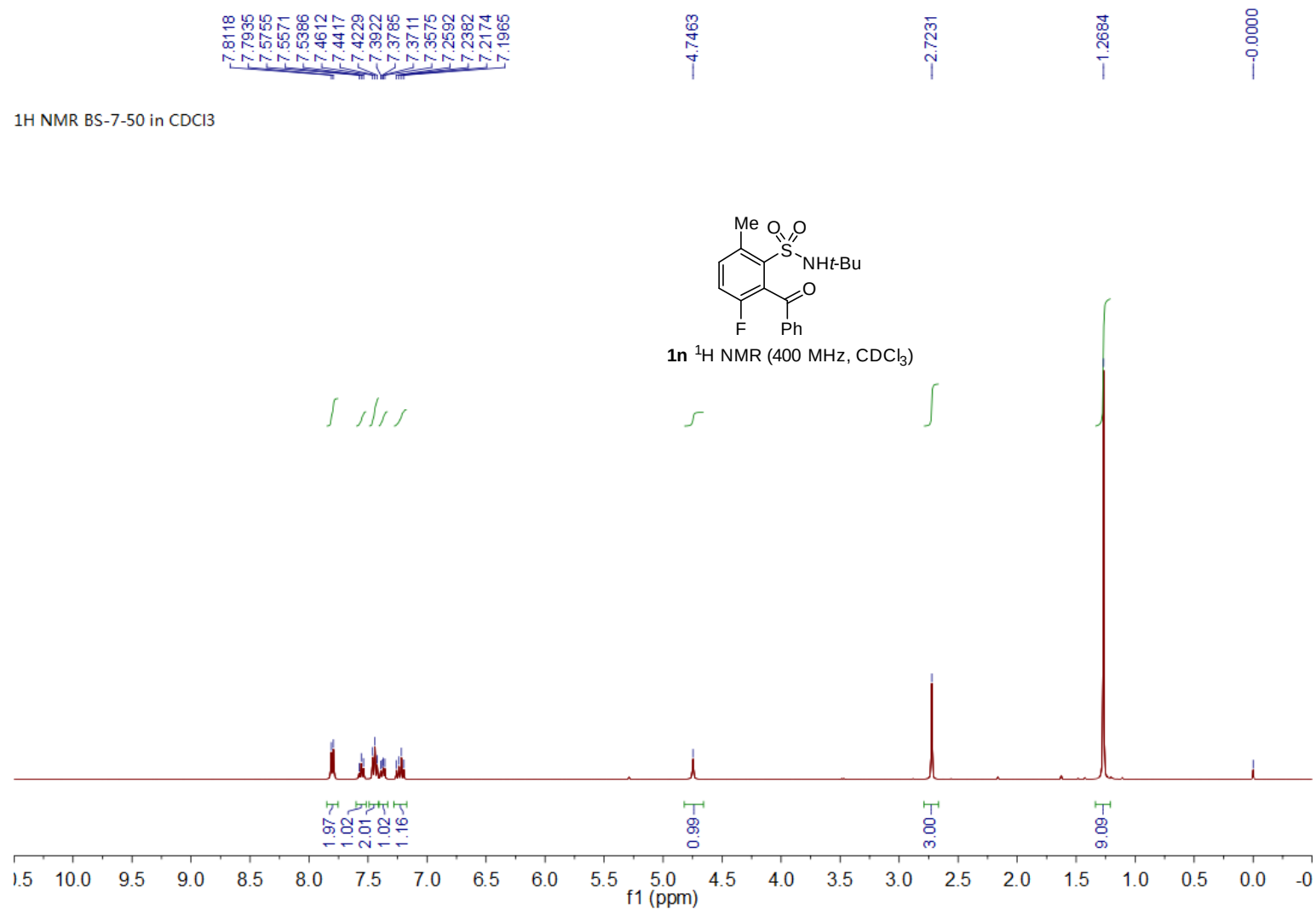


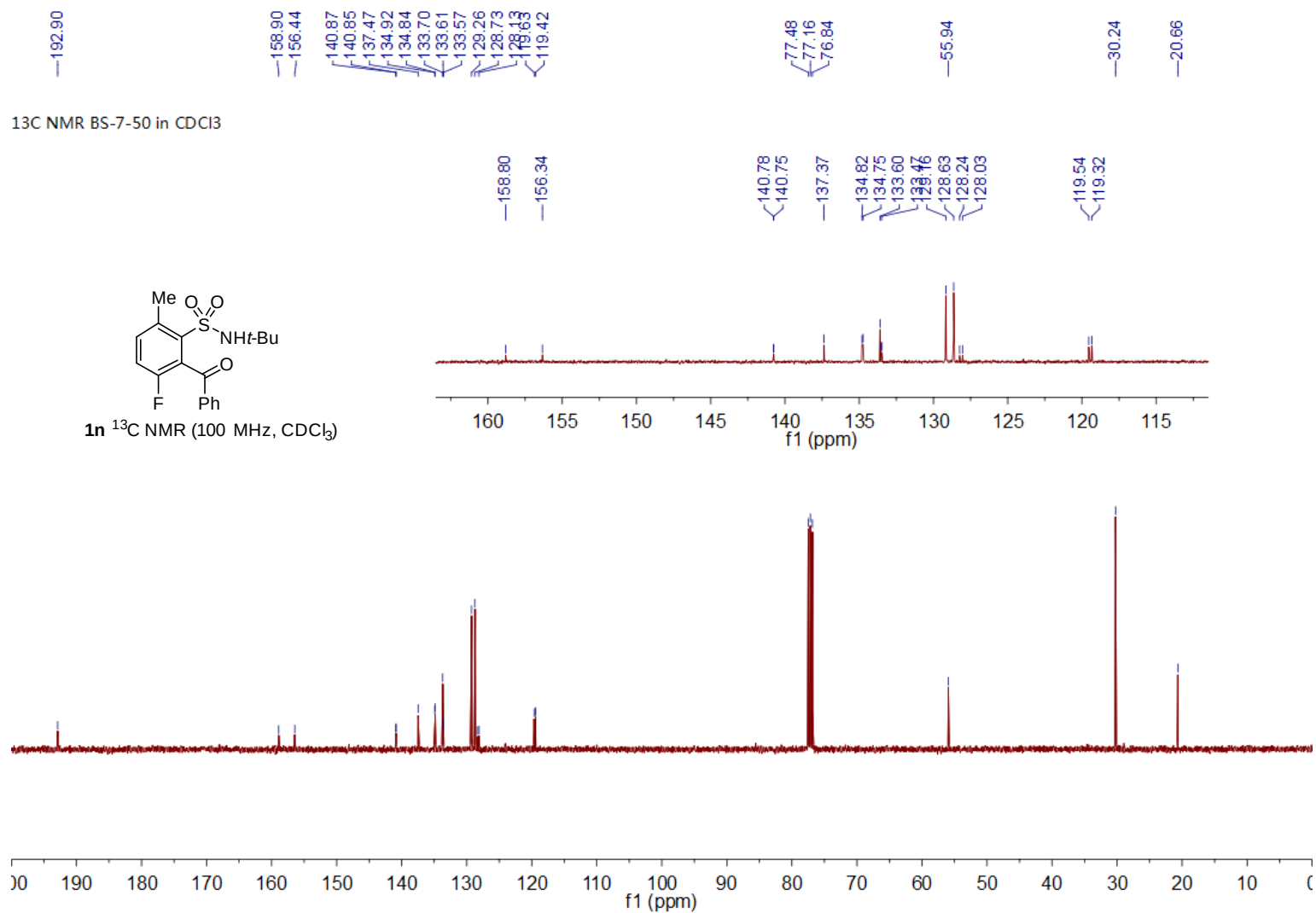




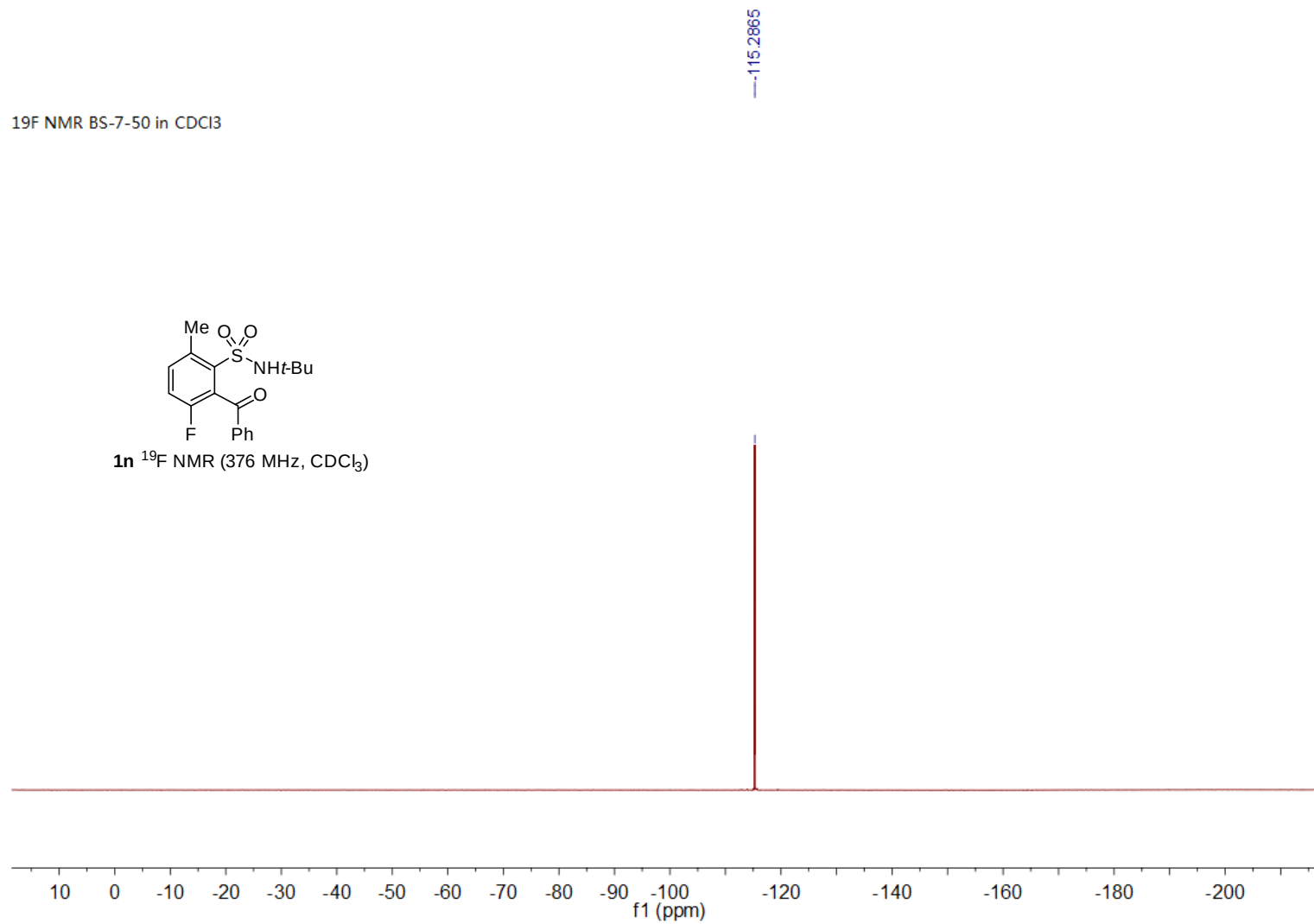
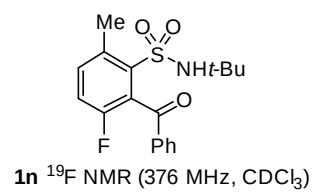




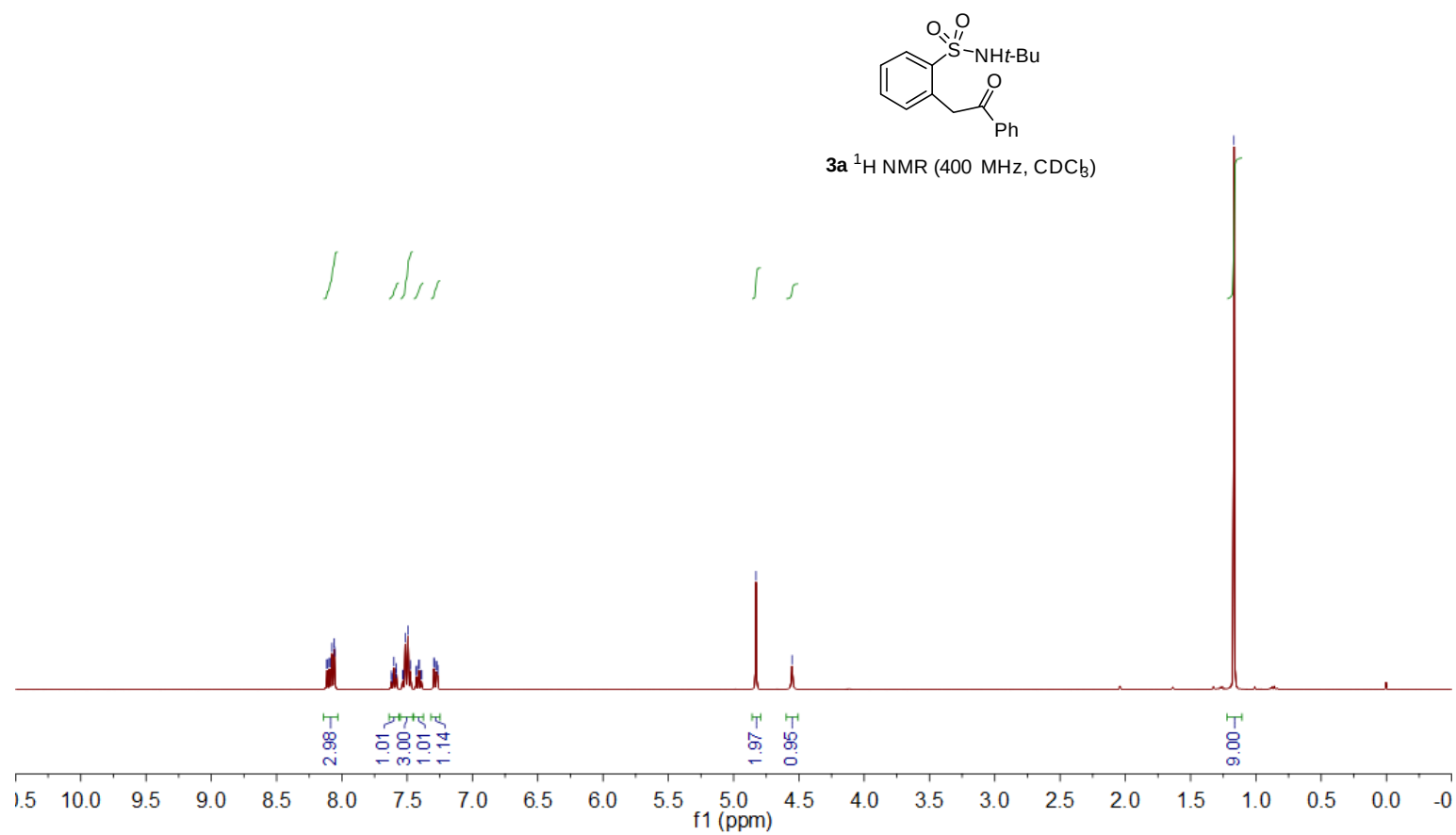


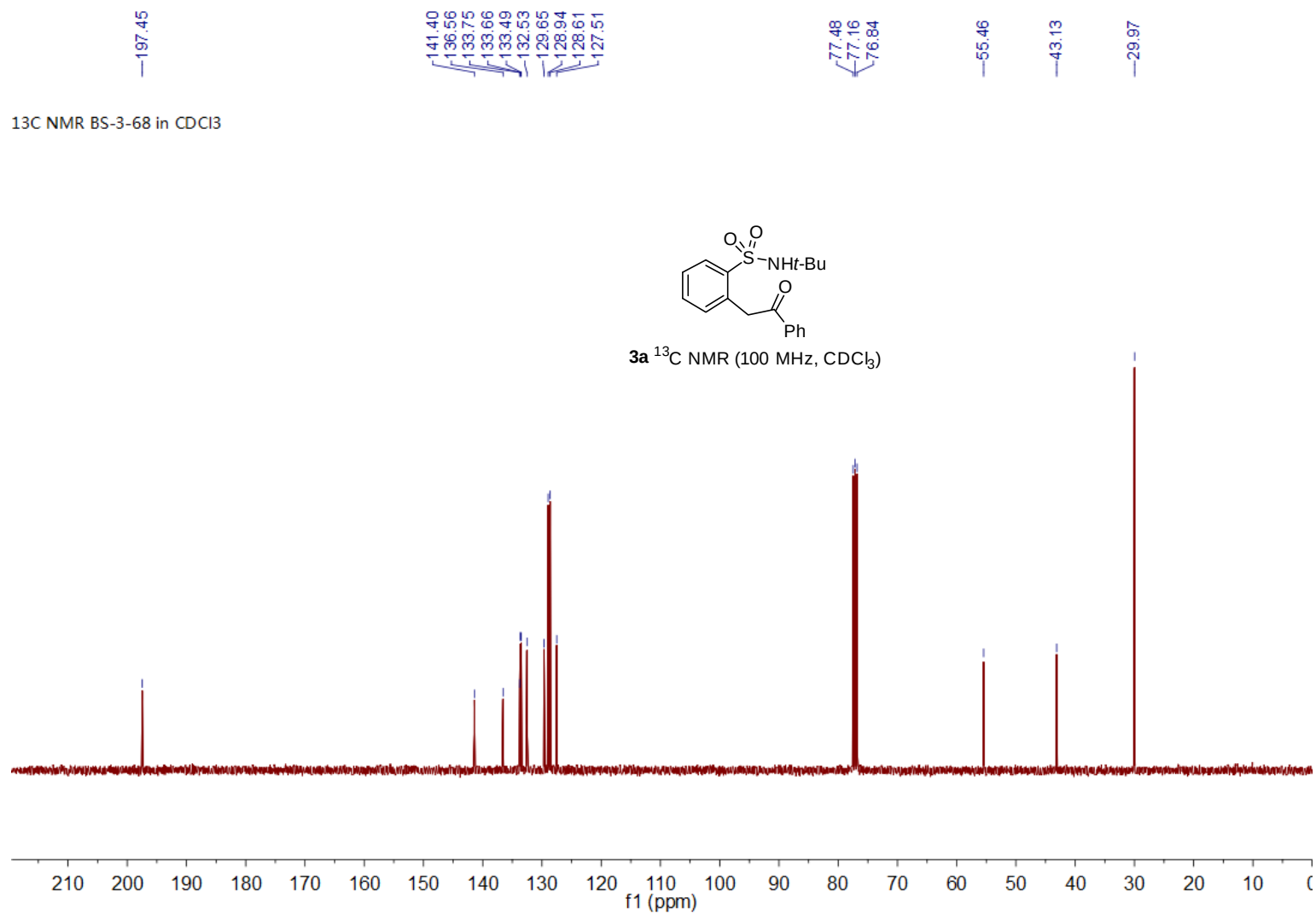


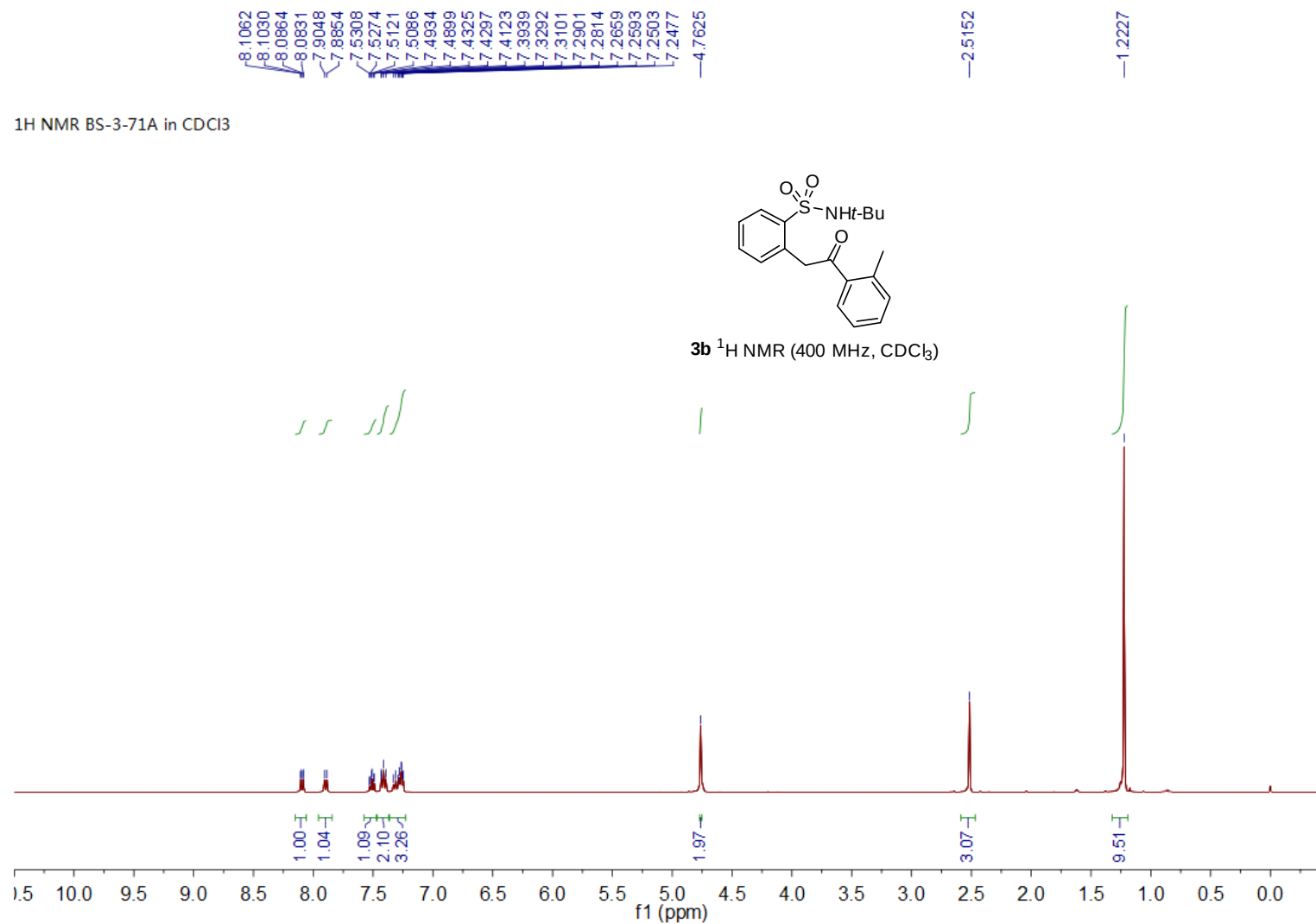
¹⁹F NMR BS-7-50 in CDCl₃

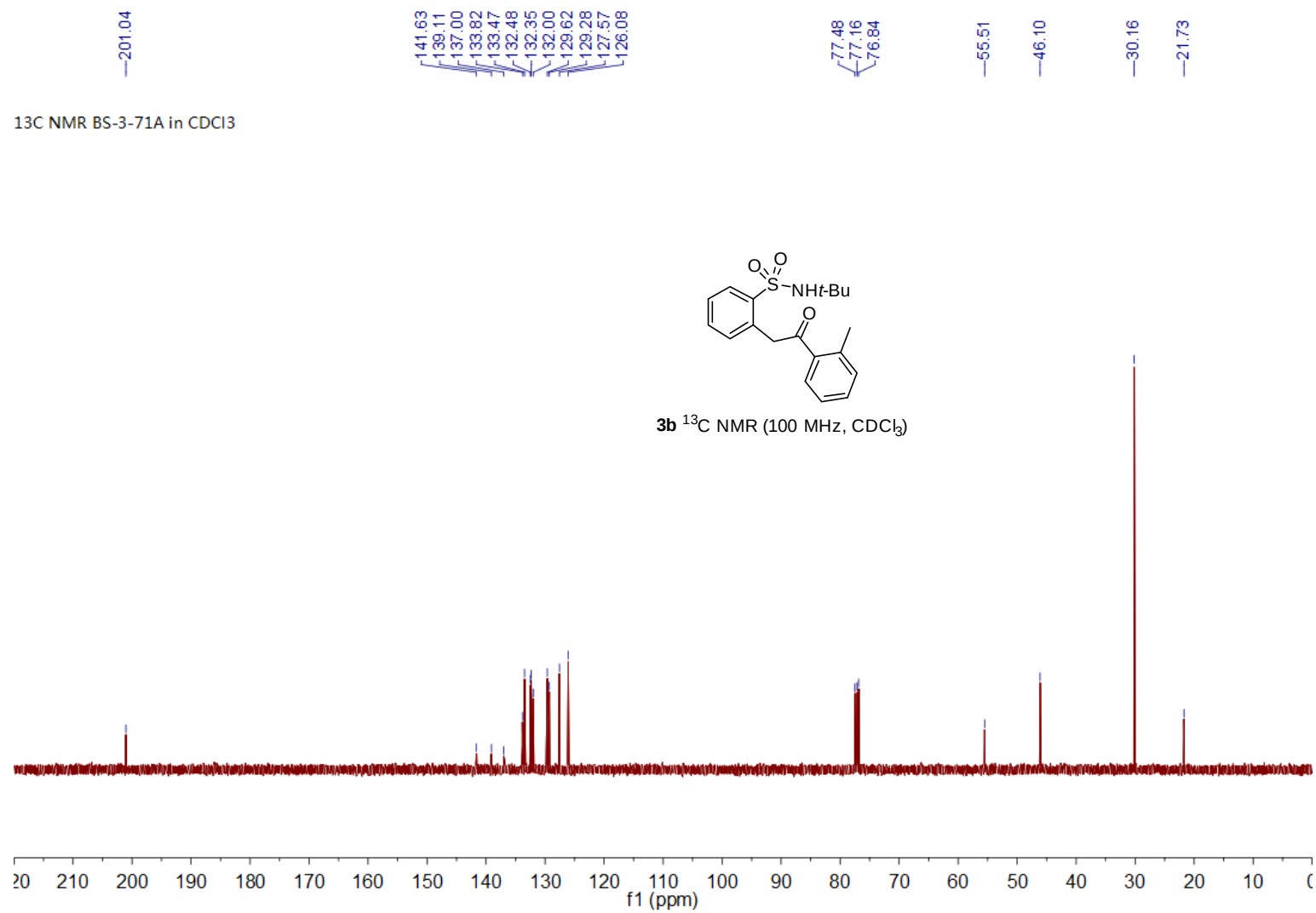


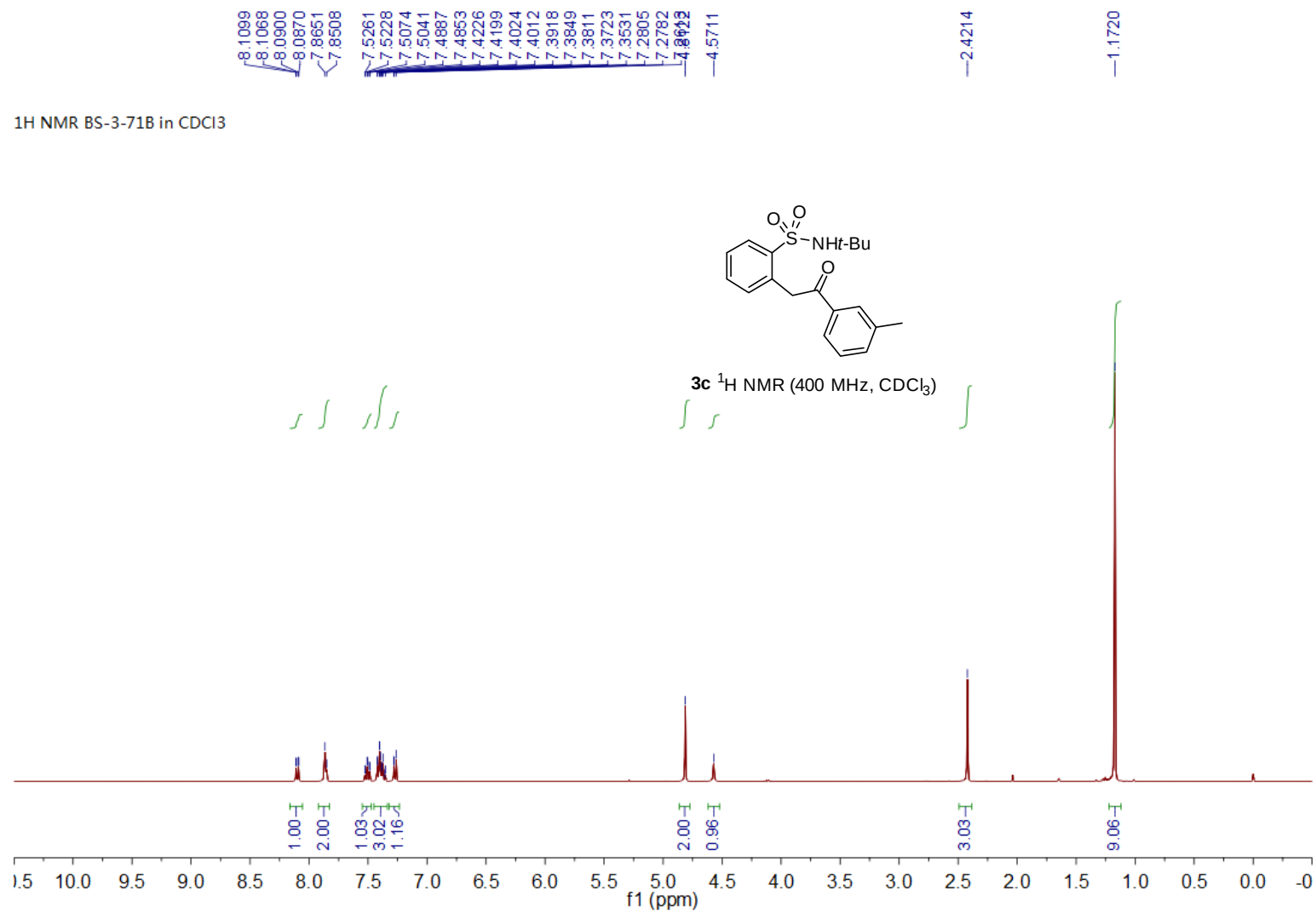
¹H NMR BS-3-68 in CDCl₃

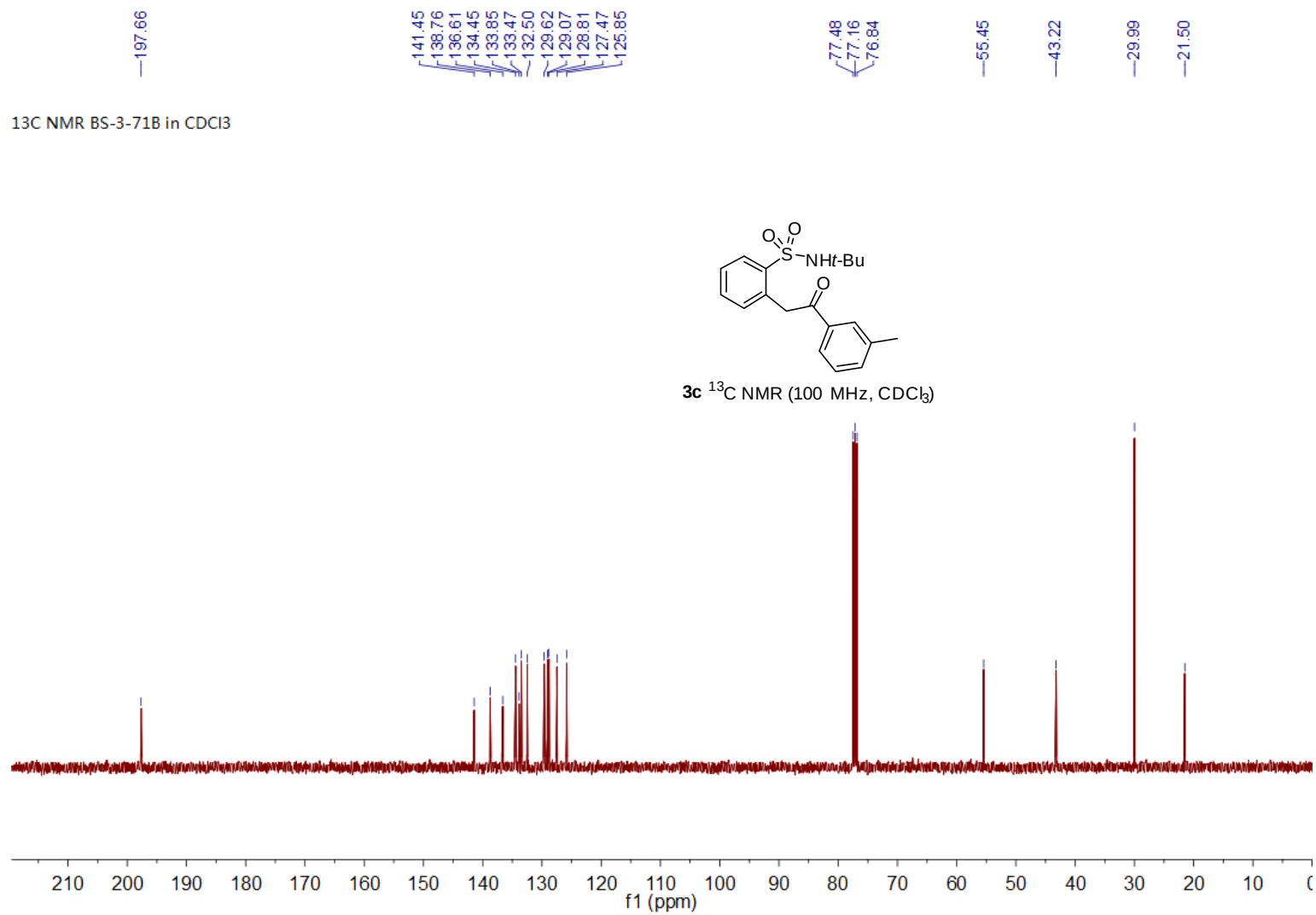


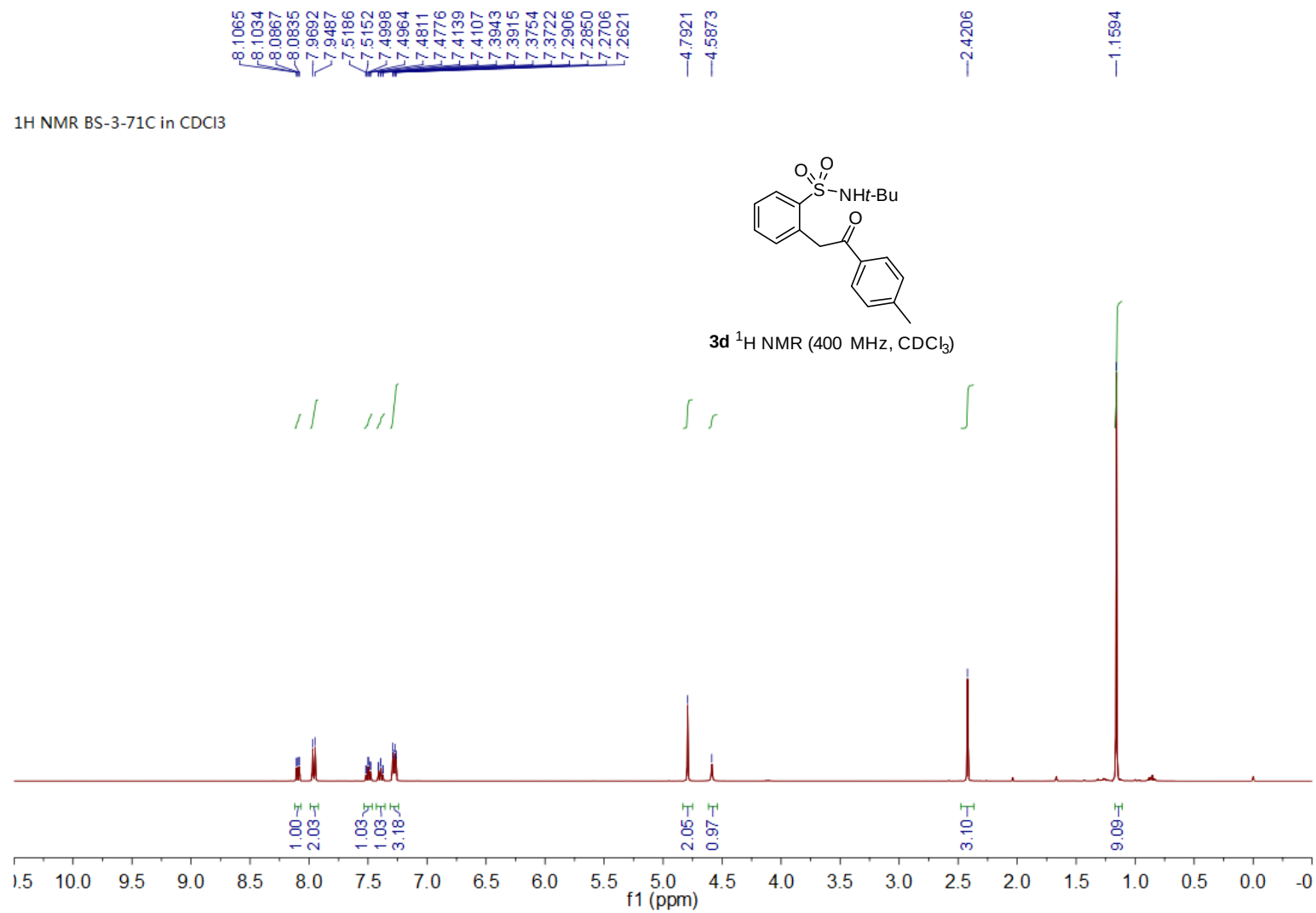


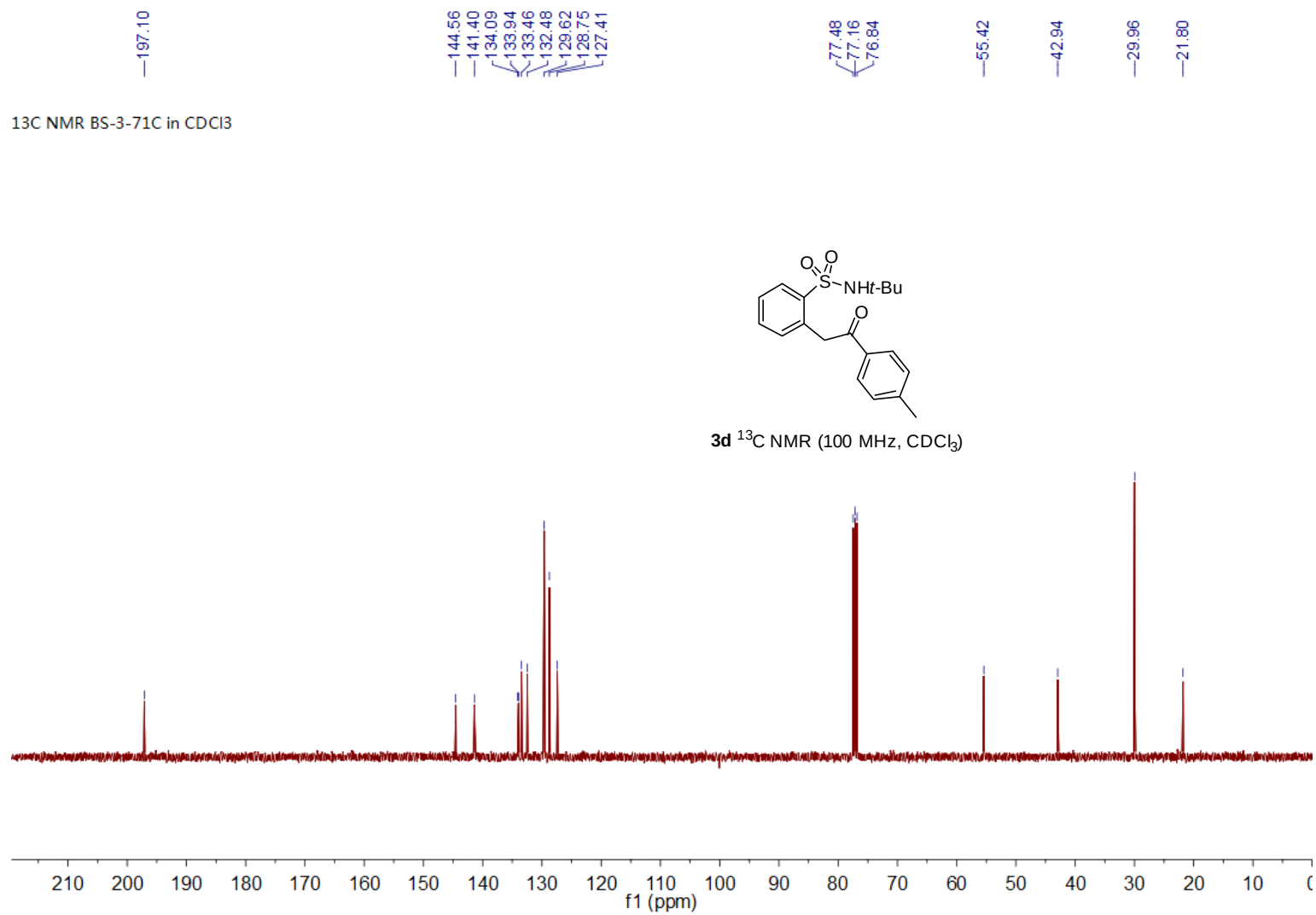


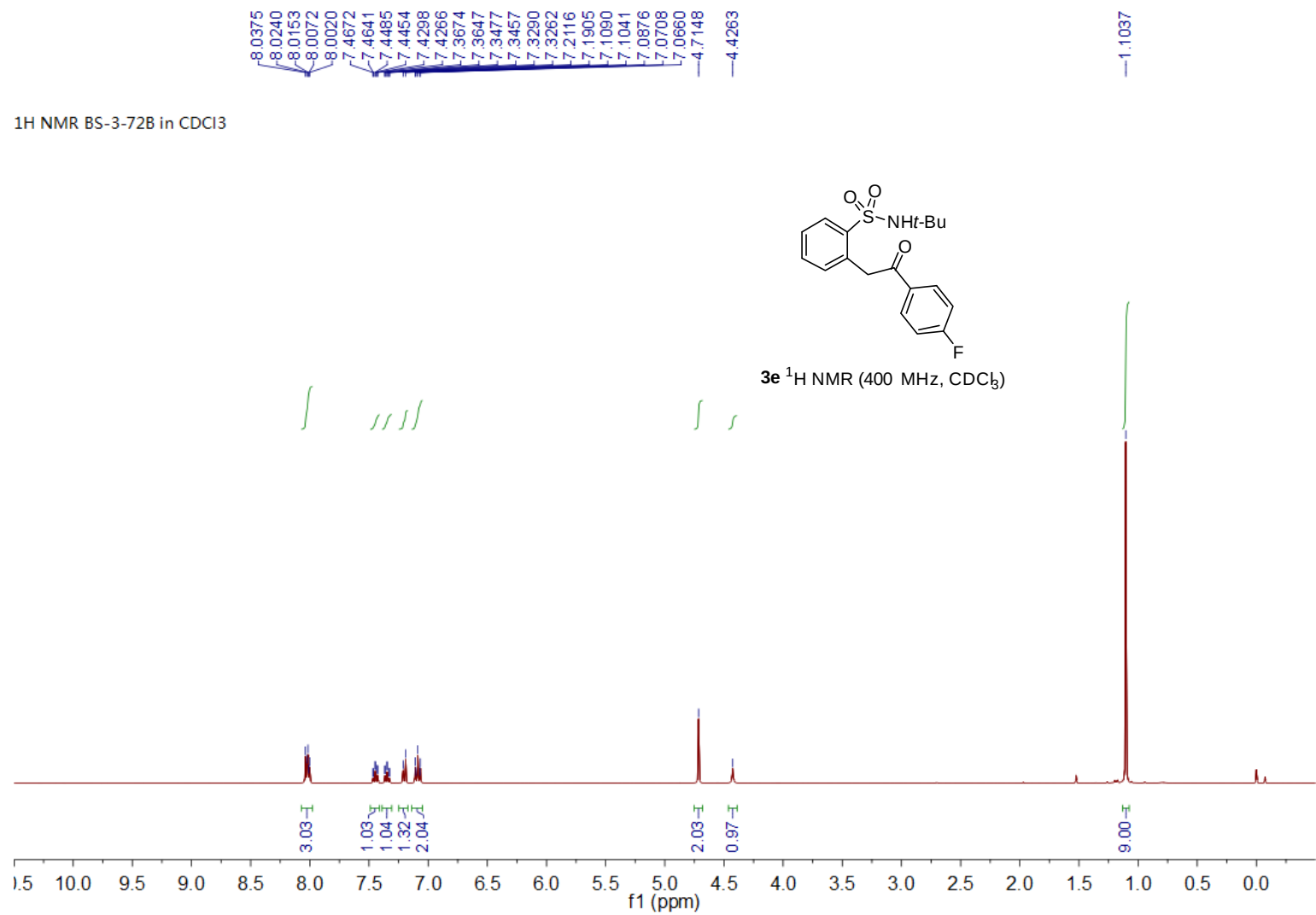




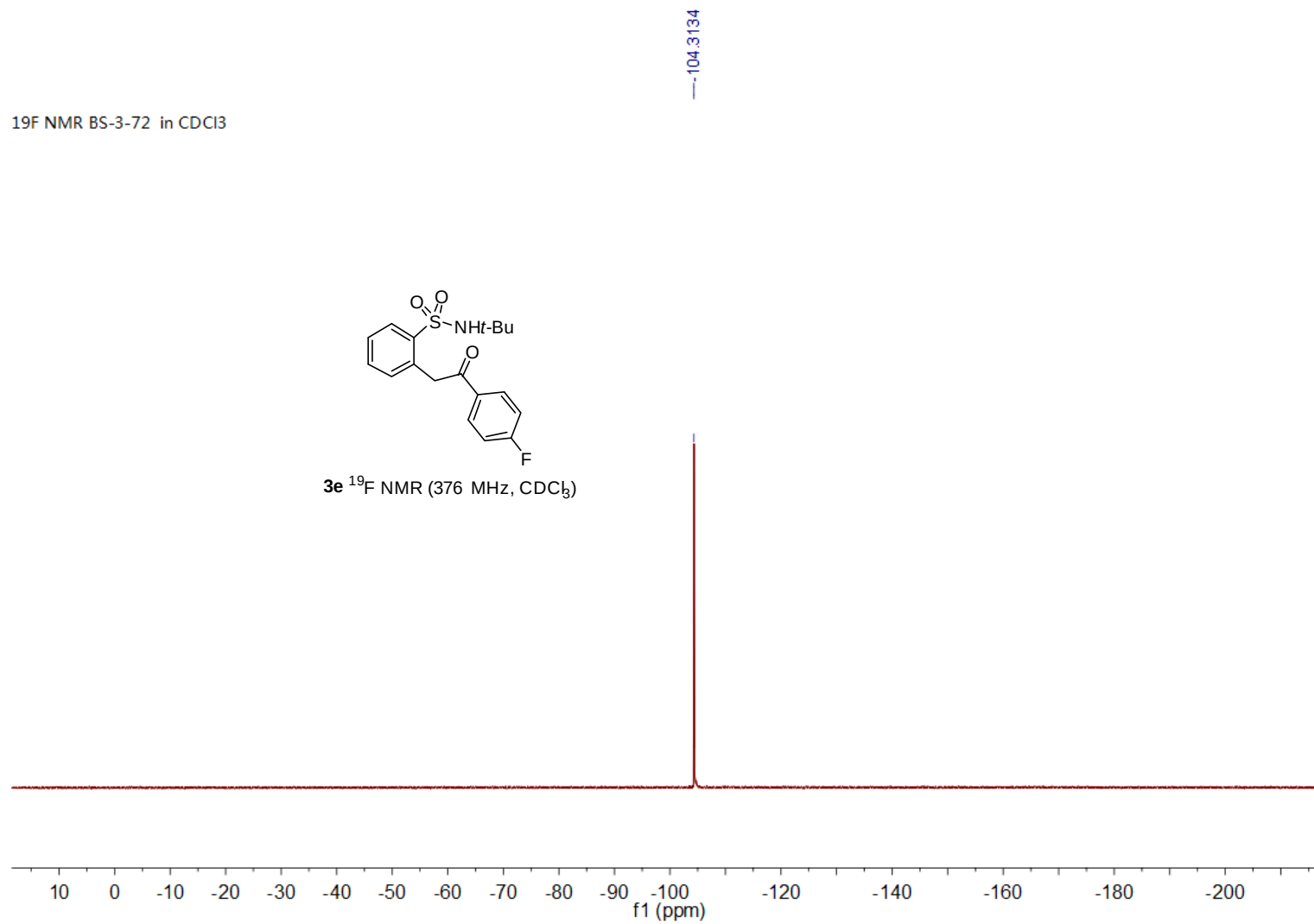
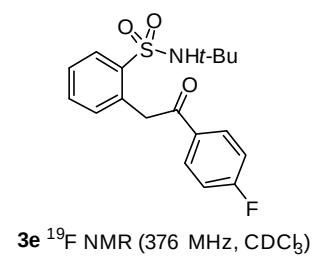


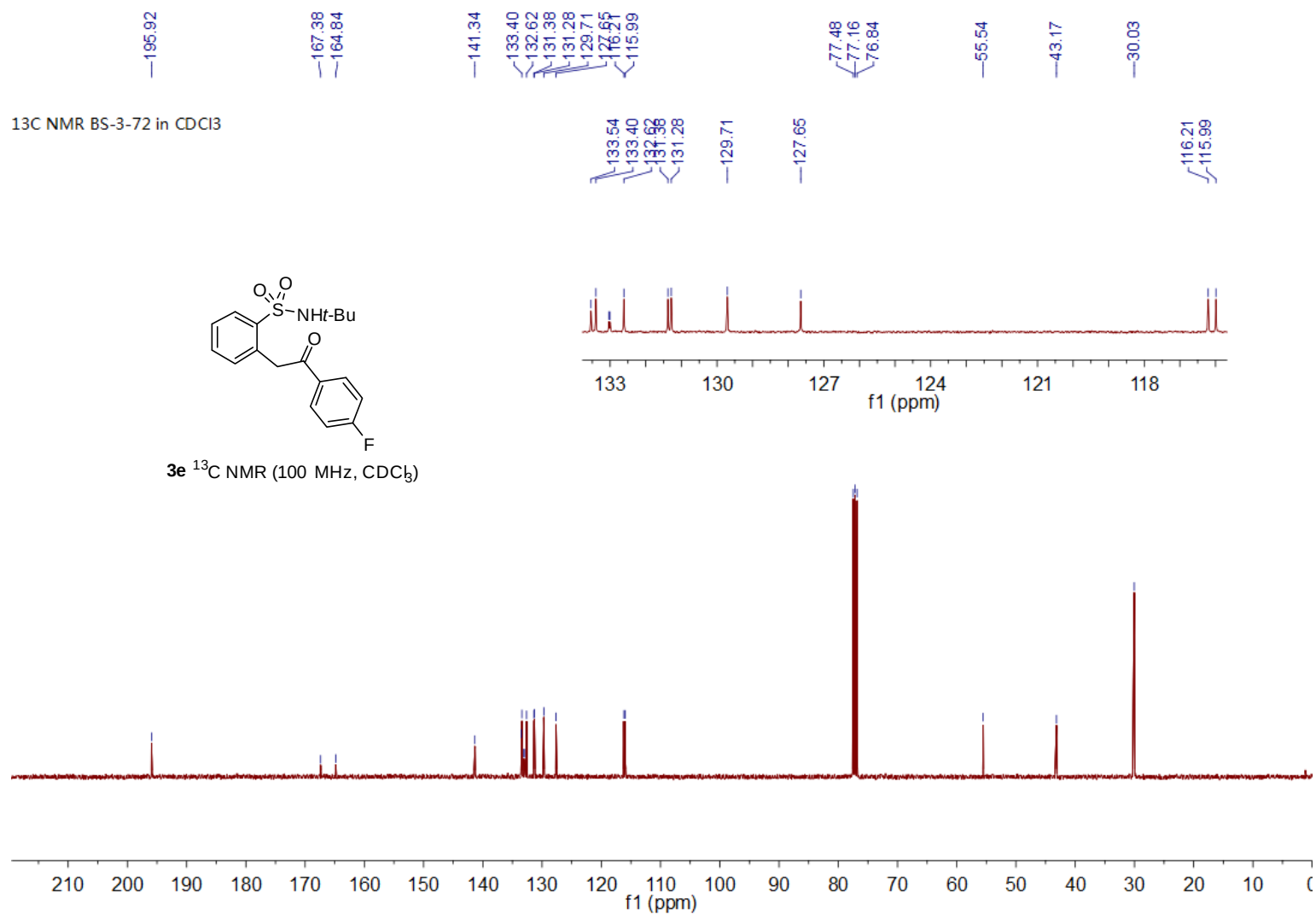




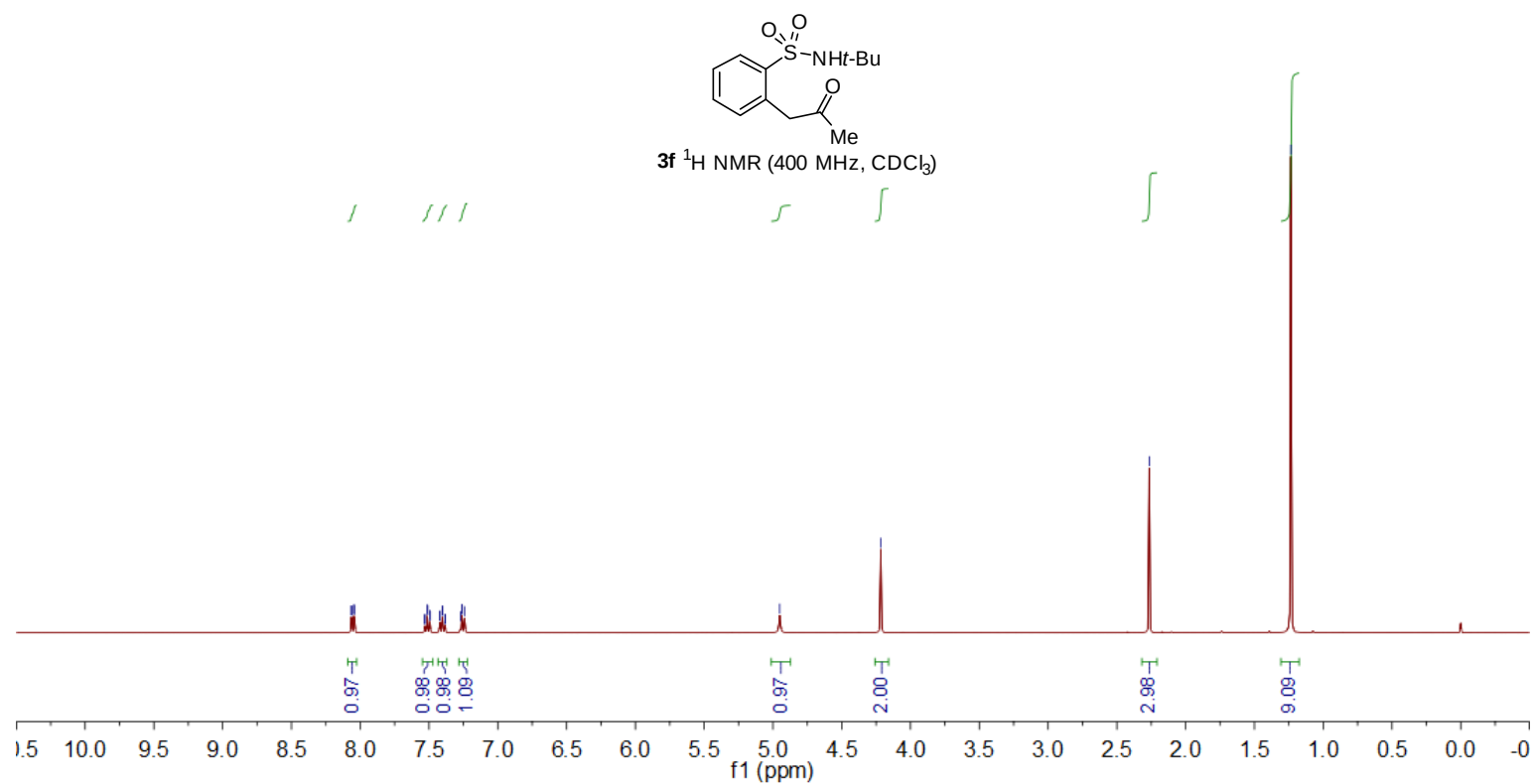


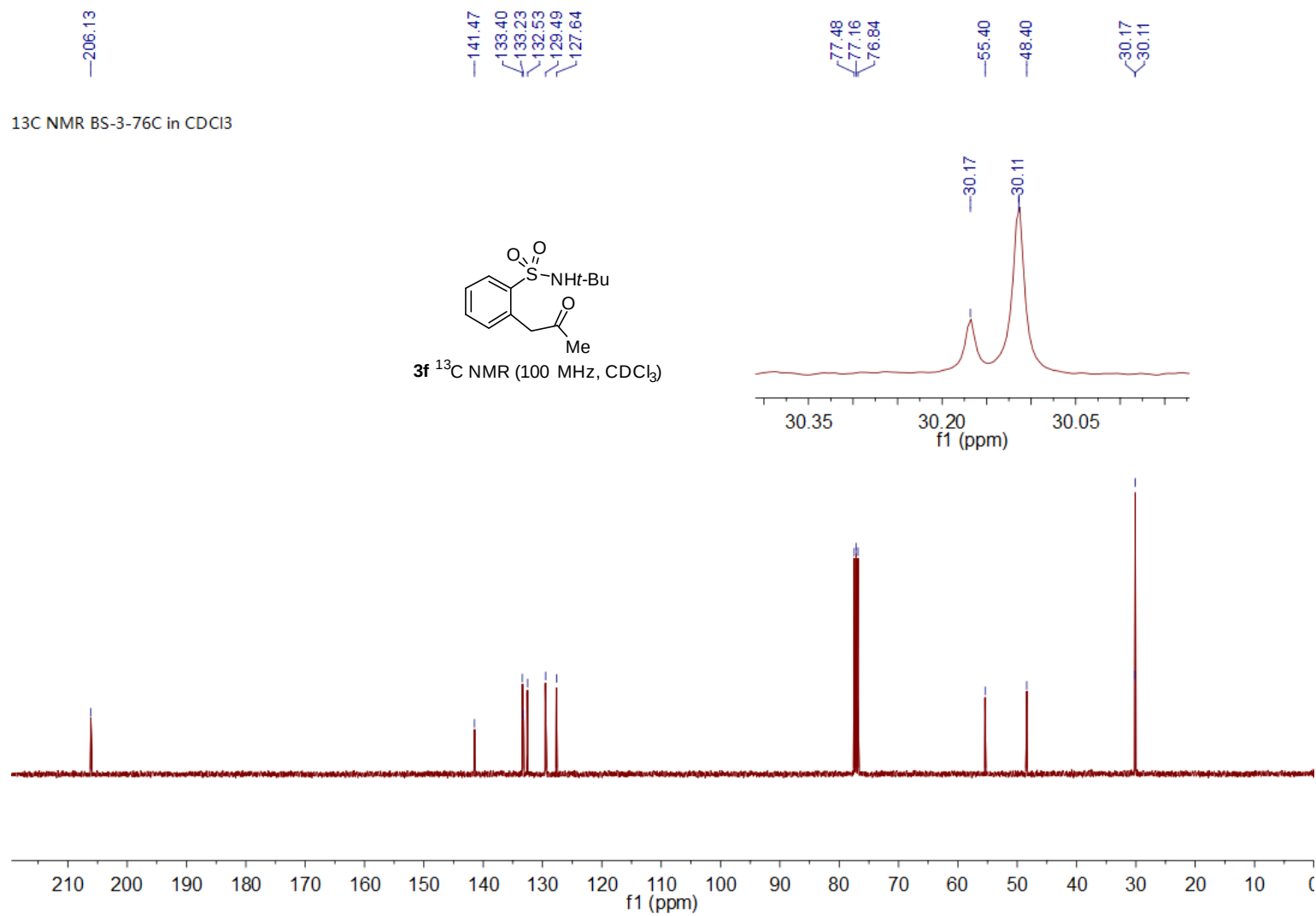
¹⁹F NMR BS-3-72 in CDCl₃



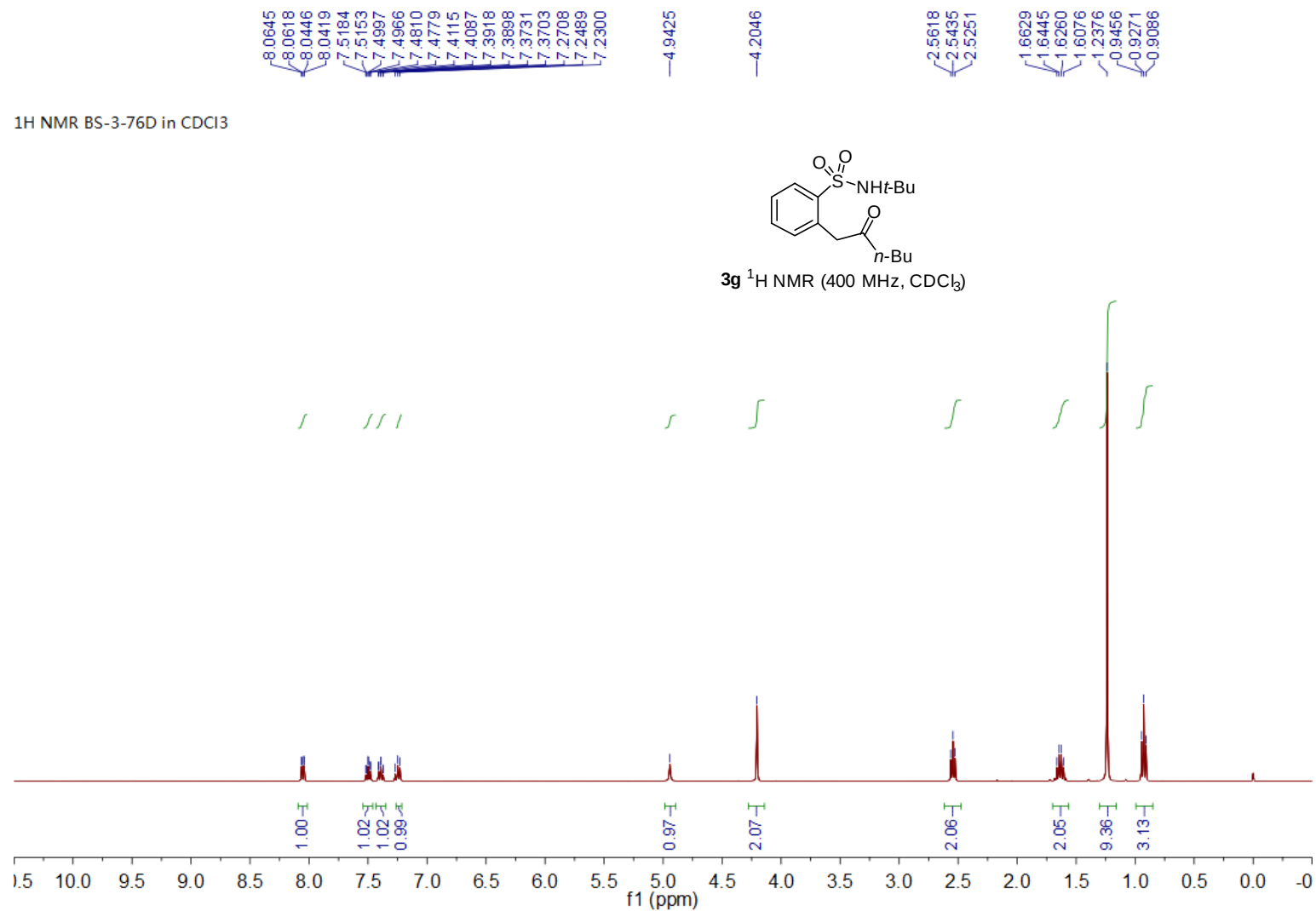
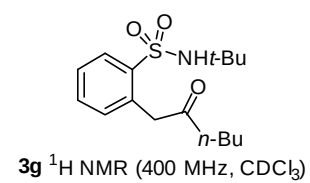


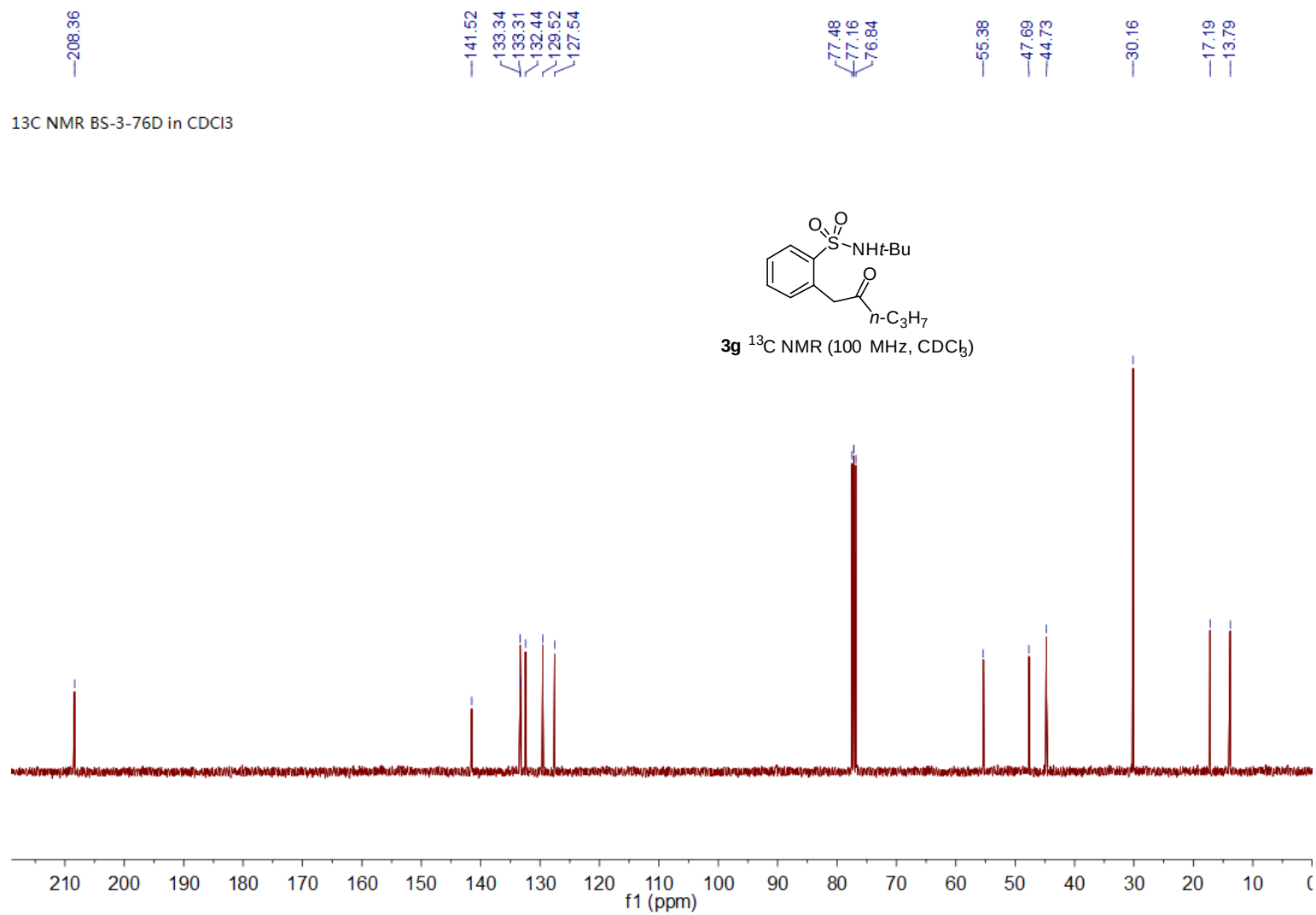
¹H NMR BS-3-76C in CDCl₃

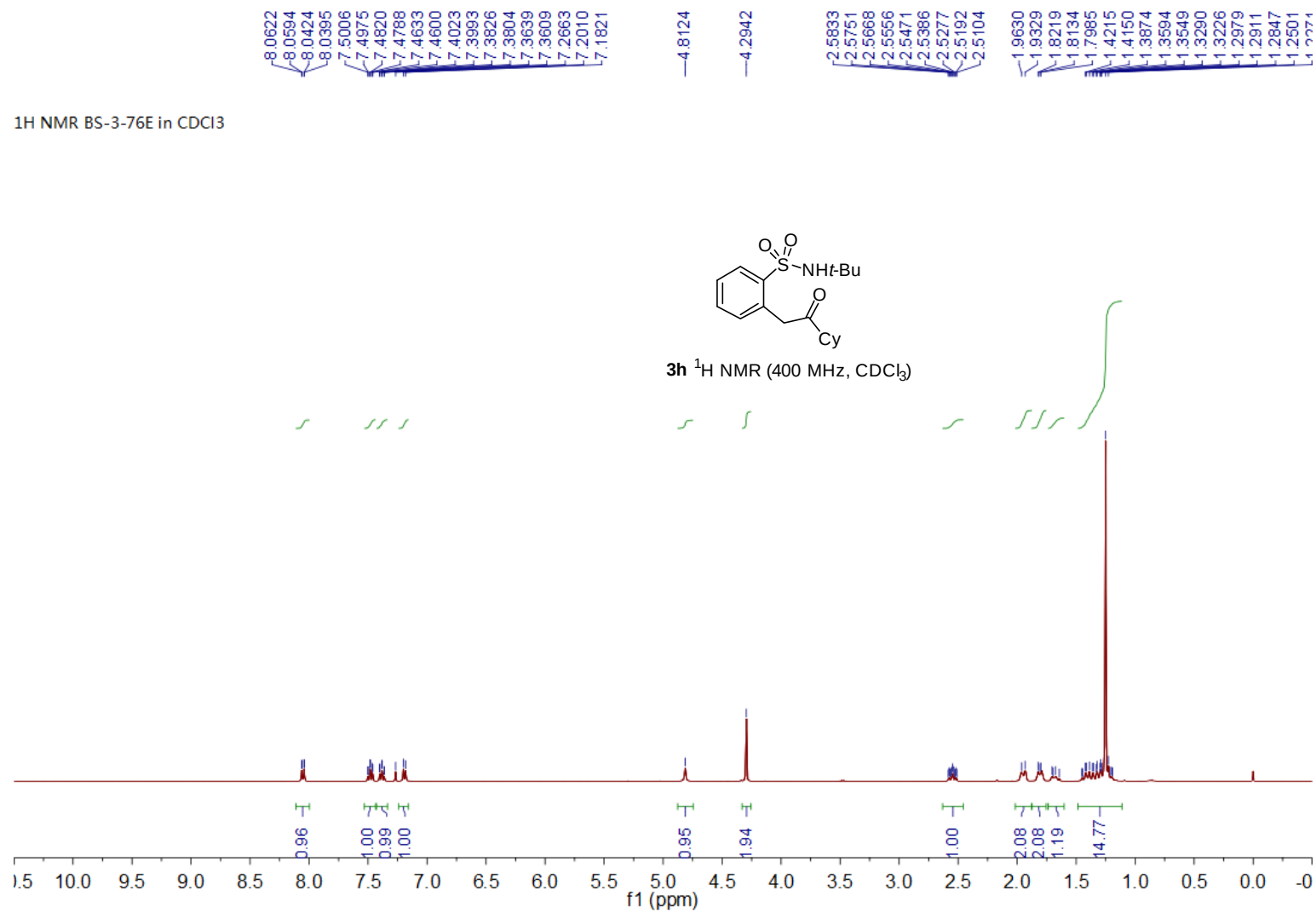


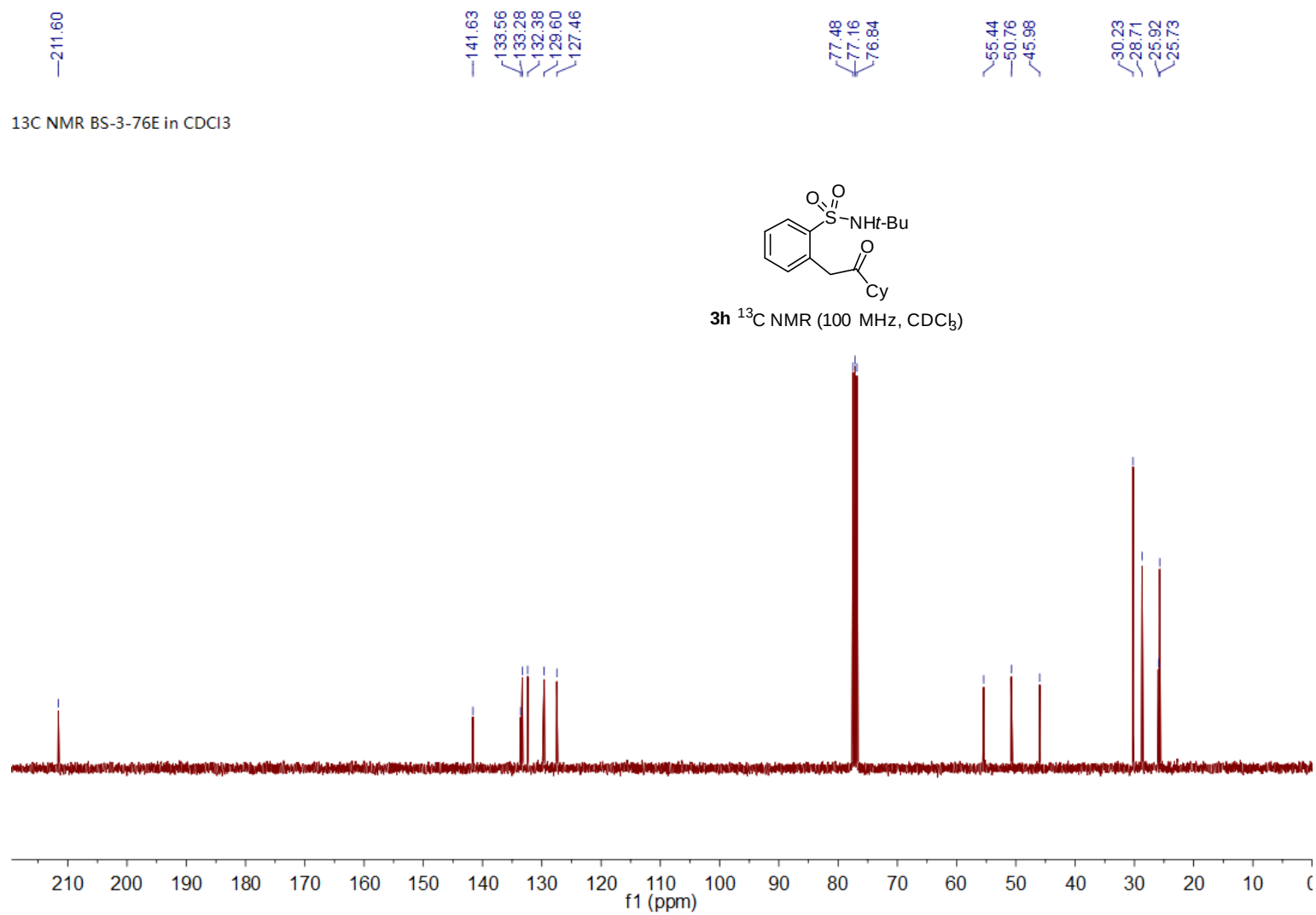


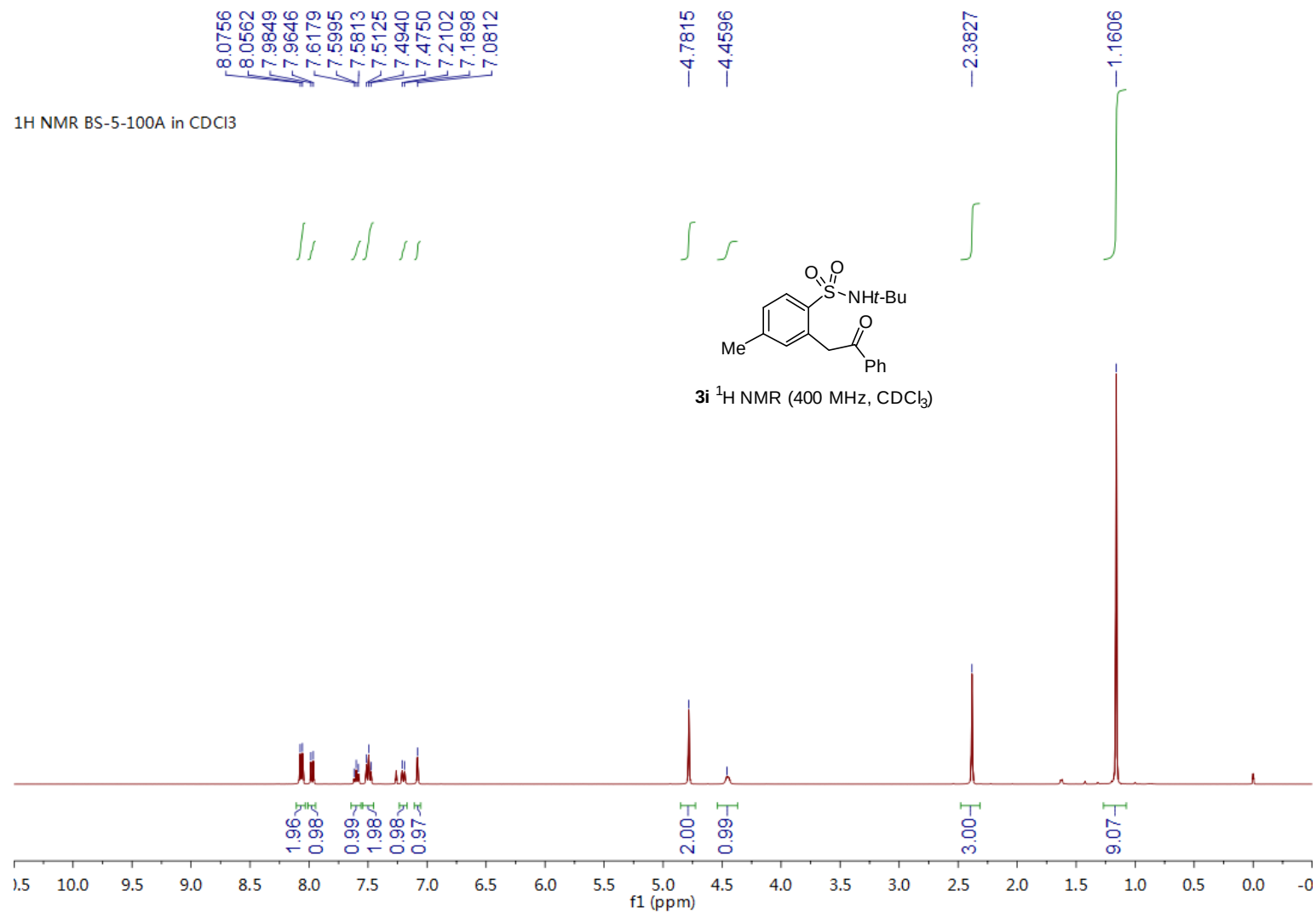
¹H NMR BS-3-76D in CDCl₃

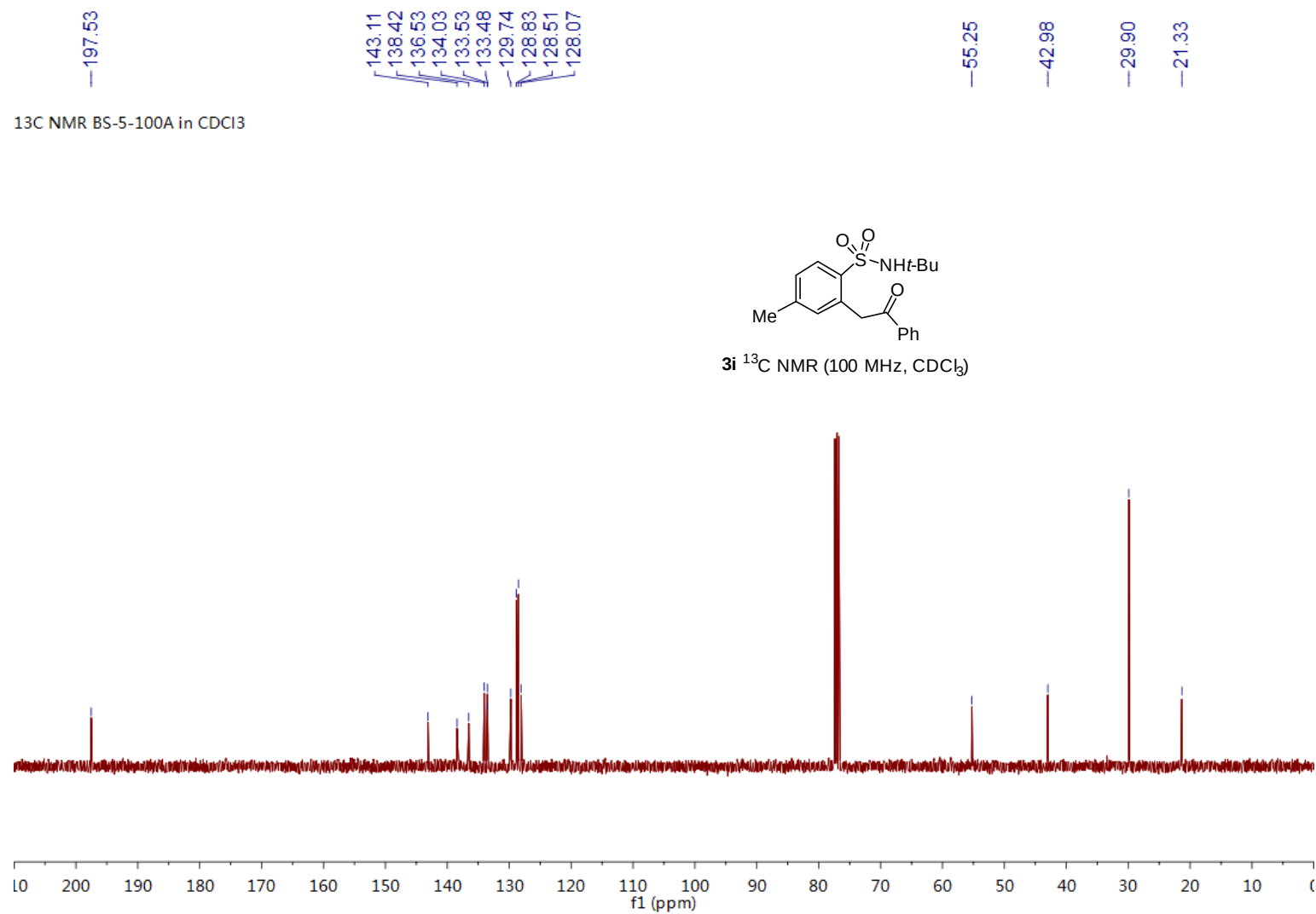


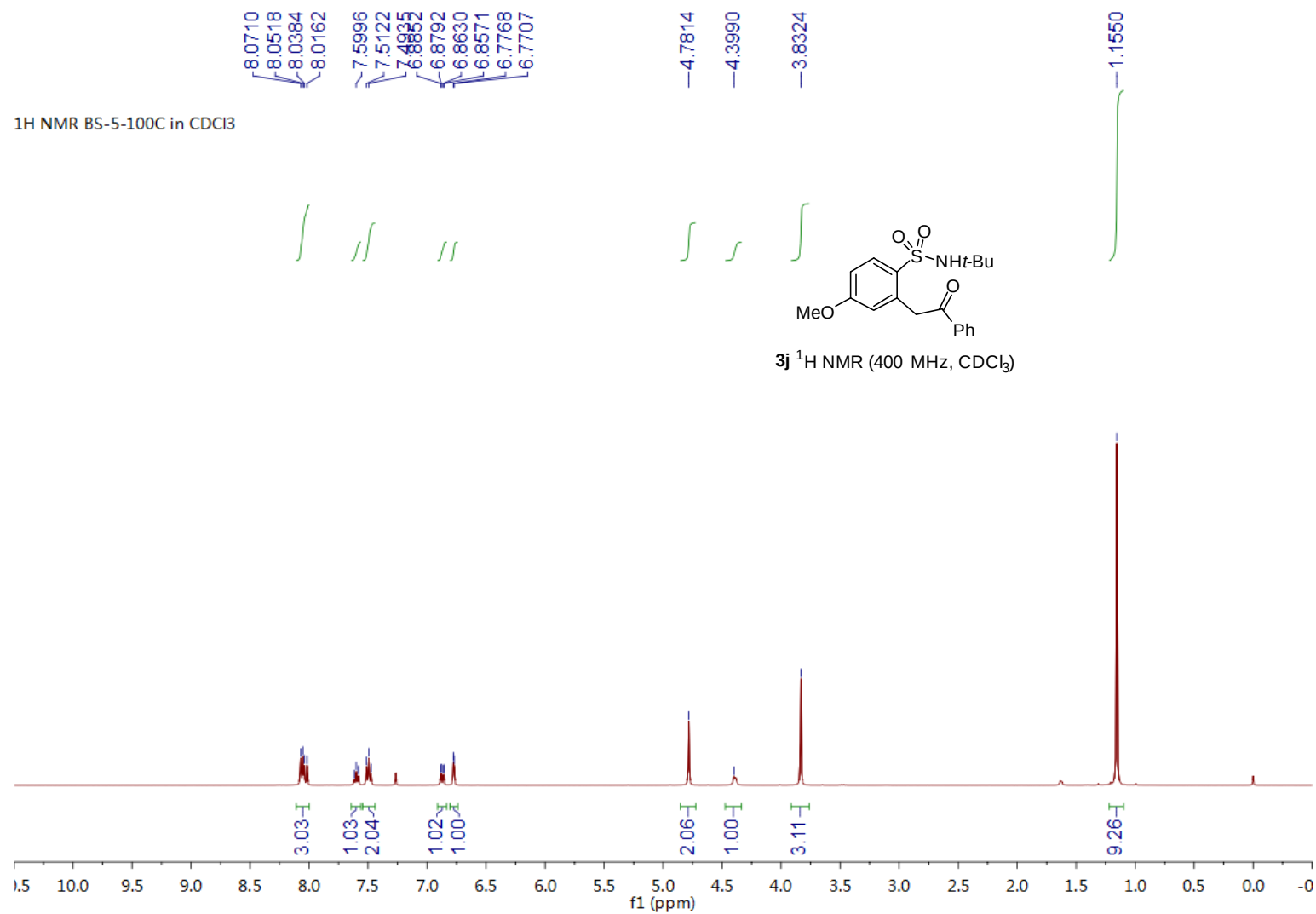


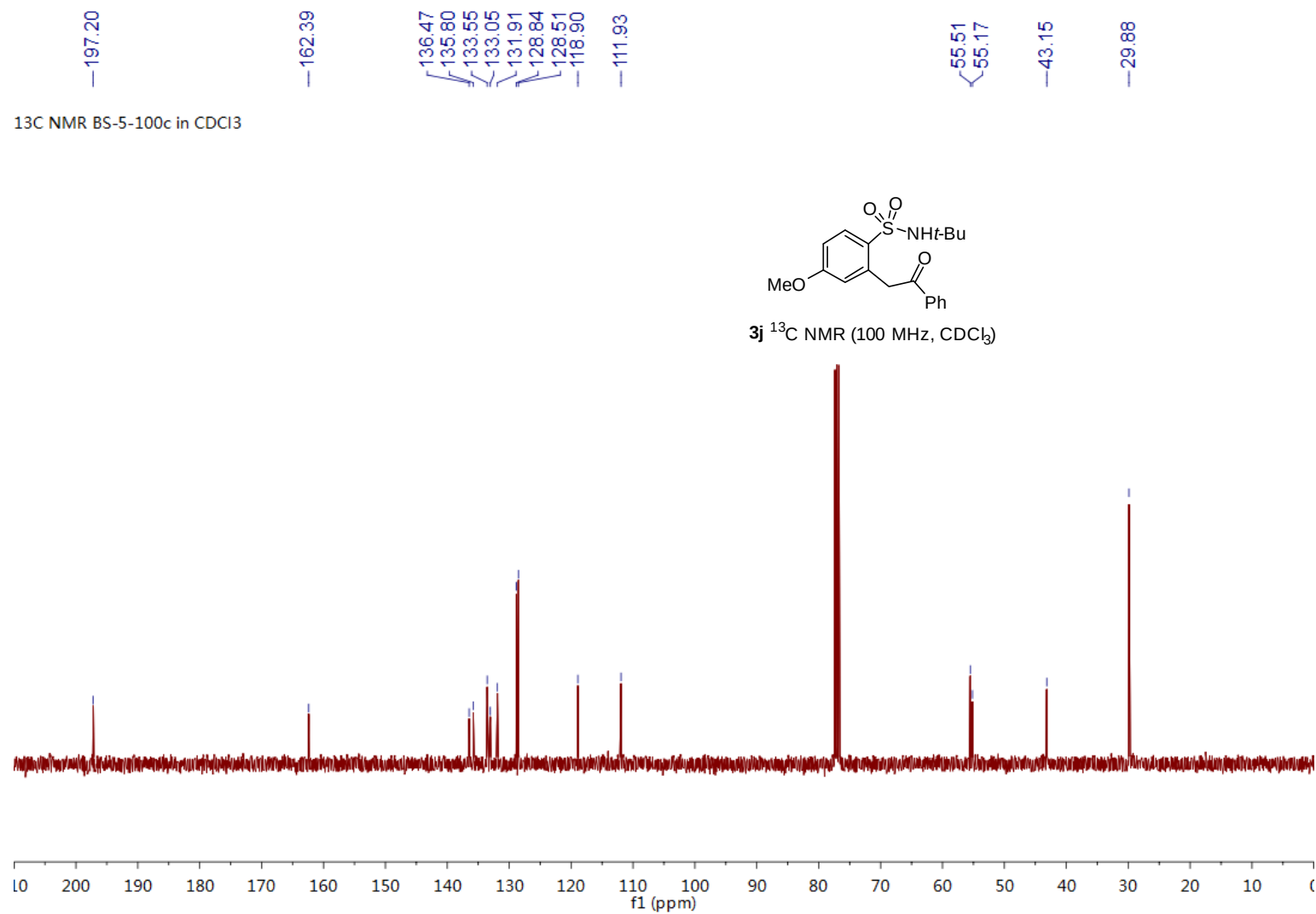


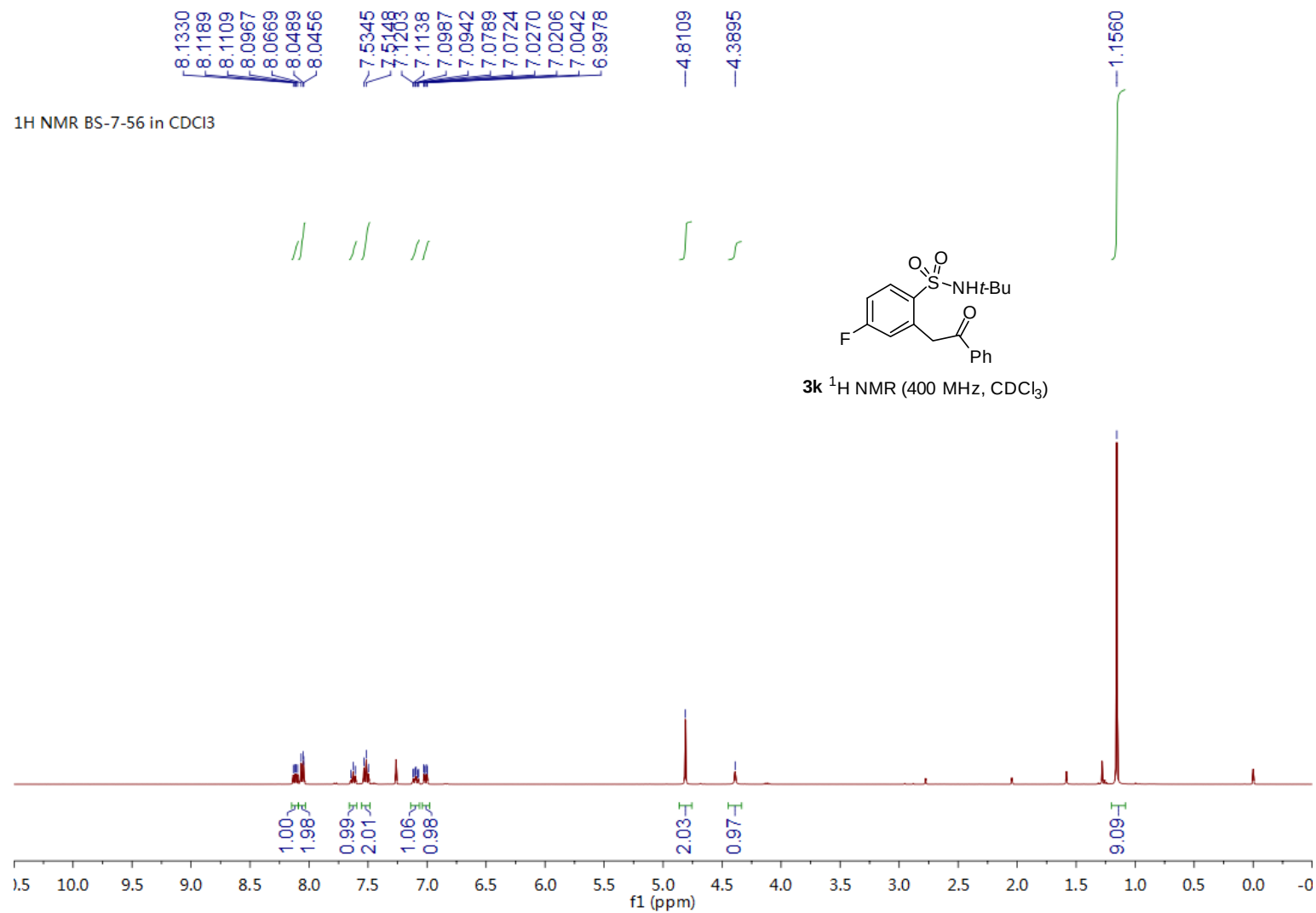


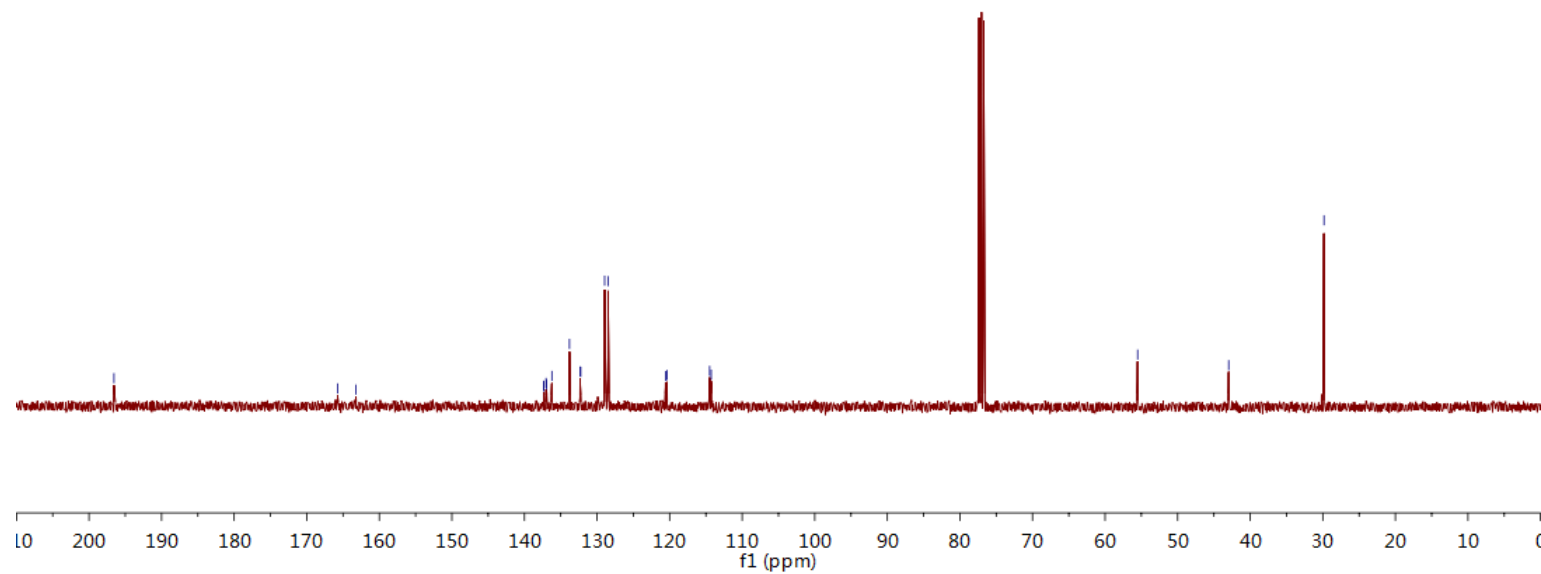
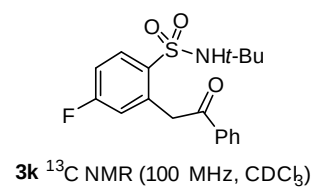
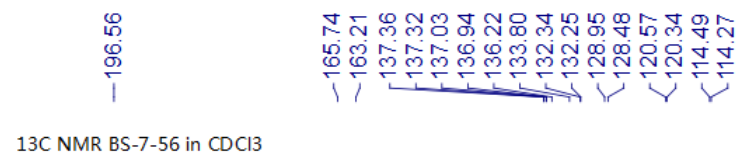






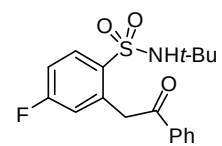




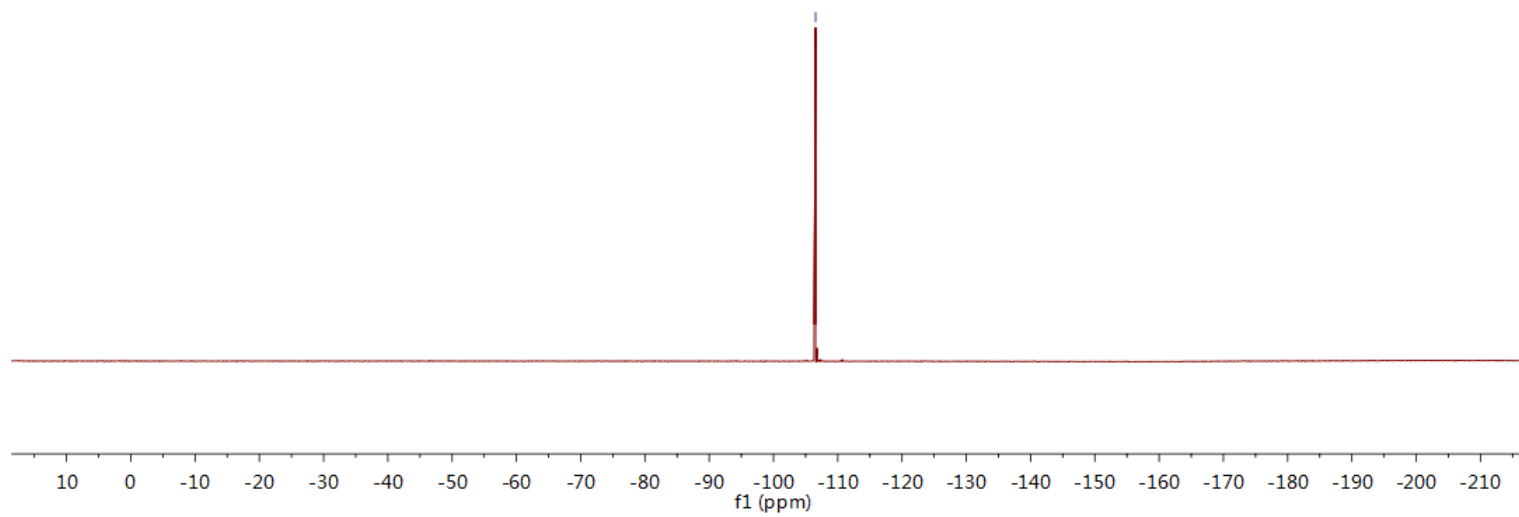


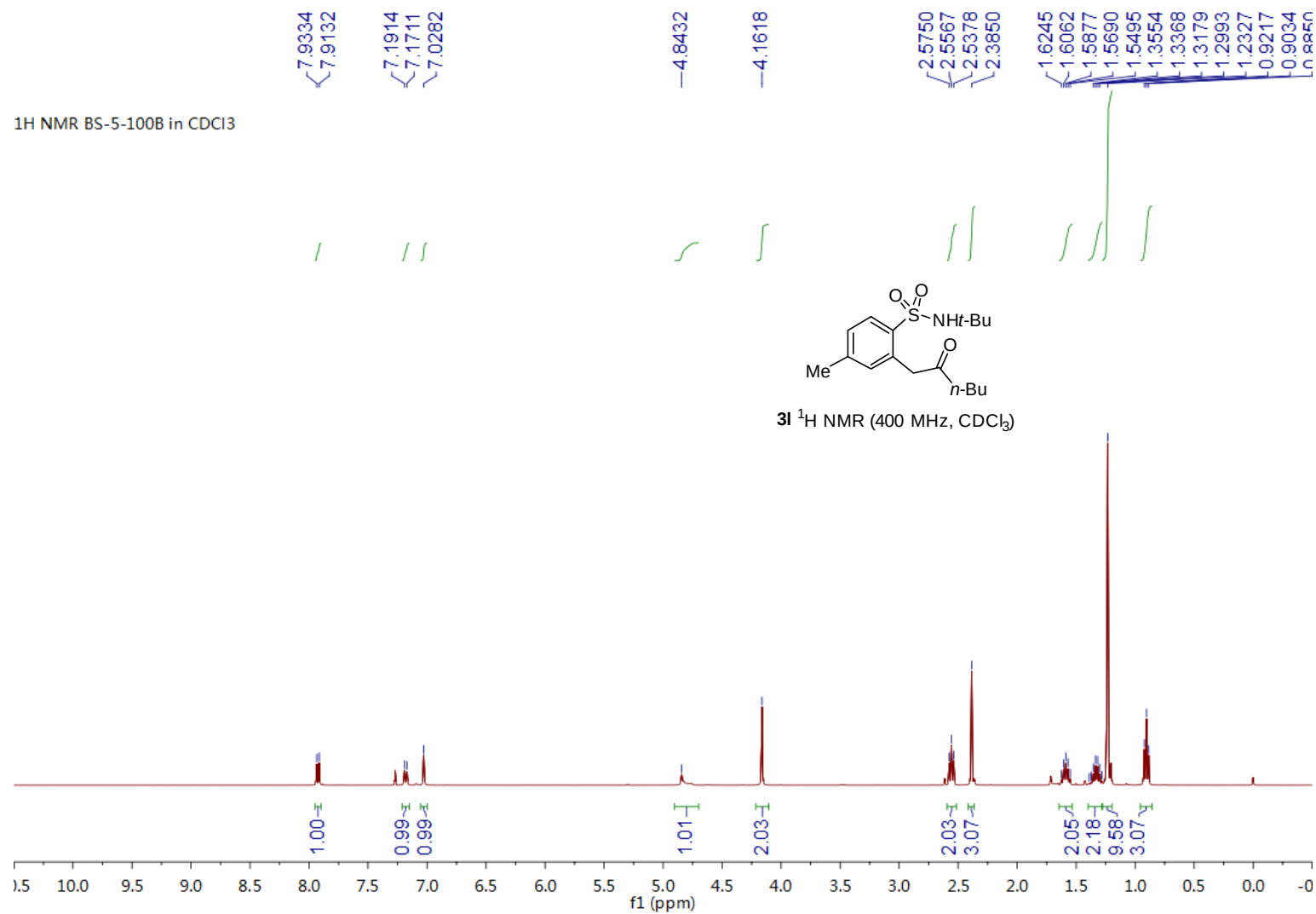
¹⁹F NMR BS-7-56 in CDCl₃

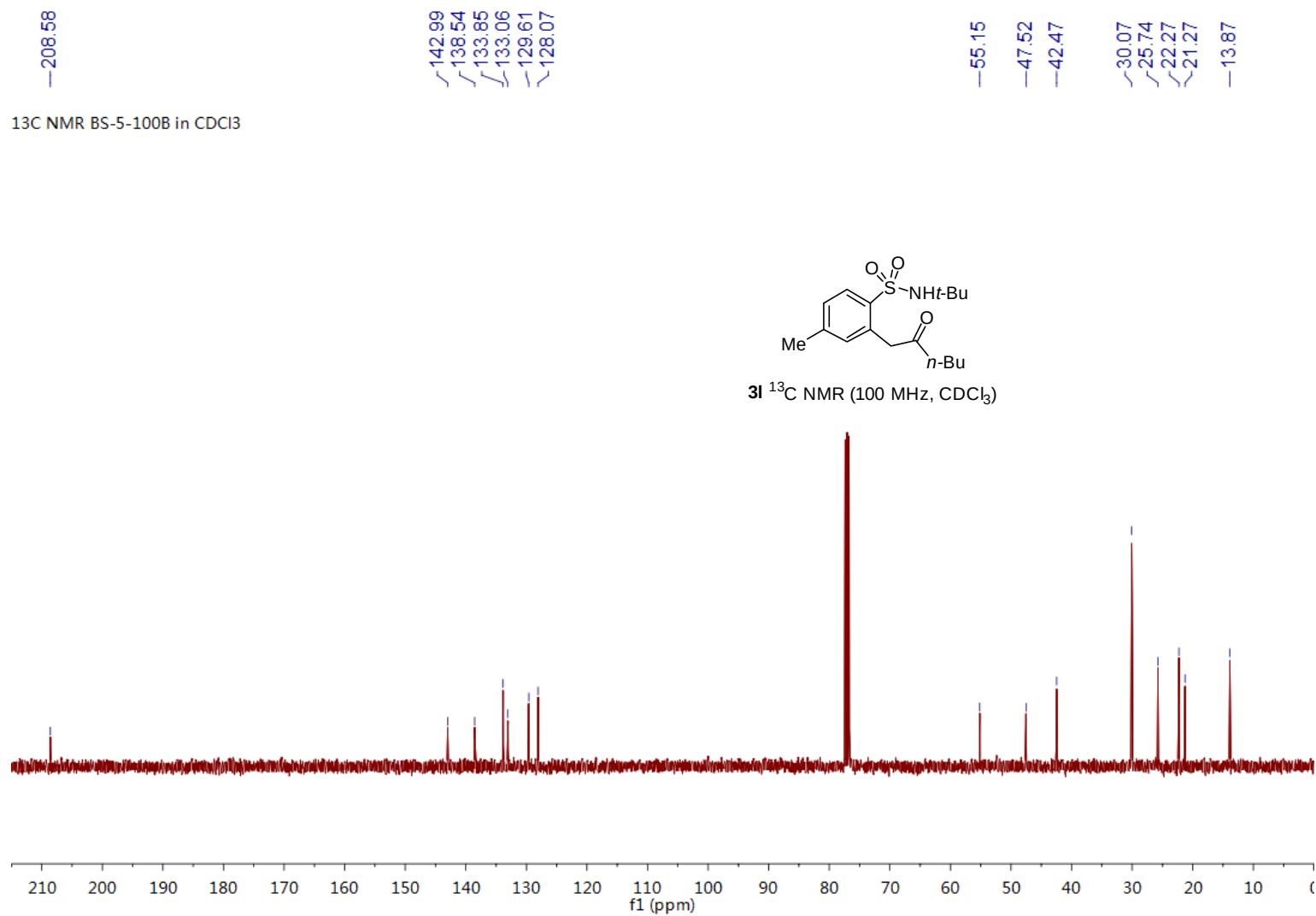
--106.5440

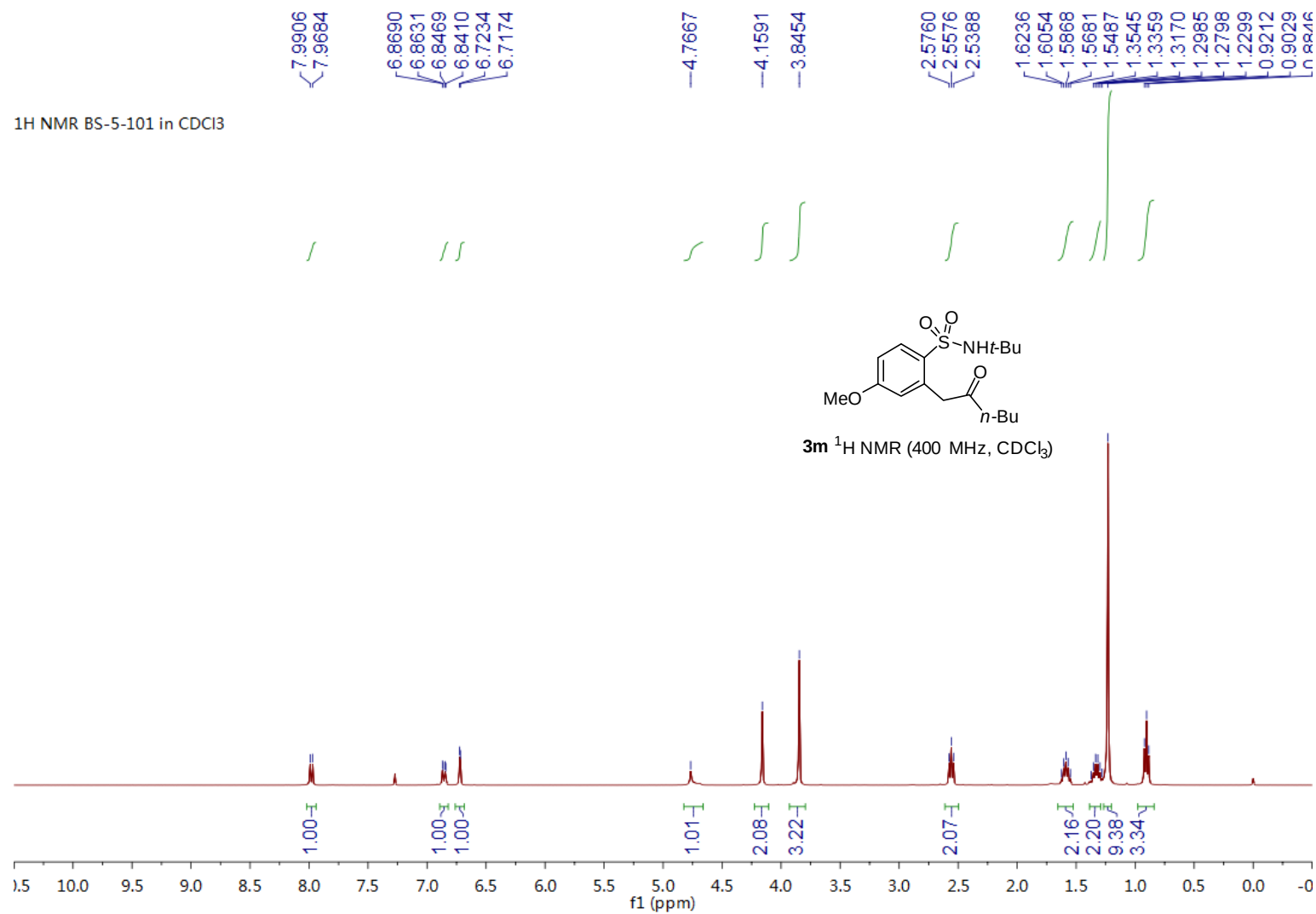


3k ¹⁹F NMR (376 MHz, CDCl₃)

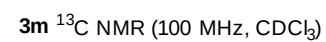


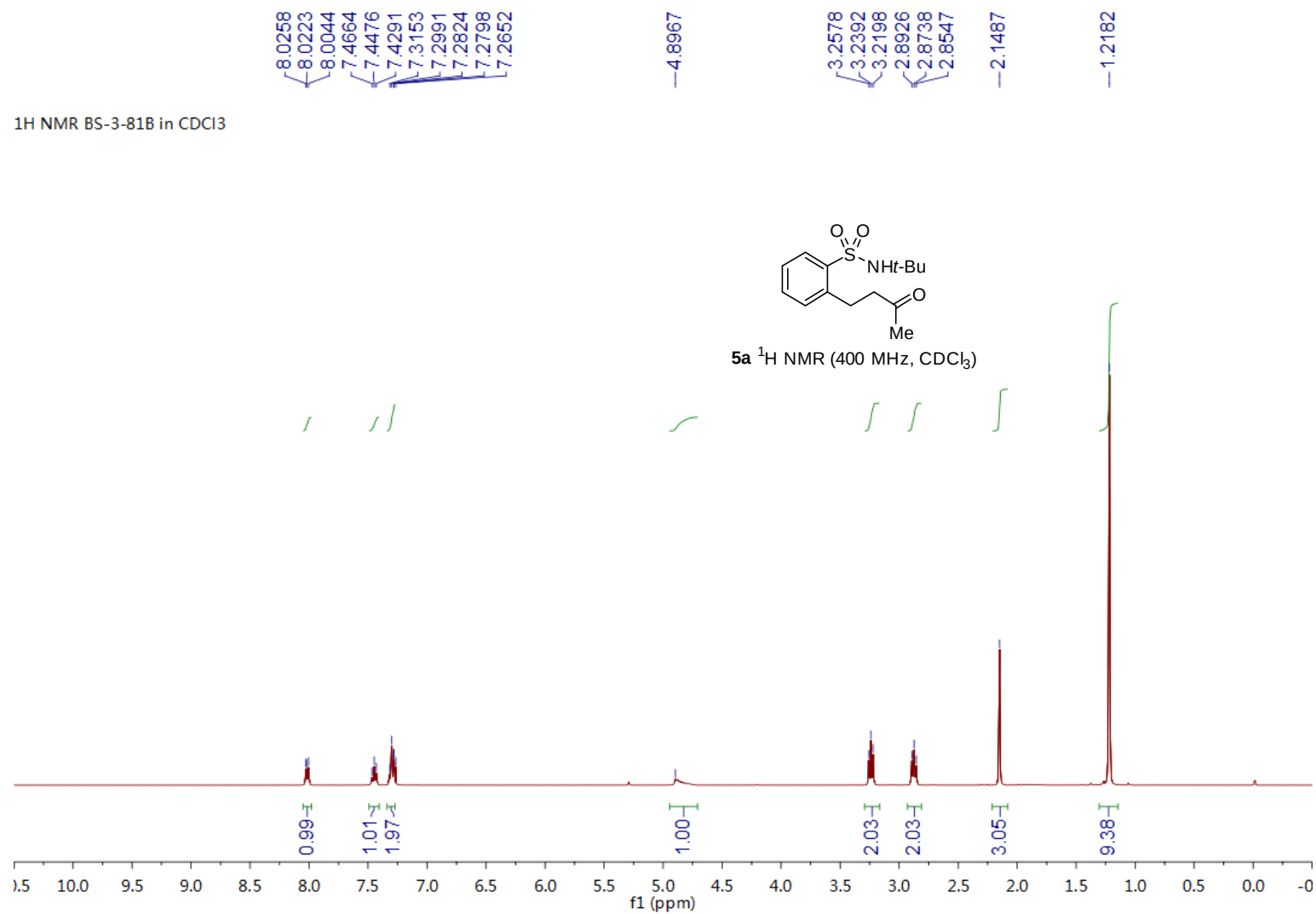


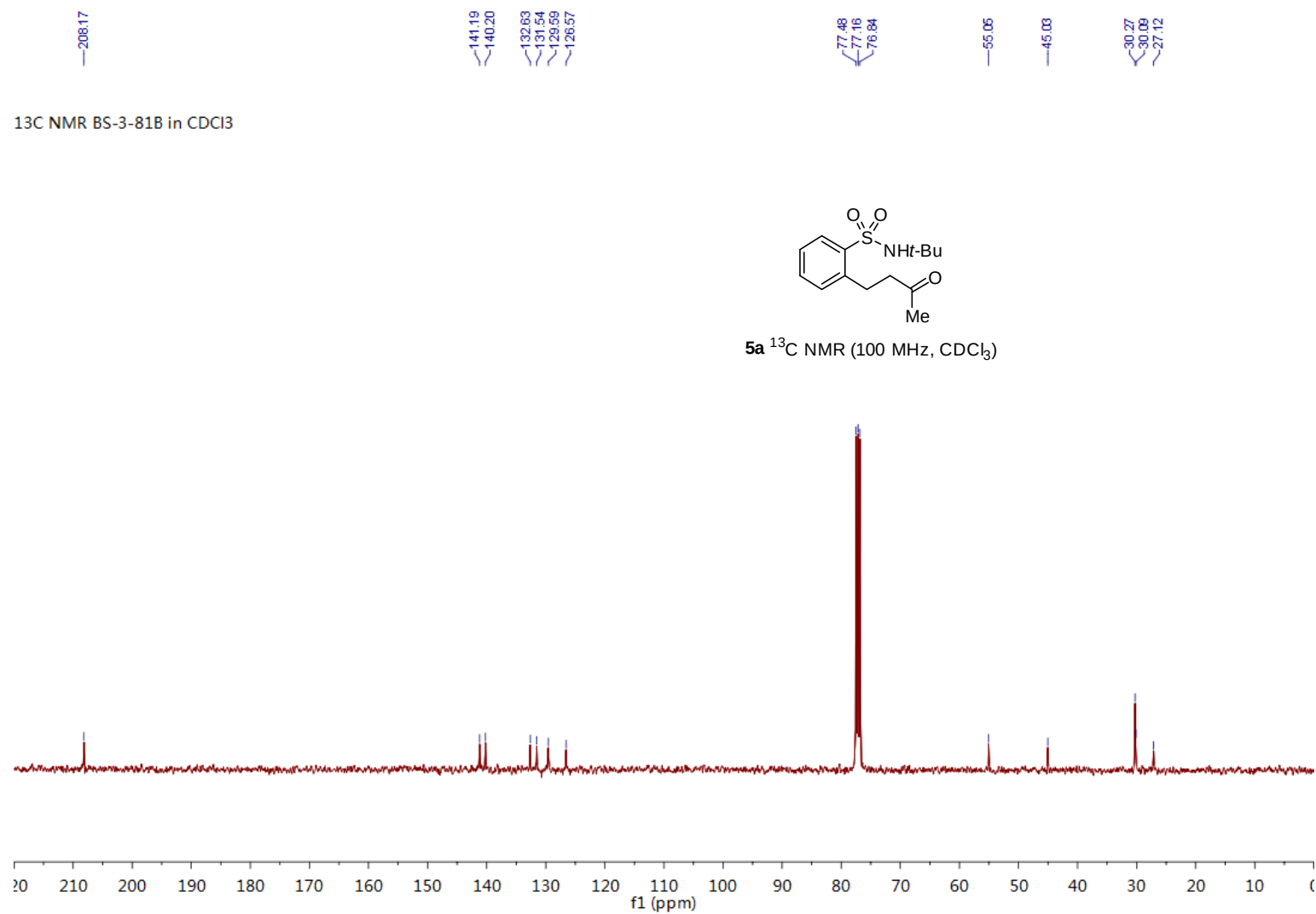


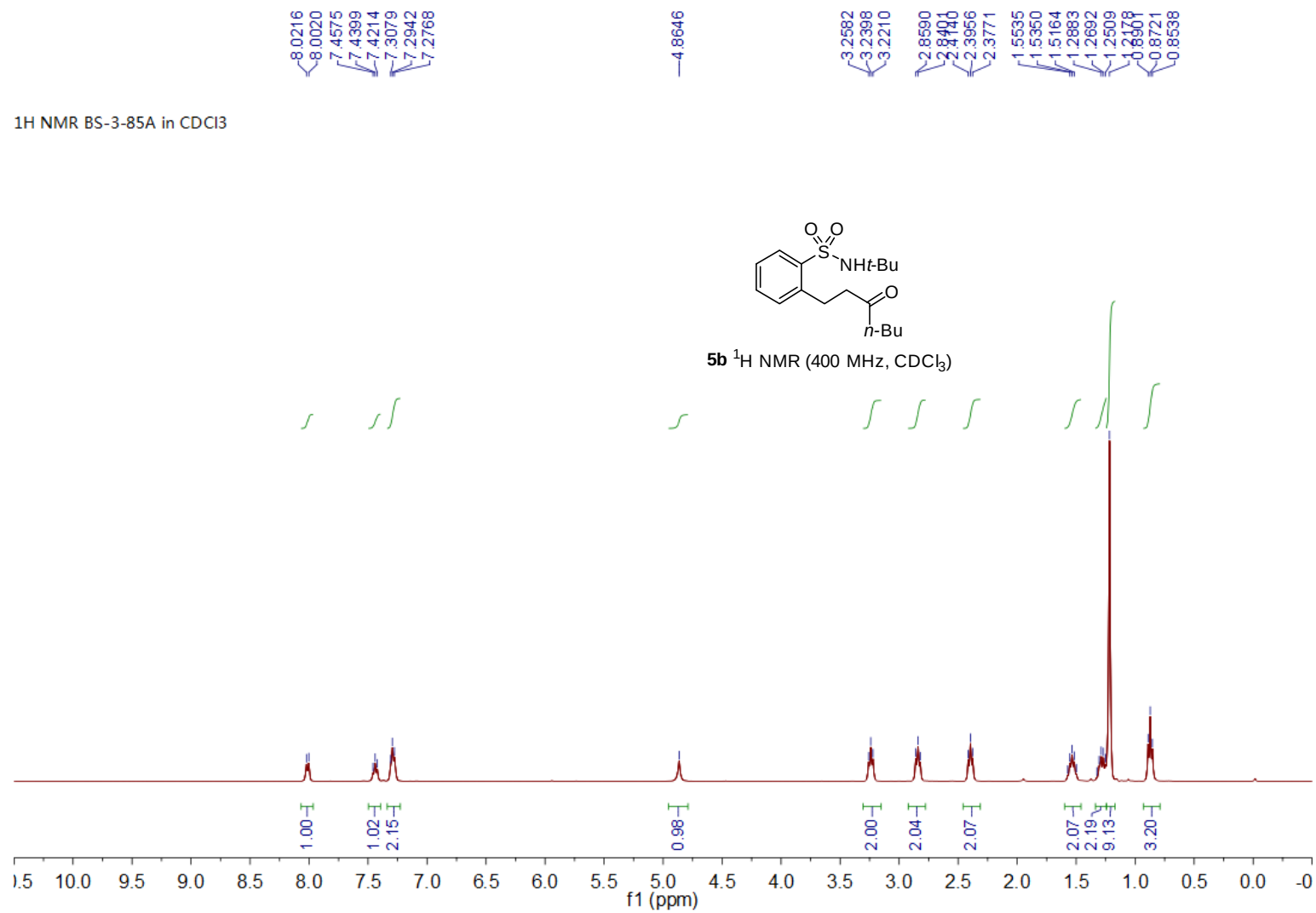


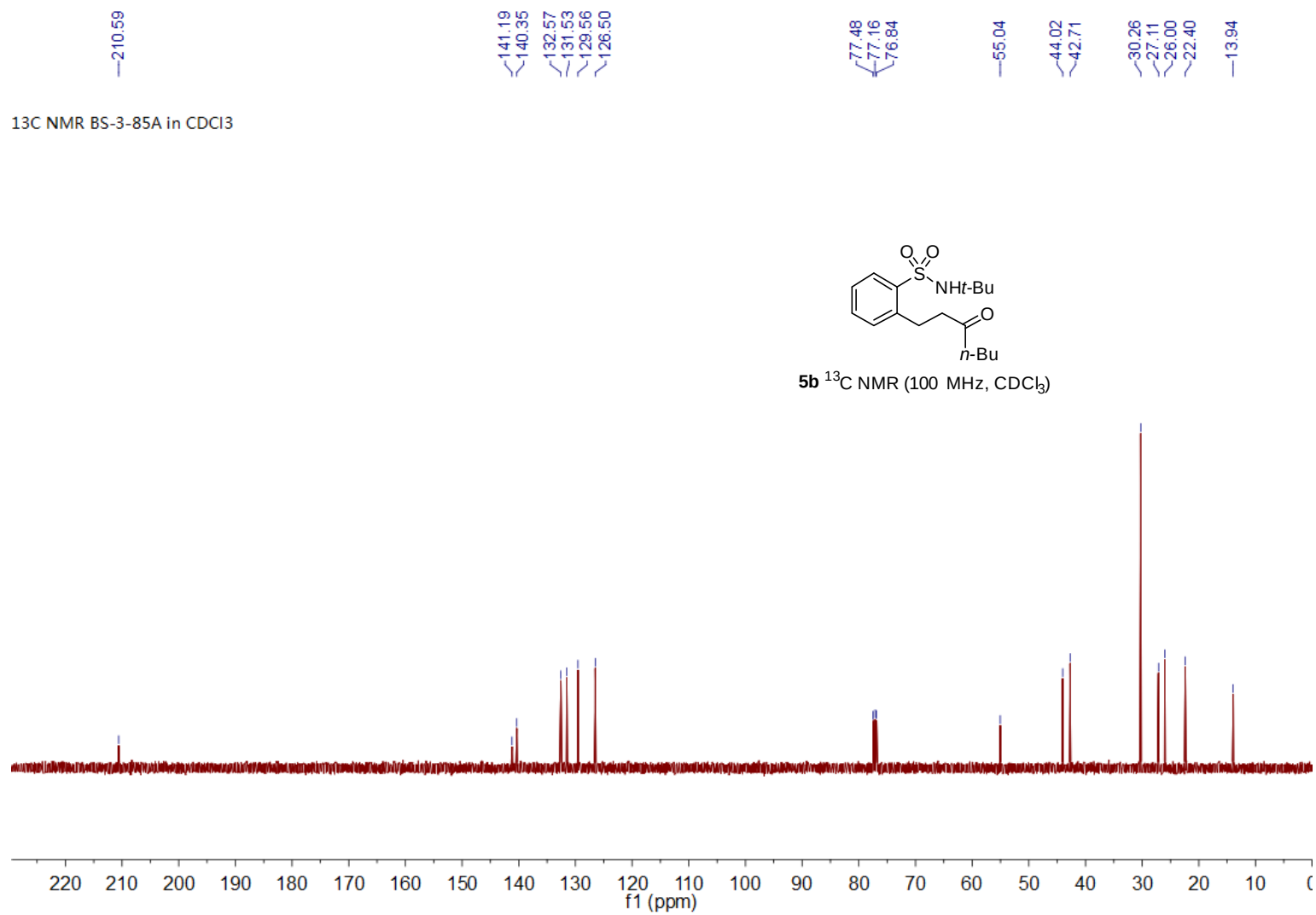
¹³C NMR BS-5-101 in CDCl₃

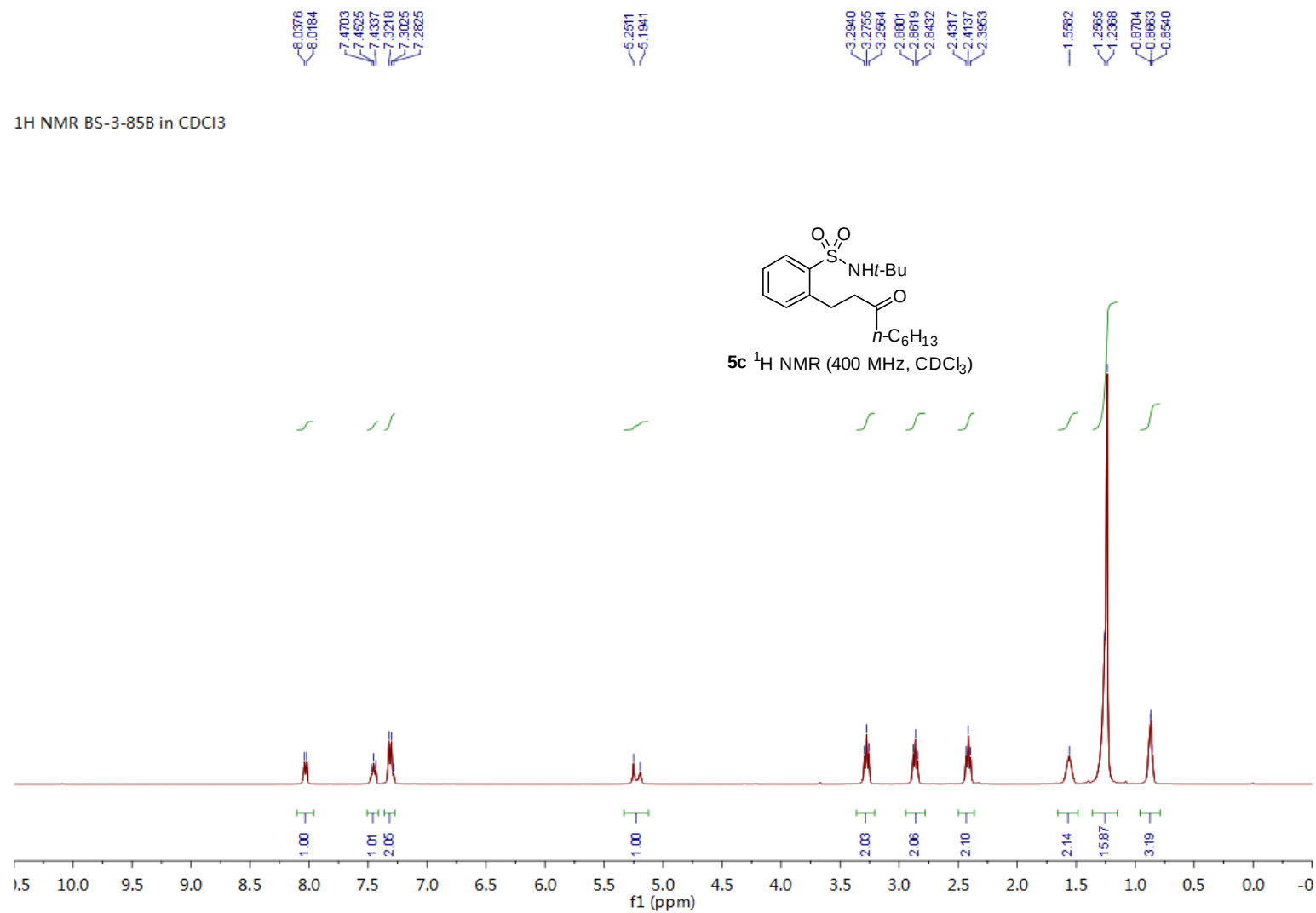


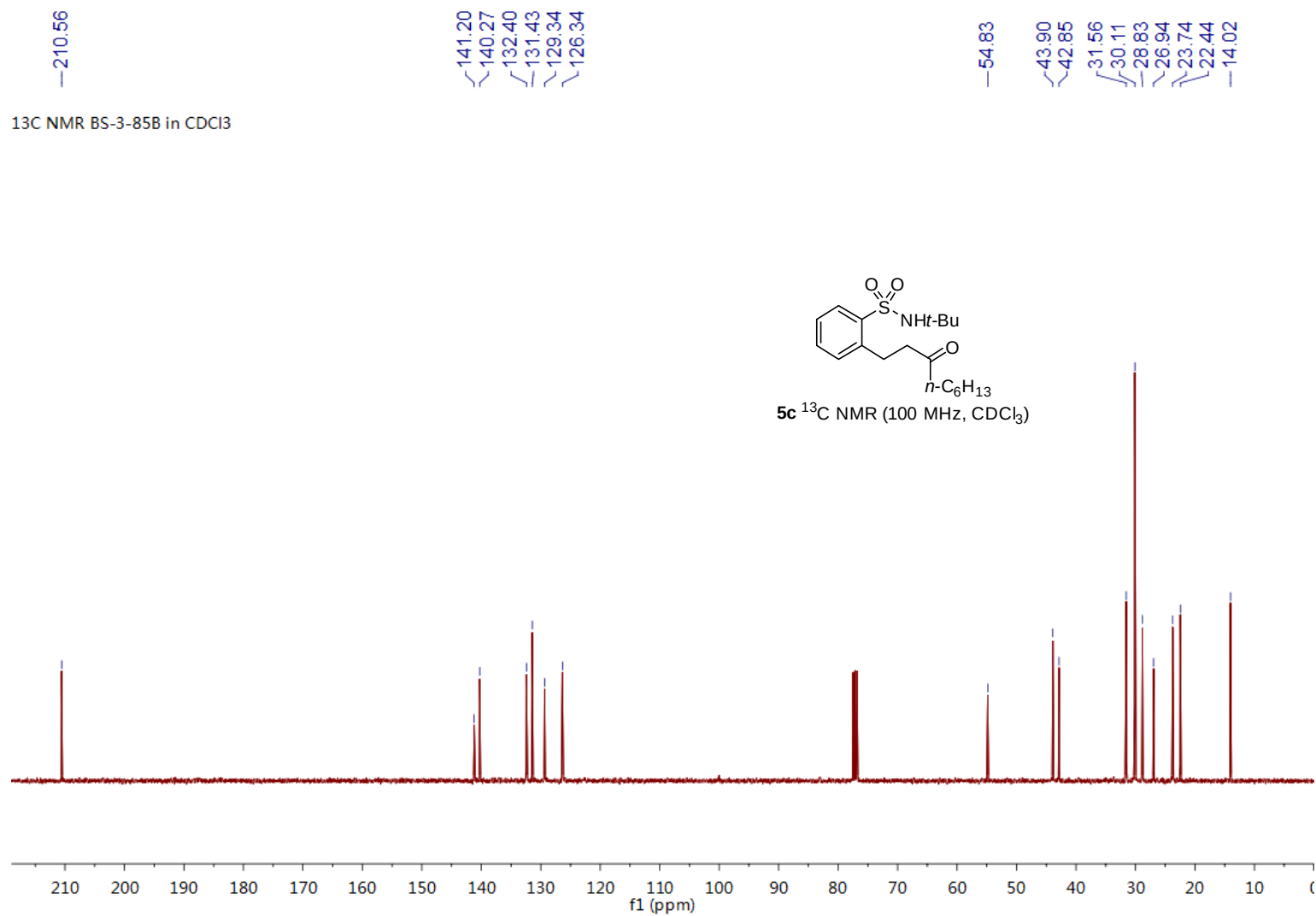


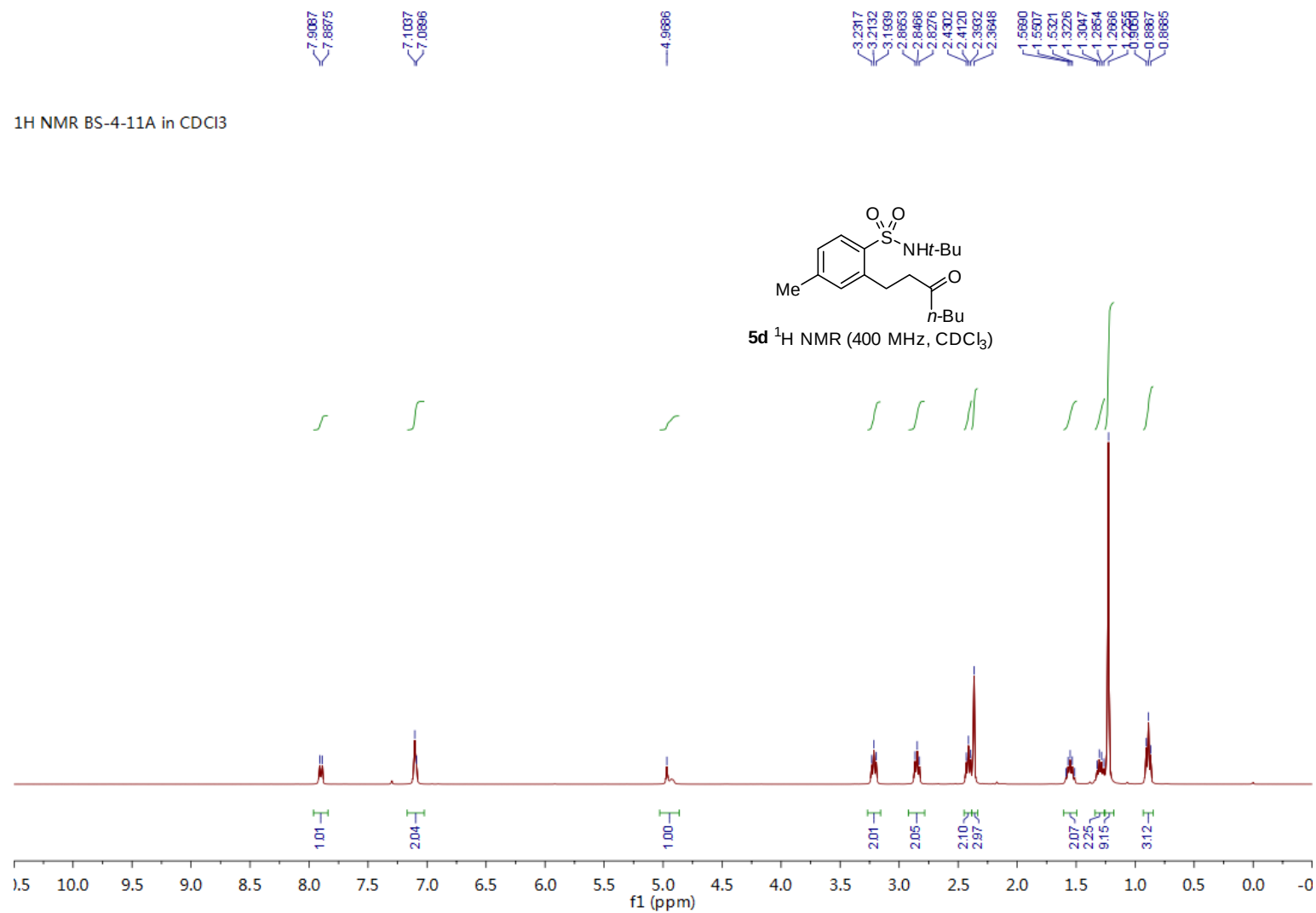


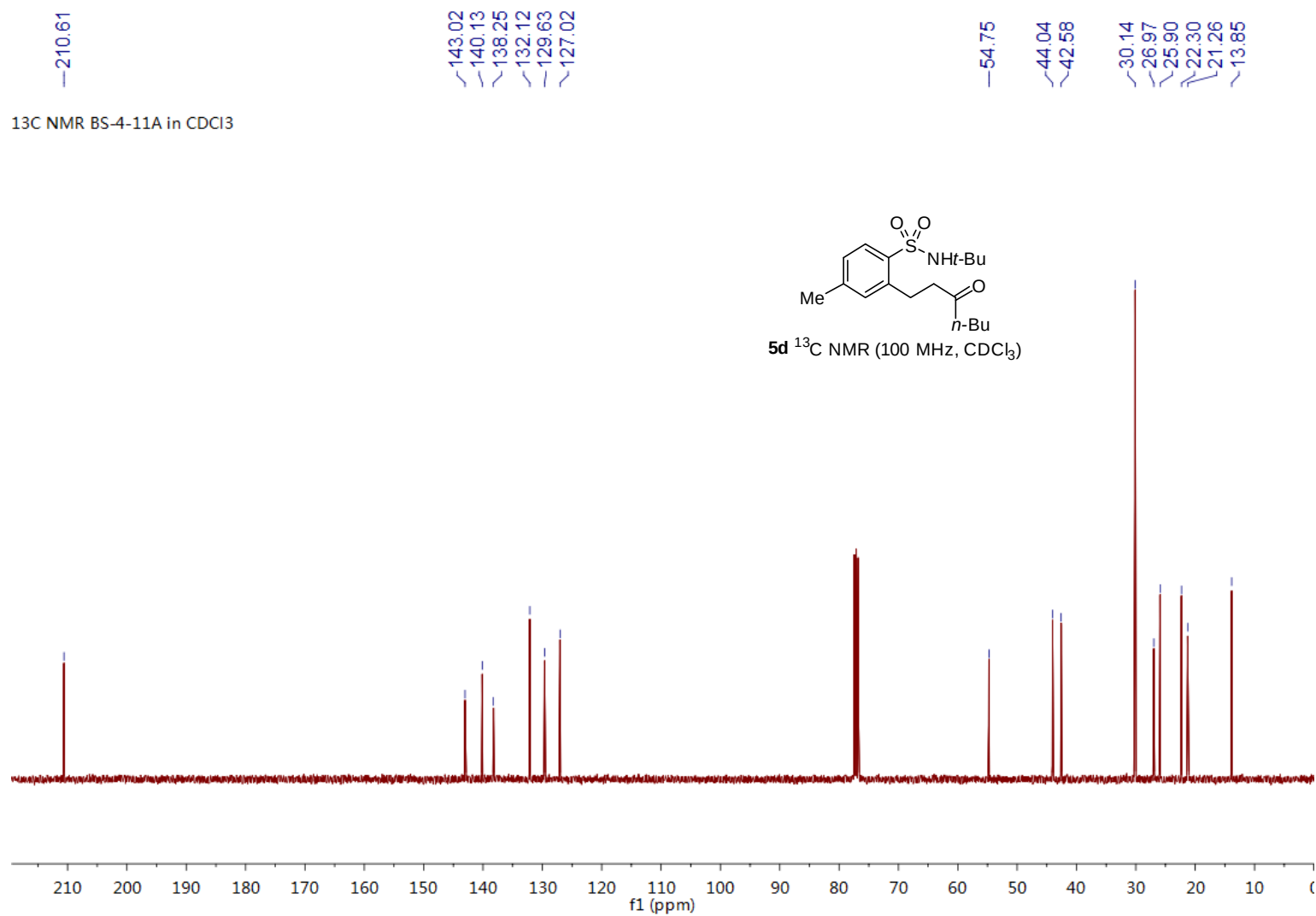


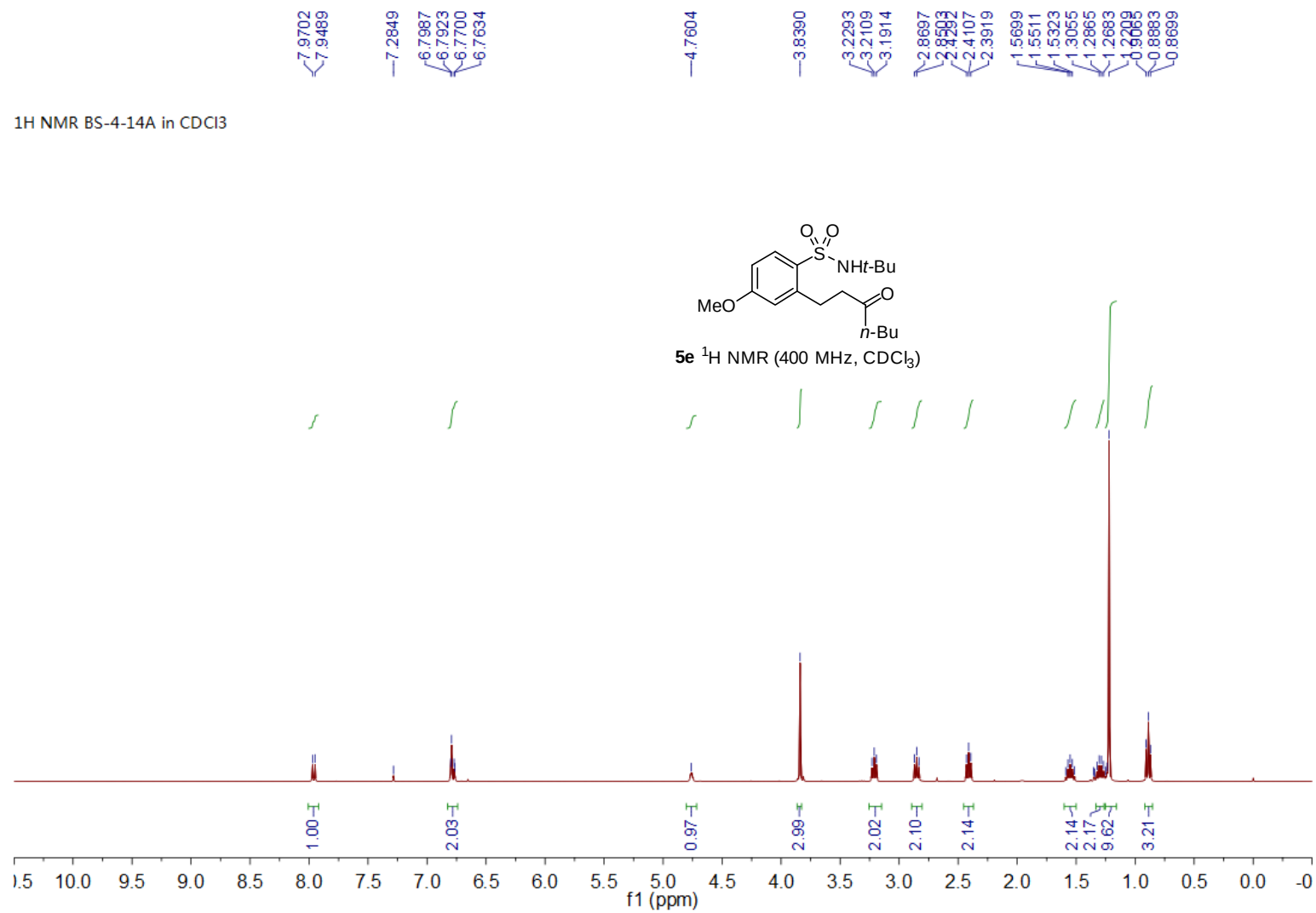


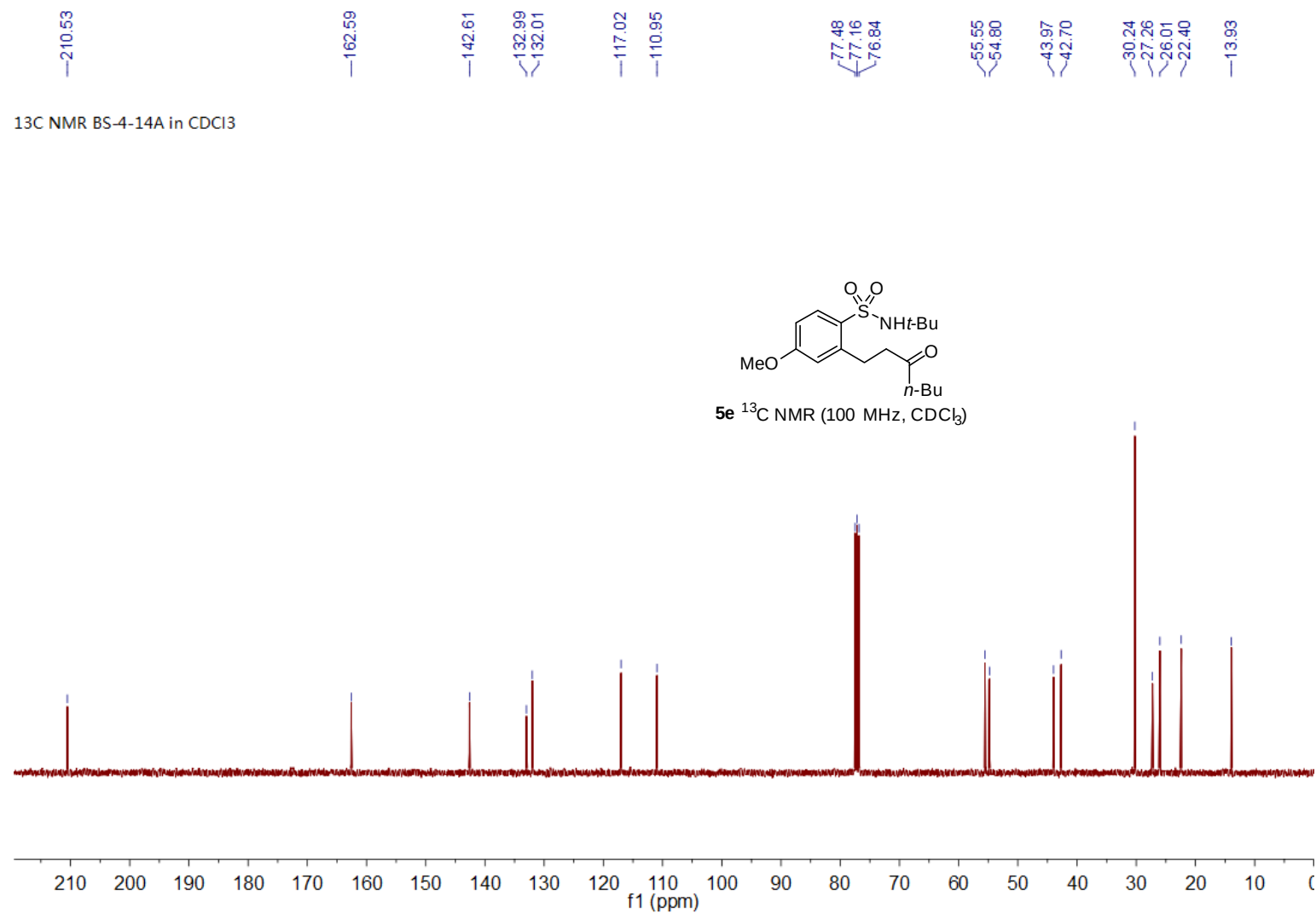


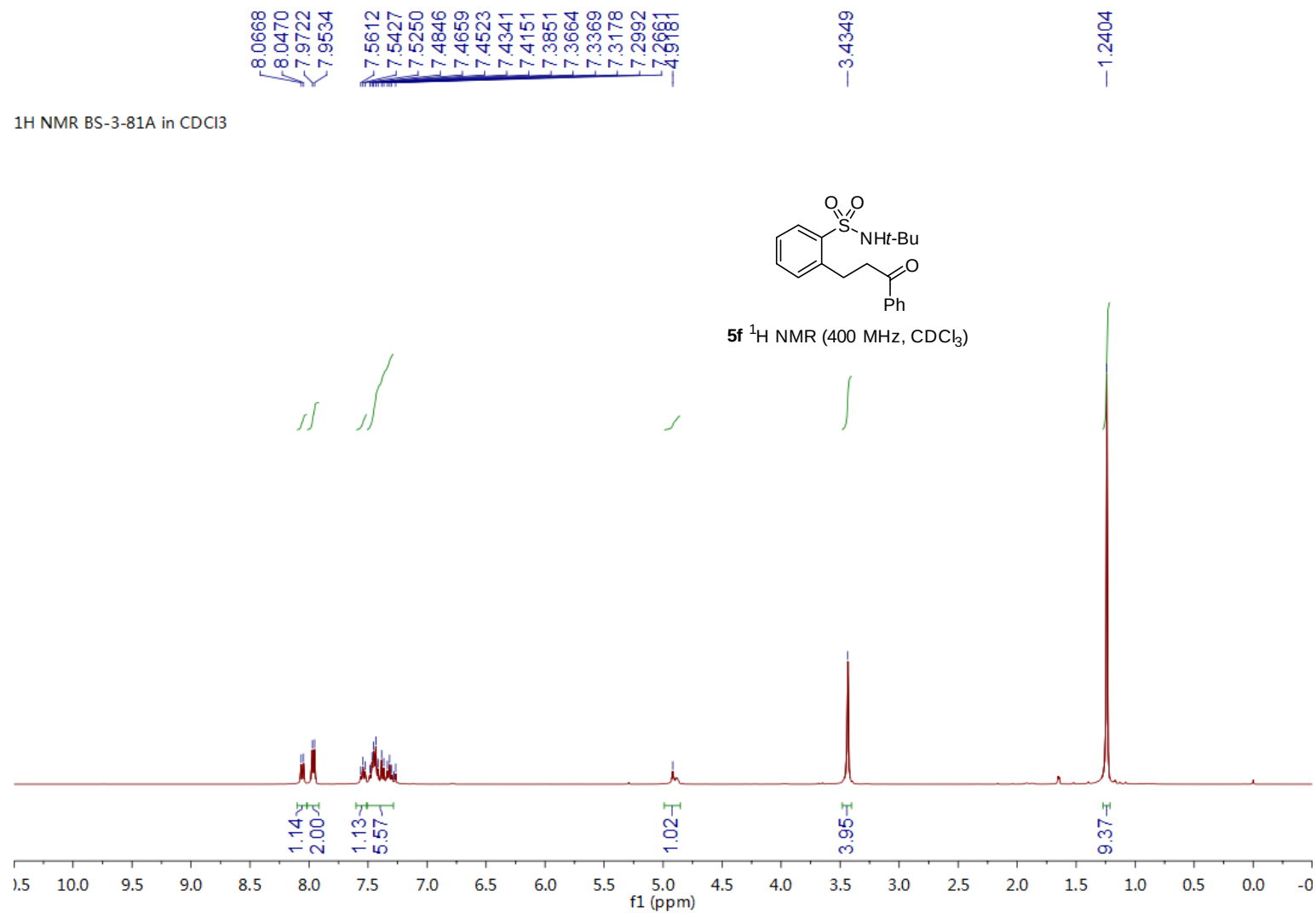


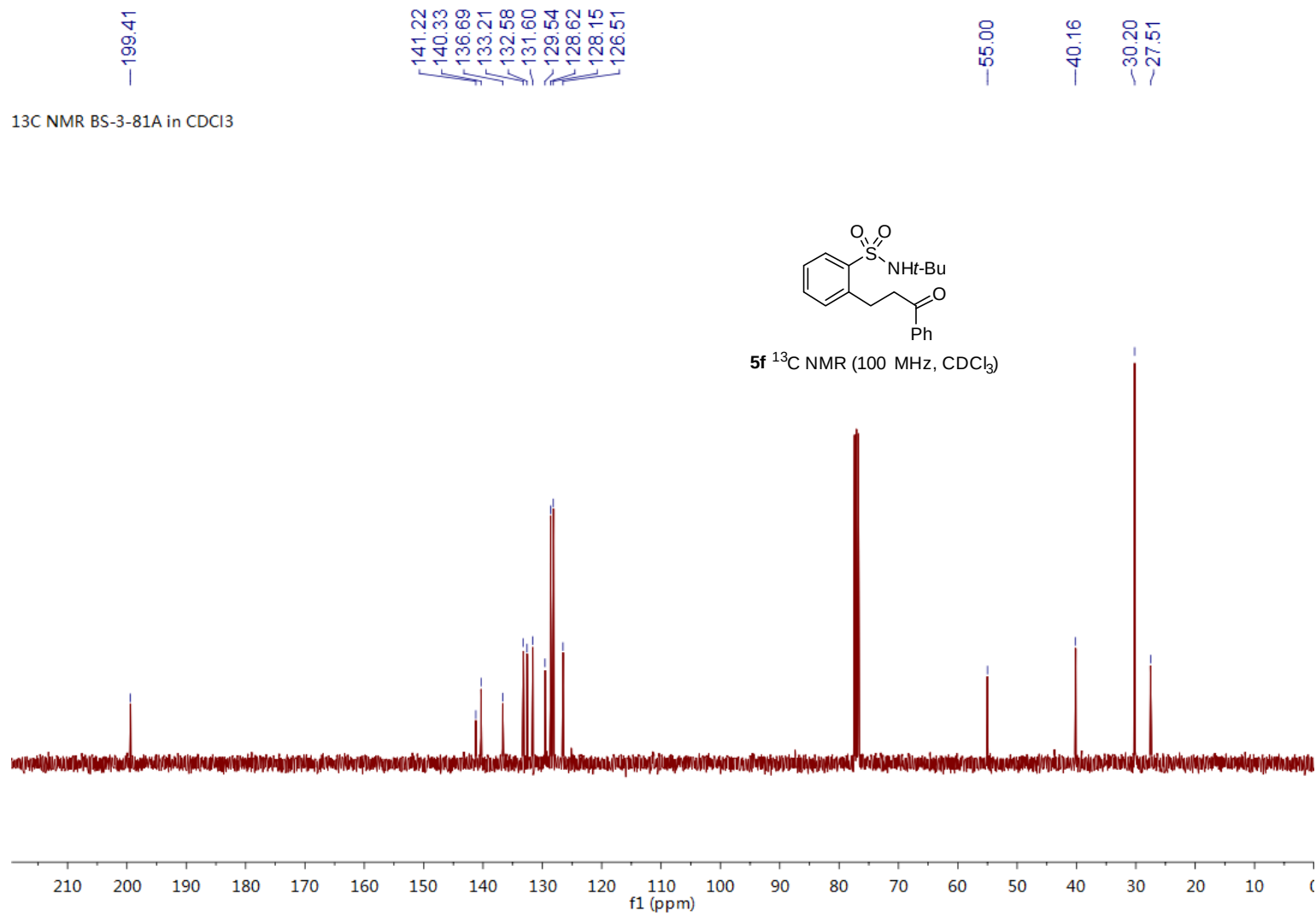


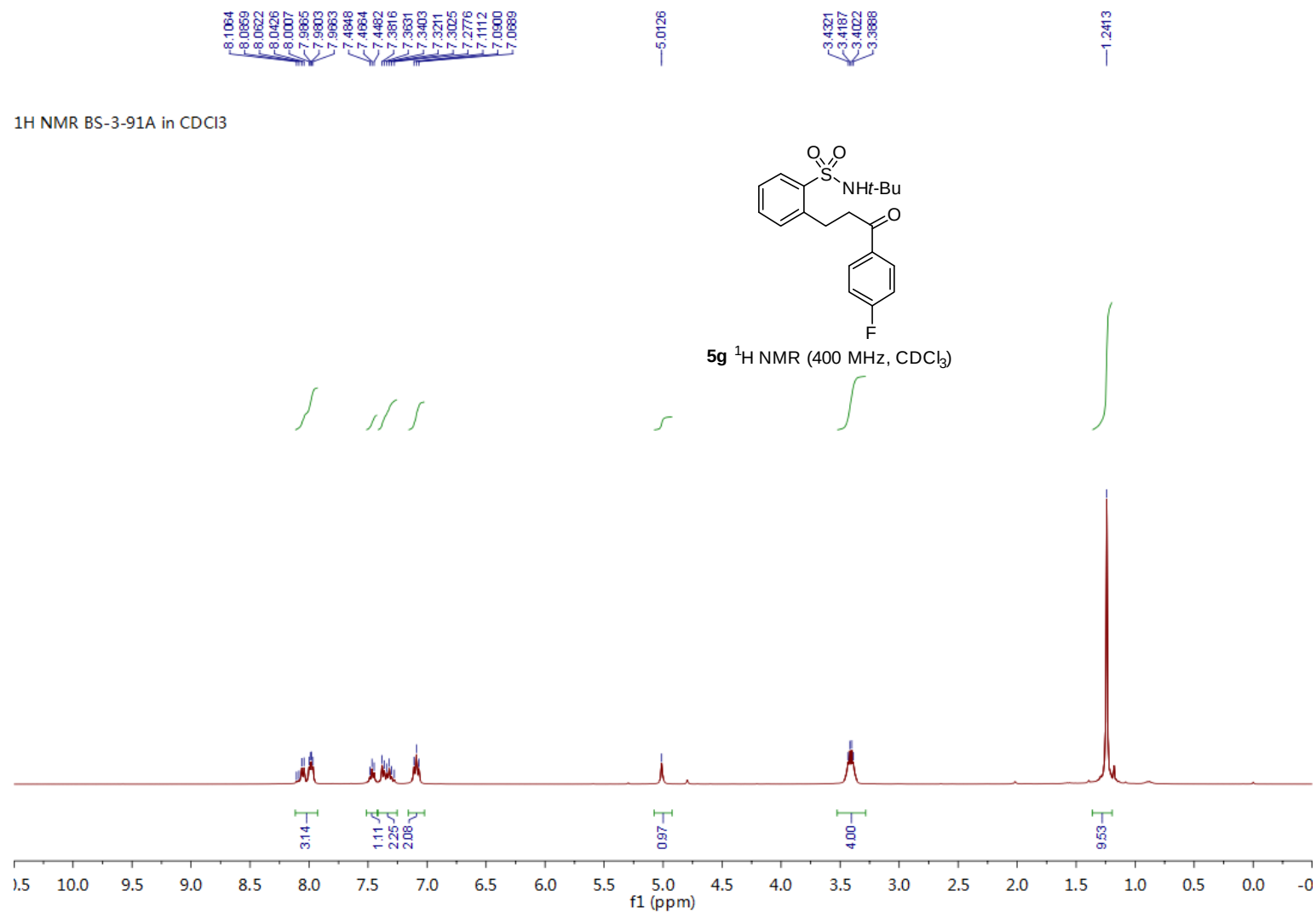






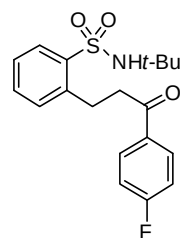




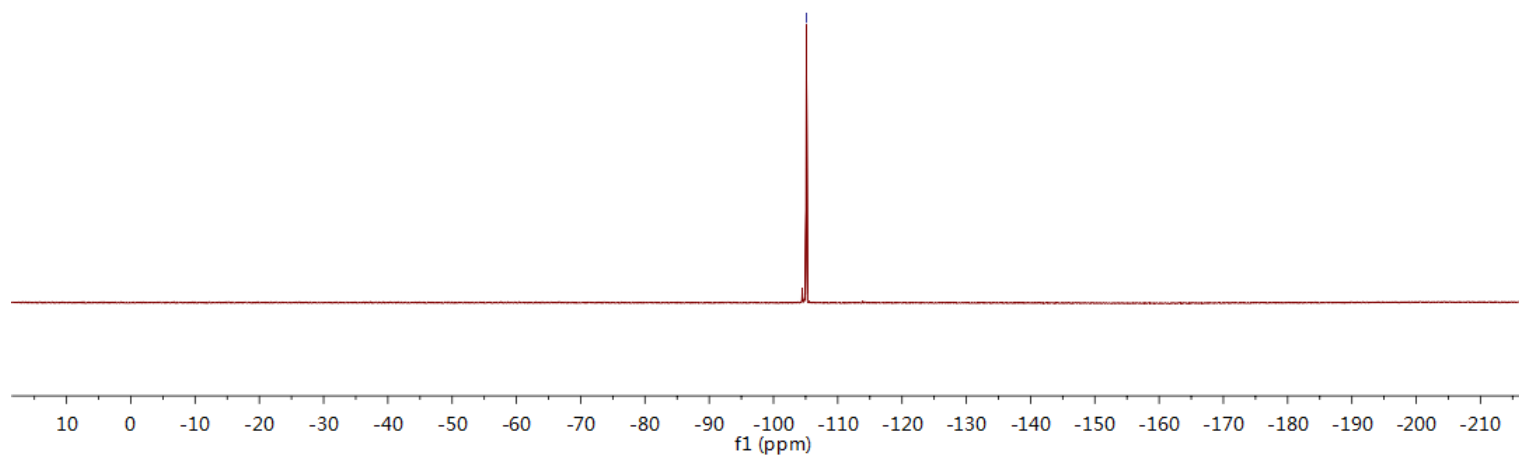


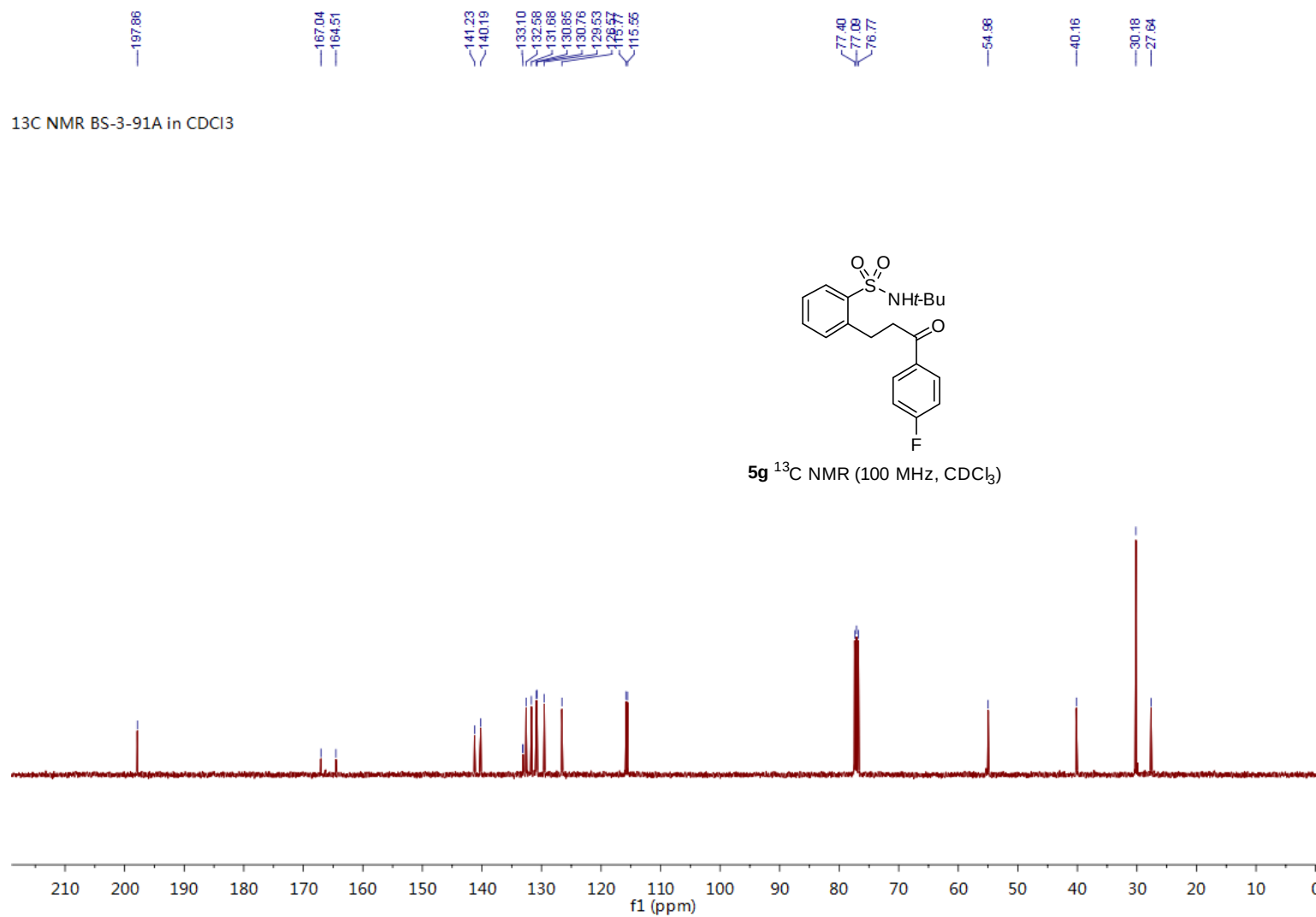
¹⁹F NMR BS-3-91 in CDCl₃

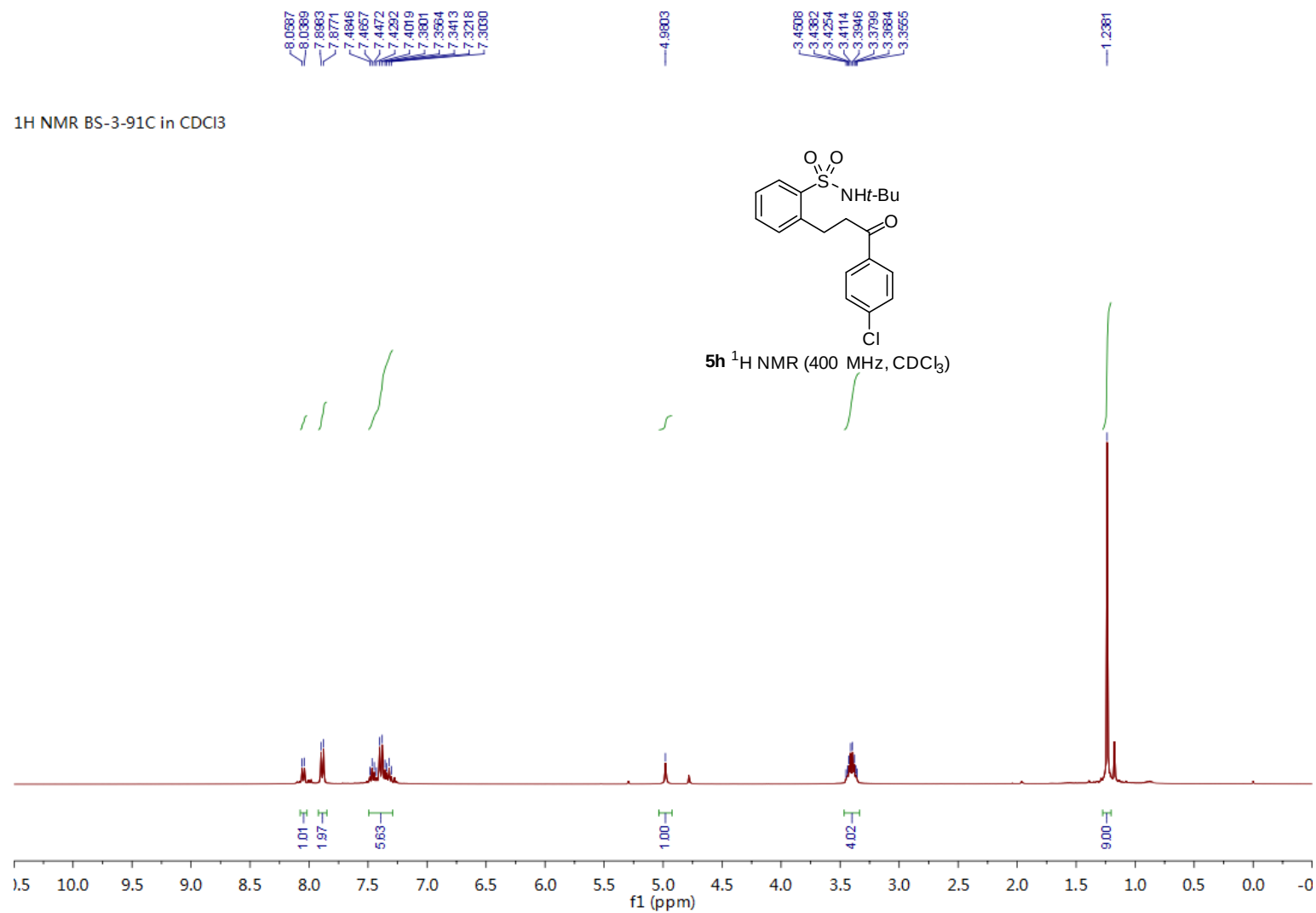
-105.1080

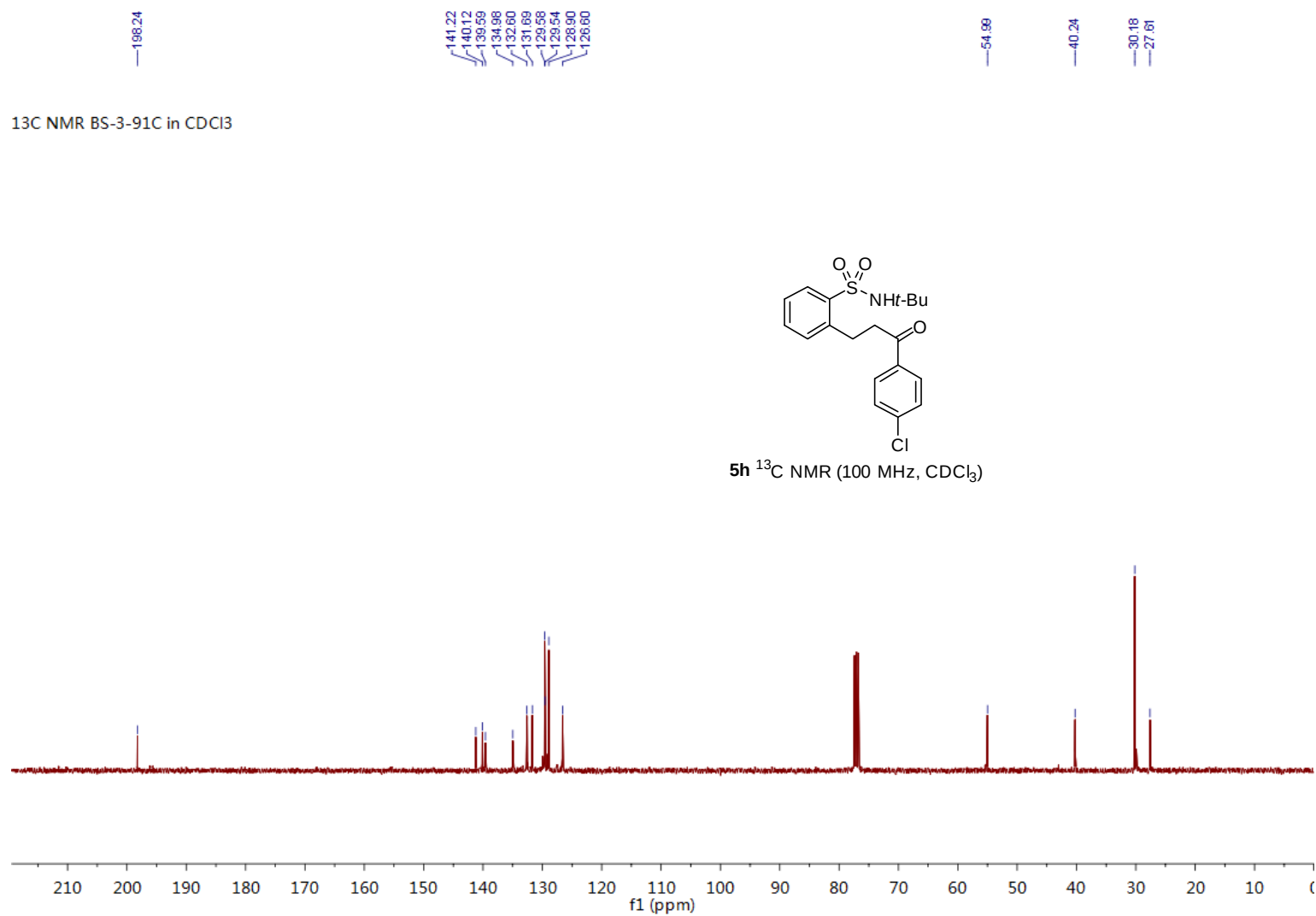


5g ¹⁹F NMR (376 MHz, CDCl₃)

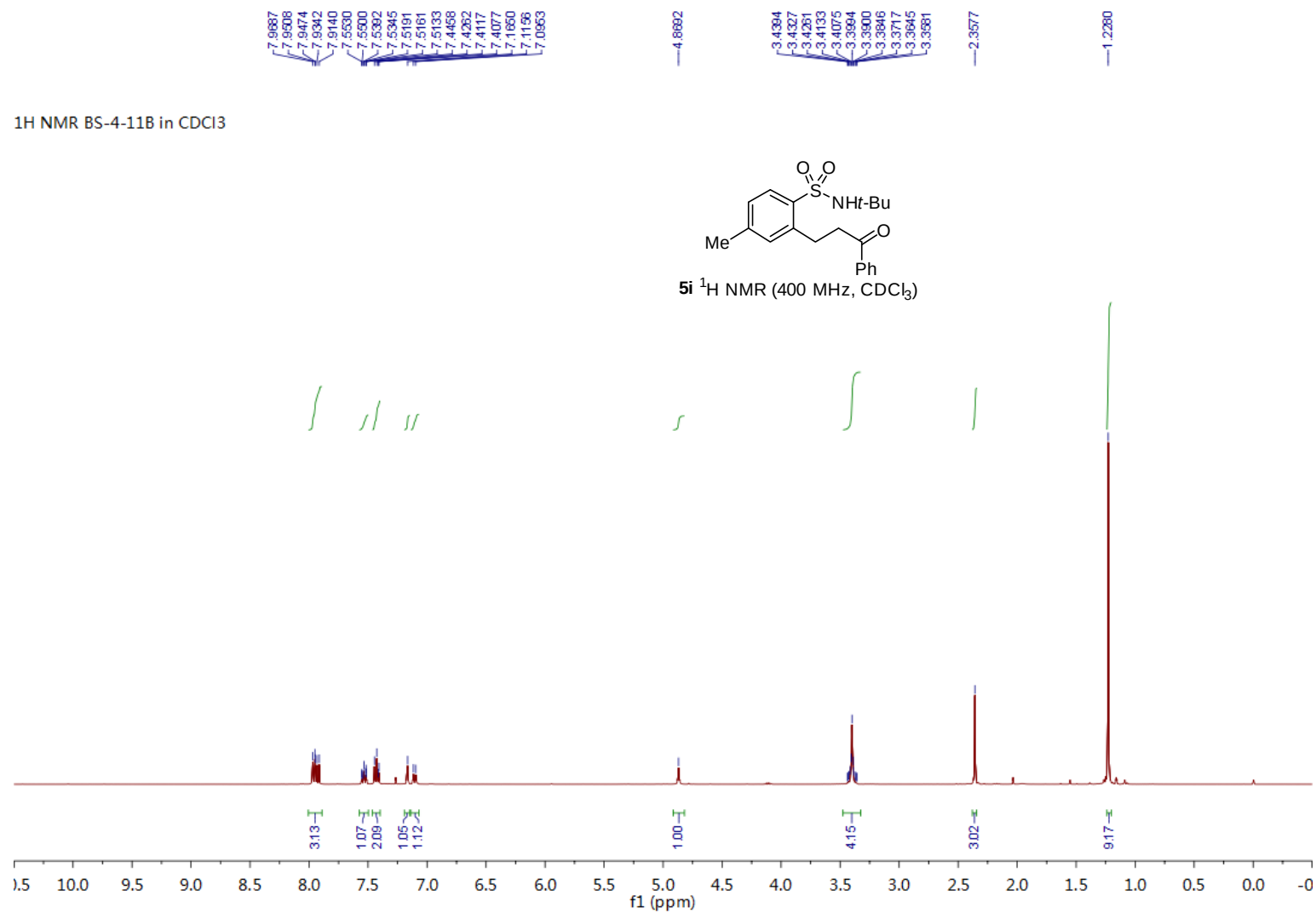
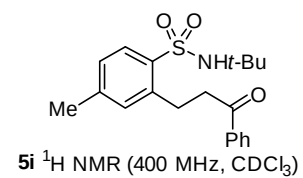


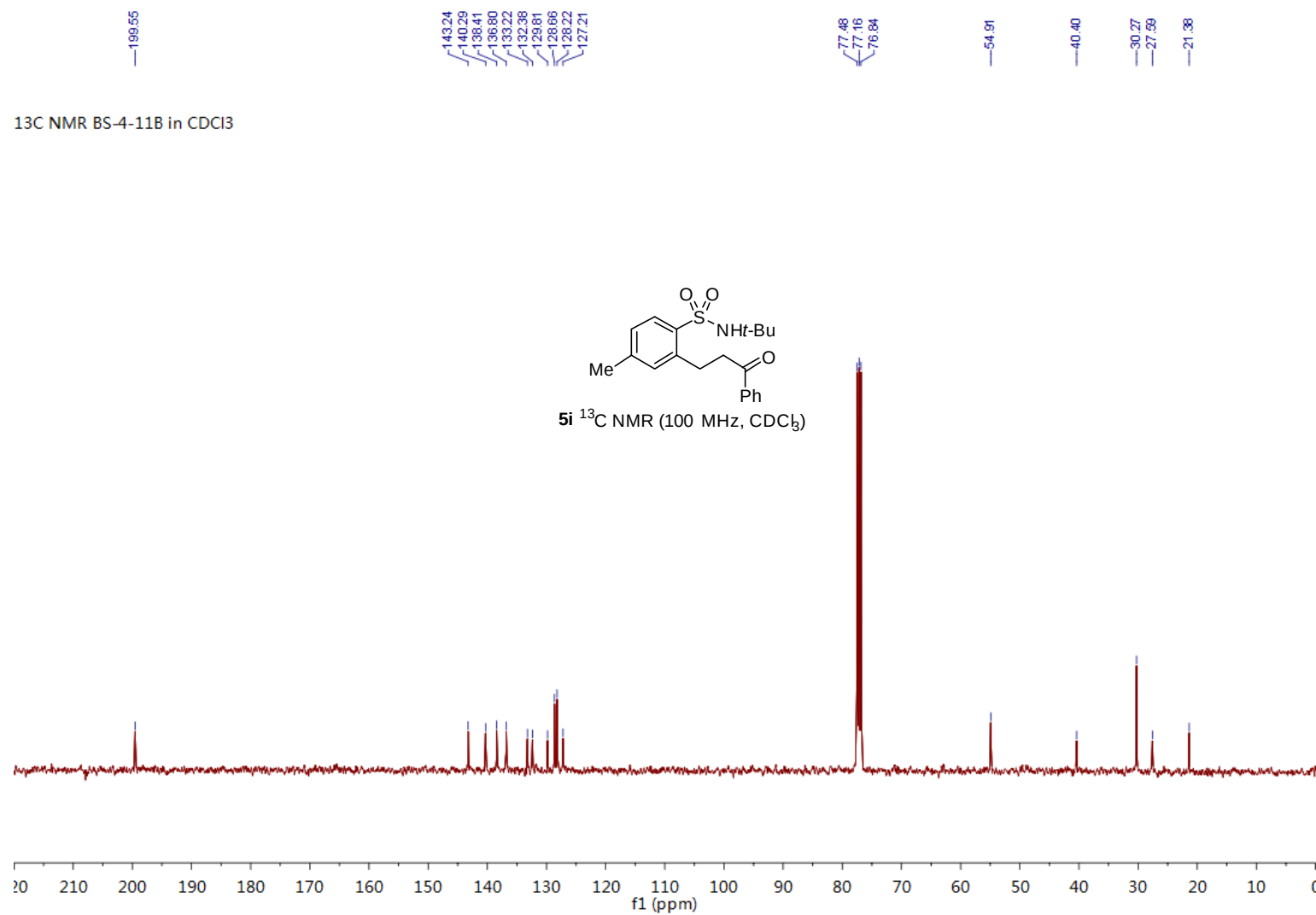




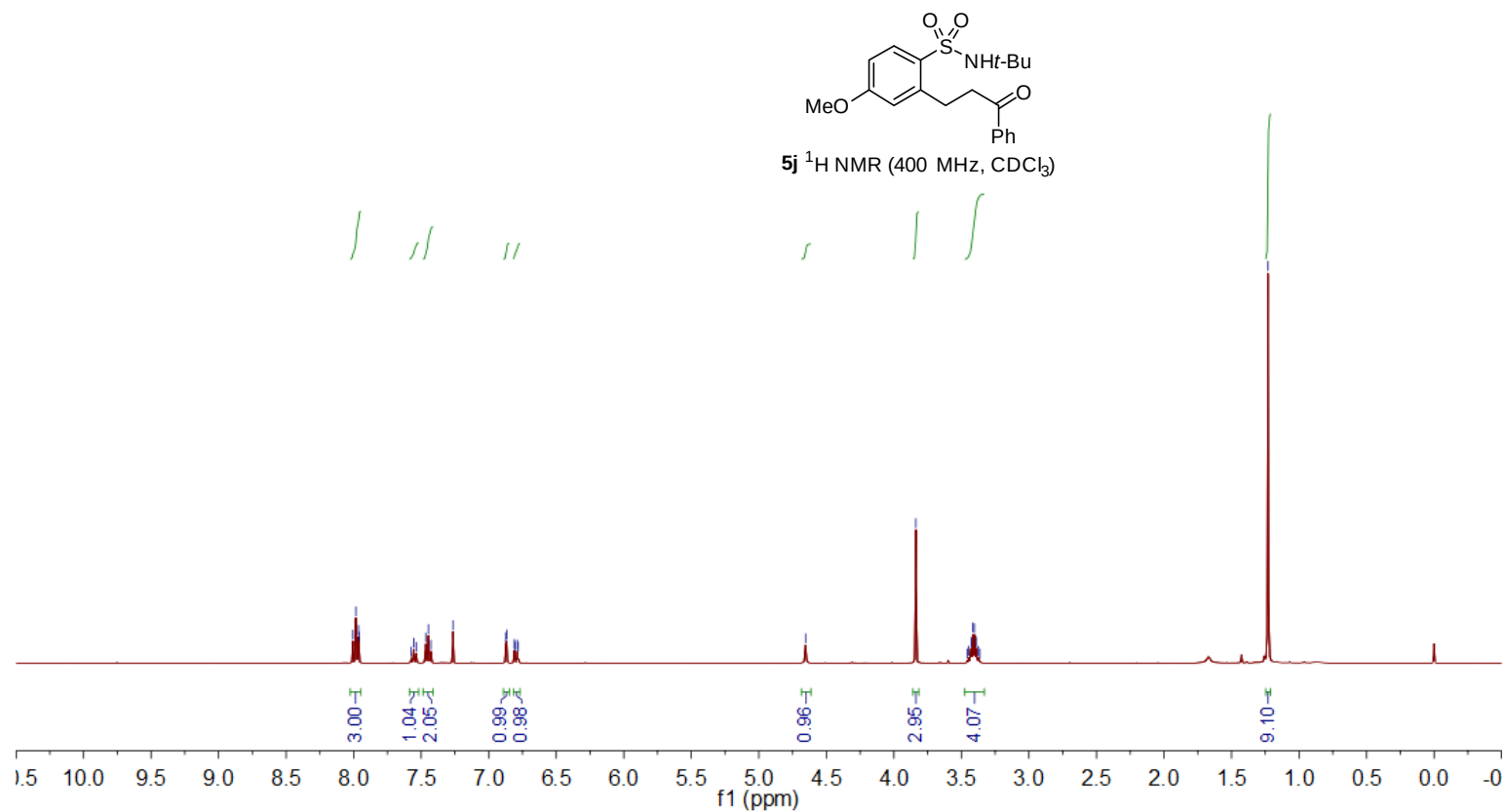


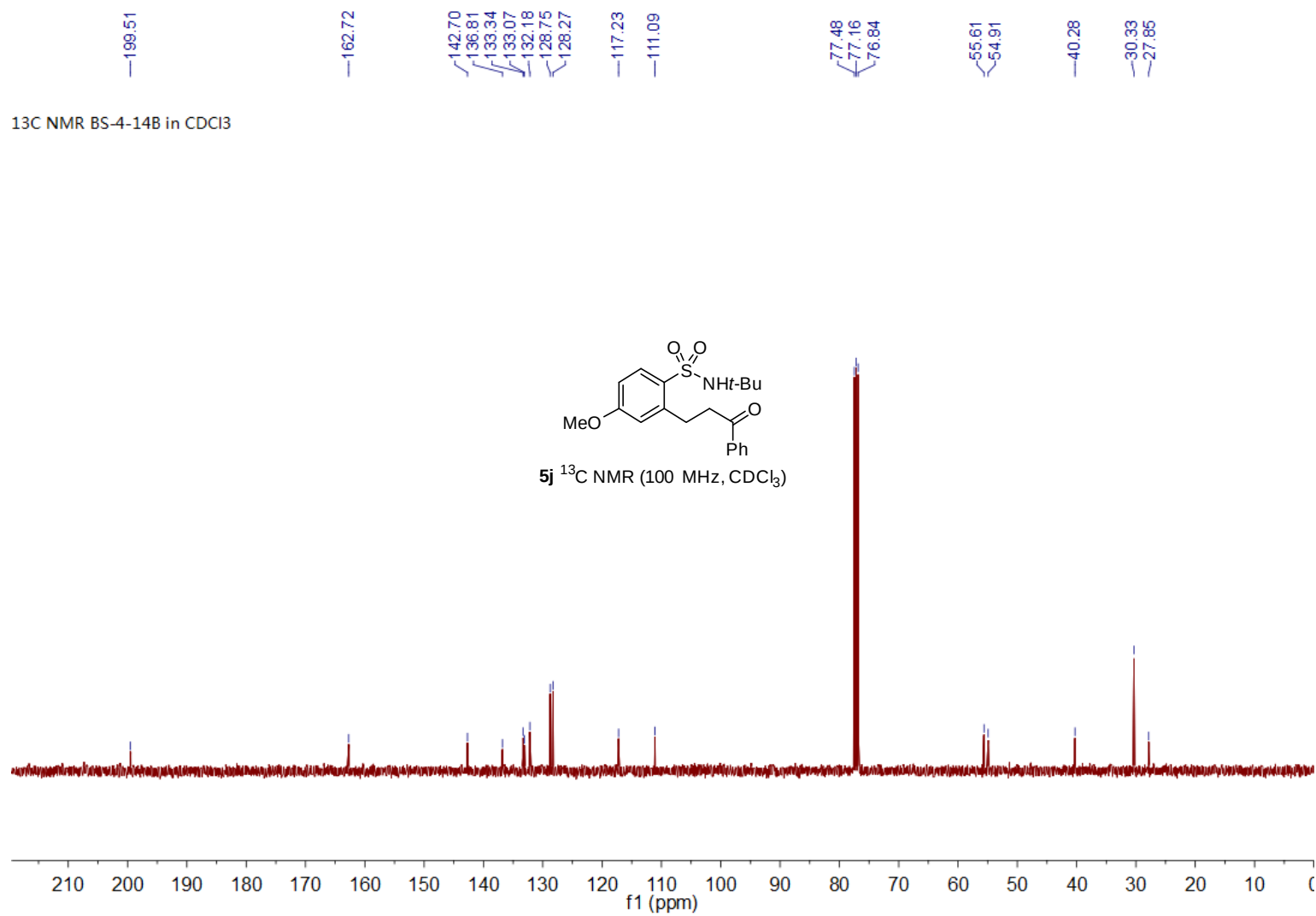
¹H NMR BS-4-11B in CDCl₃

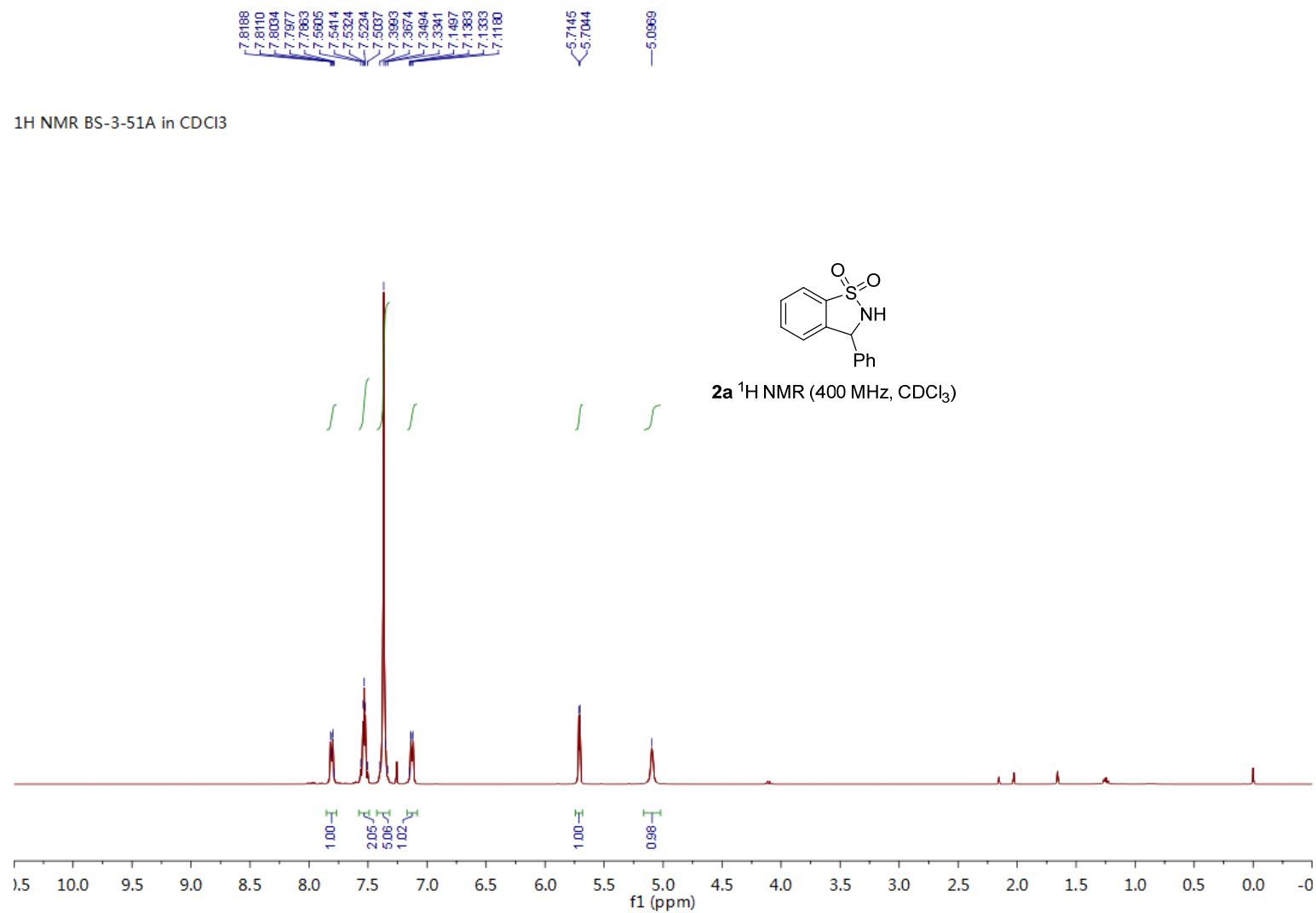




¹H NMR BS-4-14B in CDCl₃

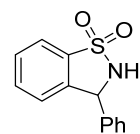




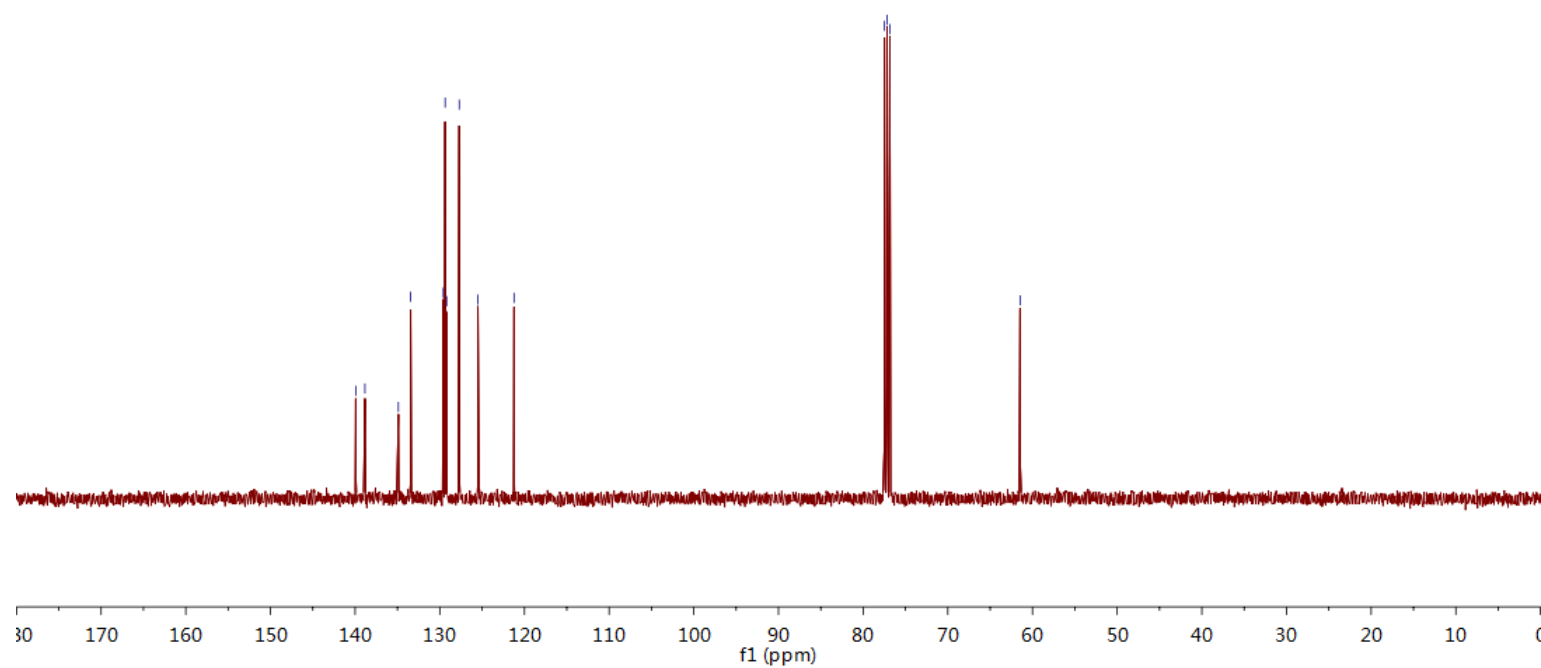


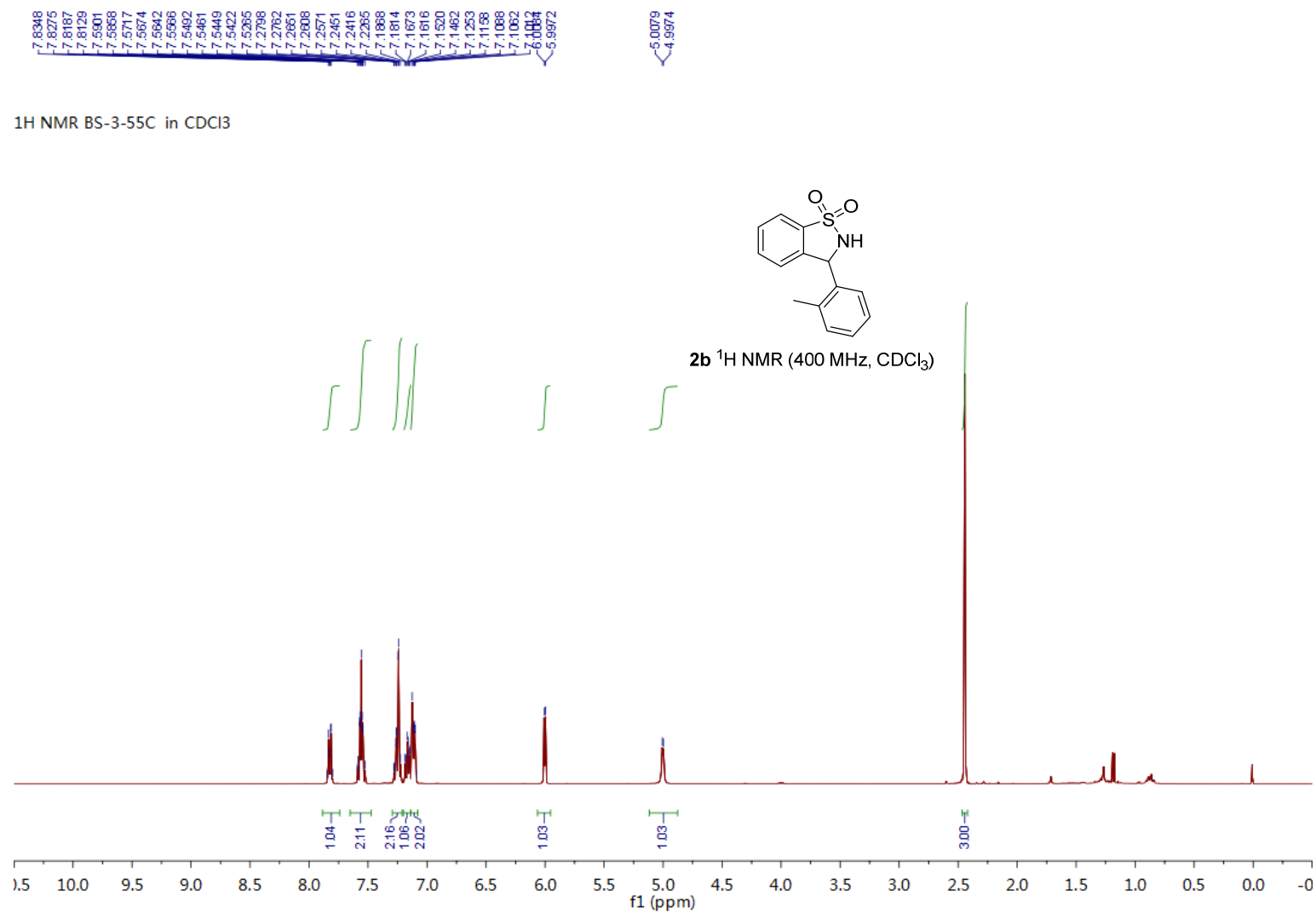


¹³C NMR BS-3-51A in CDCl₃



2a ¹³C NMR (100 MHz, CDCl₃)





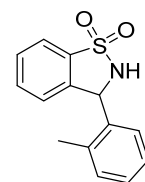
140.15
136.80
136.34
135.64
133.40
131.30
129.50
128.10
128.20
127.09
125.29
121.32

77.48
77.16
76.84

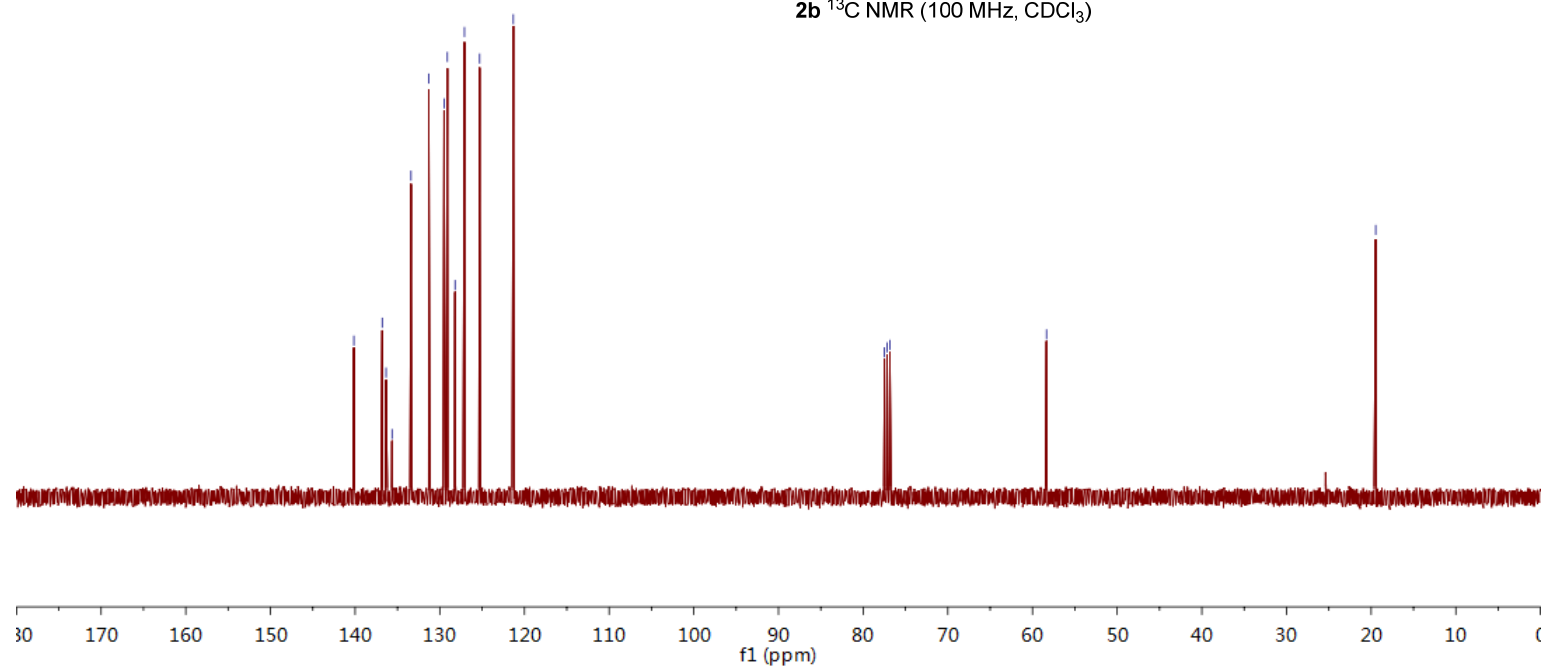
58.35

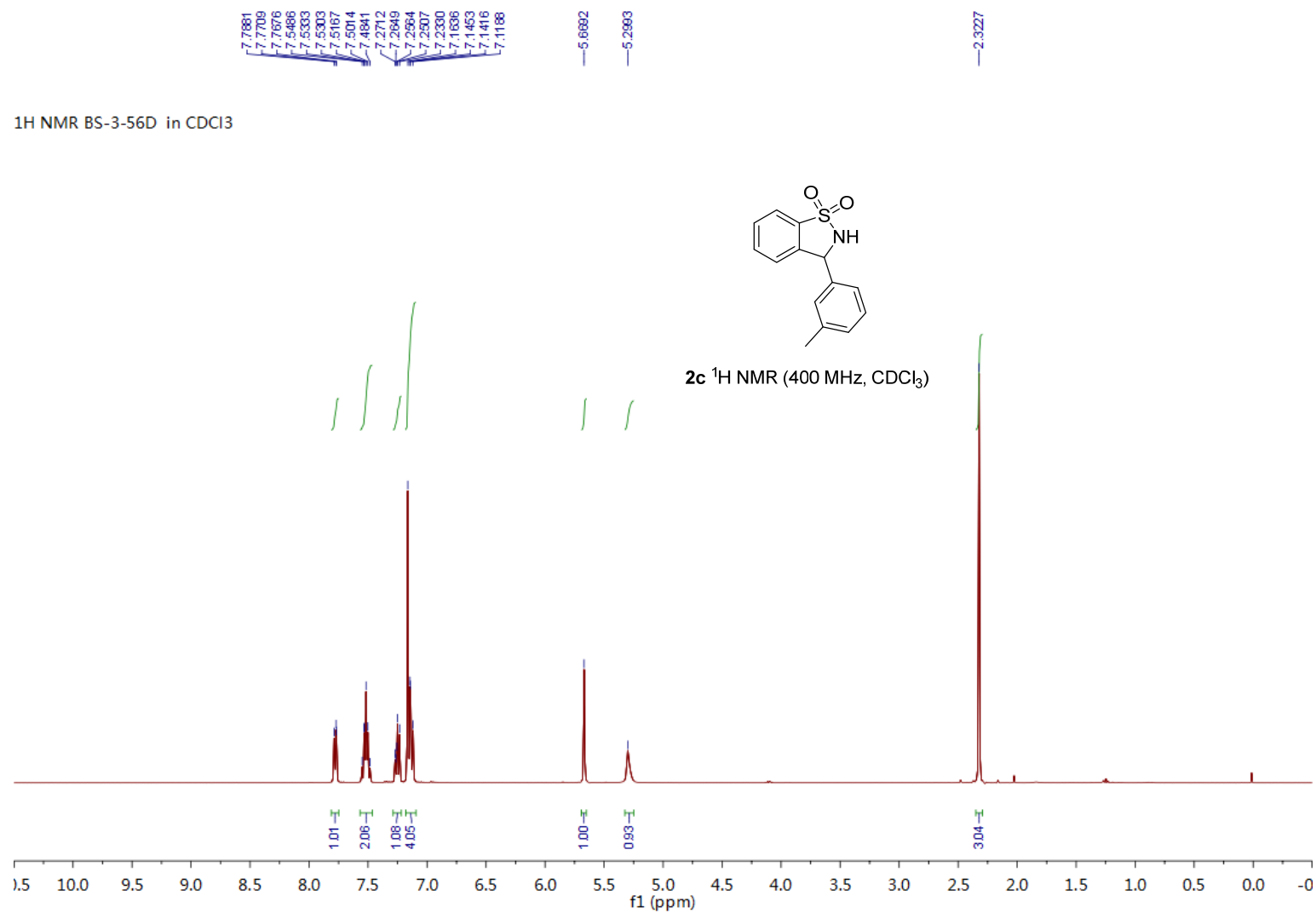
19.46

¹³C NMR BS-3-55C in CDCl₃



2b ¹³C NMR (100 MHz, CDCl₃)





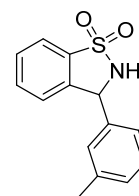
139.97
139.11
138.76
134.86
133.32
129.78
129.41
129.07
128.12
125.42
124.73
121.05

77.48
77.16
76.84

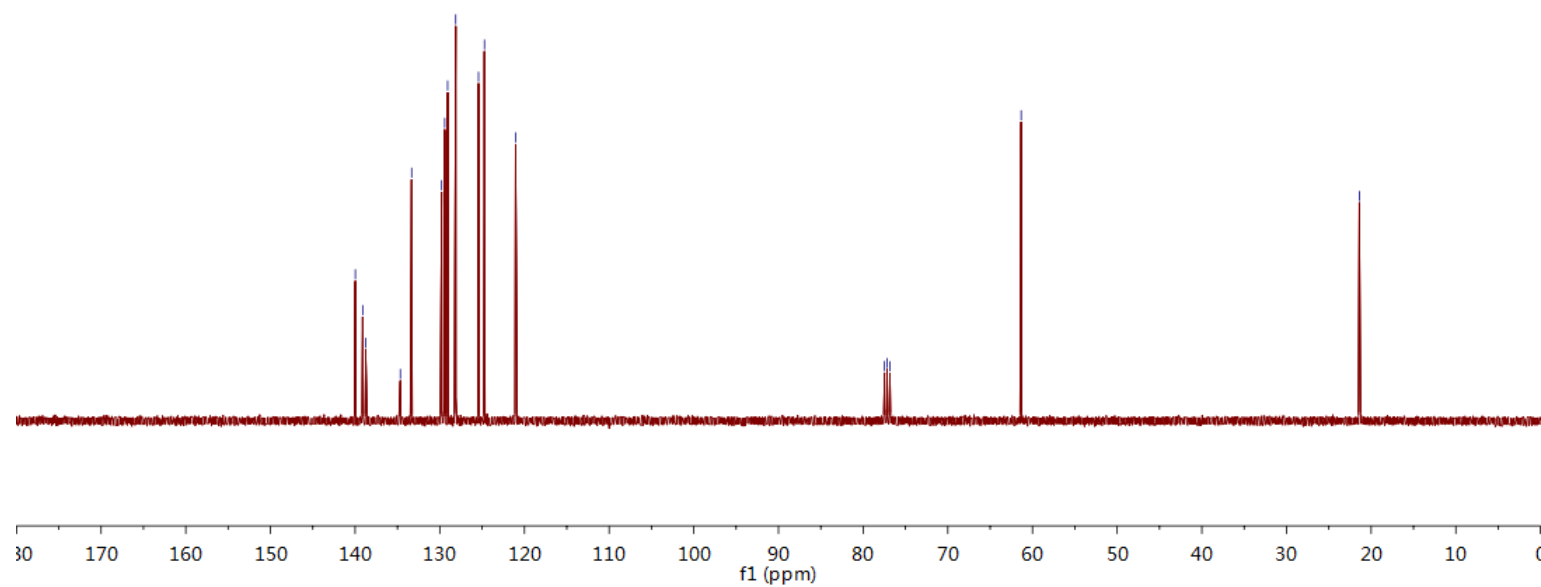
61.34

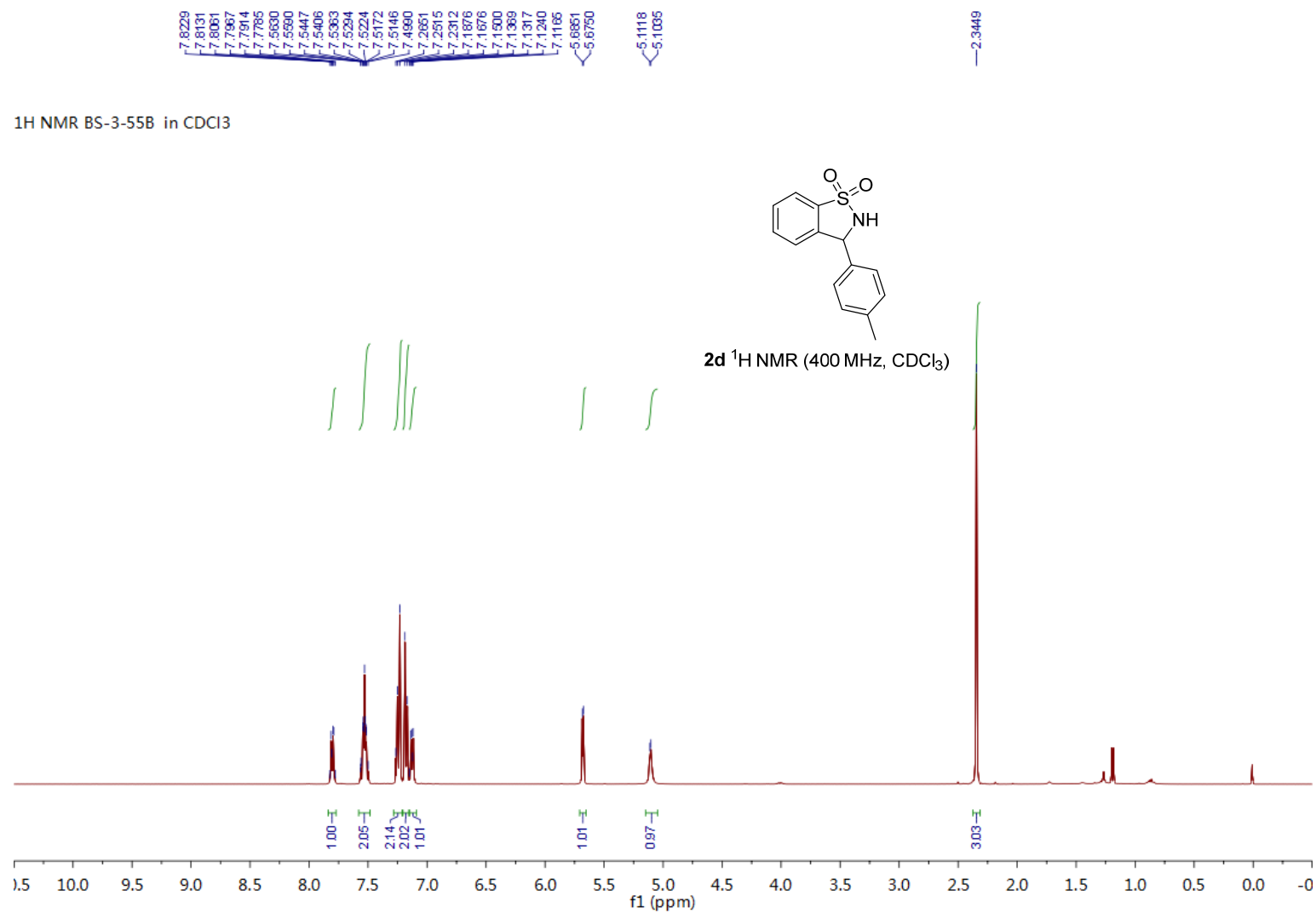
21.40

¹³C NMR BS-3-56D in CDCl₃



2c ¹³C NMR (100 MHz, CDCl₃)





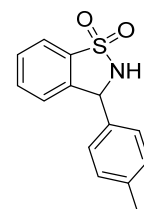
¹³C NMR BS-3-55B in CDCl₃

140.19
139.11
135.84
134.91
133.37
129.98
129.47
127.62
125.46
121.15

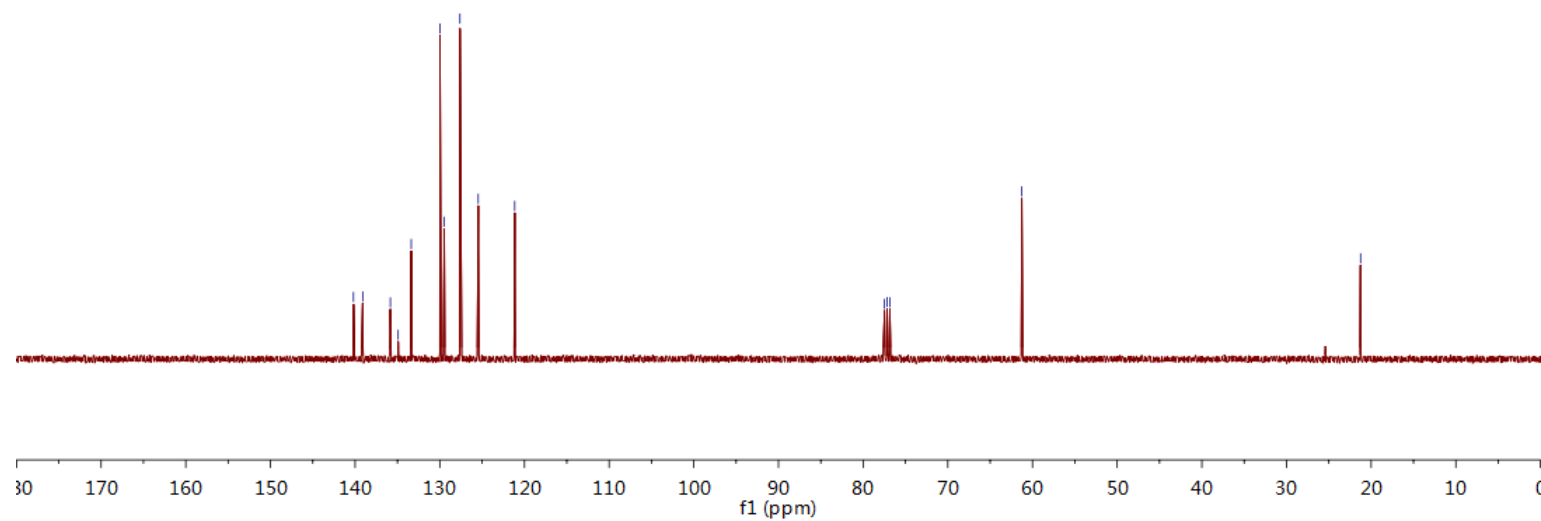
77.48
77.16
76.84

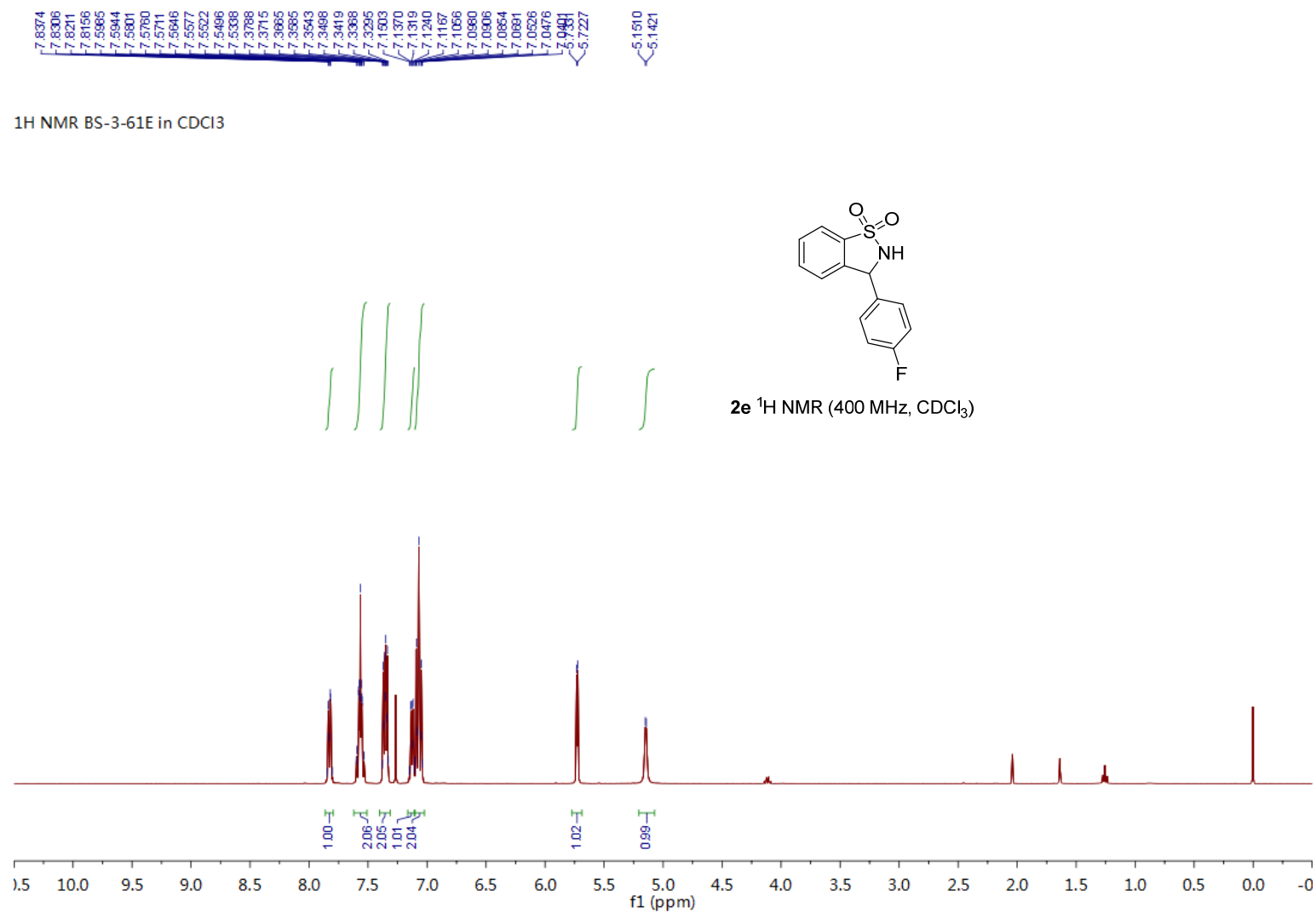
61.26

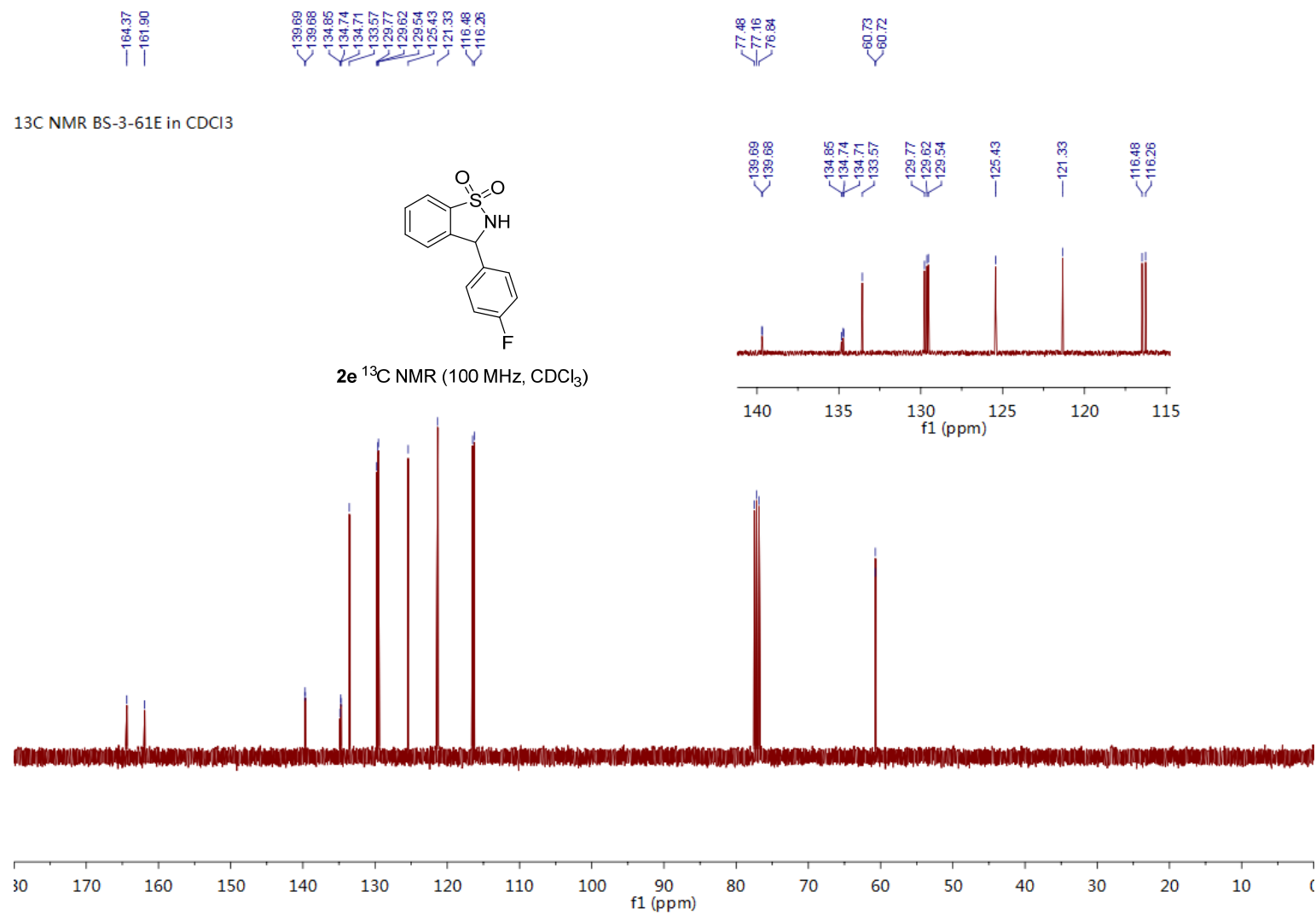
21.26



2d ¹³C NMR (100 MHz, CDCl₃)

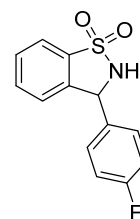




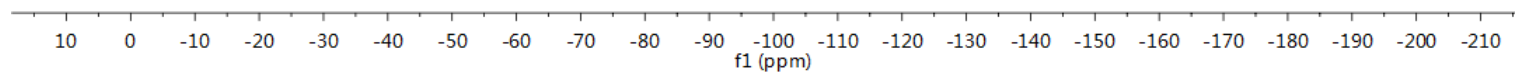


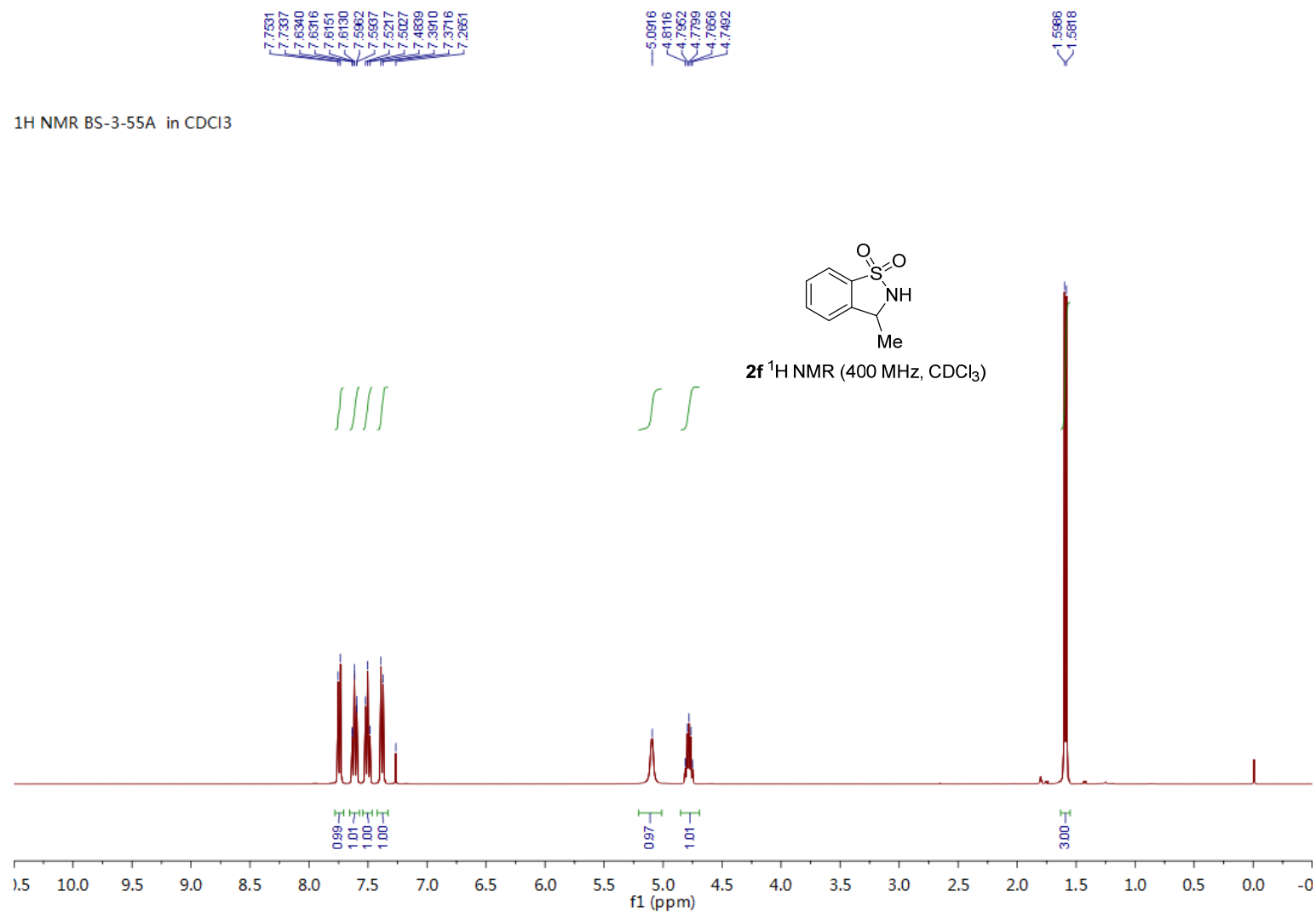
¹⁹F NMR BS-3-61E in CDCl₃

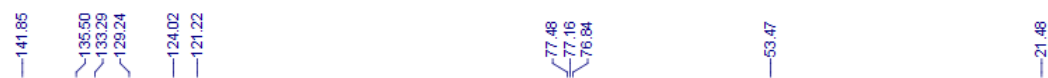
—112.2089



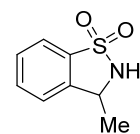
2e ¹⁹F NMR (376 MHz, CDCl₃)



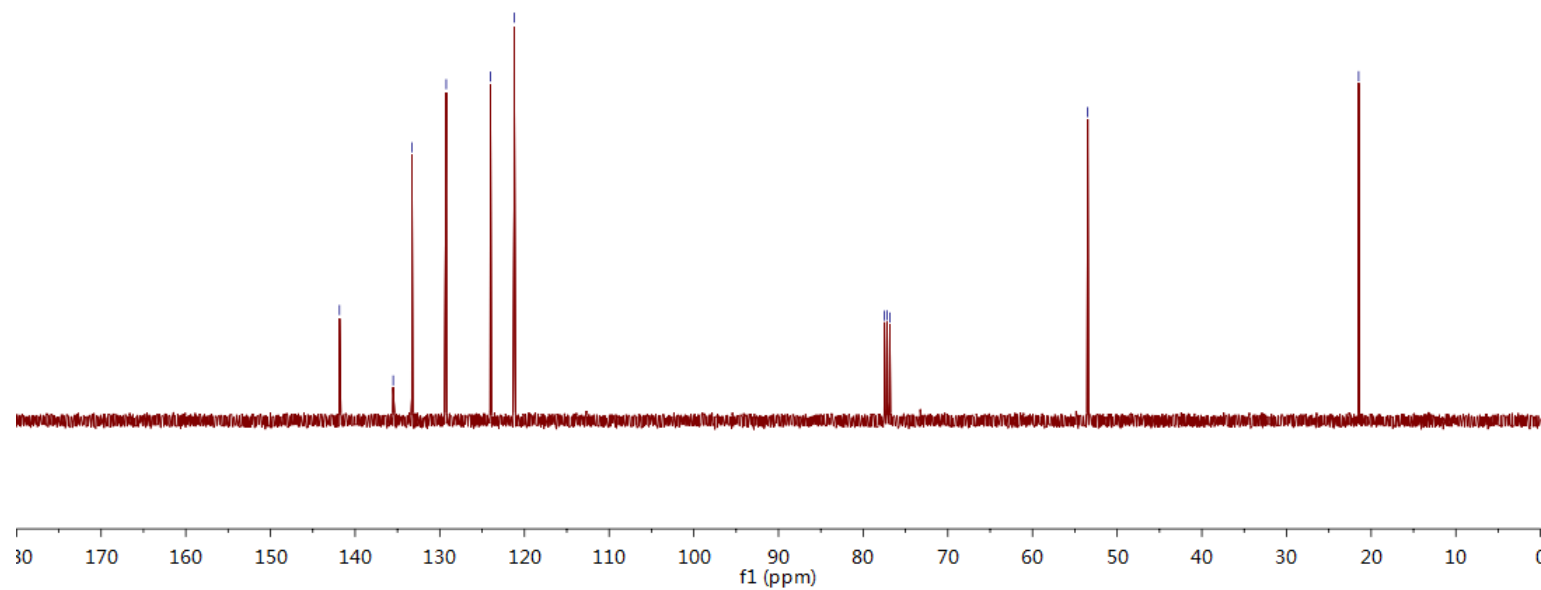


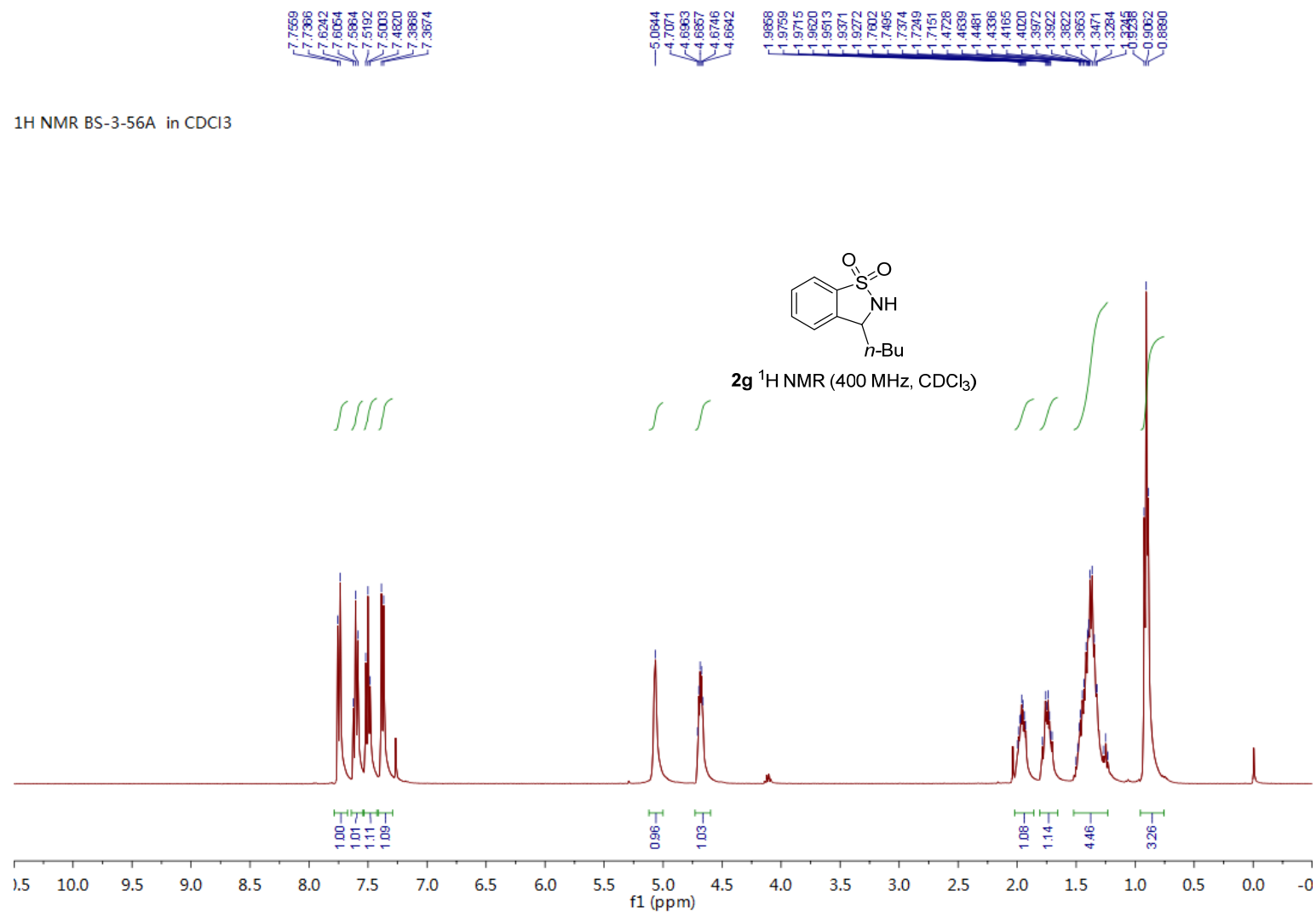


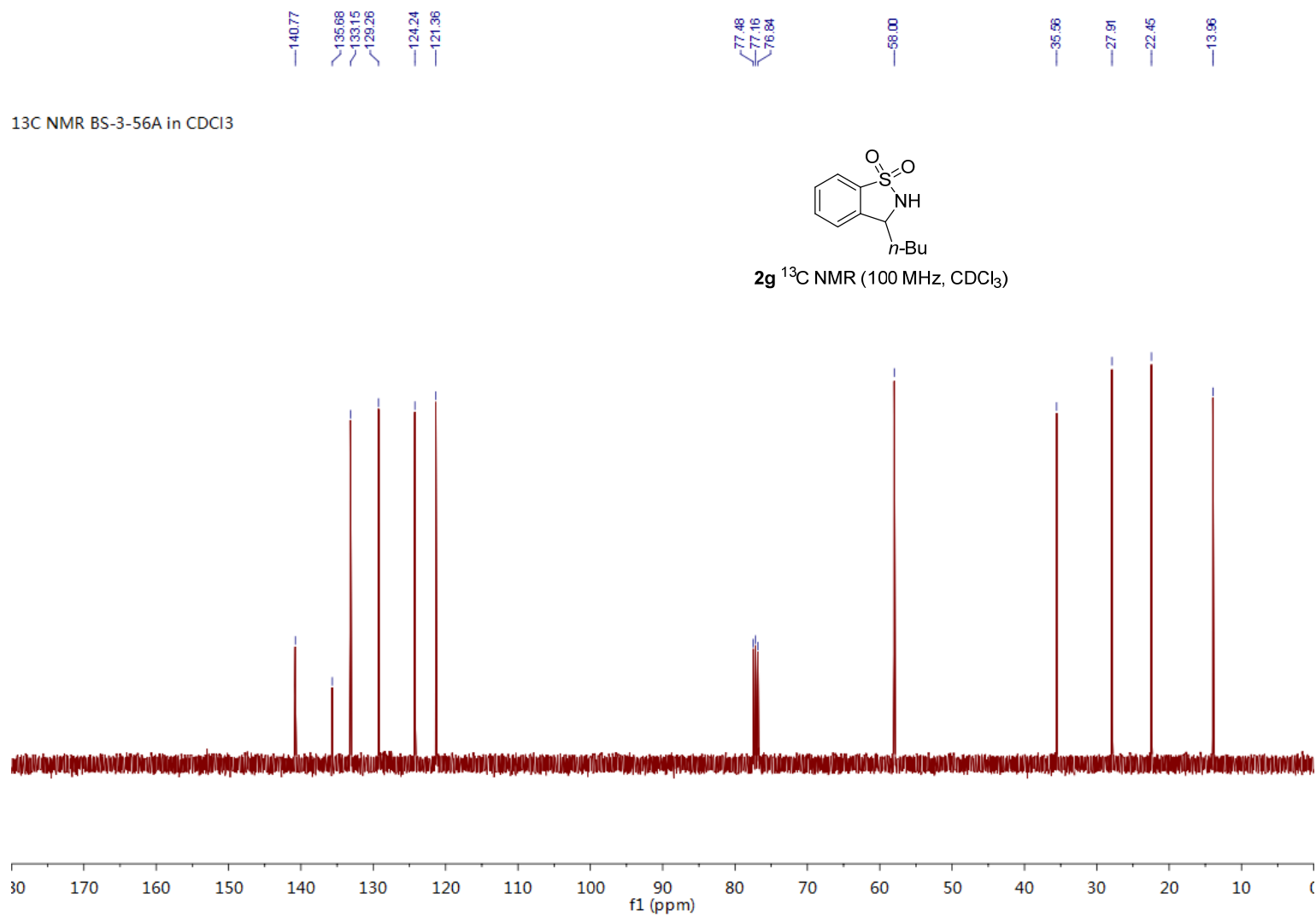
¹³C NMR BS-3-55A in CDCl₃

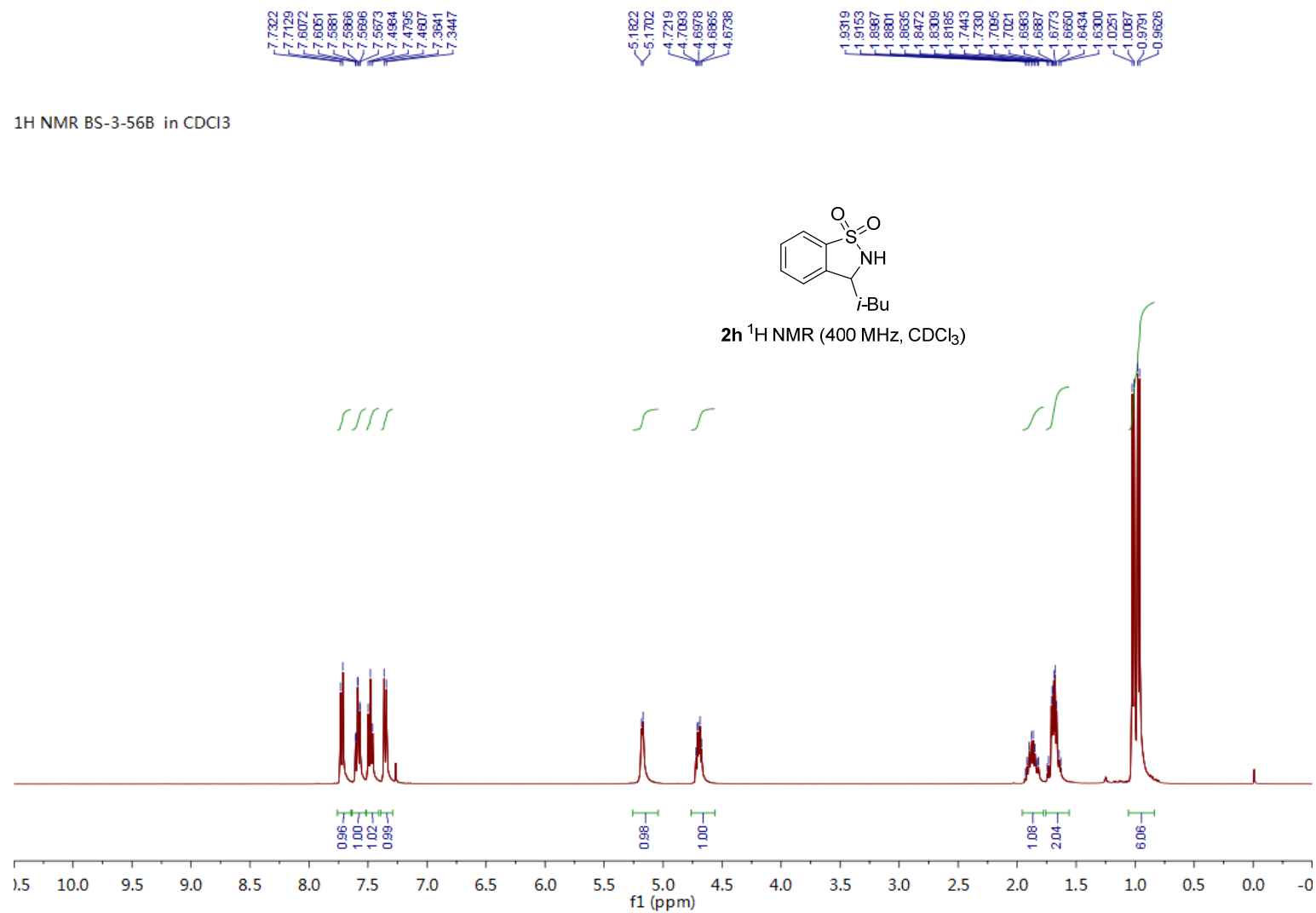


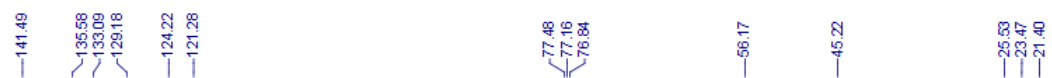
2f ¹³F NMR (100 MHz, CDCl₃)



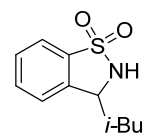




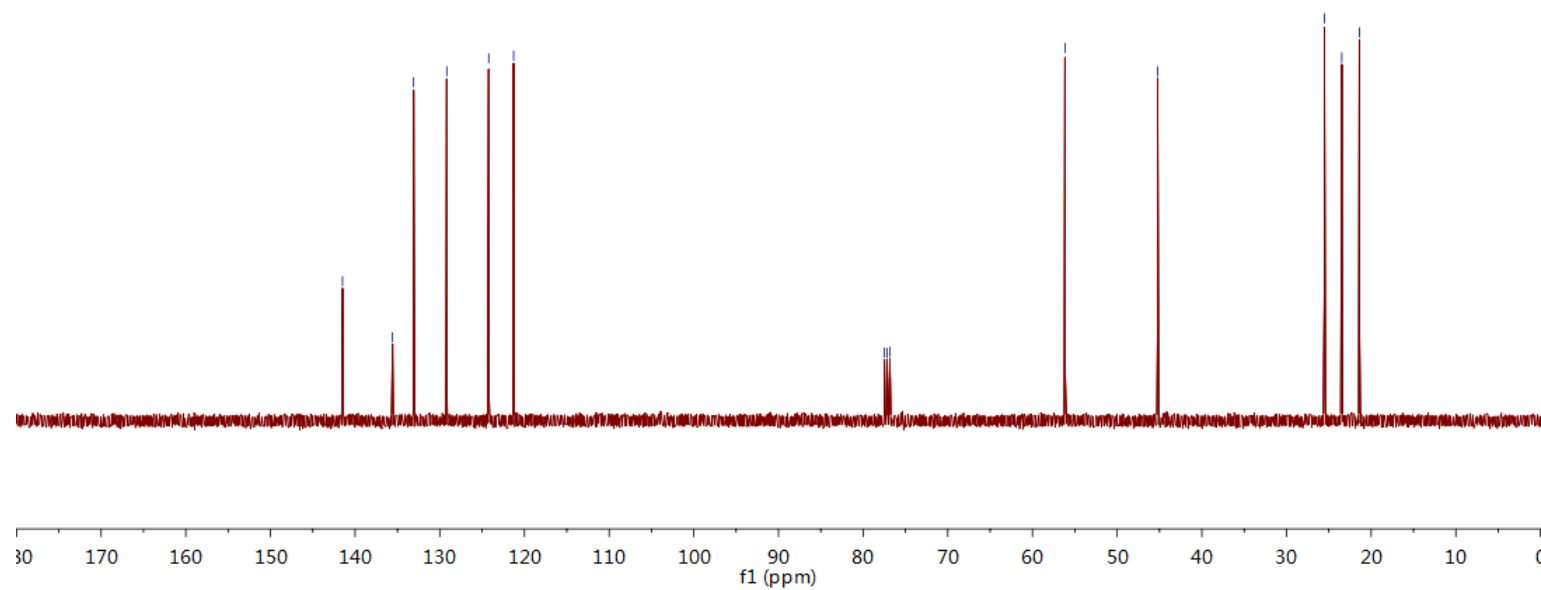


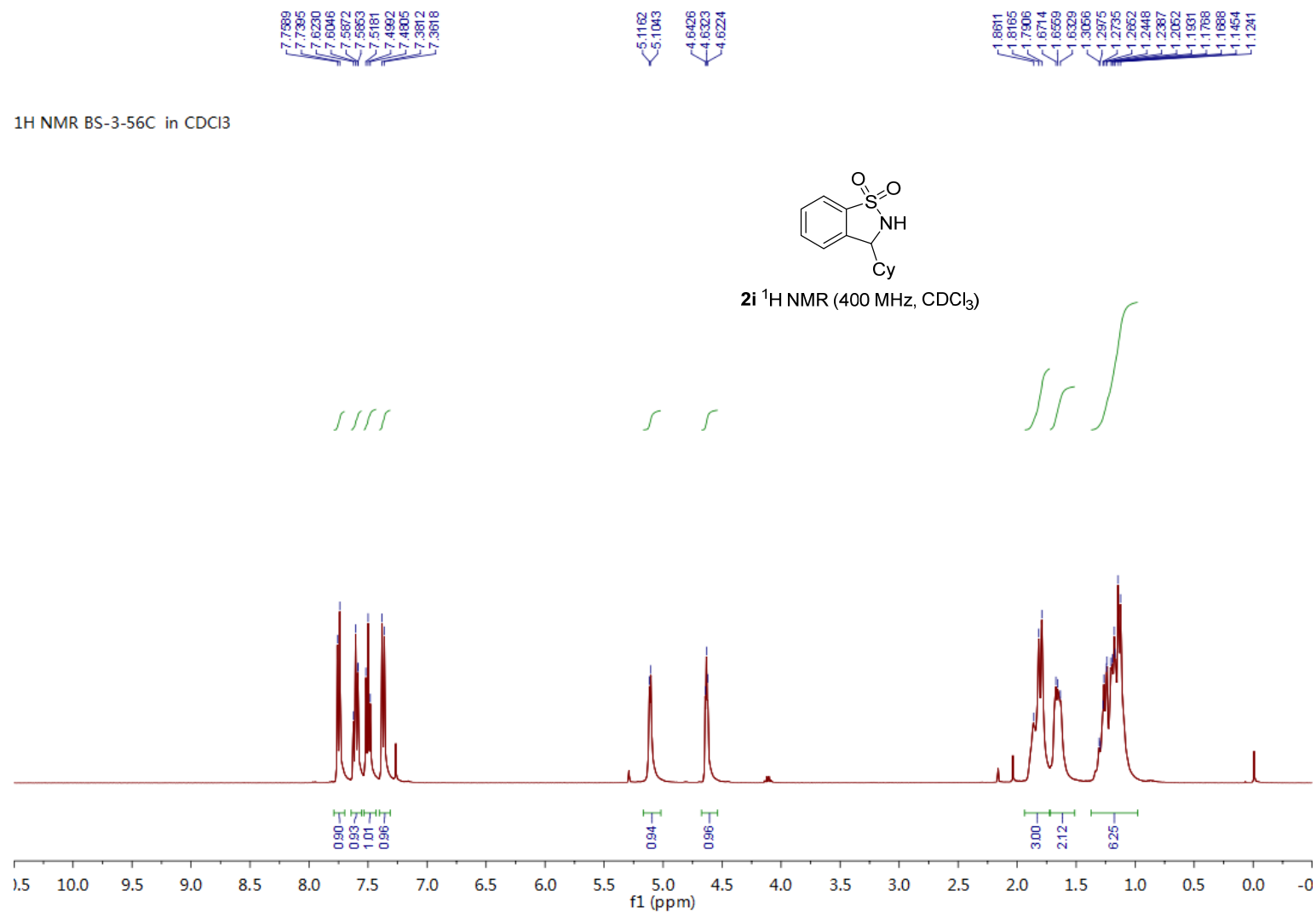


^{13}C NMR BS-3-56B in CDCl_3



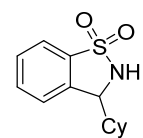
2h ^{13}C NMR (100 MHz, CDCl_3)



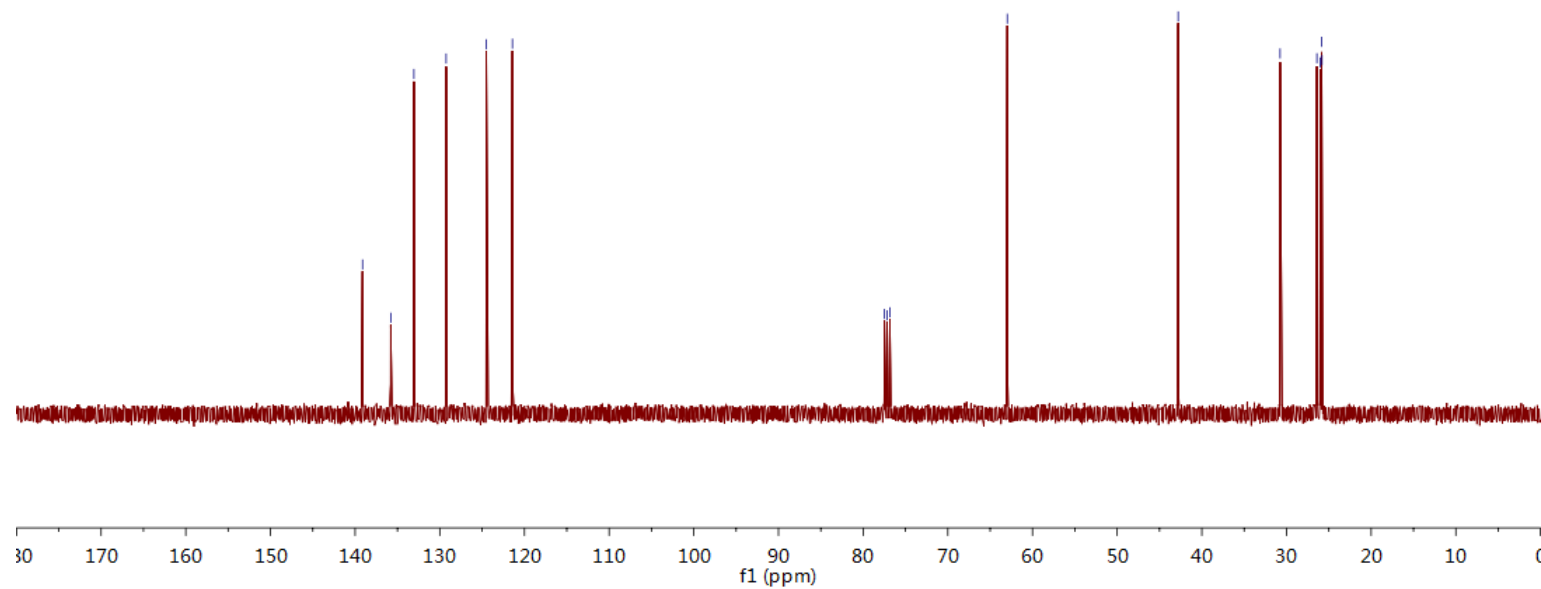


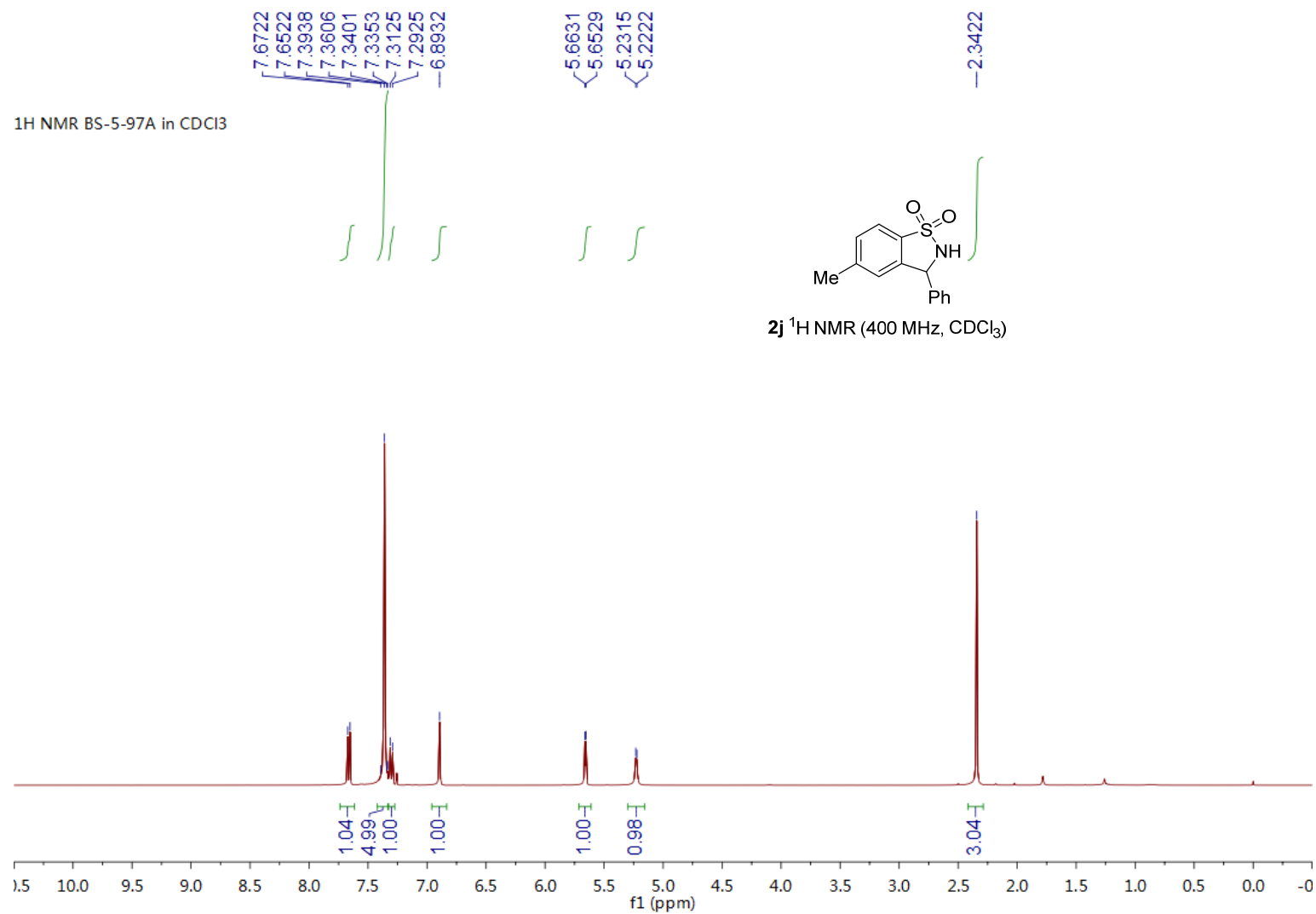


¹³C NMR BS-3-56C in CDCl₃



2i ¹³C NMR (100 MHz, CDCl₃)

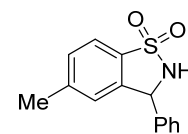




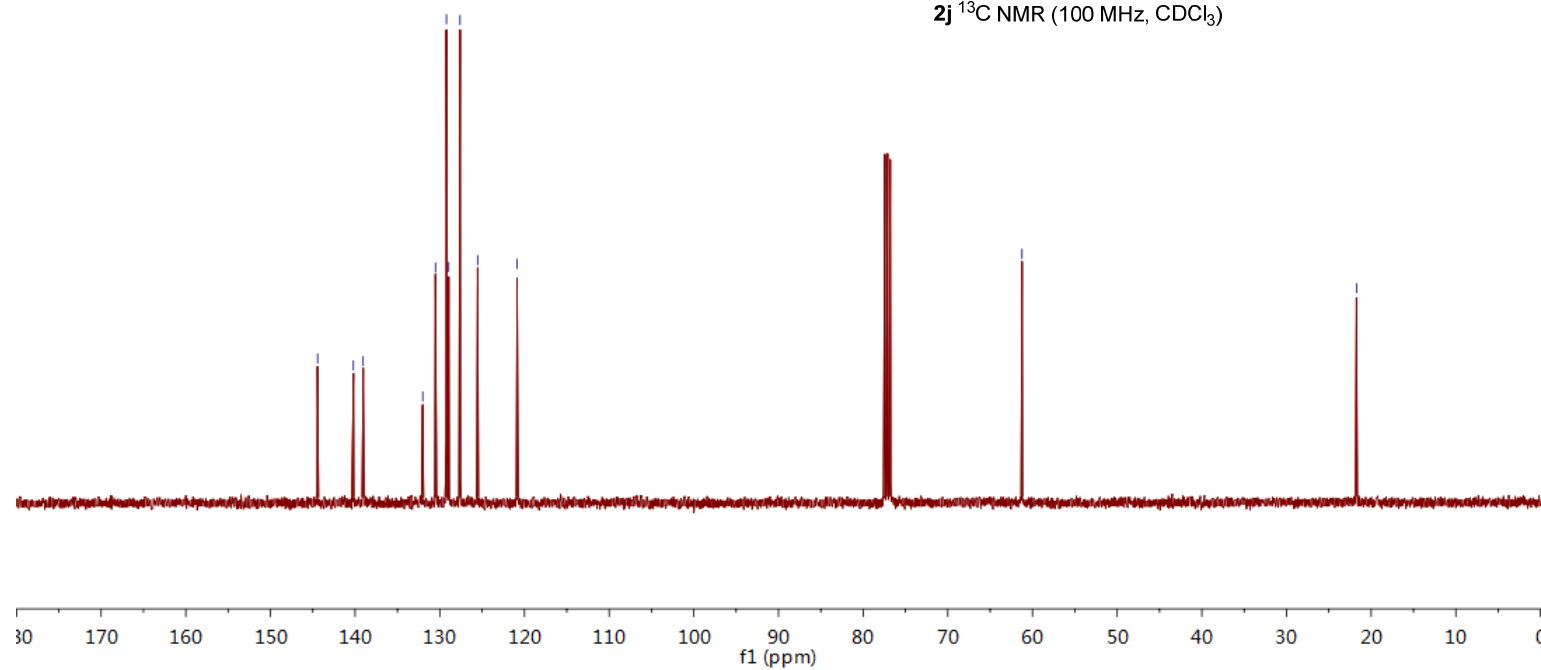
¹³C NMR BS-5-97A in CDCl₃
 144.43
 140.20
 139.05
 132.02
 130.51
 129.21
 128.95
 127.61
 125.53
 120.86

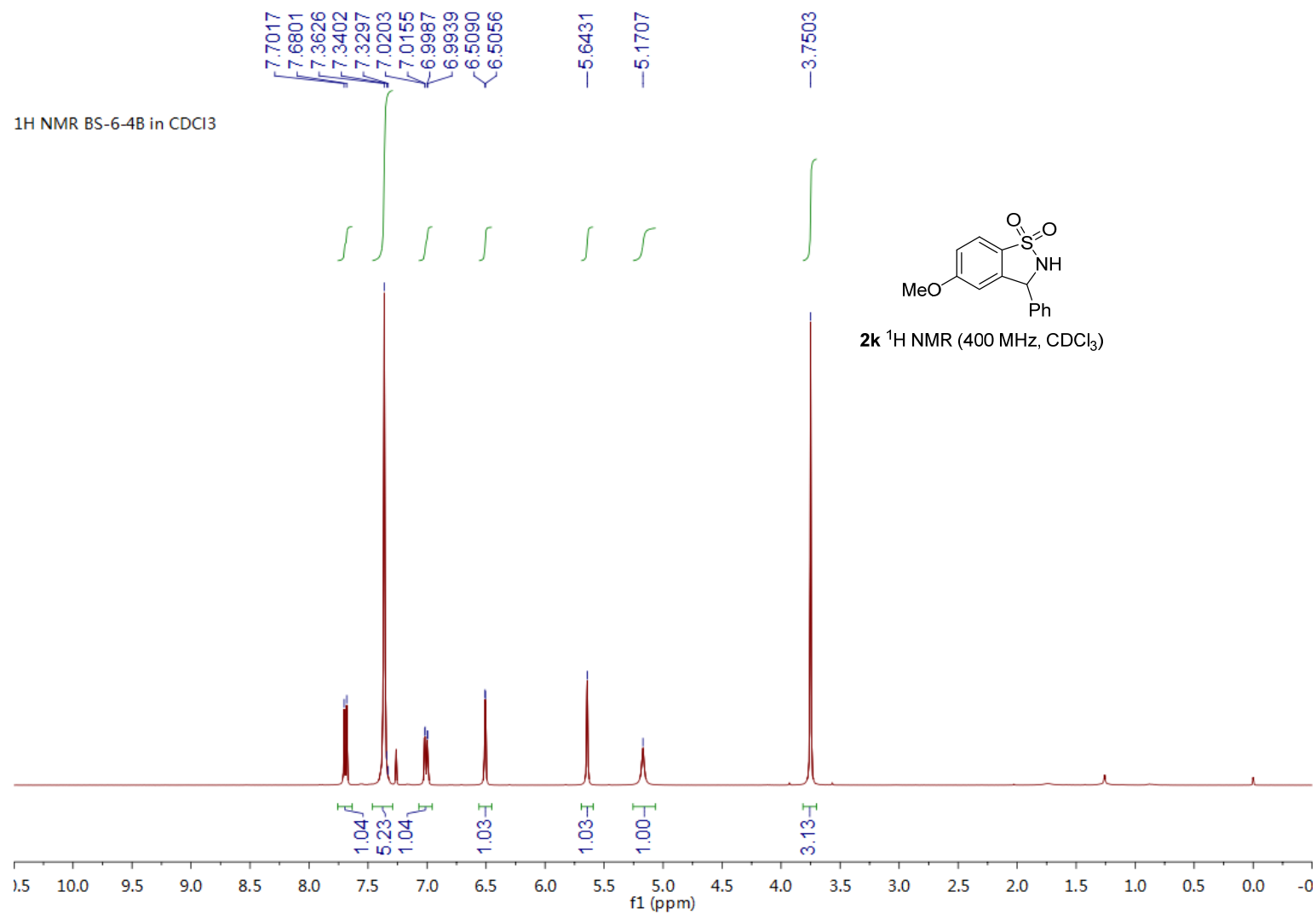
—61.24

—21.74



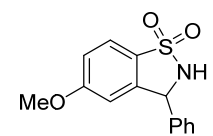
2j ¹³C NMR (100 MHz, CDCl₃)



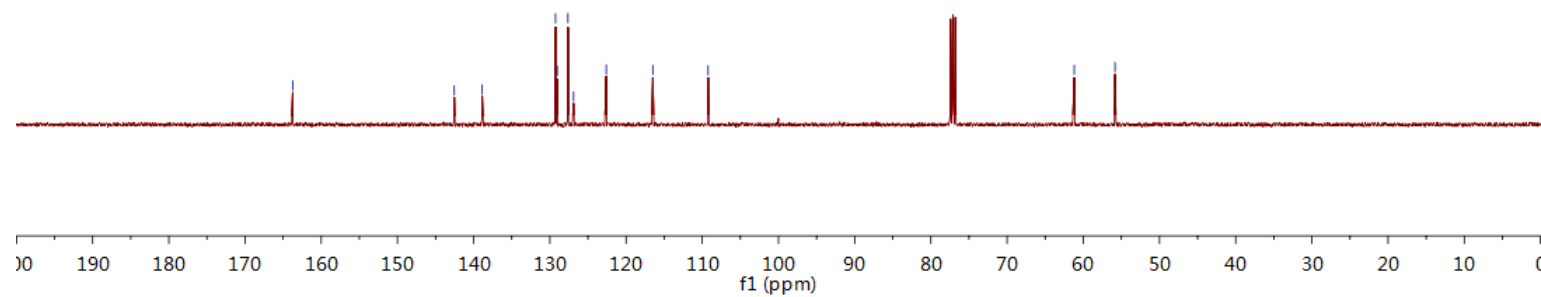


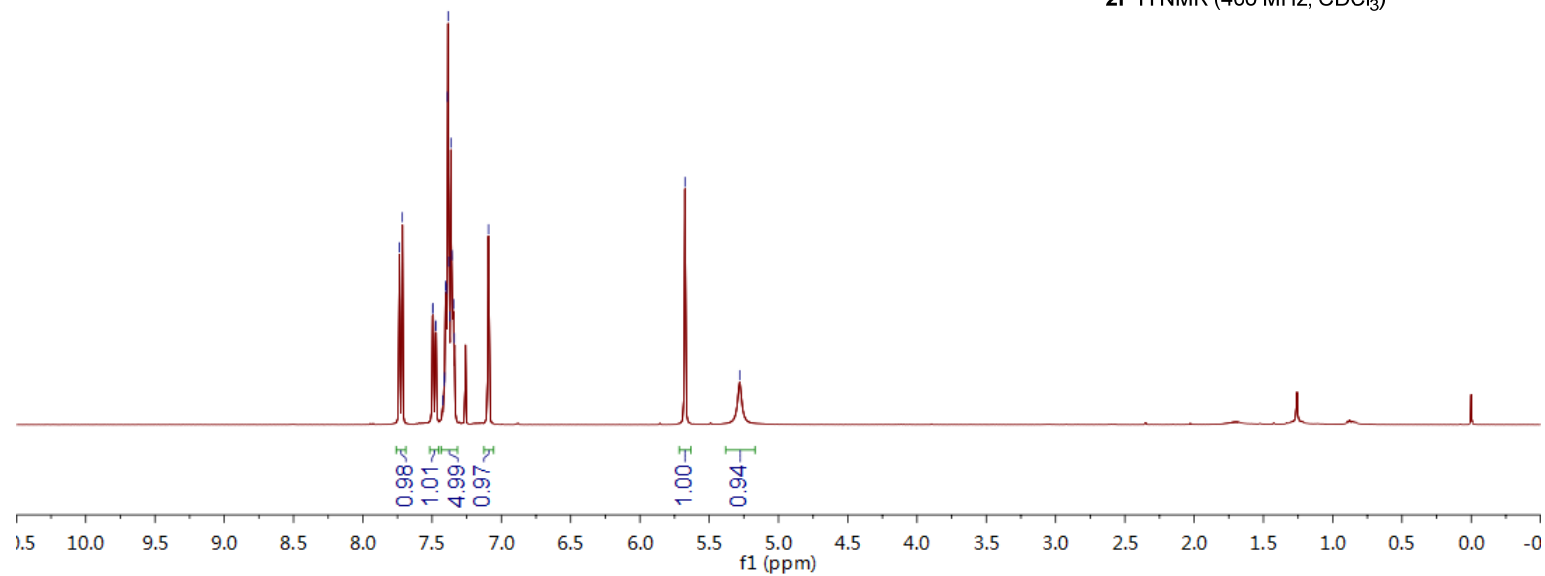
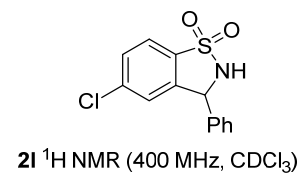
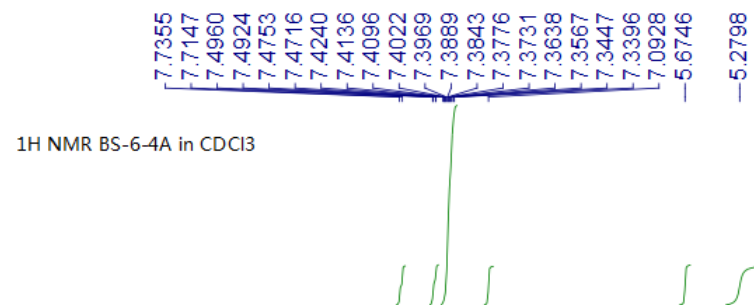
¹³C NMR BS-6-4B in CDCl₃
 —163.75
 —142.54
 —138.88
 —129.26
 —129.03
 —127.62
 —126.89
 —122.60
 —116.49
 —109.23

—61.18
 —55.84



2k ¹³C NMR (100 MHz, CDCl₃)

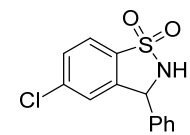




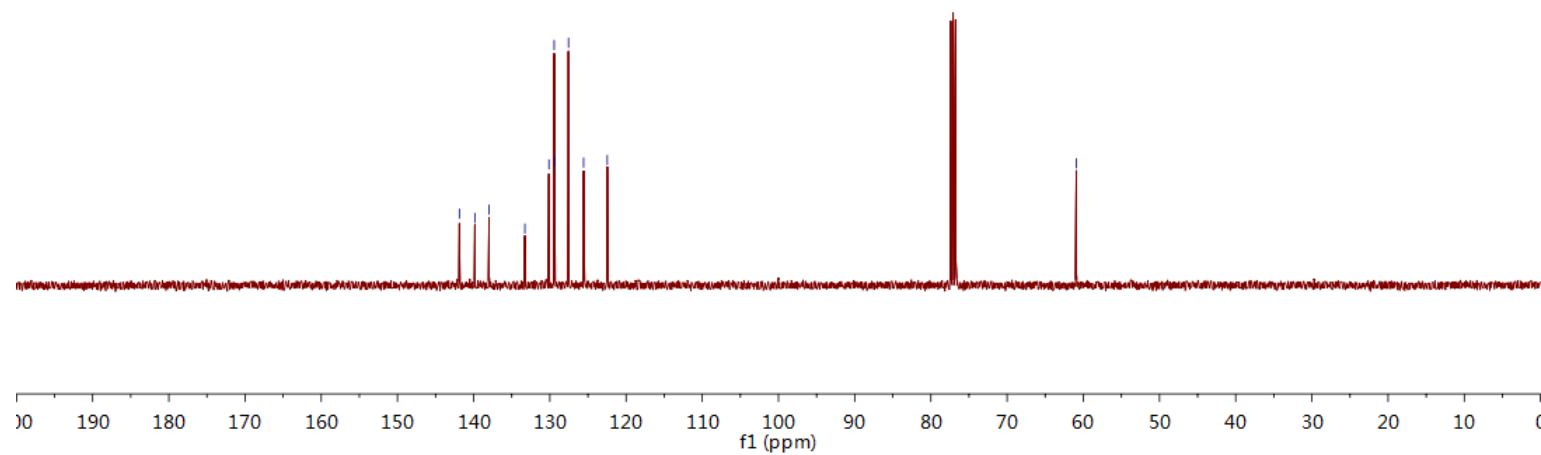
¹³C NMR BS-6-4A in CDCl₃

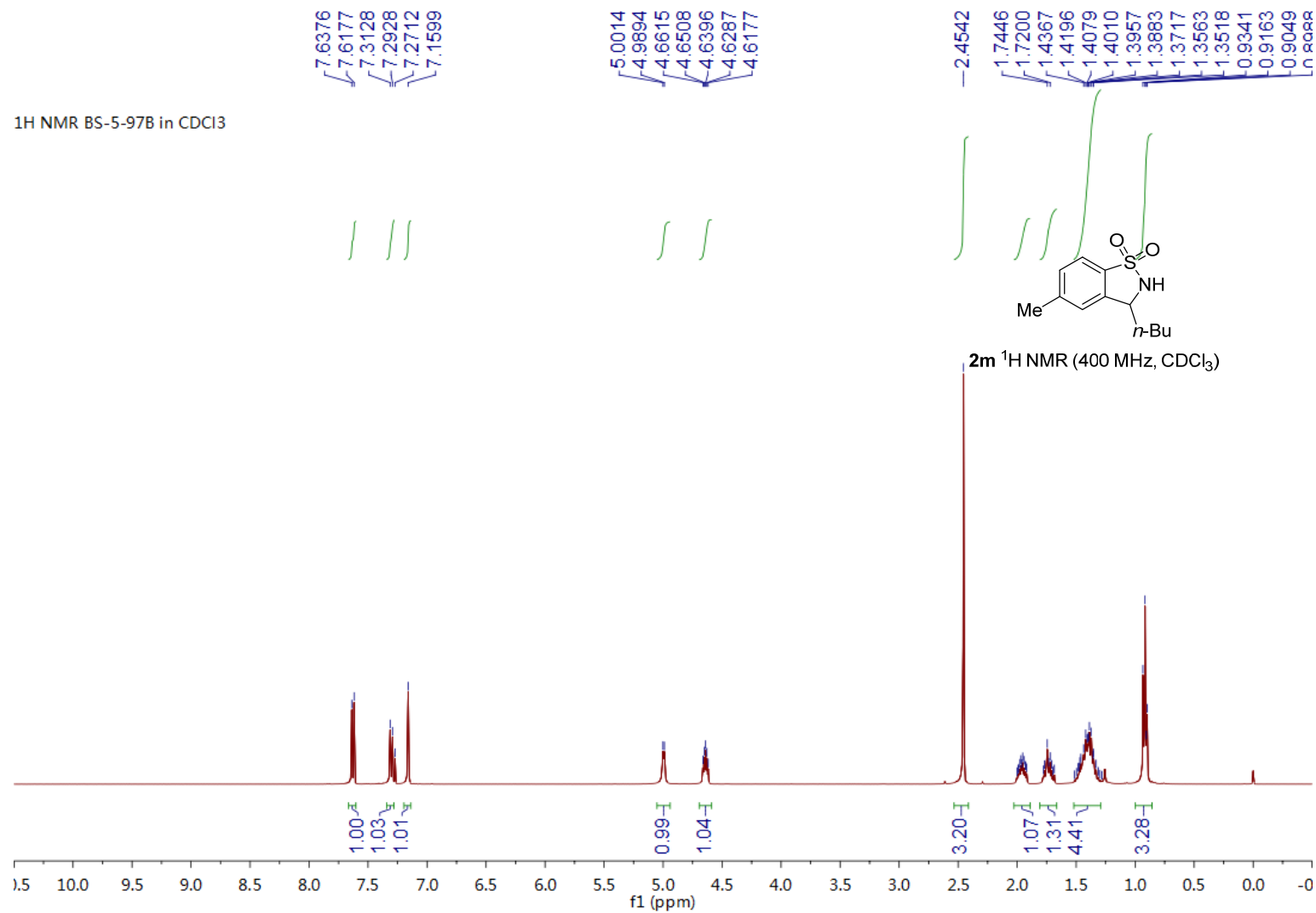
141.87
139.84
137.98
133.28
130.12
129.43
129.36
127.55
125.57
122.47

60.92



2I ¹³C NMR (100 MHz, CDCl₃)





¹³C NMR BS-5-97B in CDCl₃

~144.05
~141.07
~132.92
~130.16
~124.37
~121.02

—57.79

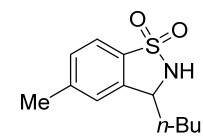
—35.55

~27.87

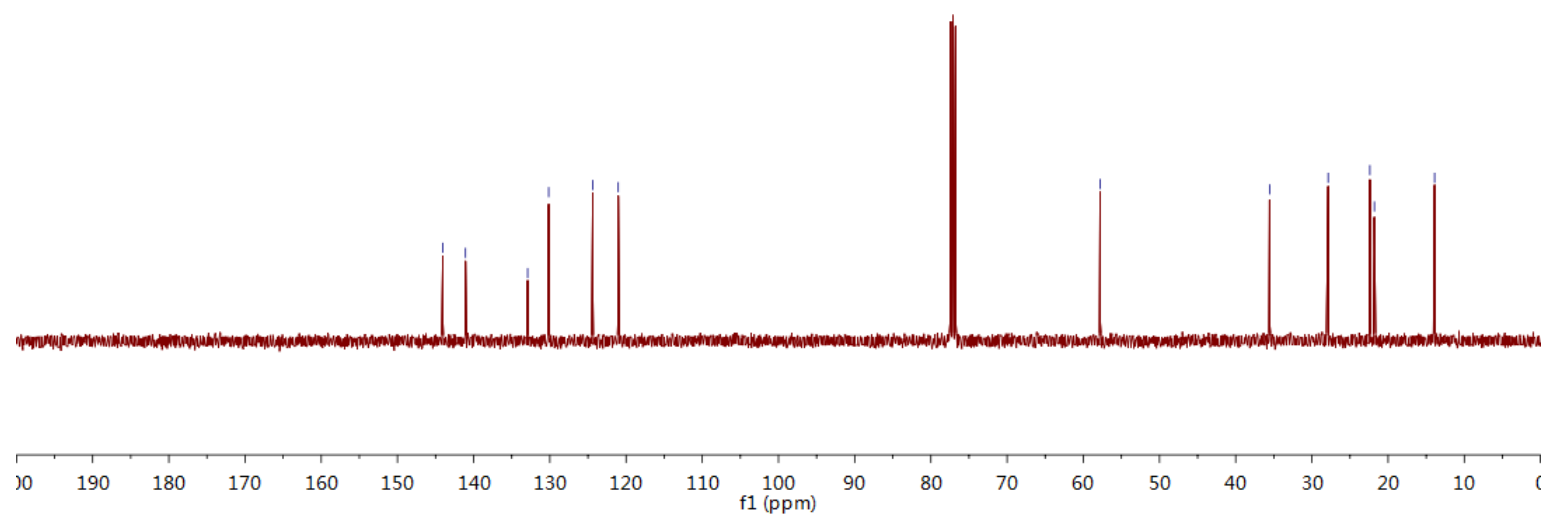
~22.40

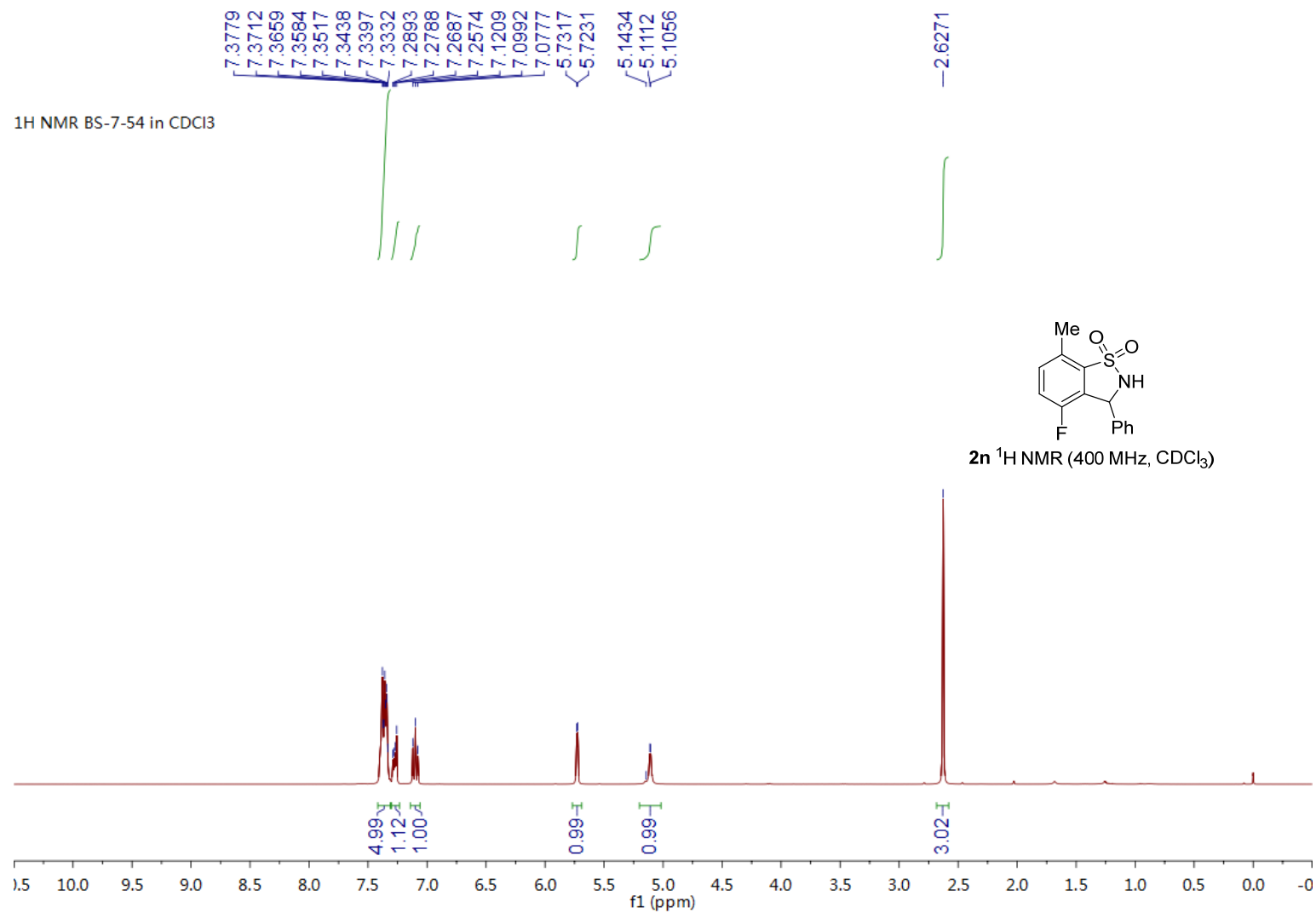
~21.80

—13.90

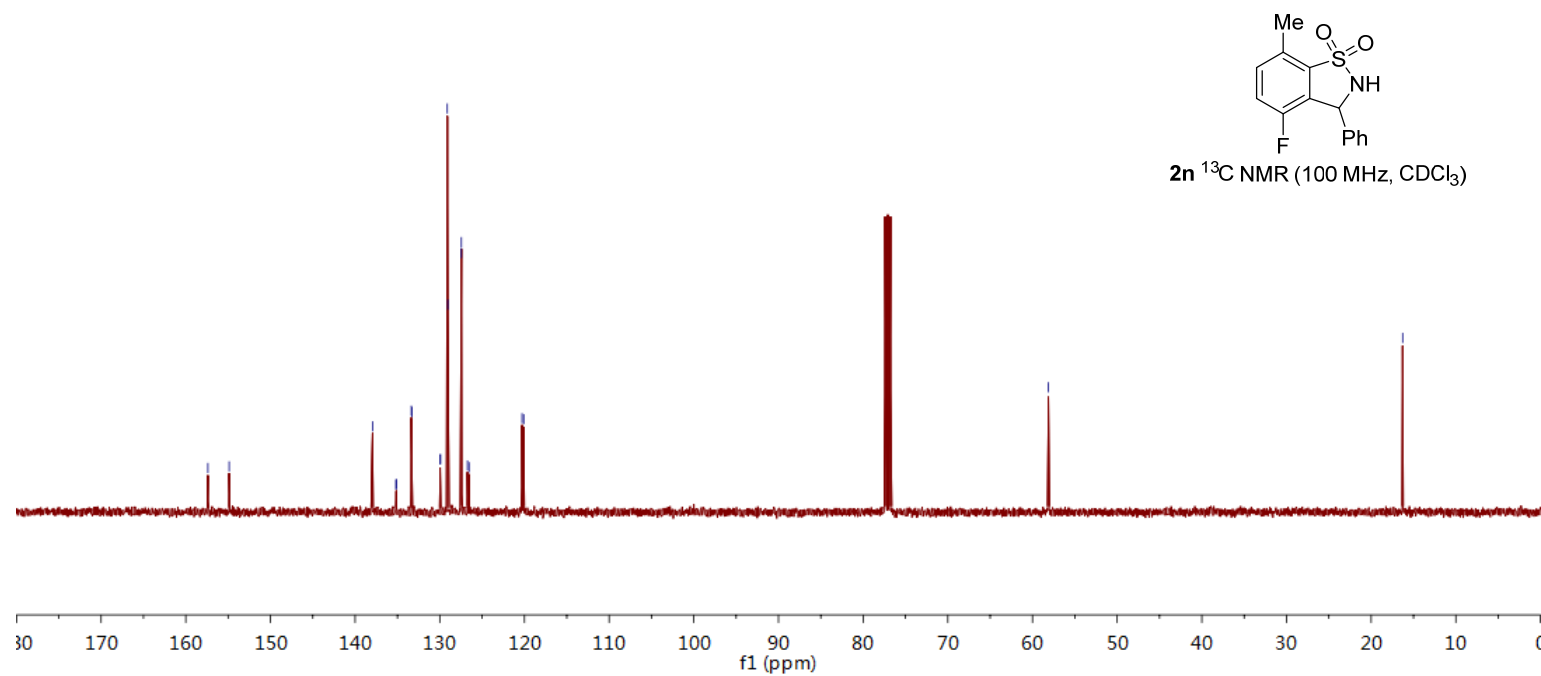


2m ¹³C NMR (100 MHz, CDCl₃)



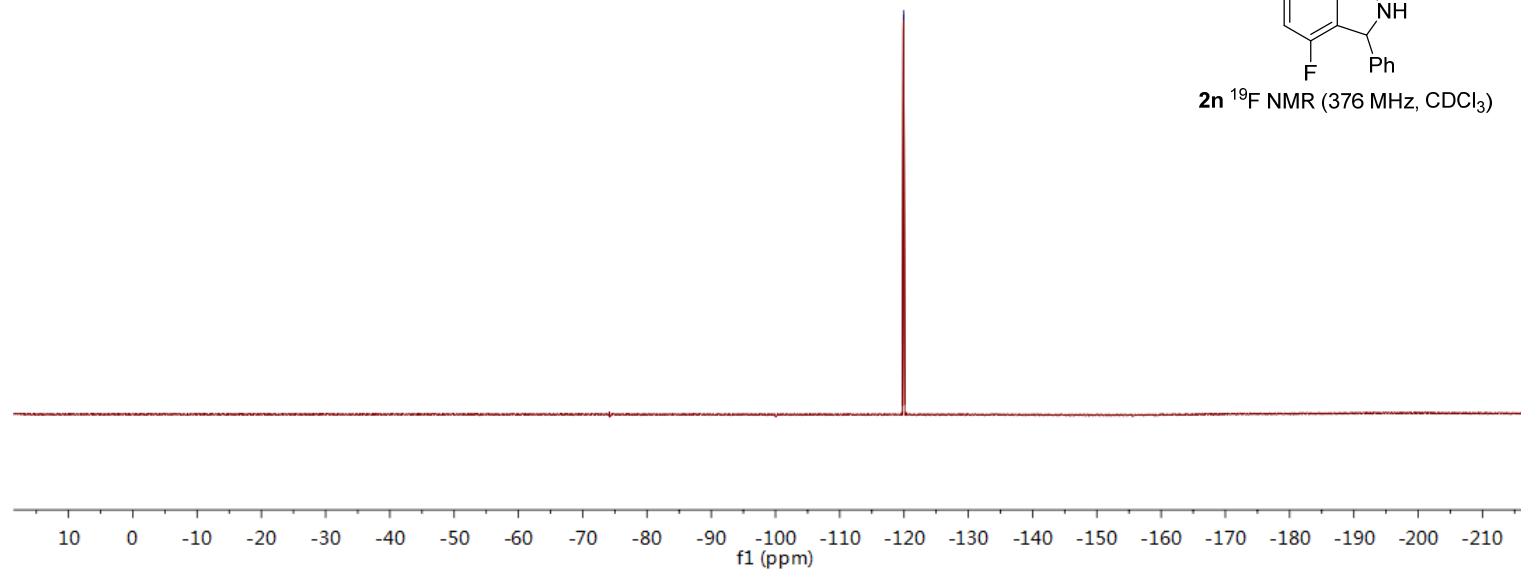
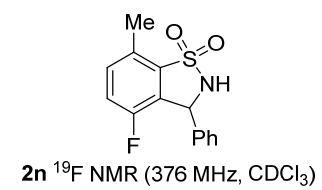


¹³C NMR BS-7-54 in CDCl₃
 157.39, 154.88, 137.94, 133.36, 133.29, 129.96, 129.91, 129.11, 129.05, 127.46, 127.45, 120.48, 120.09, 58.12, 16.28

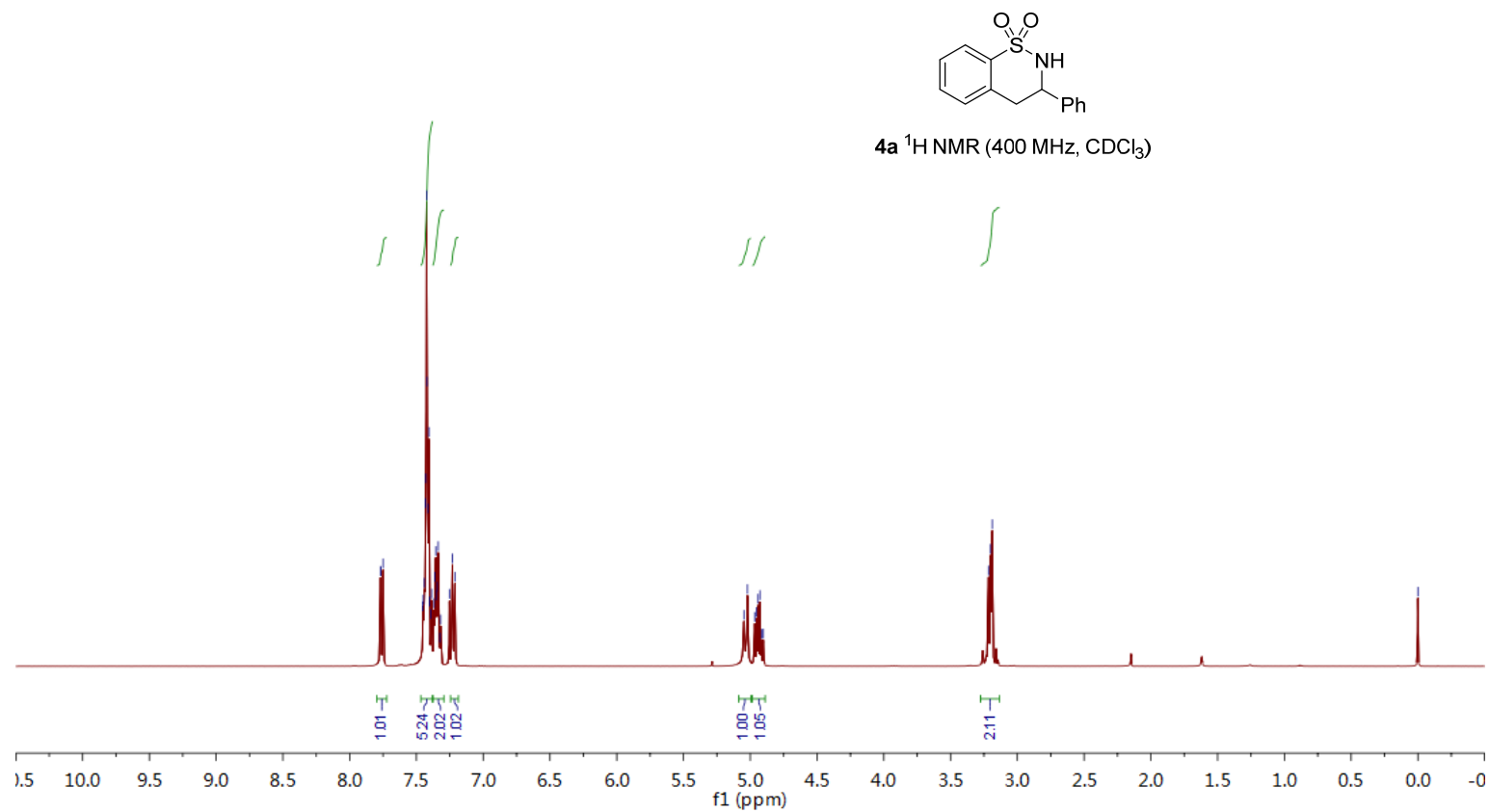


¹⁹F NMR BS-7-54 in CDCl₃

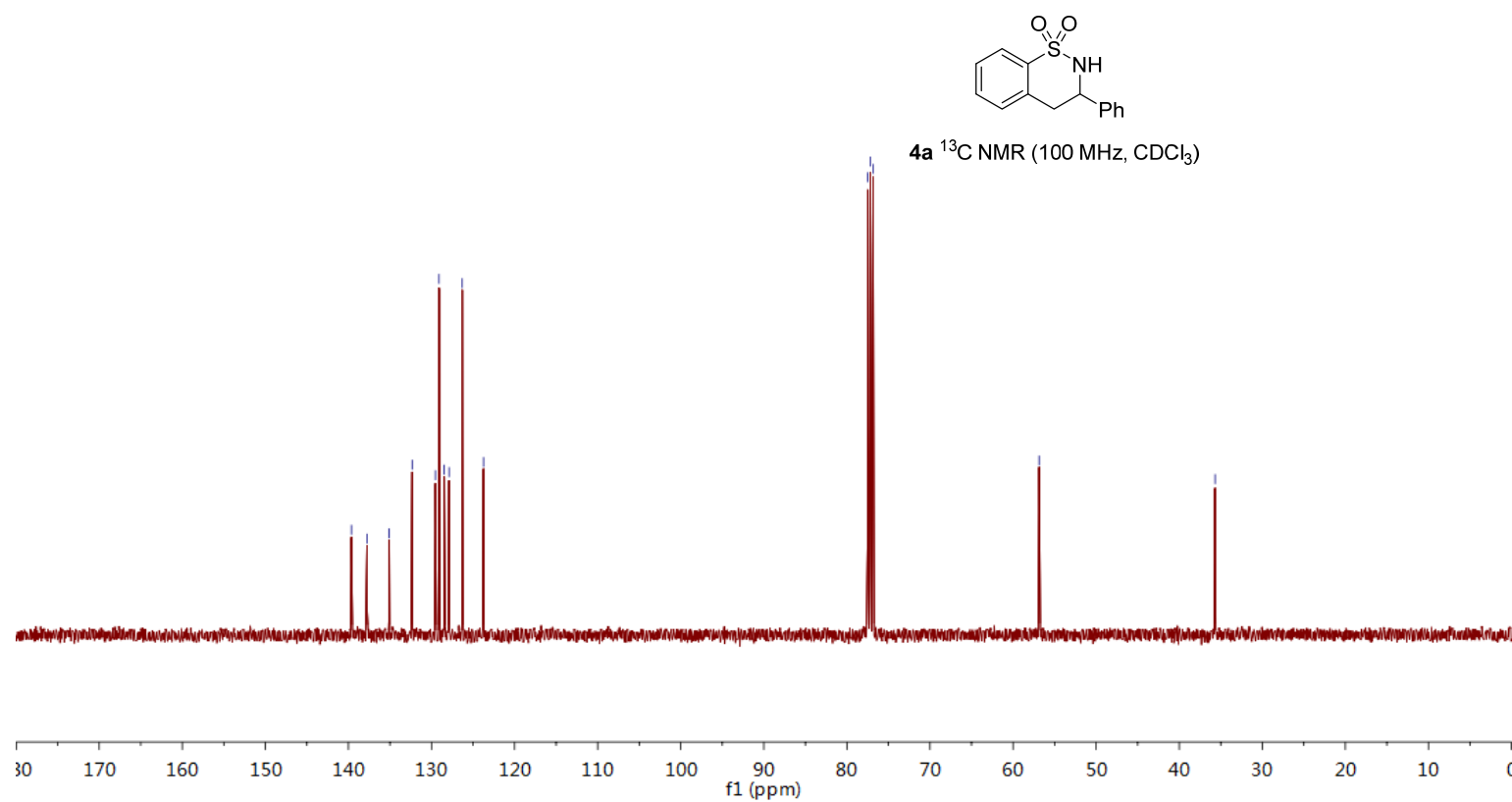
—119.9244

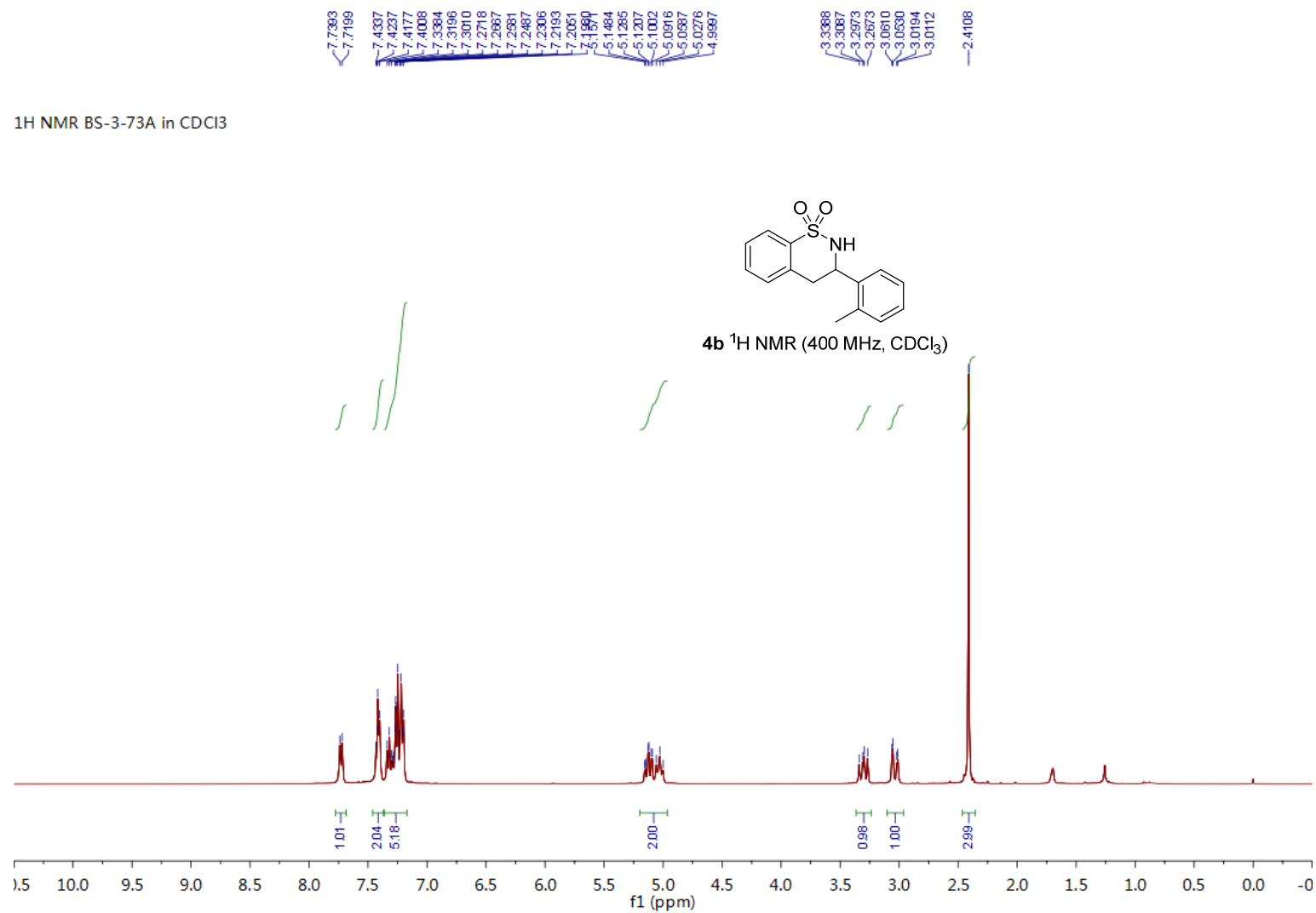


¹H NMR BS-3-70 in CDCl₃



¹³C NMR BS-3-70 in CDCl₃





137.52
137.19
136.18
135.28
132.25
131.07
129.57
128.42
127.86
126.72
125.21
123.88

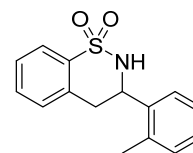
77.48
77.16
76.84

53.42

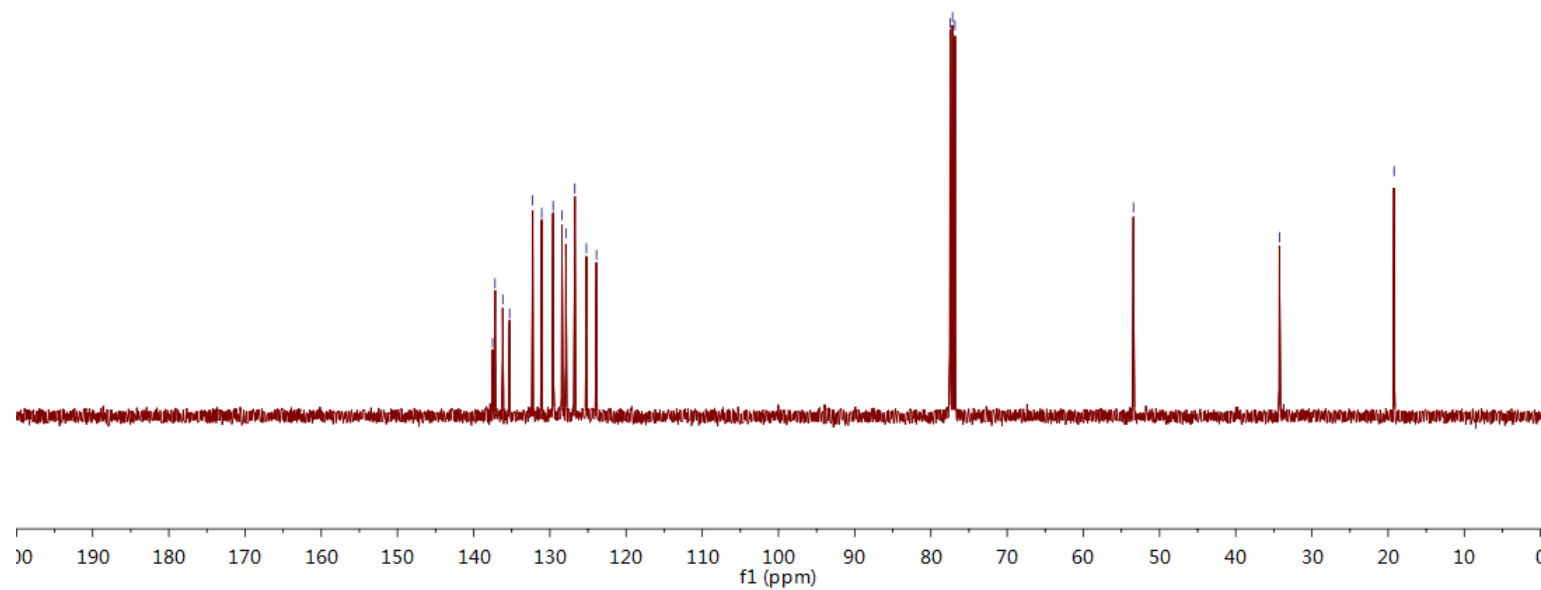
34.27

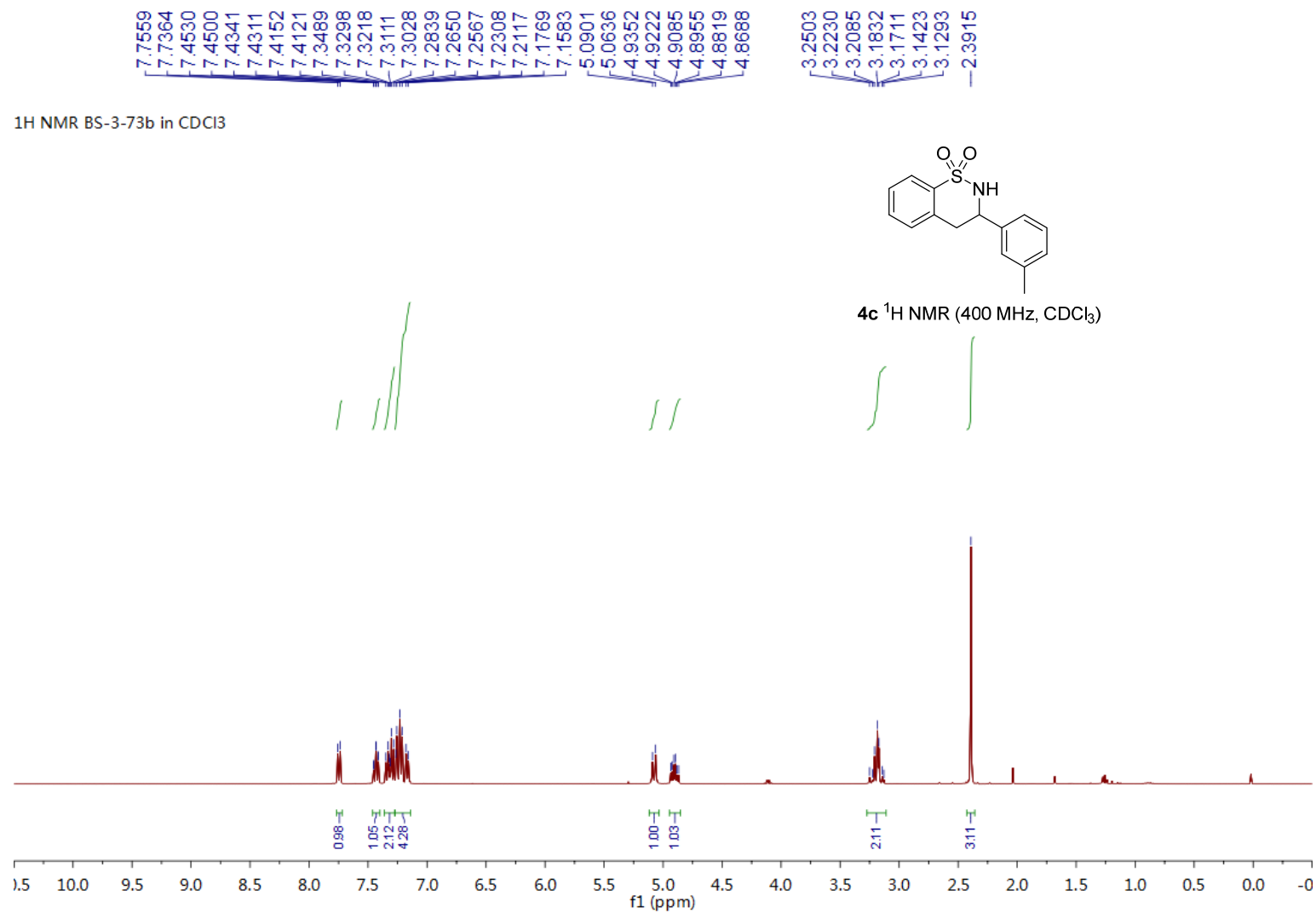
19.22

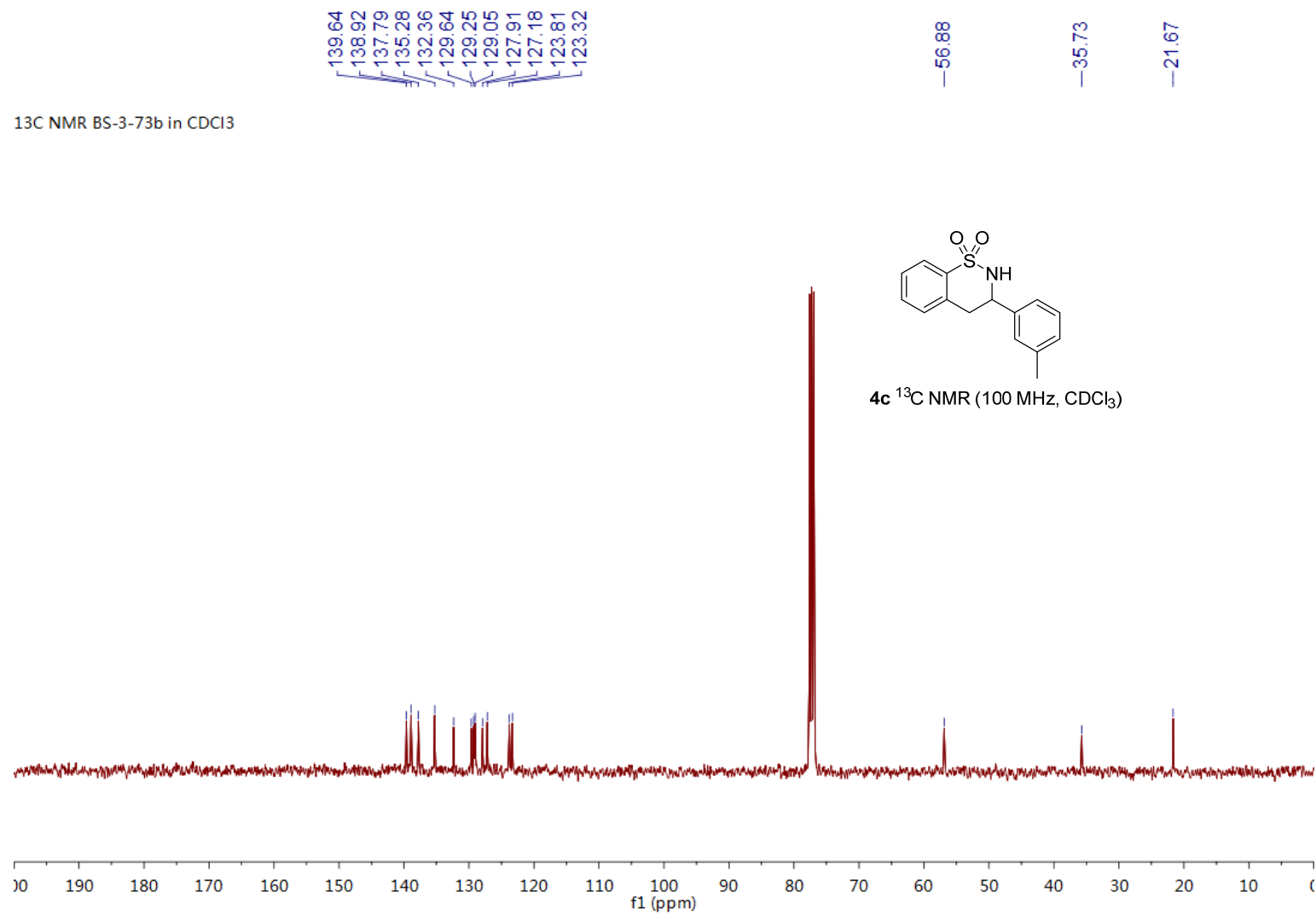
¹³C NMR BS-3-73A in CDCl₃

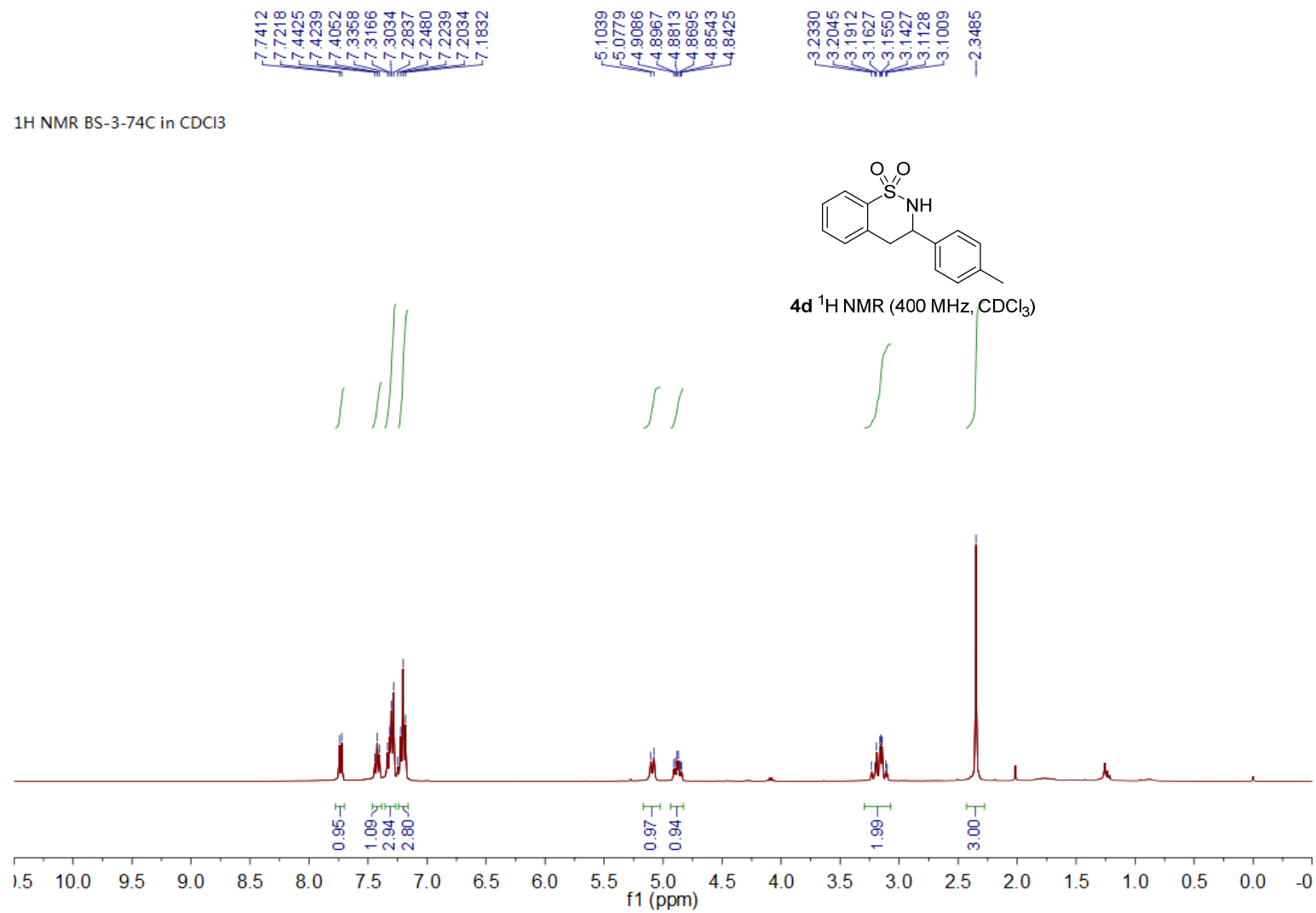


4b ¹³C NMR (100 MHz, CDCl₃)









¹³C NMR BS-3-74C in CDCl₃

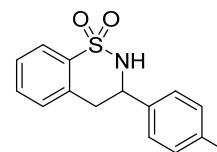
136.18
137.70
136.69
135.22
132.24
129.66
129.56
127.78
126.21
123.75

77.48
77.16
76.84

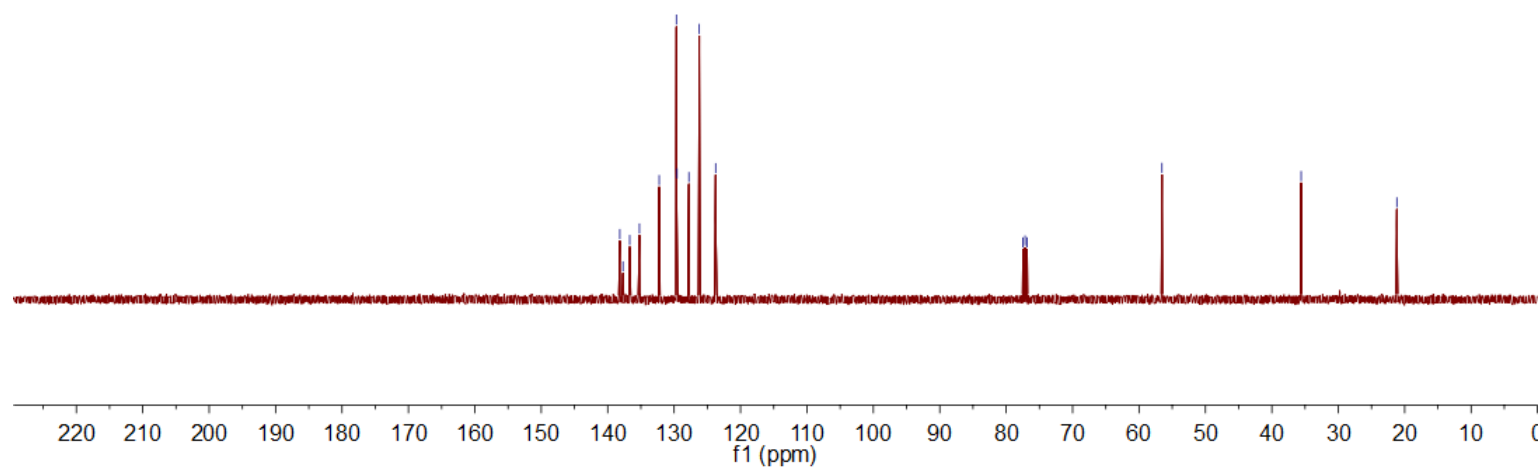
56.54

35.59

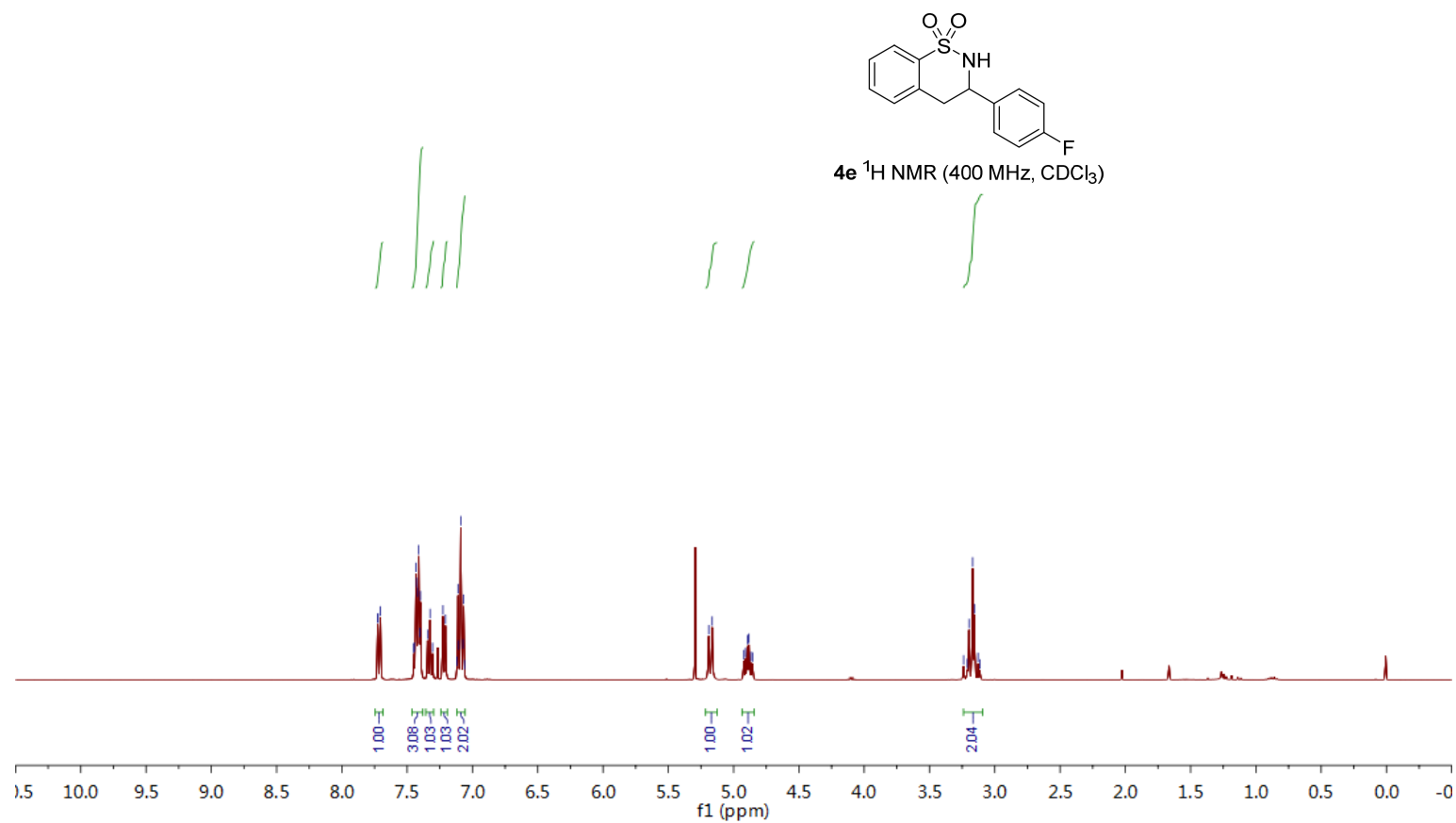
21.18



4d ¹³C NMR (100 MHz, CDCl₃)

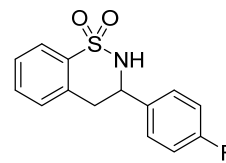


¹H NMR BS-3-74D in CDCl₃

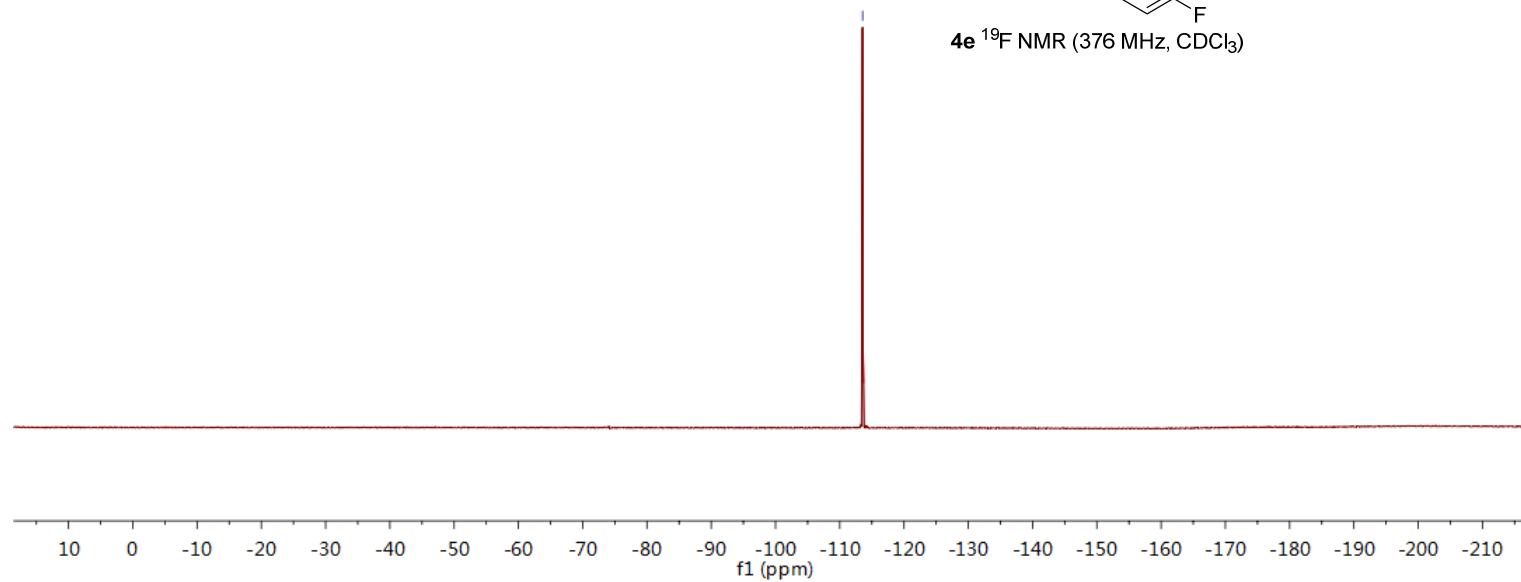


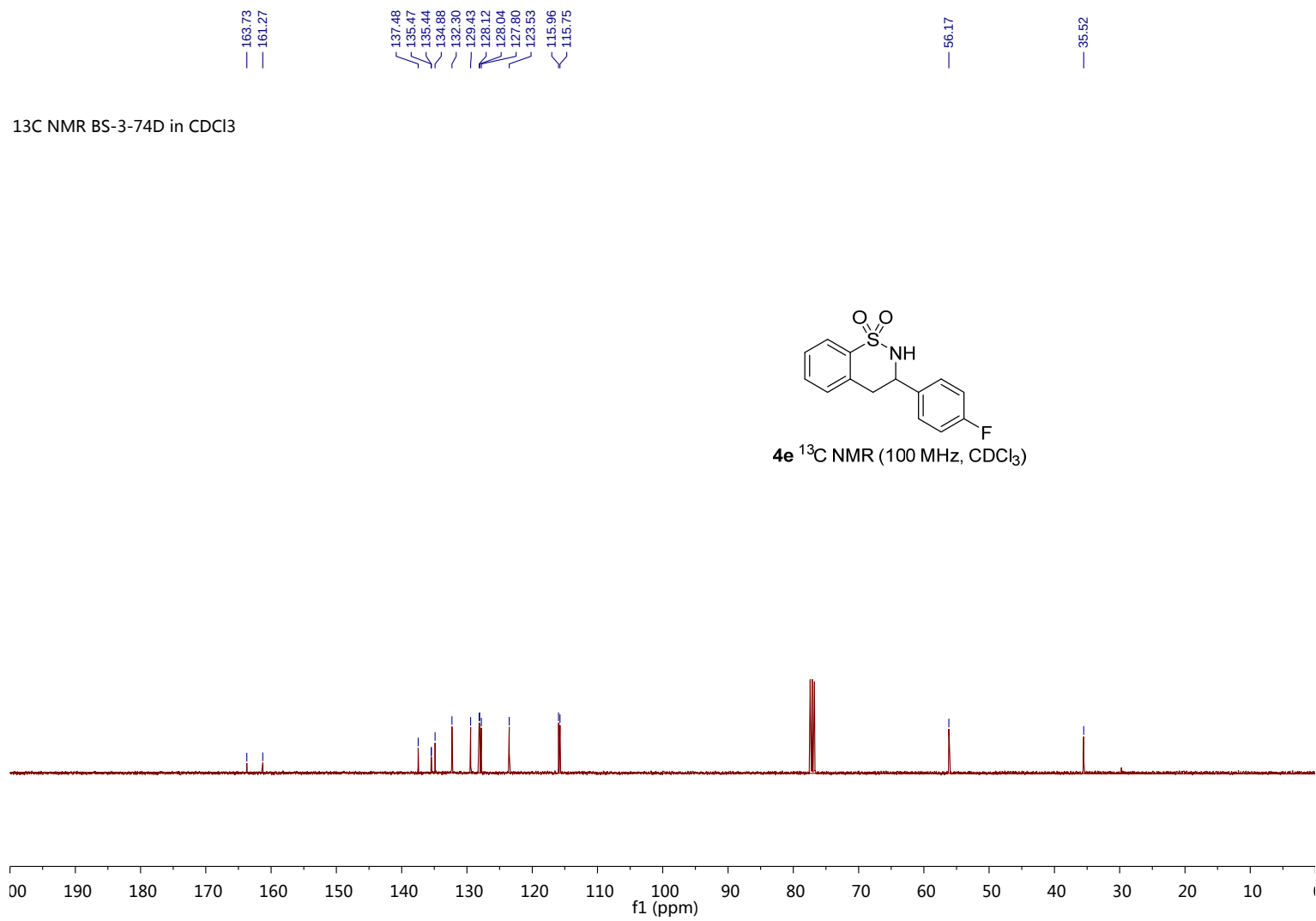
¹⁹F BS-3-74D in CDCl₃

—113.5576

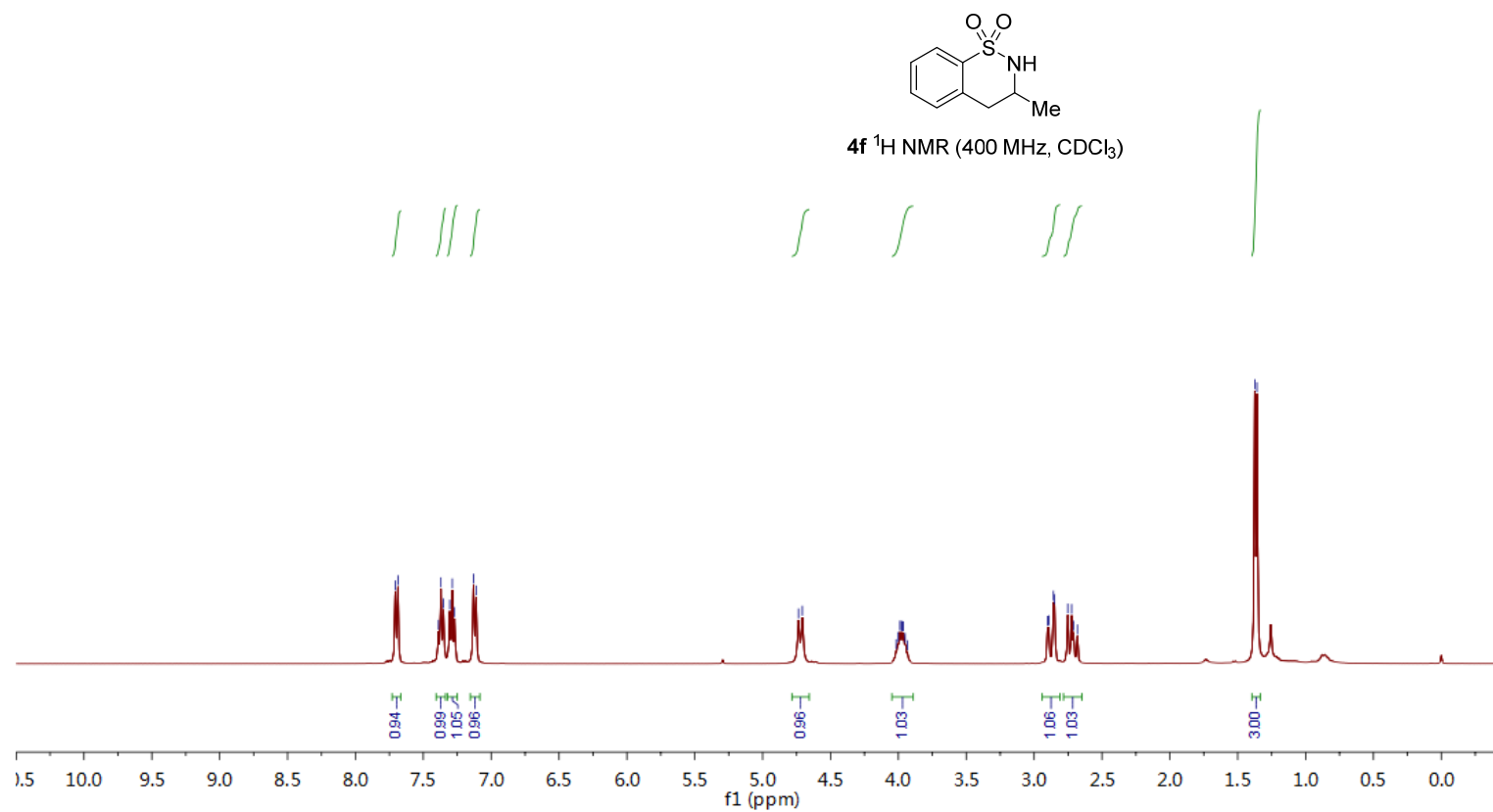


4e ¹⁹F NMR (376 MHz, CDCl₃)





¹H NMR BS-3-78C in CDCl₃



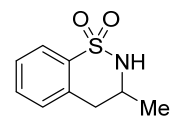
137.01
135.29
131.99
129.44
127.57
123.83

49.42

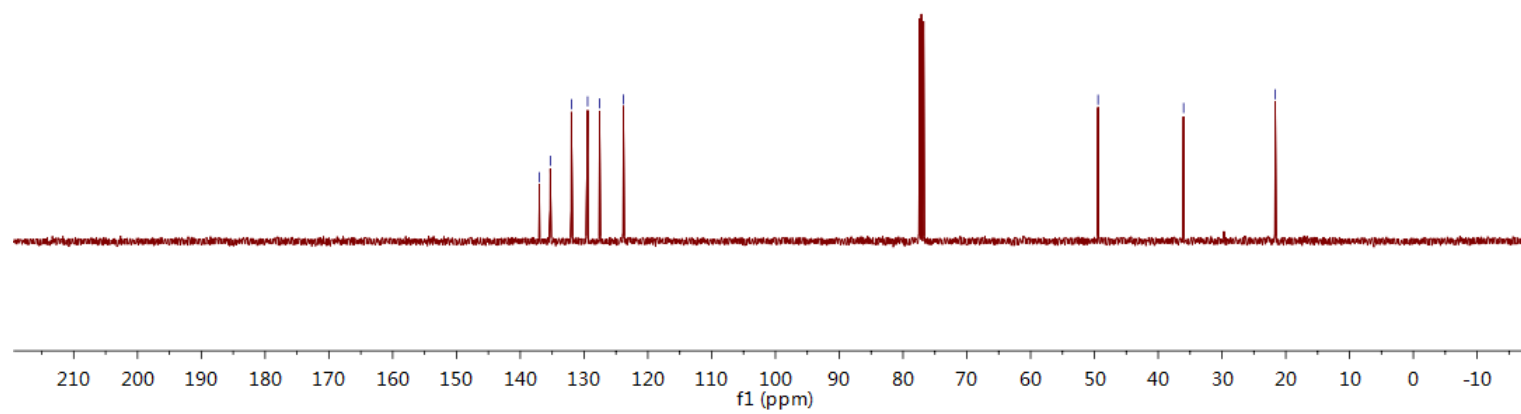
36.02

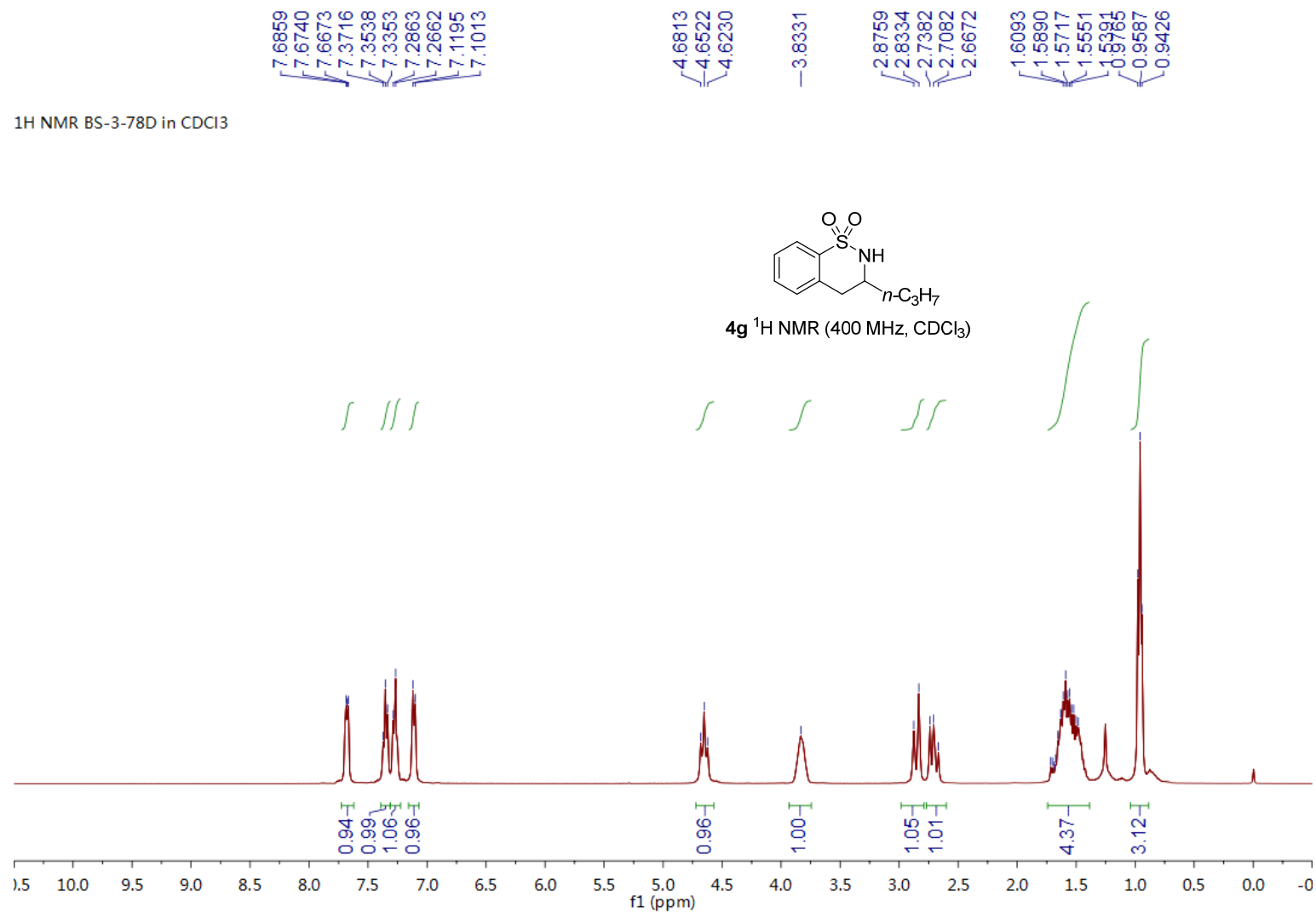
21.85

^{13}C NMR BS-3-78C in CDCl_3



4f ^{13}C NMR (100 MHz, CDCl_3)





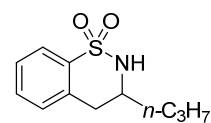
¹³C NMR BS-3-78D in CDCl₃

137.36
135.37
131.92
129.51
127.49
123.82

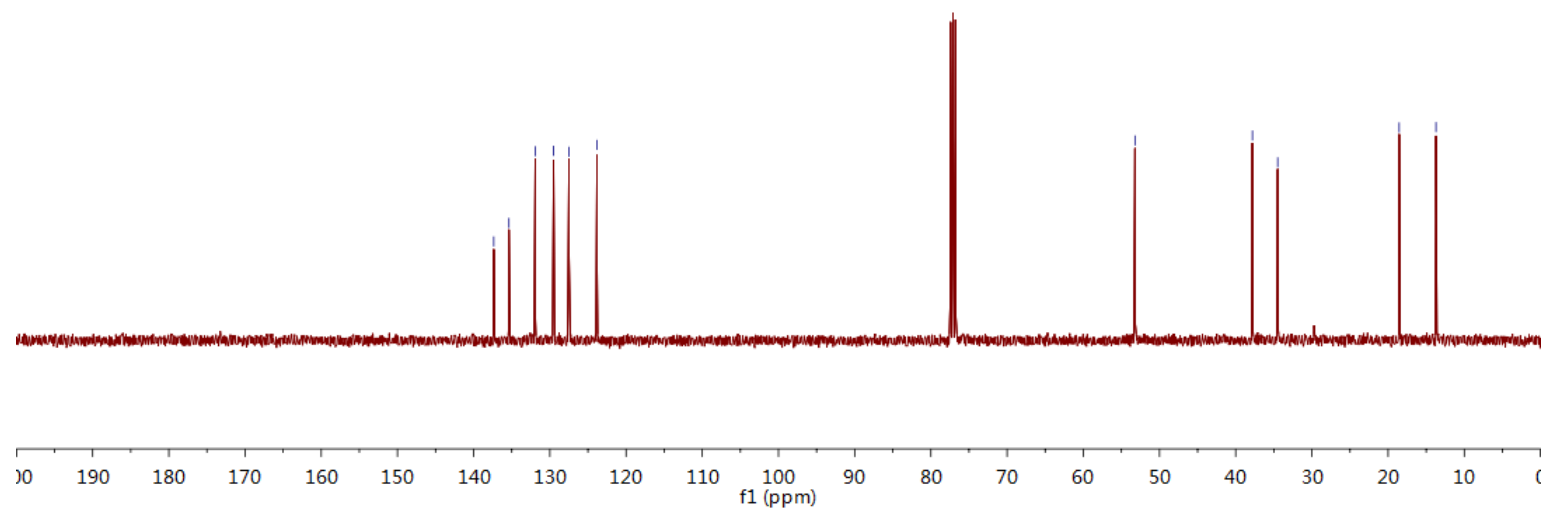
53.22

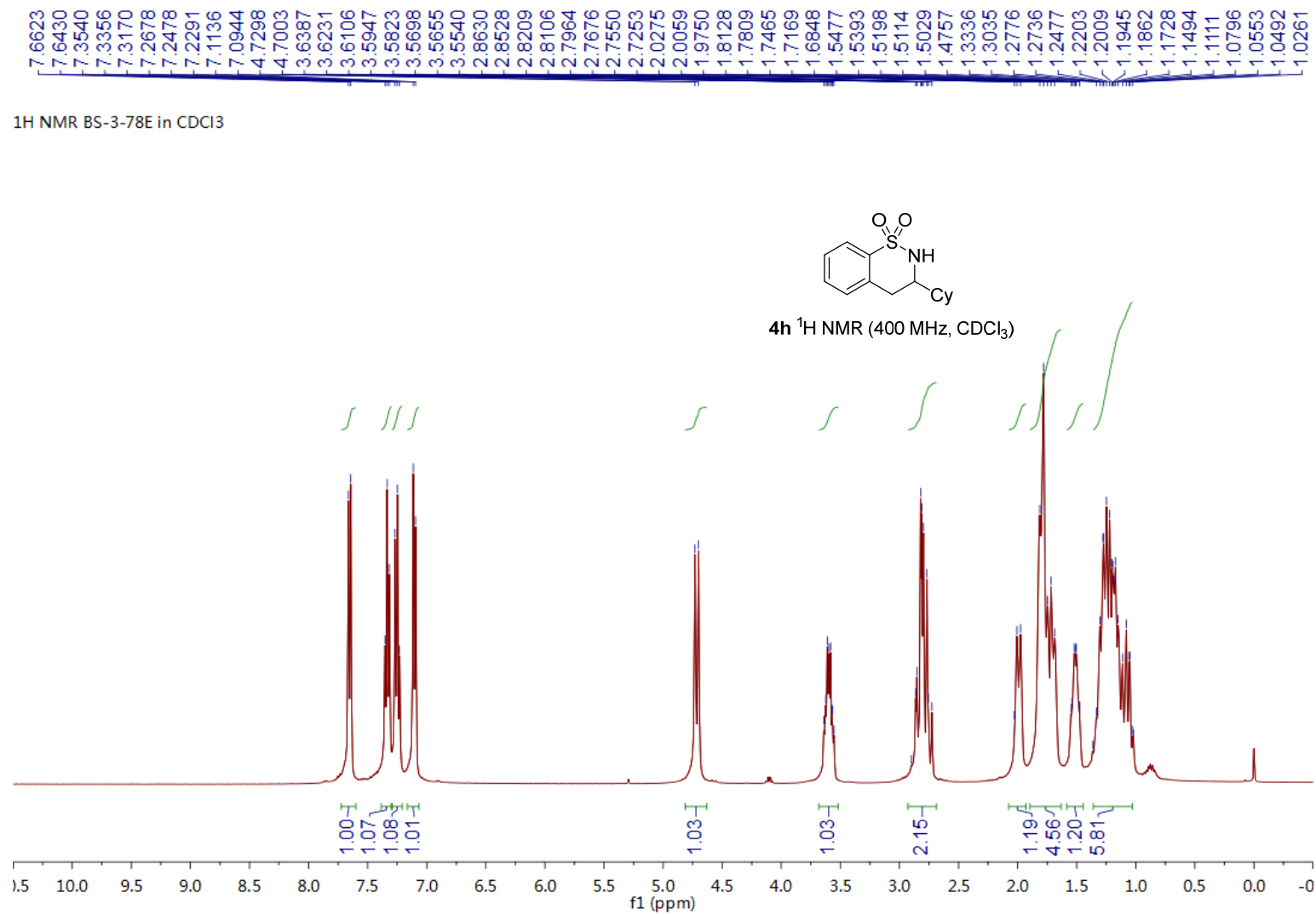
37.81
34.49

18.55
13.72

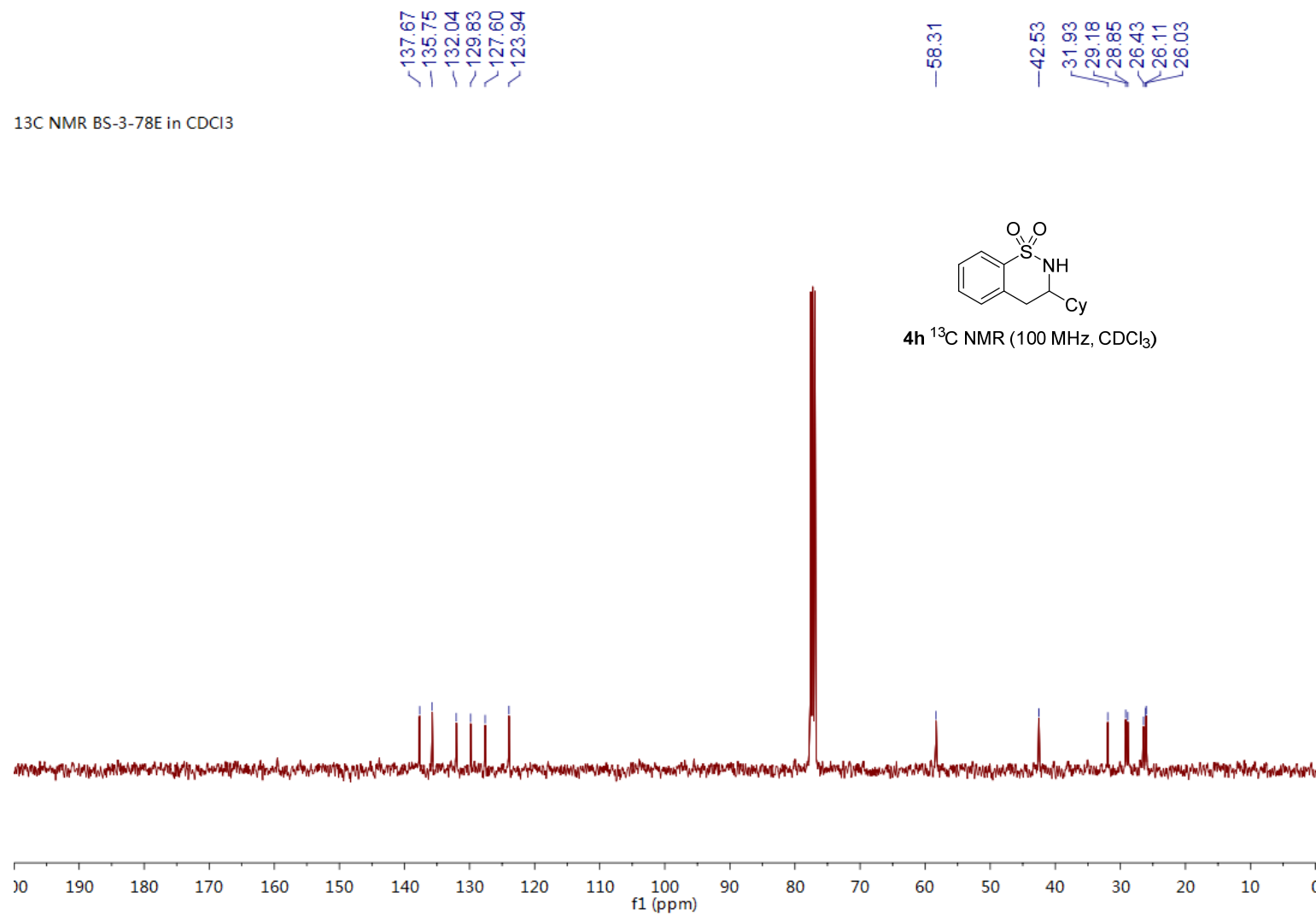


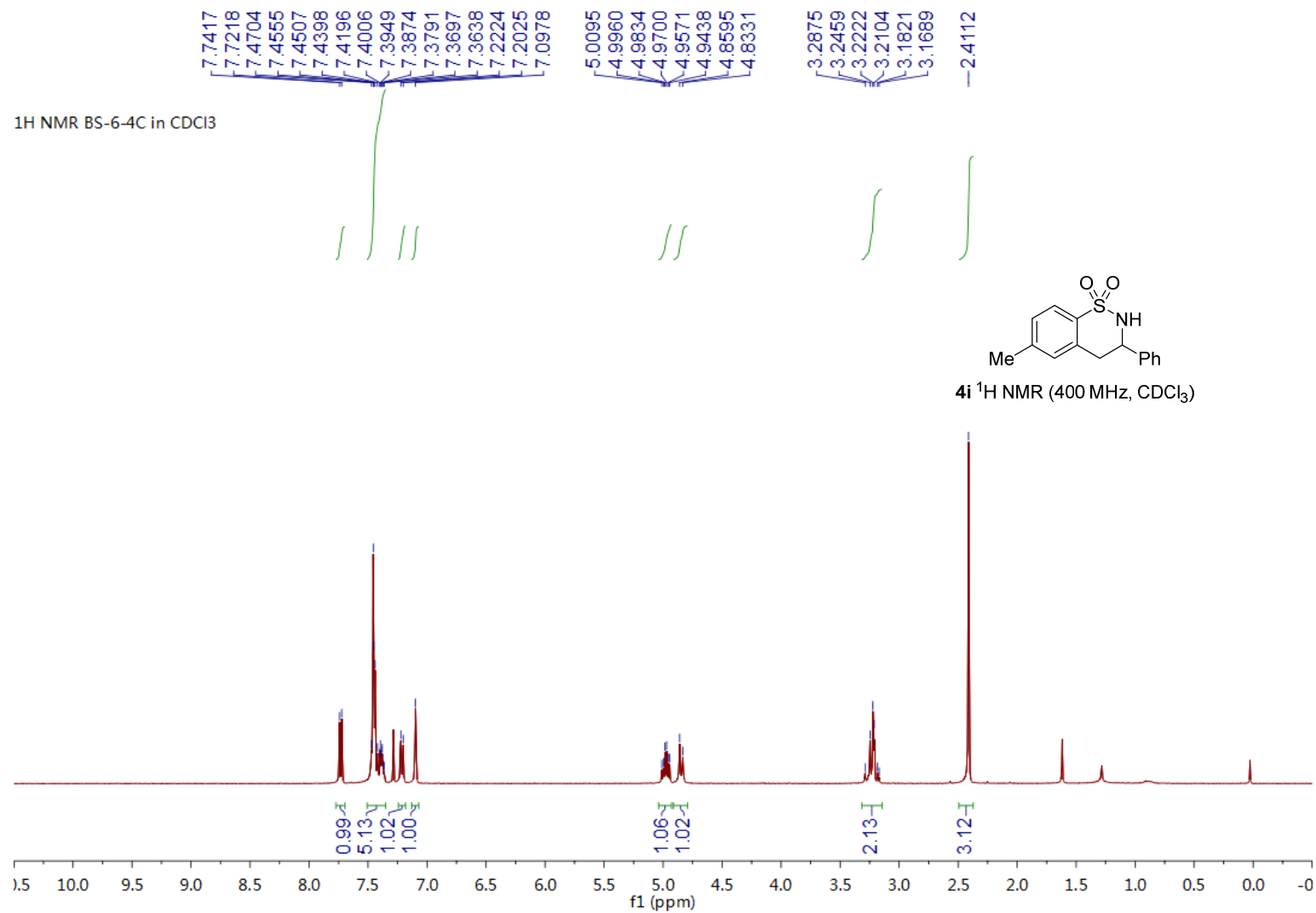
4g ¹³C NMR (100 MHz, CDCl₃)





¹³C NMR BS-3-78E in CDCl₃





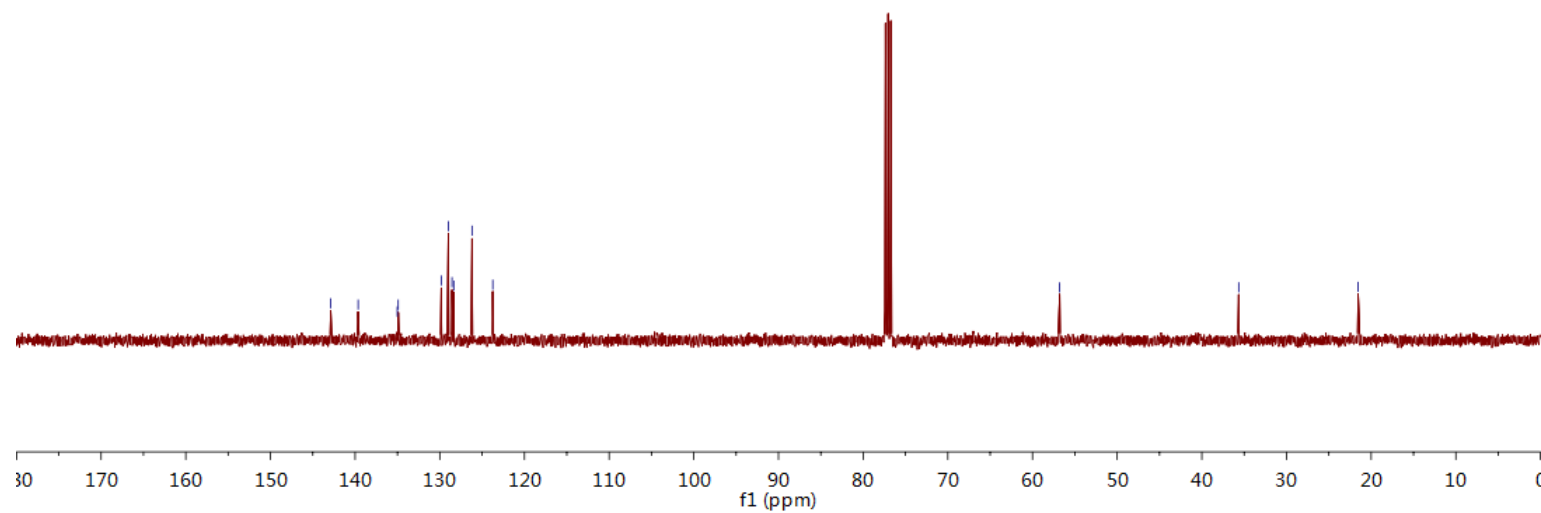
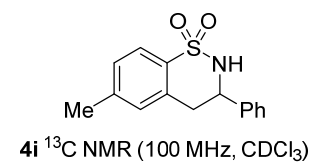
13C NMR BS-6-4C in CDCl₃

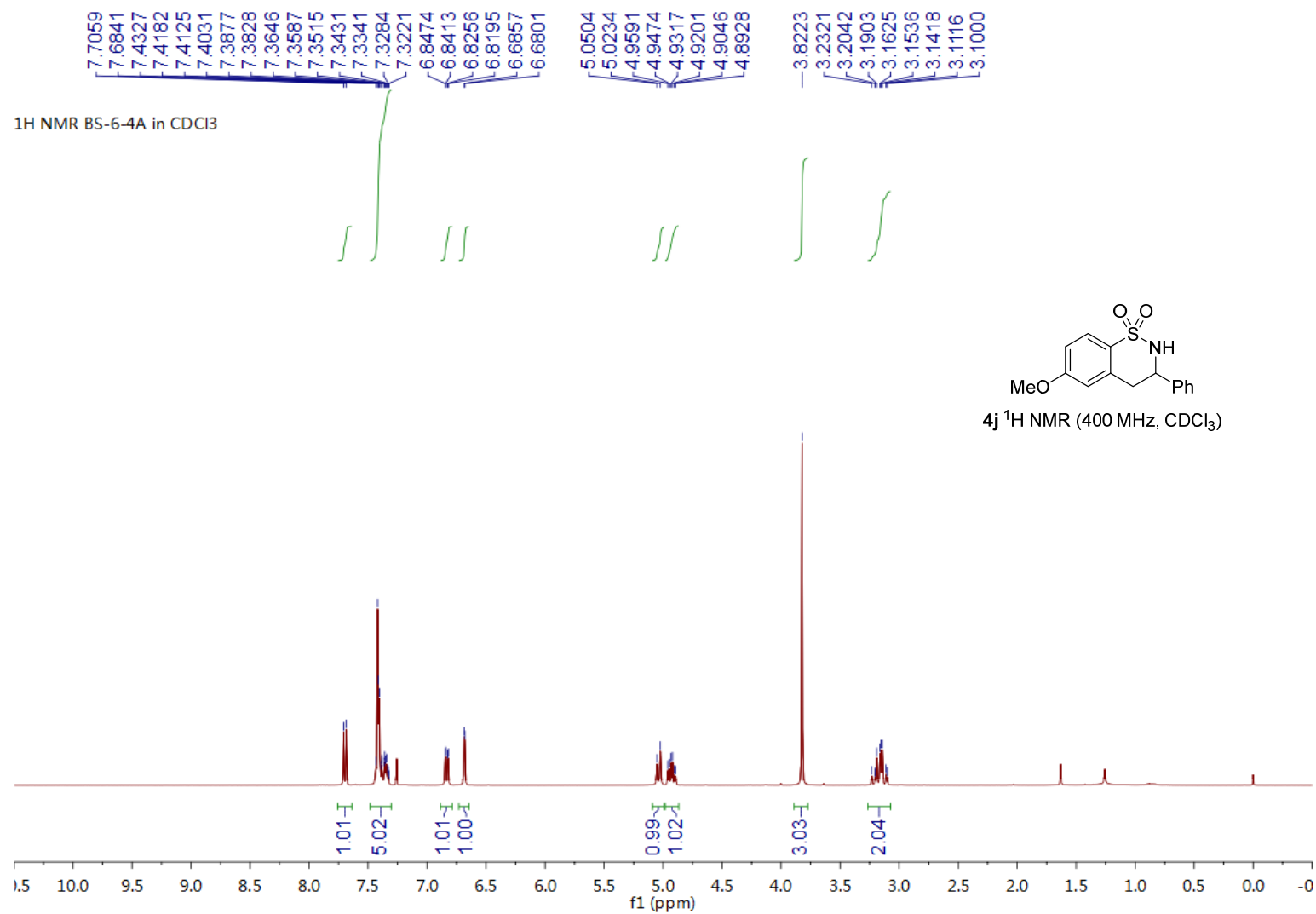
142.88
139.64
135.02
134.91
129.83
128.99
128.56
128.35
126.19
123.74

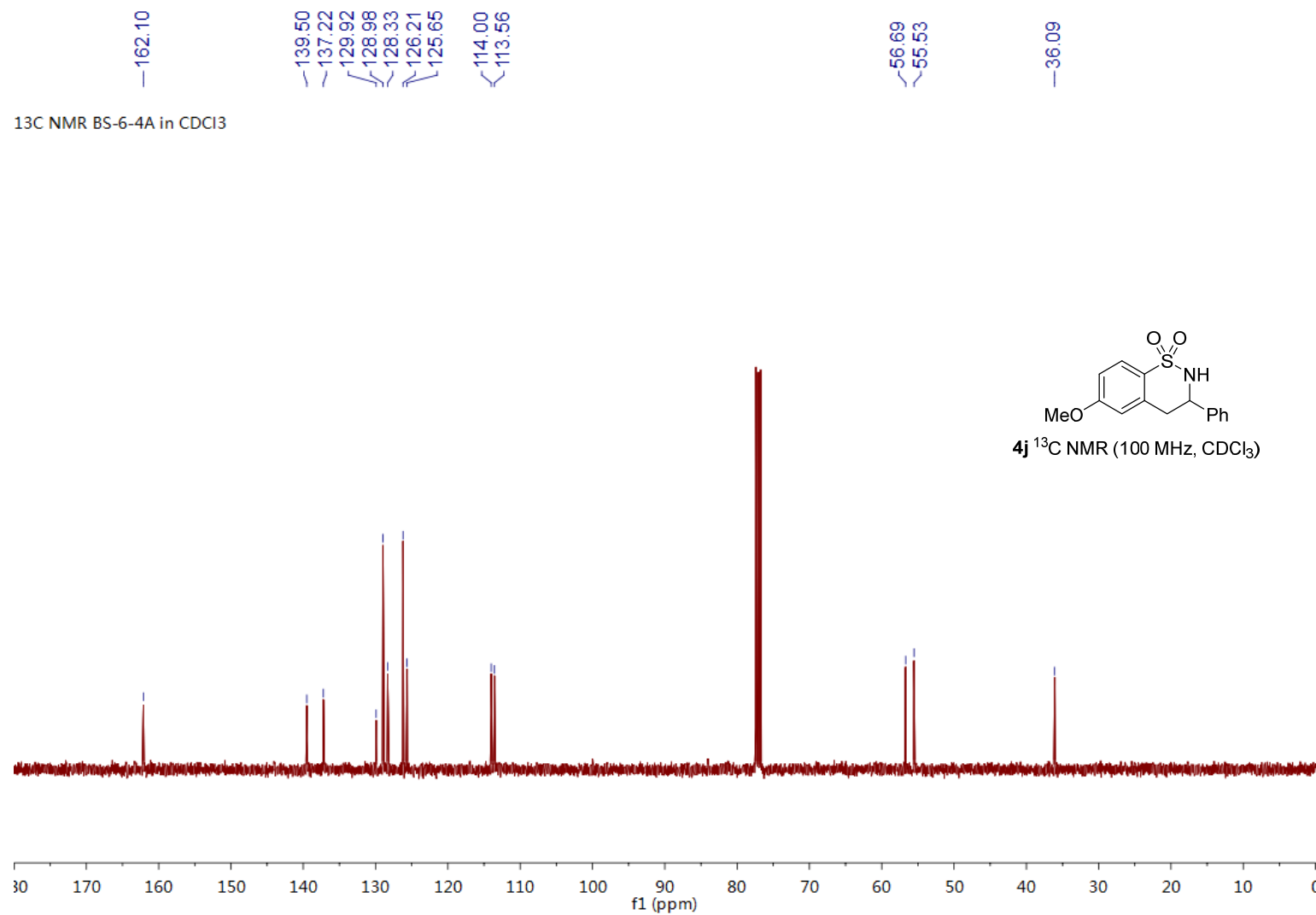
—56.79

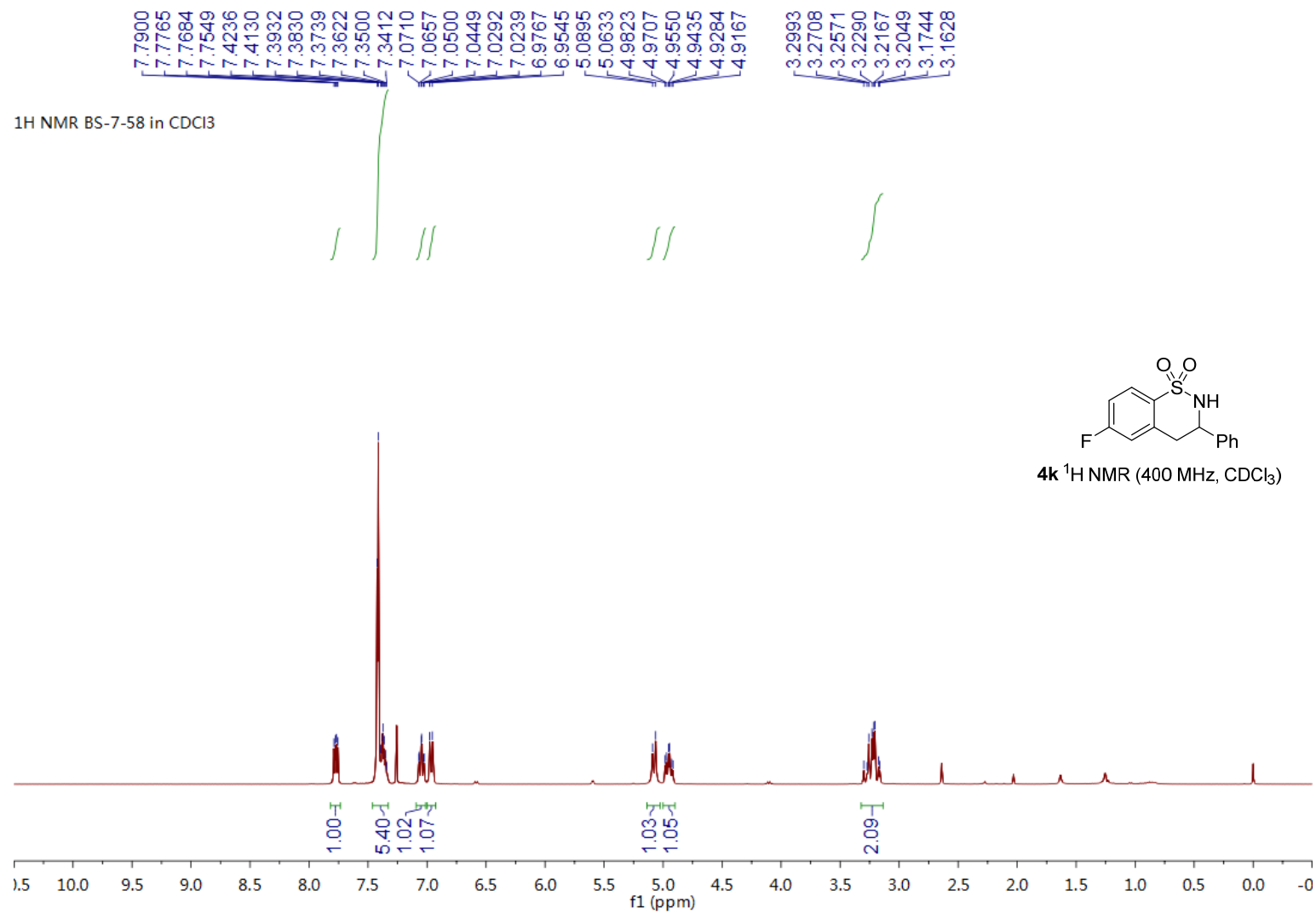
—35.65

—21.54







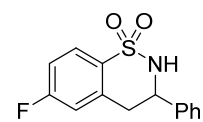


165.53
 163.00
 139.08
 138.29
 138.20
 133.92
 133.89
 129.09
 128.55
 126.49
 126.40
 126.19
 116.34
 116.12
 115.46
 115.24

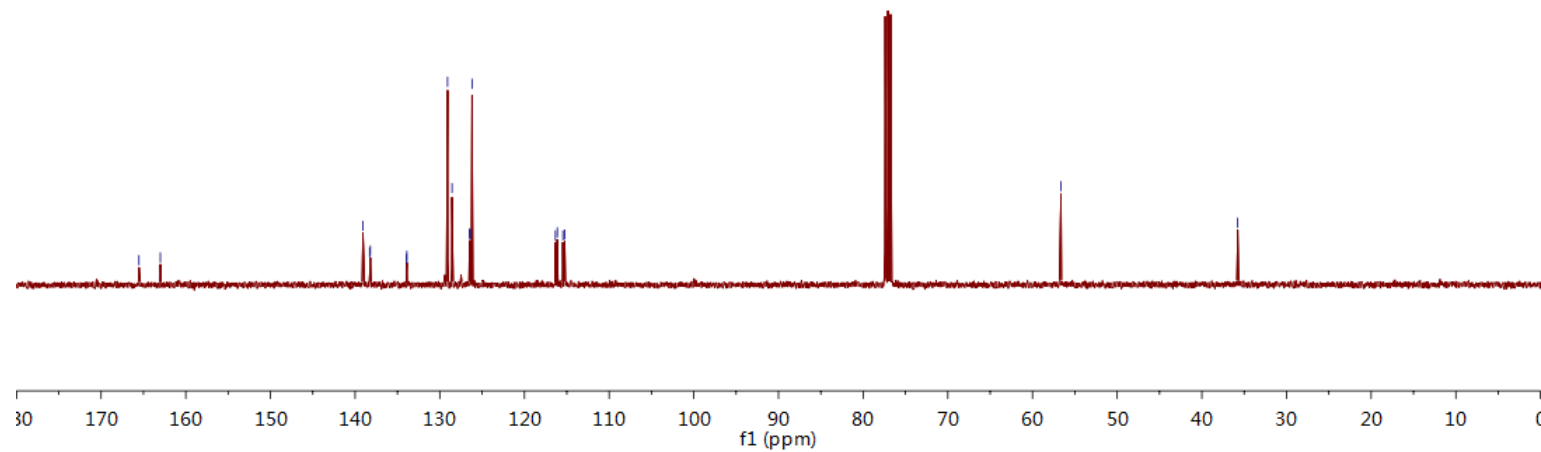
56.64

35.76

¹³C NMR BS-7-58 in CDCl₃

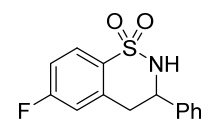


4k ¹³C NMR (100 MHz, CDCl₃)

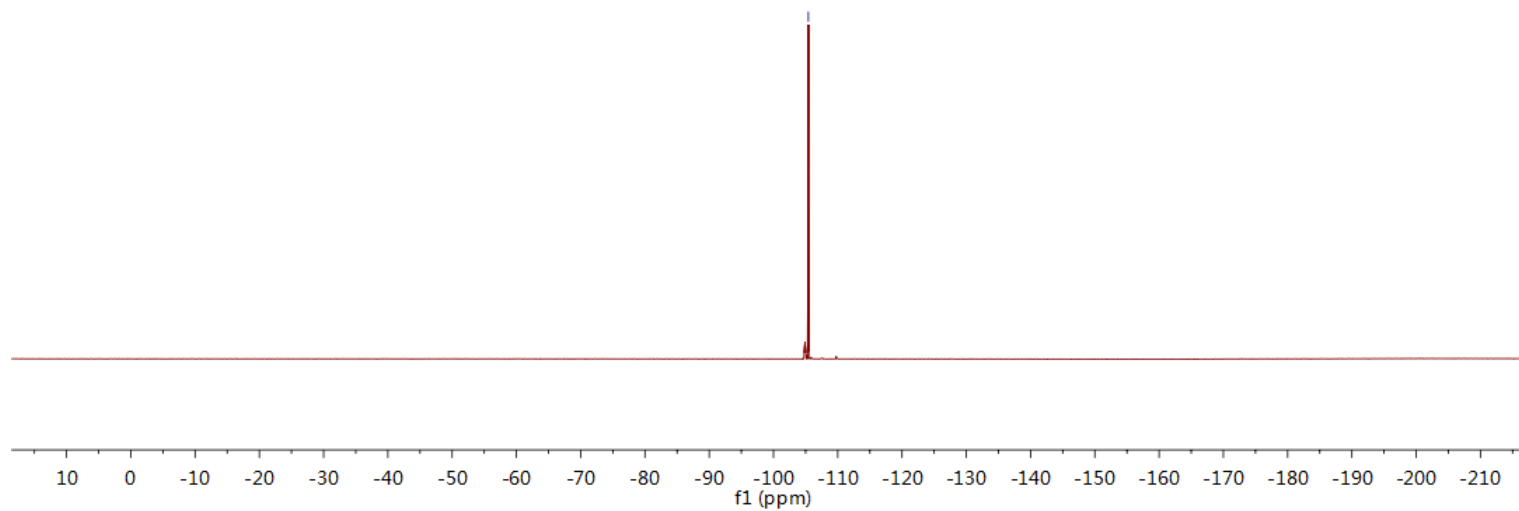


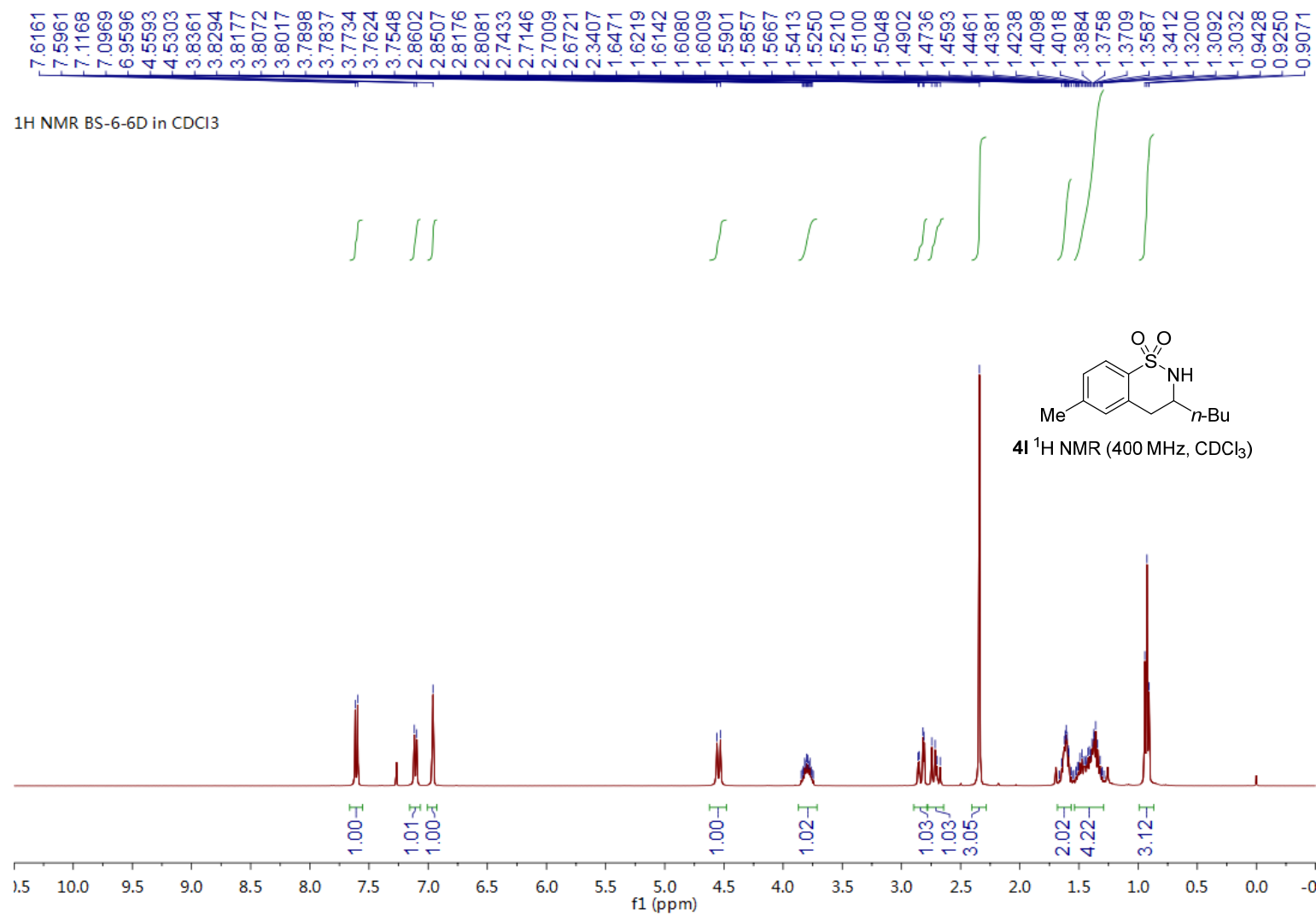
¹⁹F NMR BS-7-58 in CDCl₃

—105.4253



4k ¹⁹F NMR (376 MHz, CDCl₃)





¹³C NMR BS-6-4D in CDCl₃

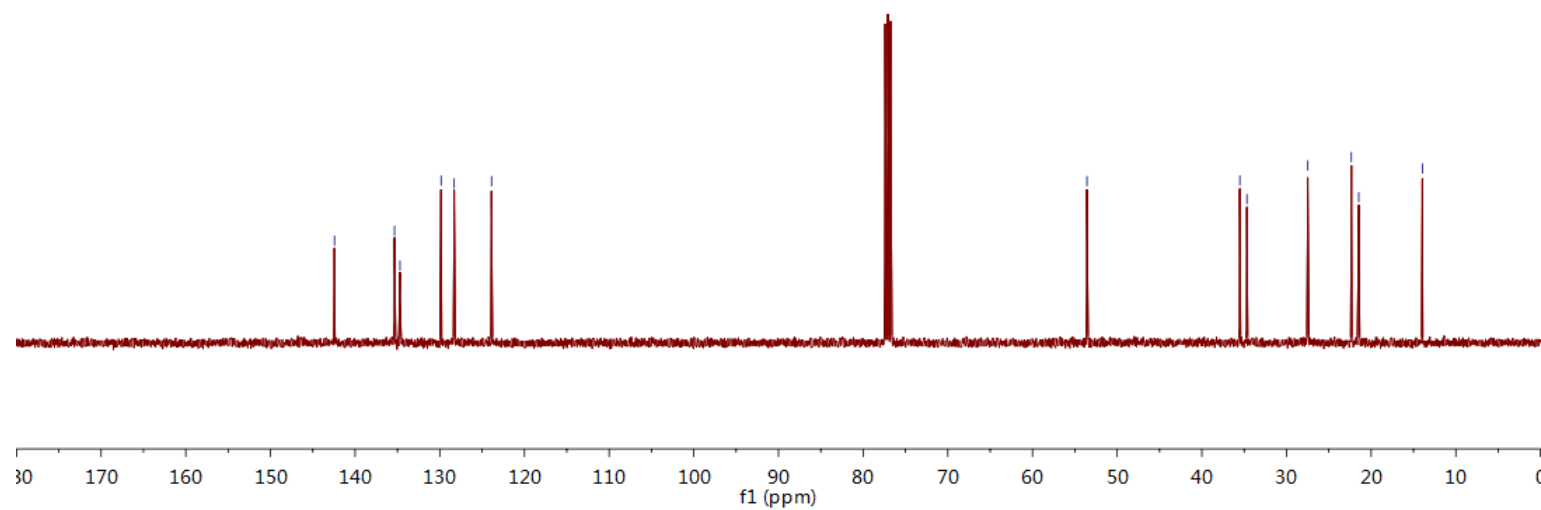
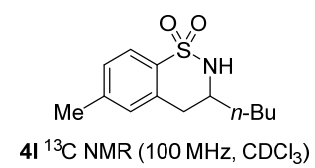
—142.44
 ~135.35
 ~134.70
 ~129.85
 ~128.30
 ~123.89

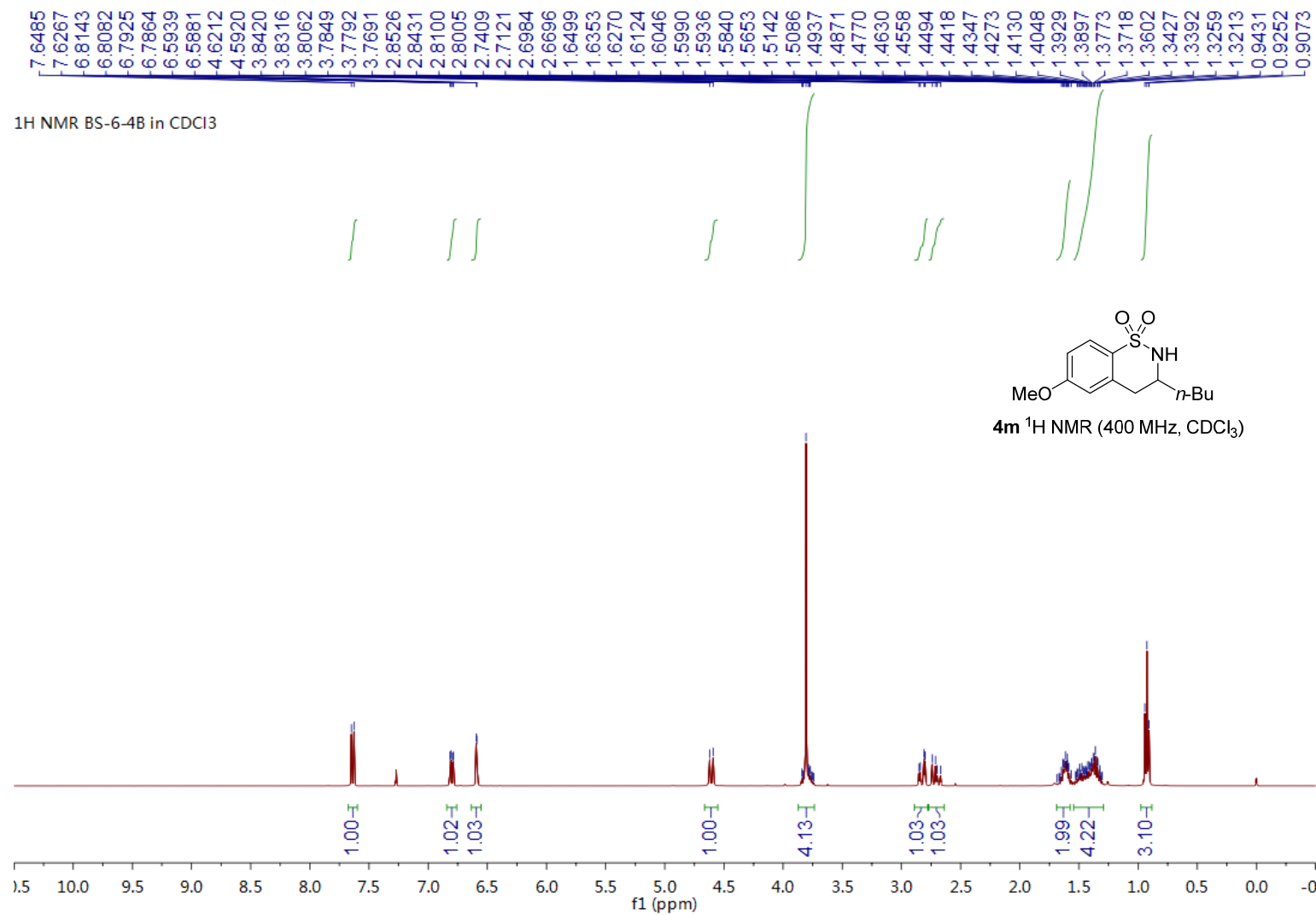
—53.57

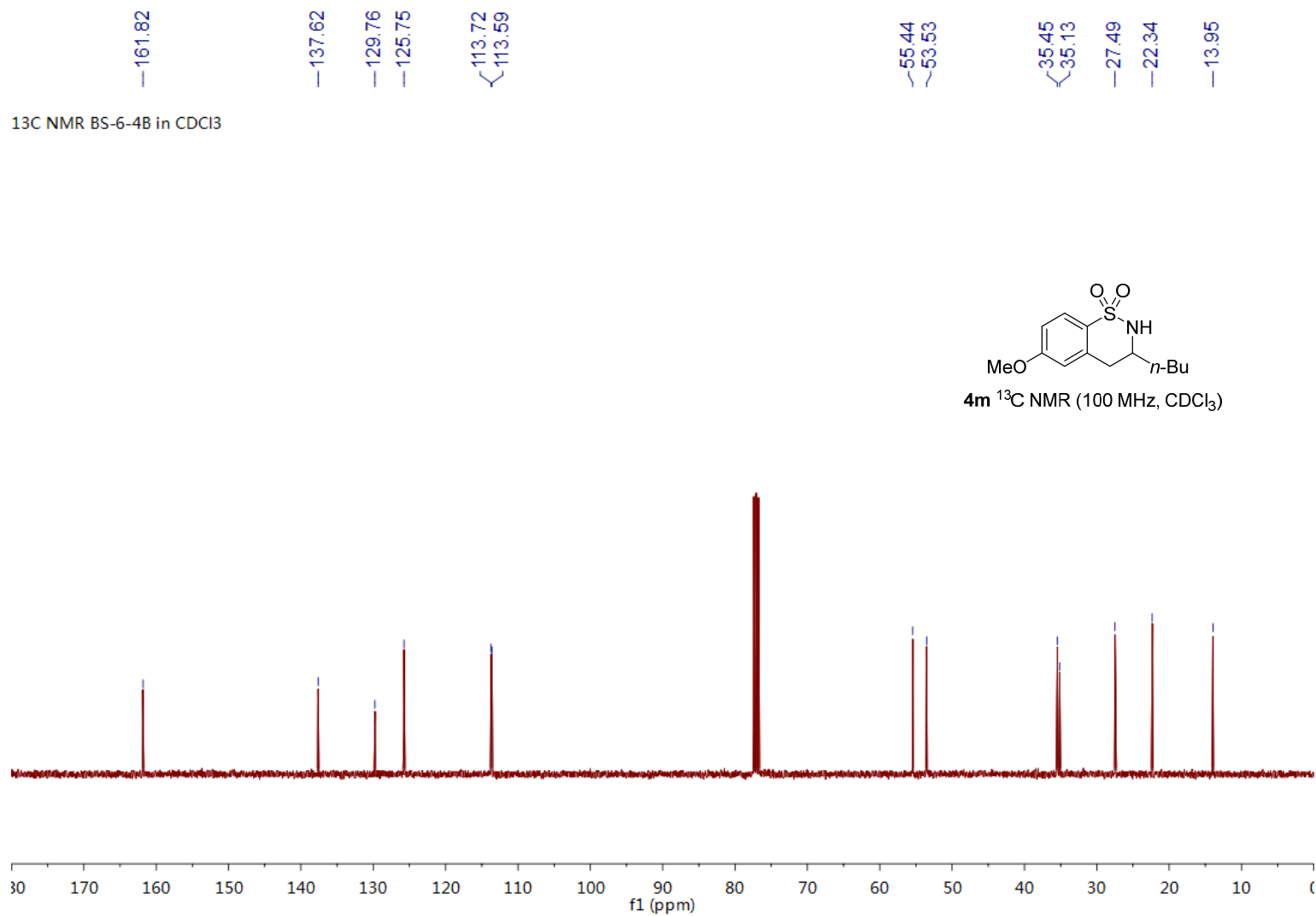
~35.51
 ~34.68

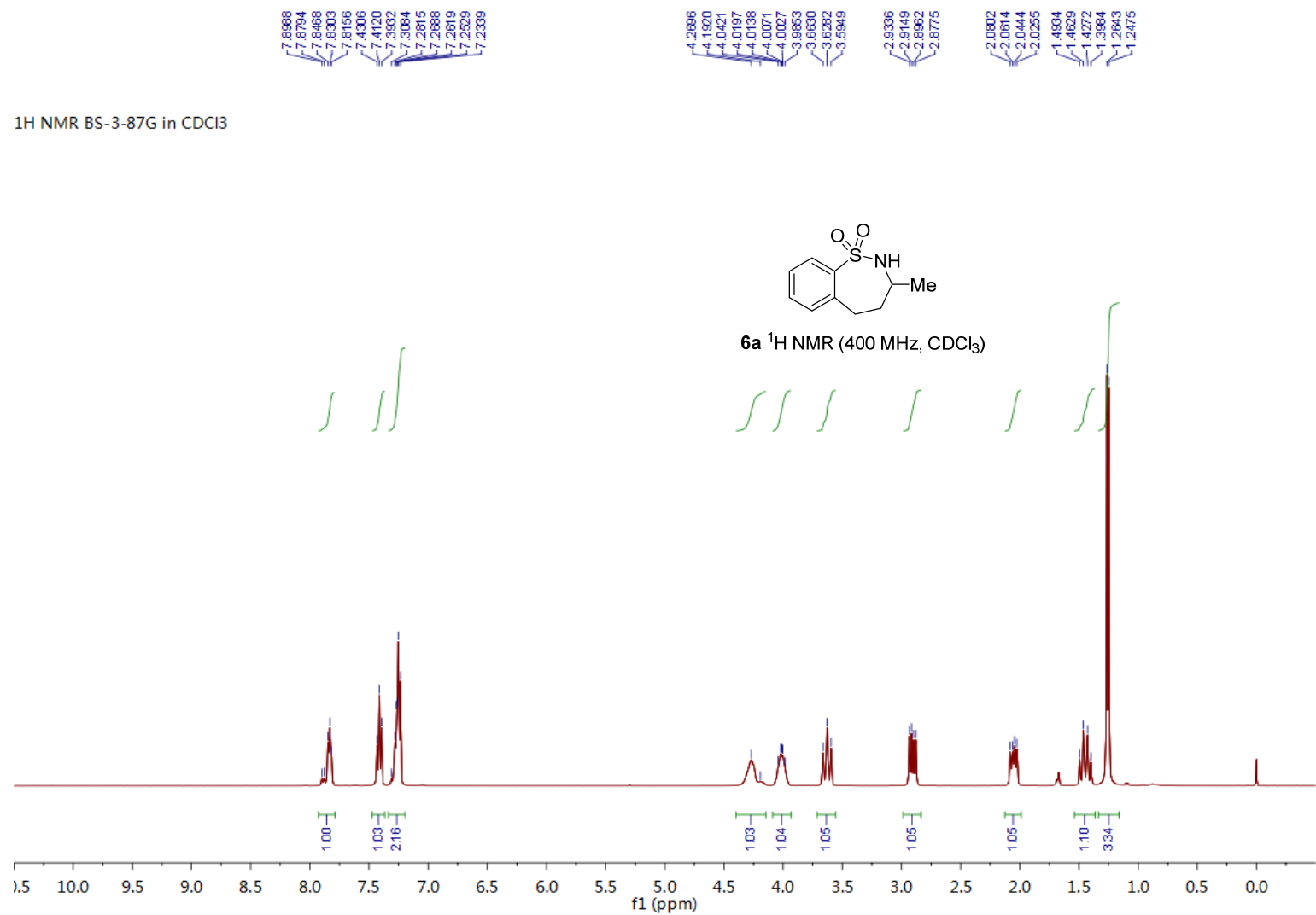
—27.48
 ~22.34
 ~21.47

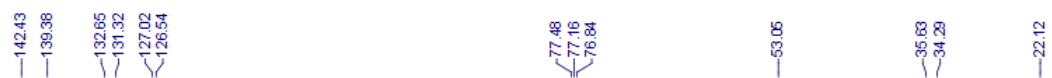
—13.96



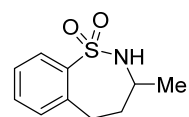




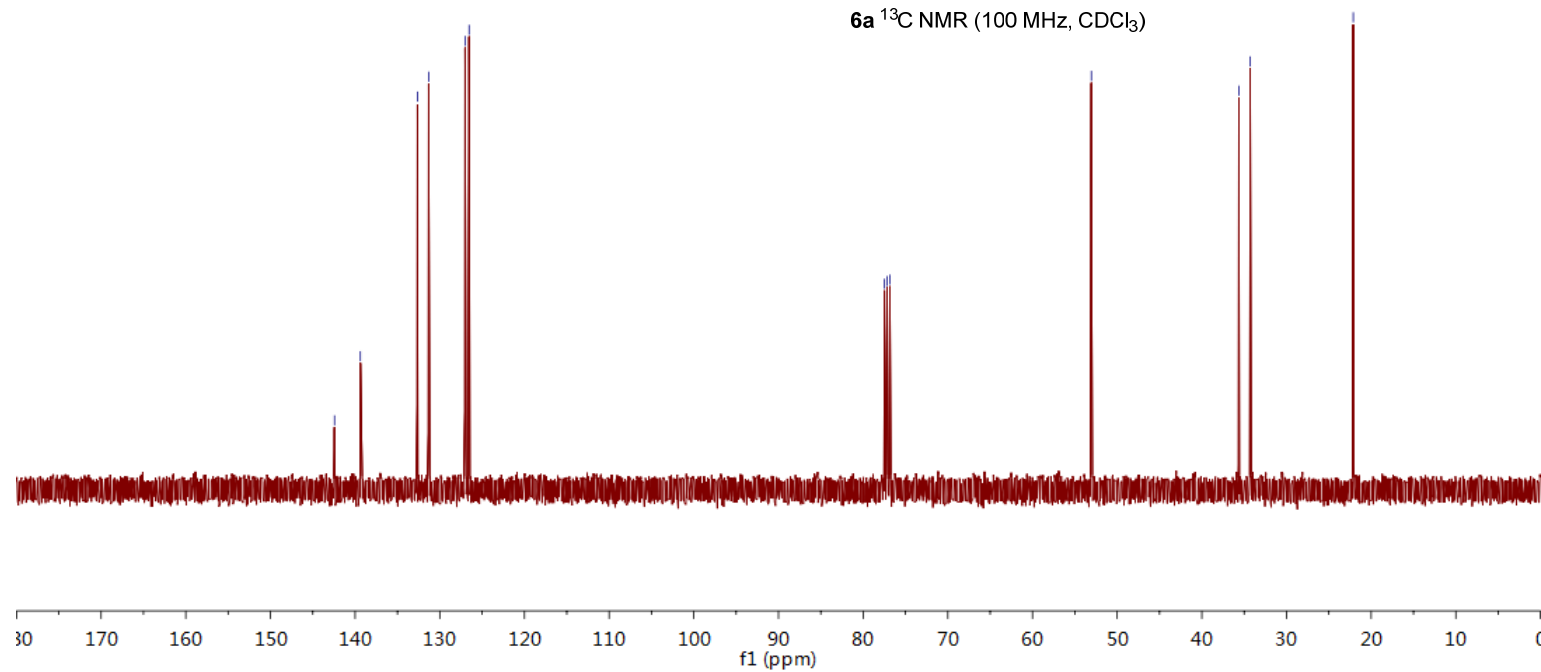




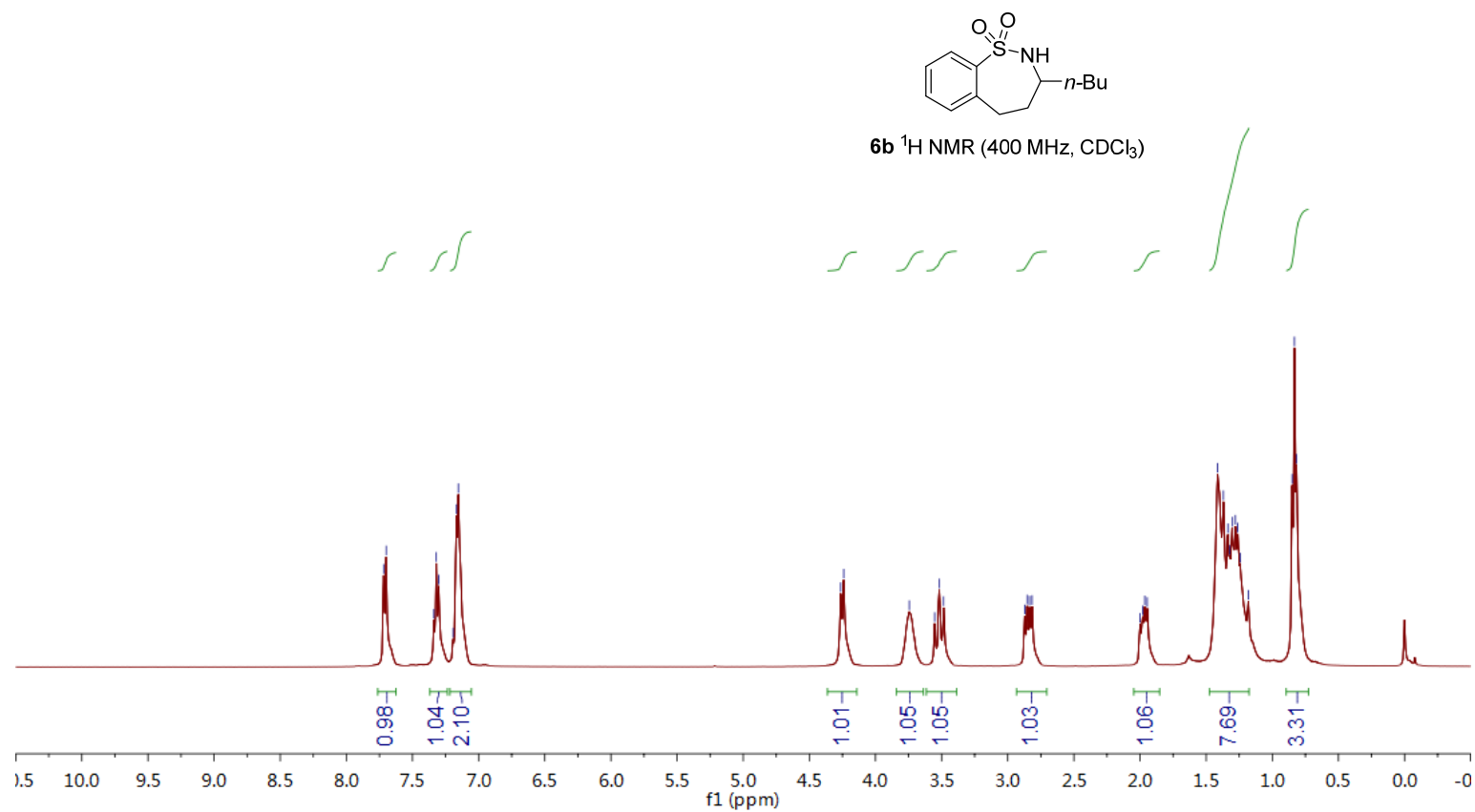
13C NMR BS-3-87G in CDCl₃



6a ¹³C NMR (100 MHz, CDCl₃)



¹H NMR BS-3-88A in CDCl₃

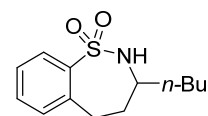


¹³C NMR BS-3-88A in CDCl₃

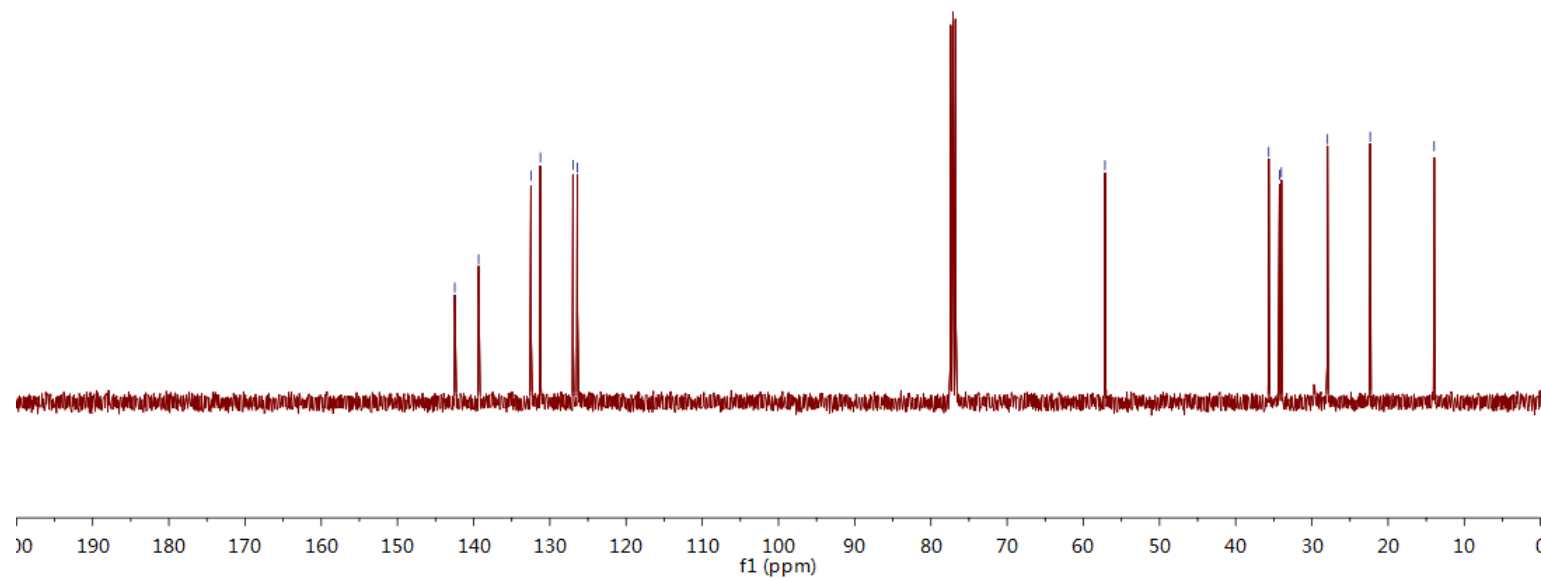
—142.48
—139.35
—132.48
—131.24
—126.97
—126.40

—57.15

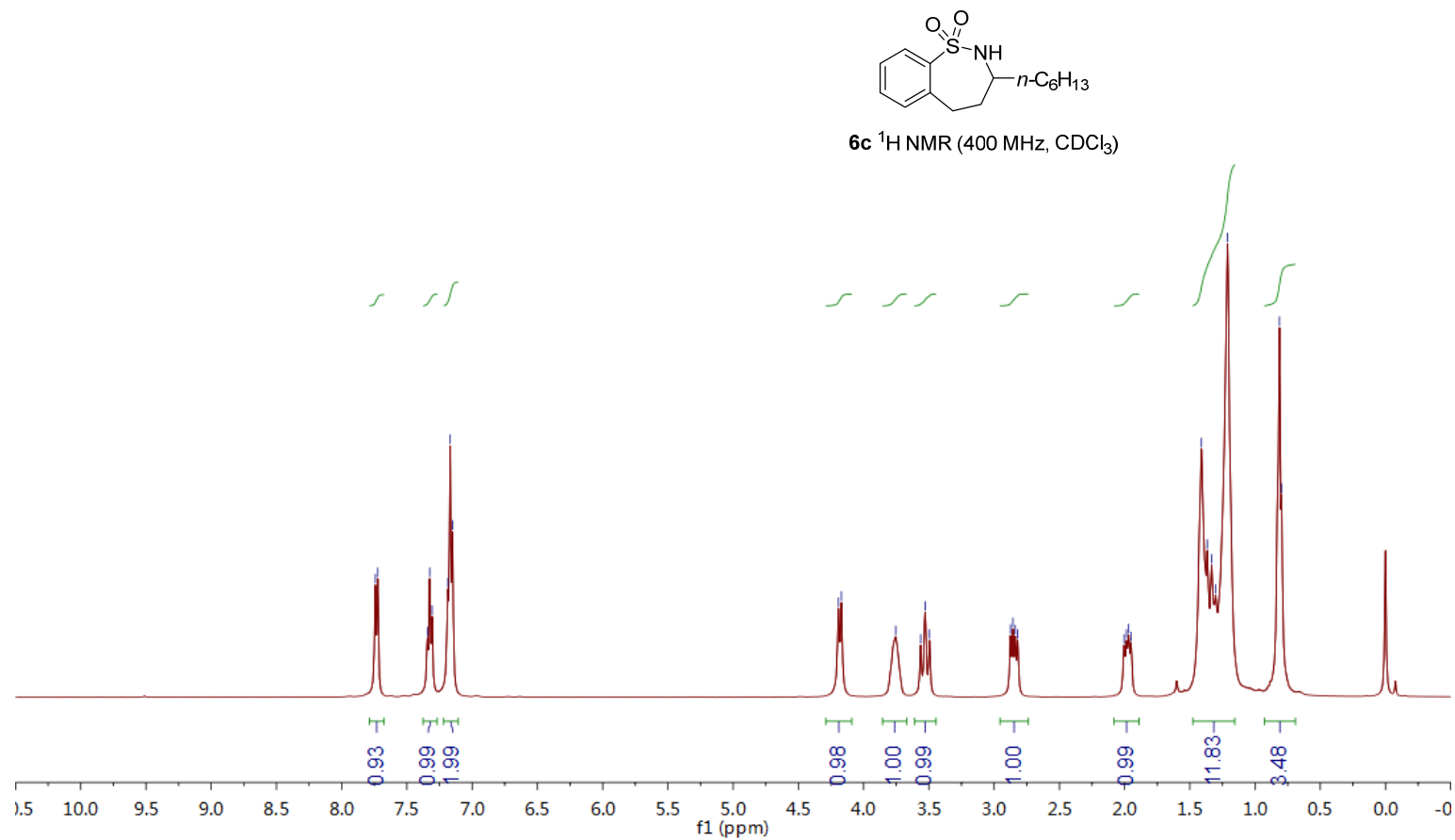
—35.67
—34.26
—34.01
—27.96
—22.35
—13.96



6b ¹³C NMR (100 MHz, CDCl₃)



¹H NMR BS-3-88B in CDCl₃

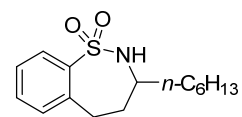


¹³C NMR BS-3-88B in CDCl₃

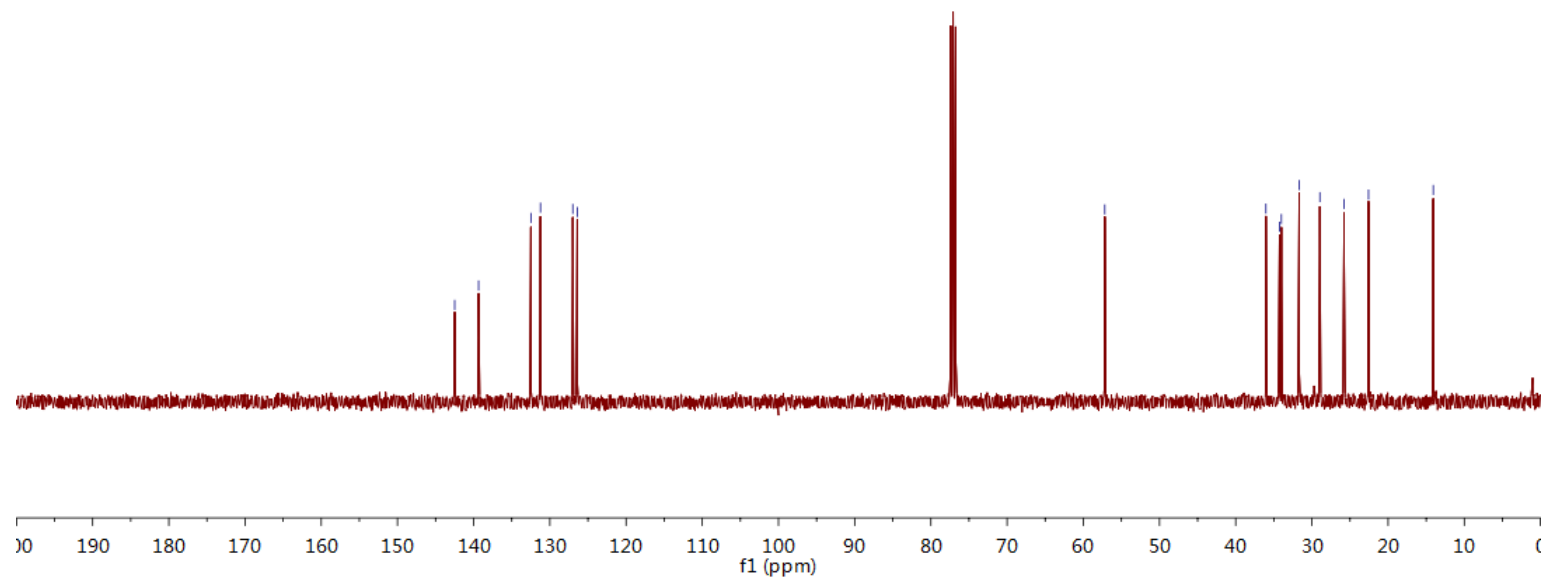
142.49
139.35
132.49
131.23
126.99
126.41

57.18

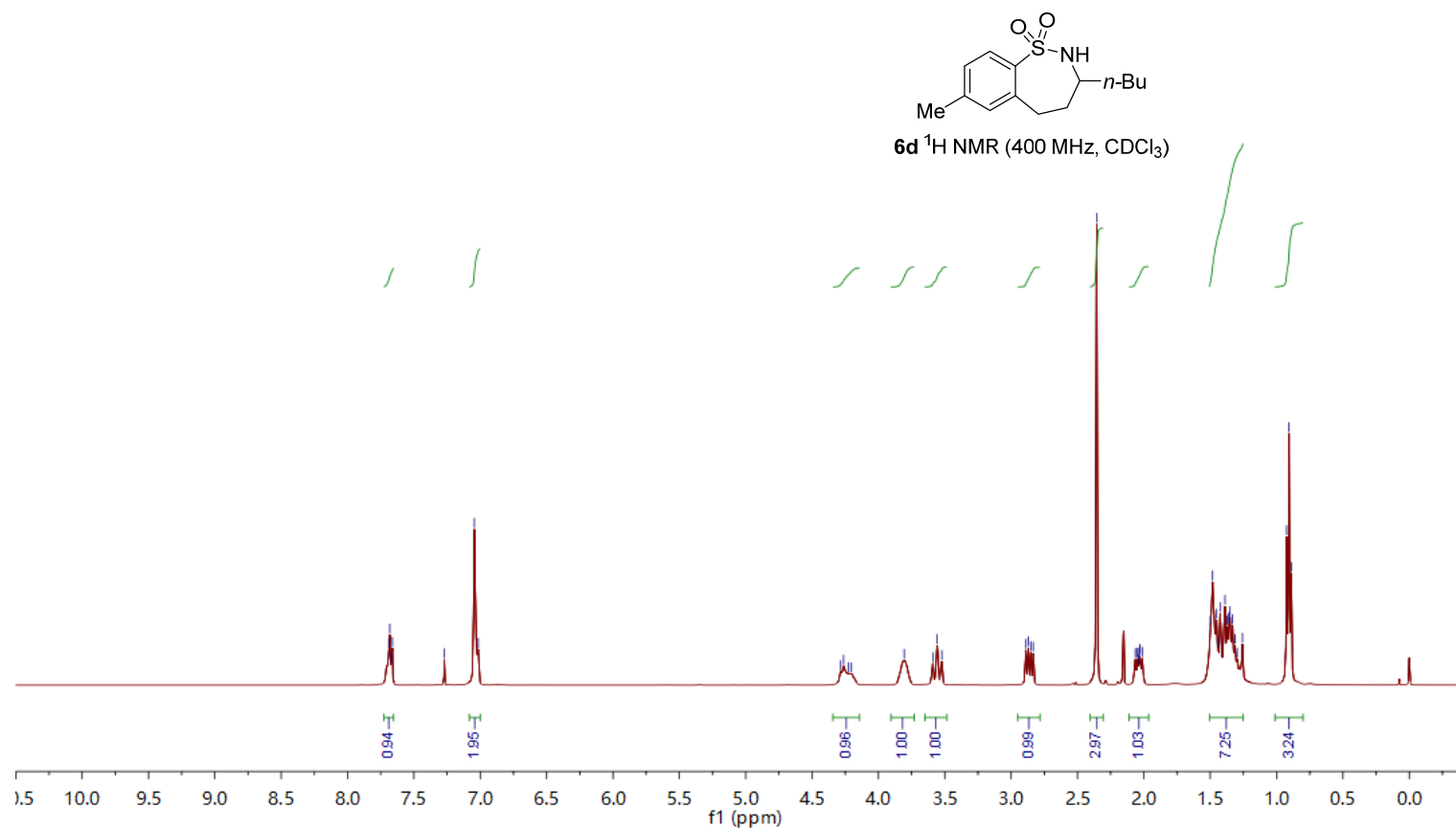
36.03
34.26
34.00
31.68
28.96
25.78
22.57
14.08



6c ¹³C NMR (100 MHz, CDCl₃)



¹H NMR BS-4-12A in CDCl₃



143.09
139.75
139.26
132.07
127.20
126.85

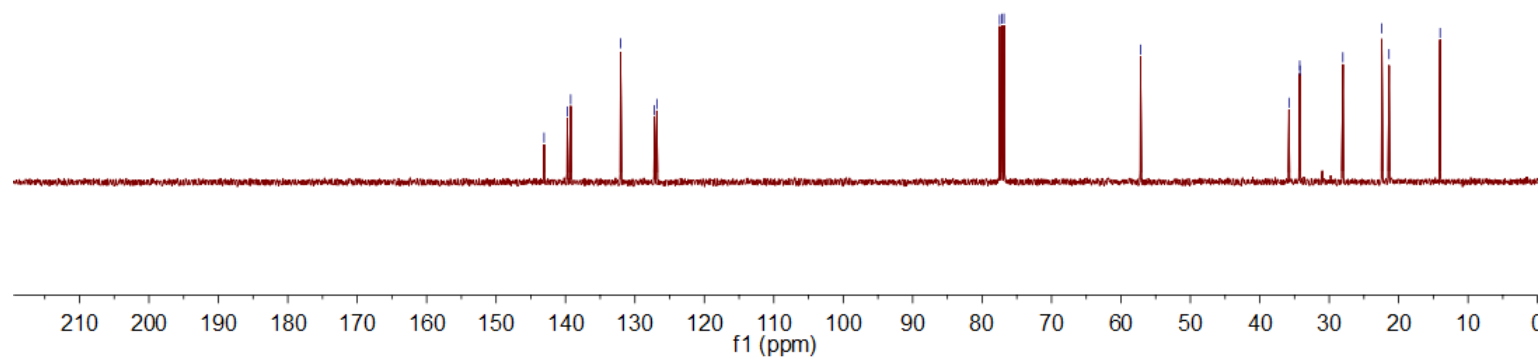
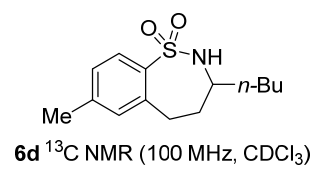
77.48
77.16
76.84

57.18

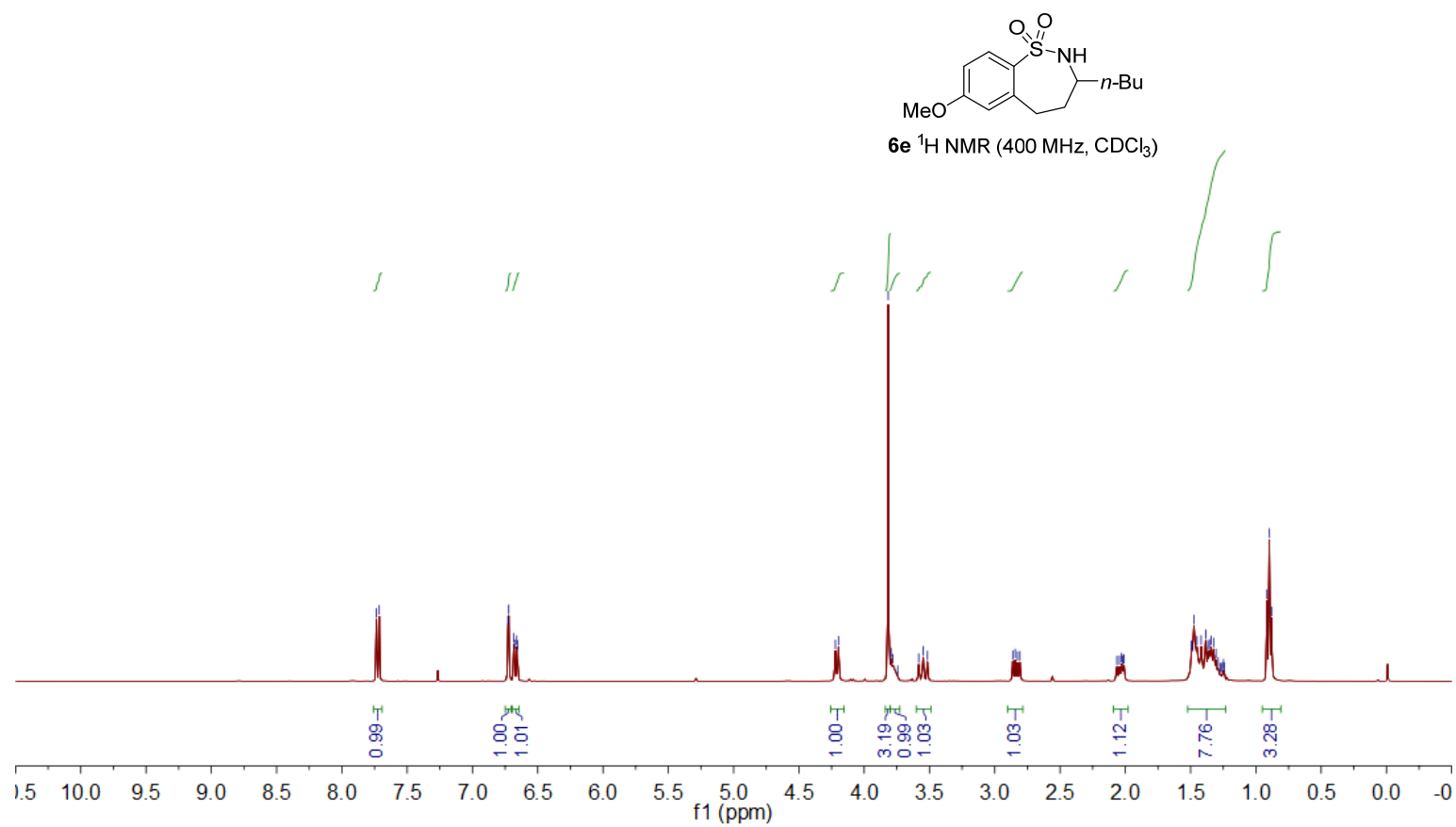
35.79
34.26
34.22
28.04
22.43
21.39

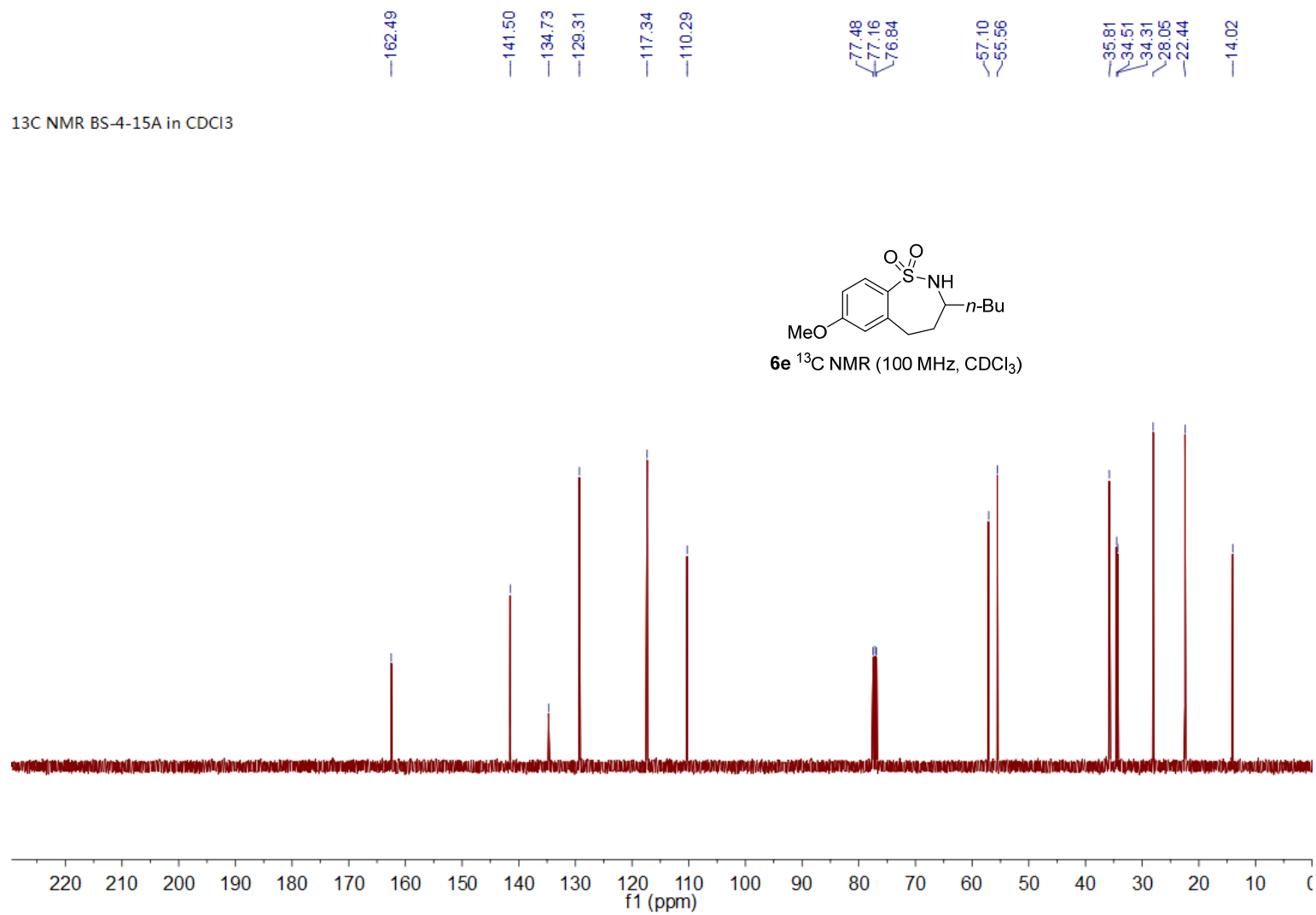
14.03

¹³C NMR BS-4-12A in CDCl₃

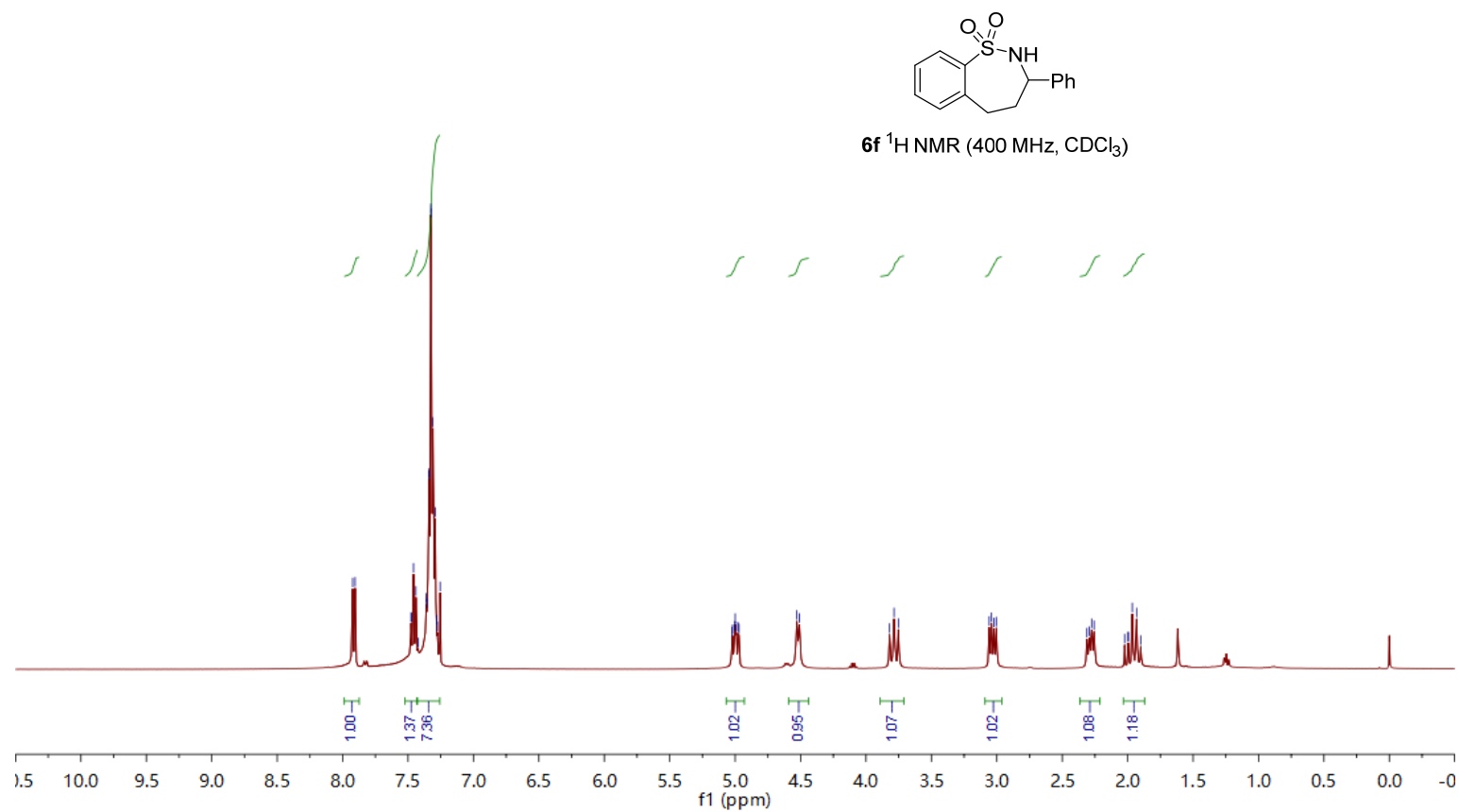


¹H NMR BS-3-15A in CDCl₃





¹H NMR BS-3-100 in CDCl₃

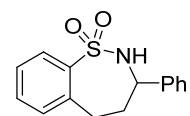


142.11
140.50
138.87
132.76
131.58
128.99
128.53
126.89
126.68
126.63

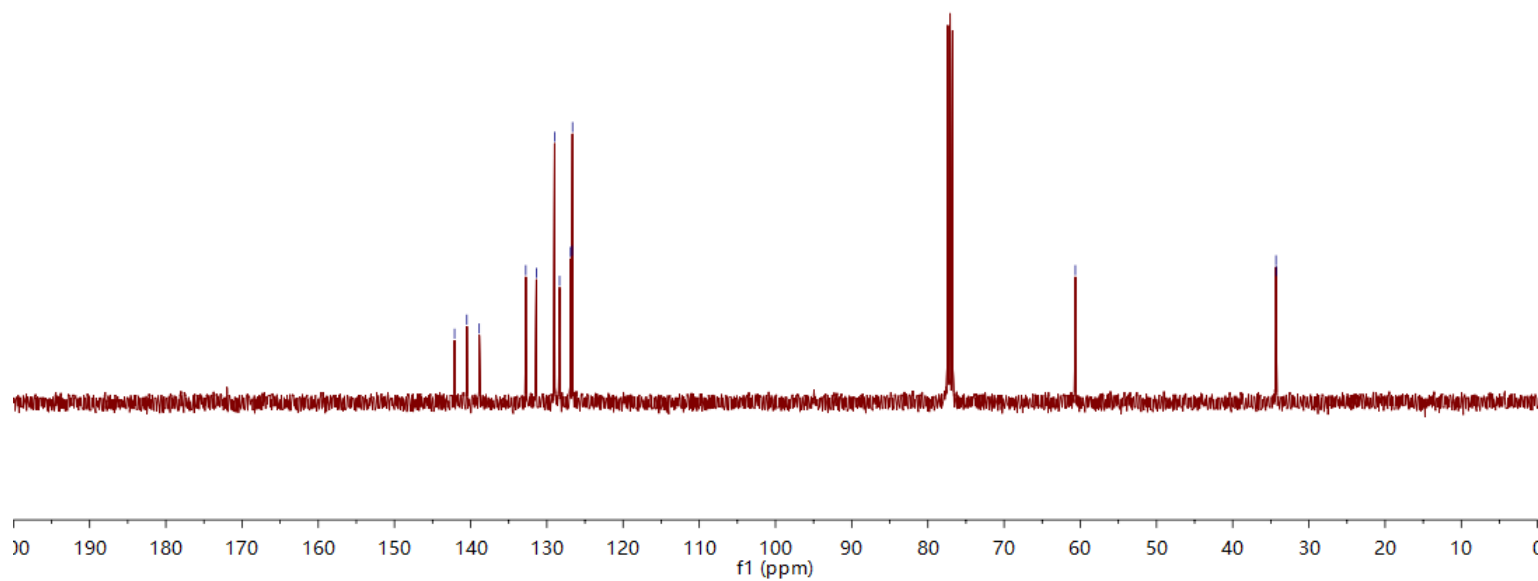
60.65

34.32
34.30

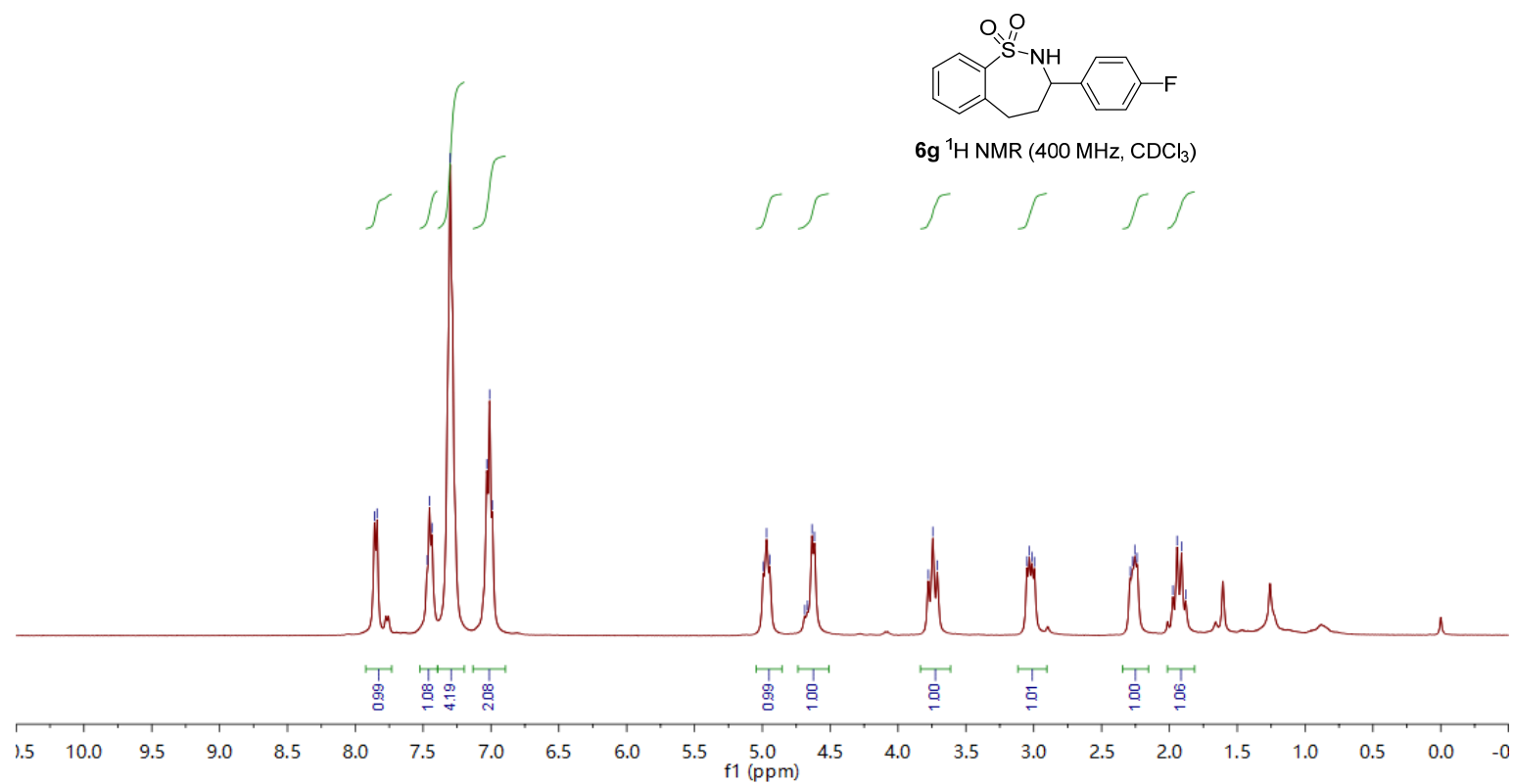
^{13}C NMR BS-3-100 in CDCl_3



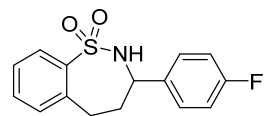
6f ^{13}C NMR (100 MHz, CDCl_3)



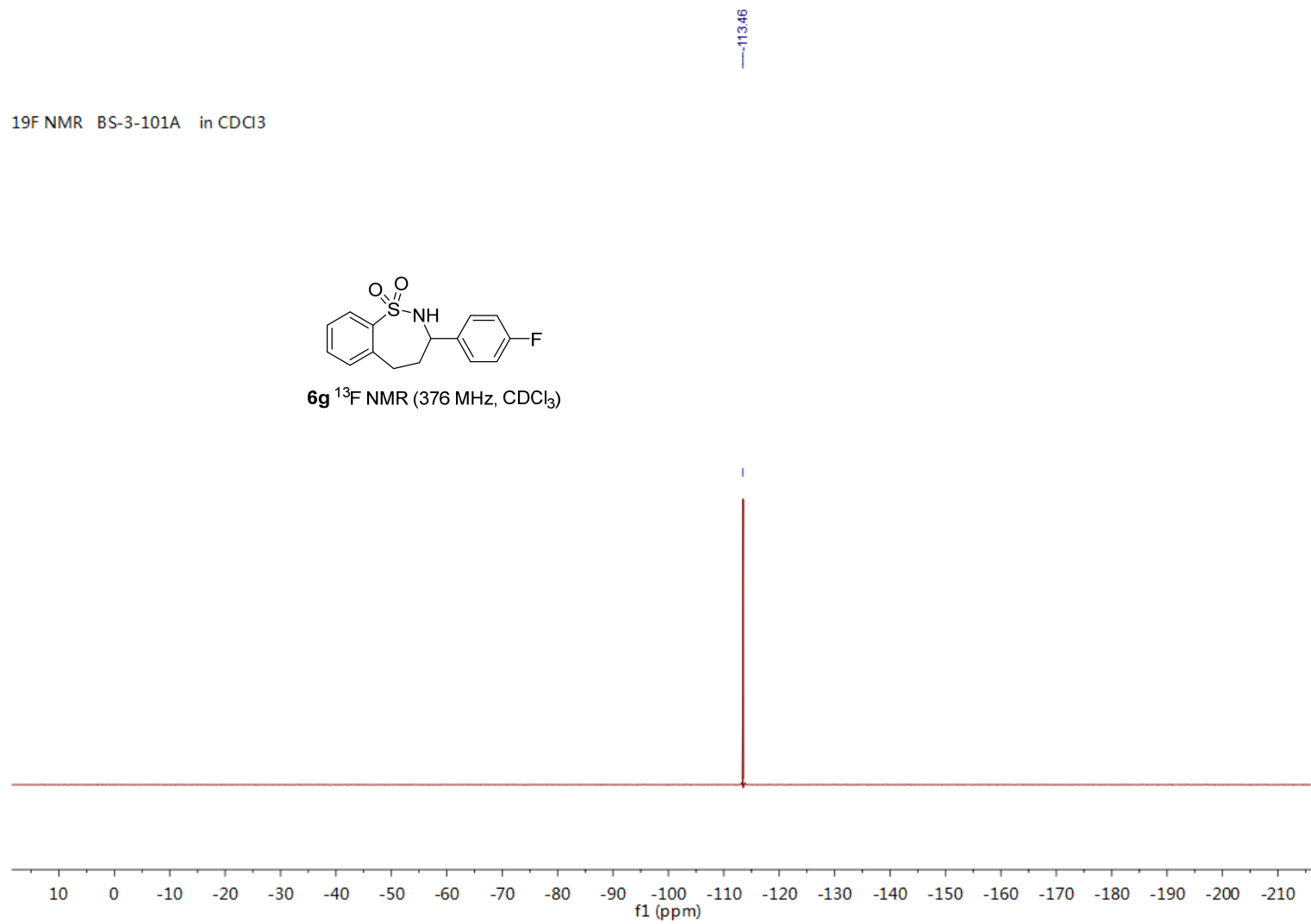
¹H NMR BS-3-101A in CDCl₃

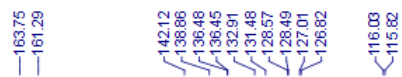


¹⁹F NMR BS-3-101A in CDCl₃

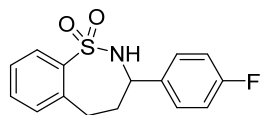


6g ¹³F NMR (376 MHz, CDCl₃)

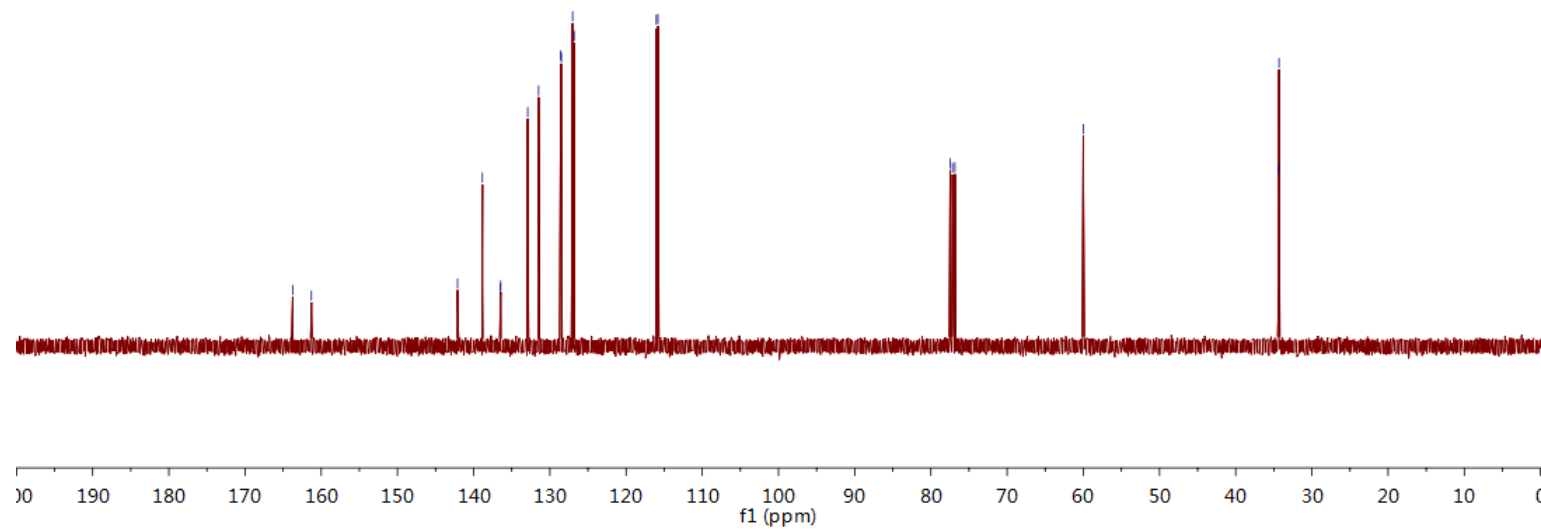
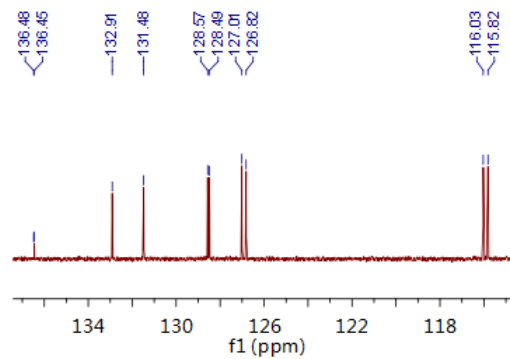




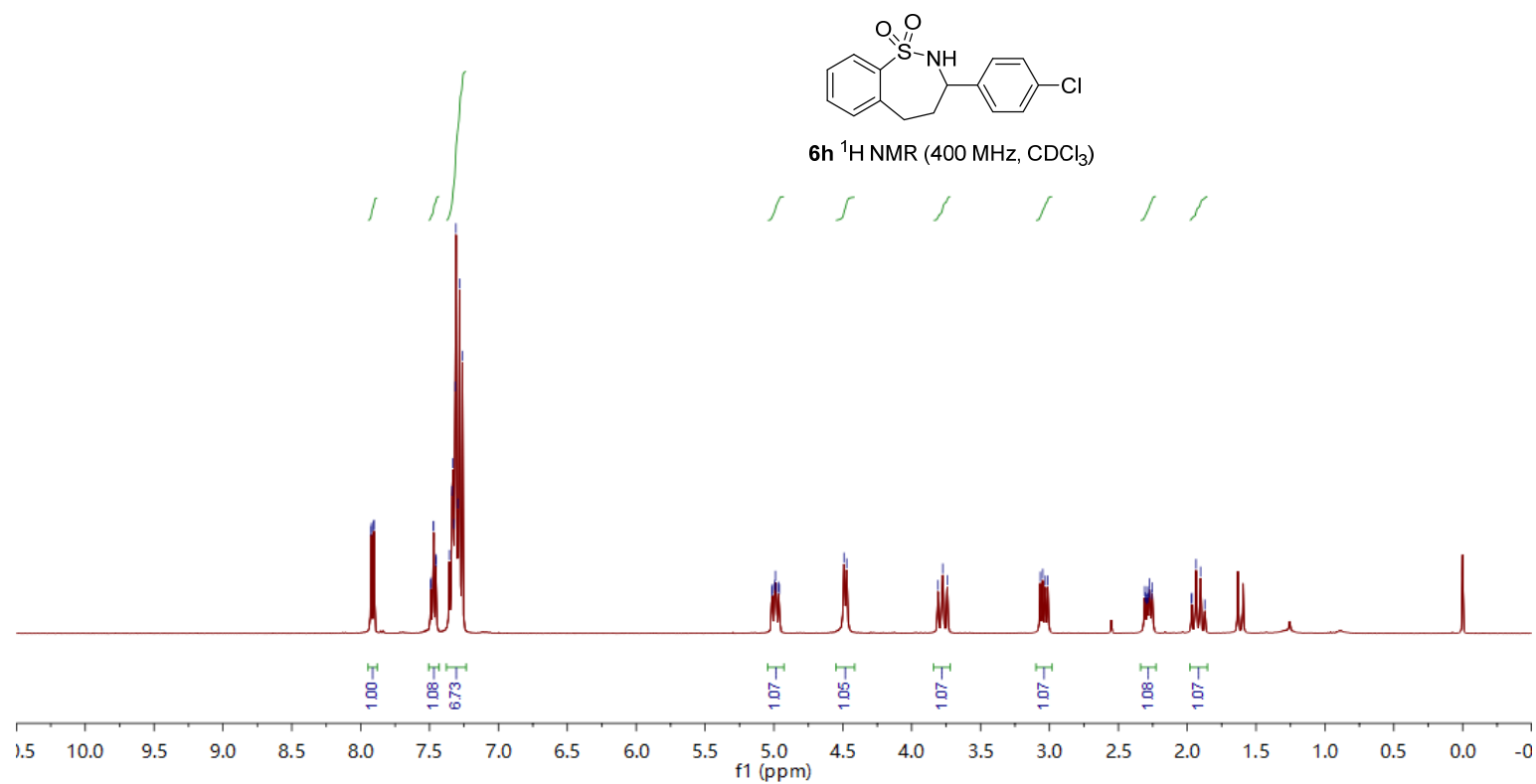
¹³C NMR BS-3-101A in CDCl₃



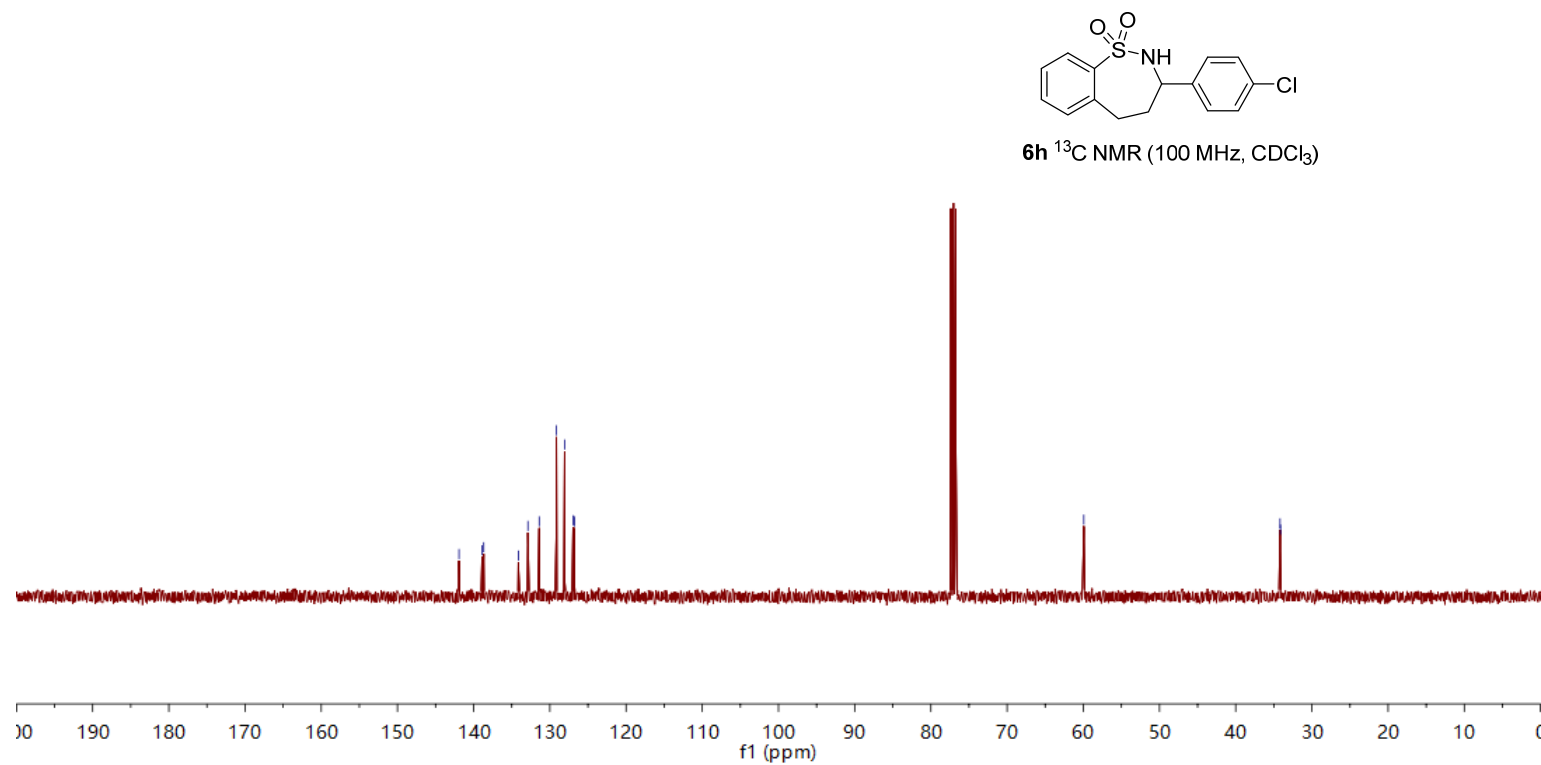
6g ¹³C NMR (100 MHz, CDCl₃)



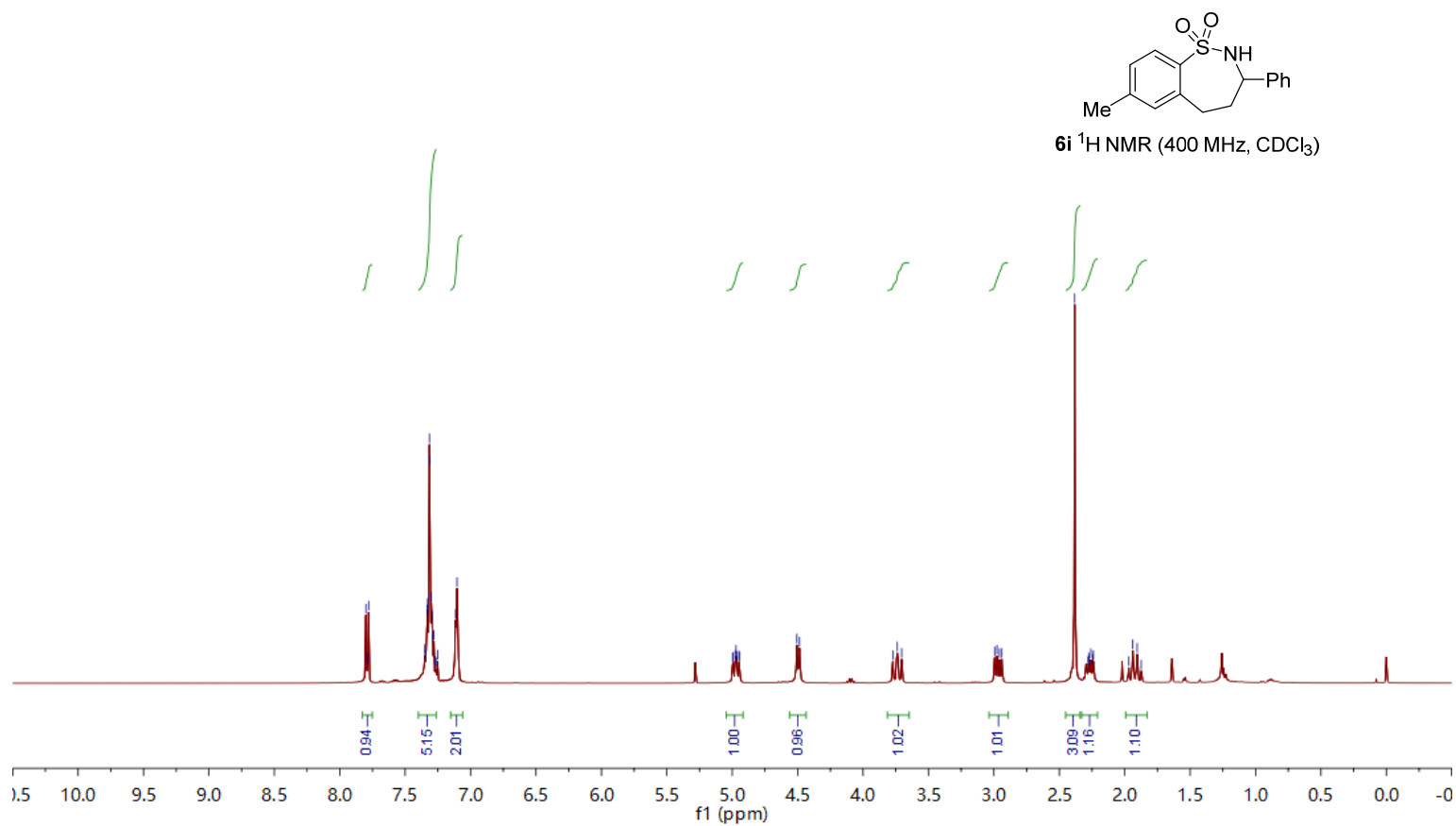
¹H NMR BS-3-93B in CDCl₃



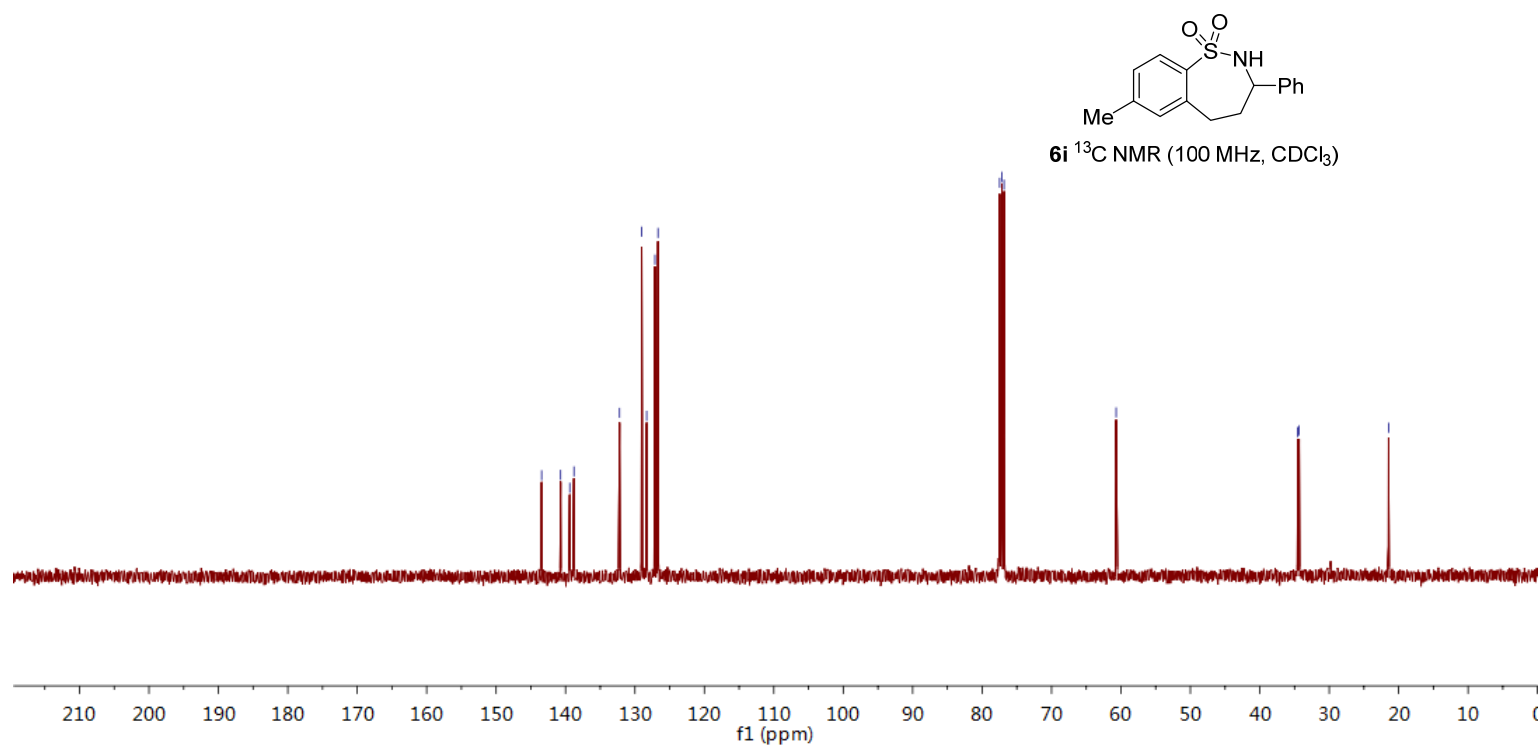
¹³C NMR BS-39B in CDCl₃

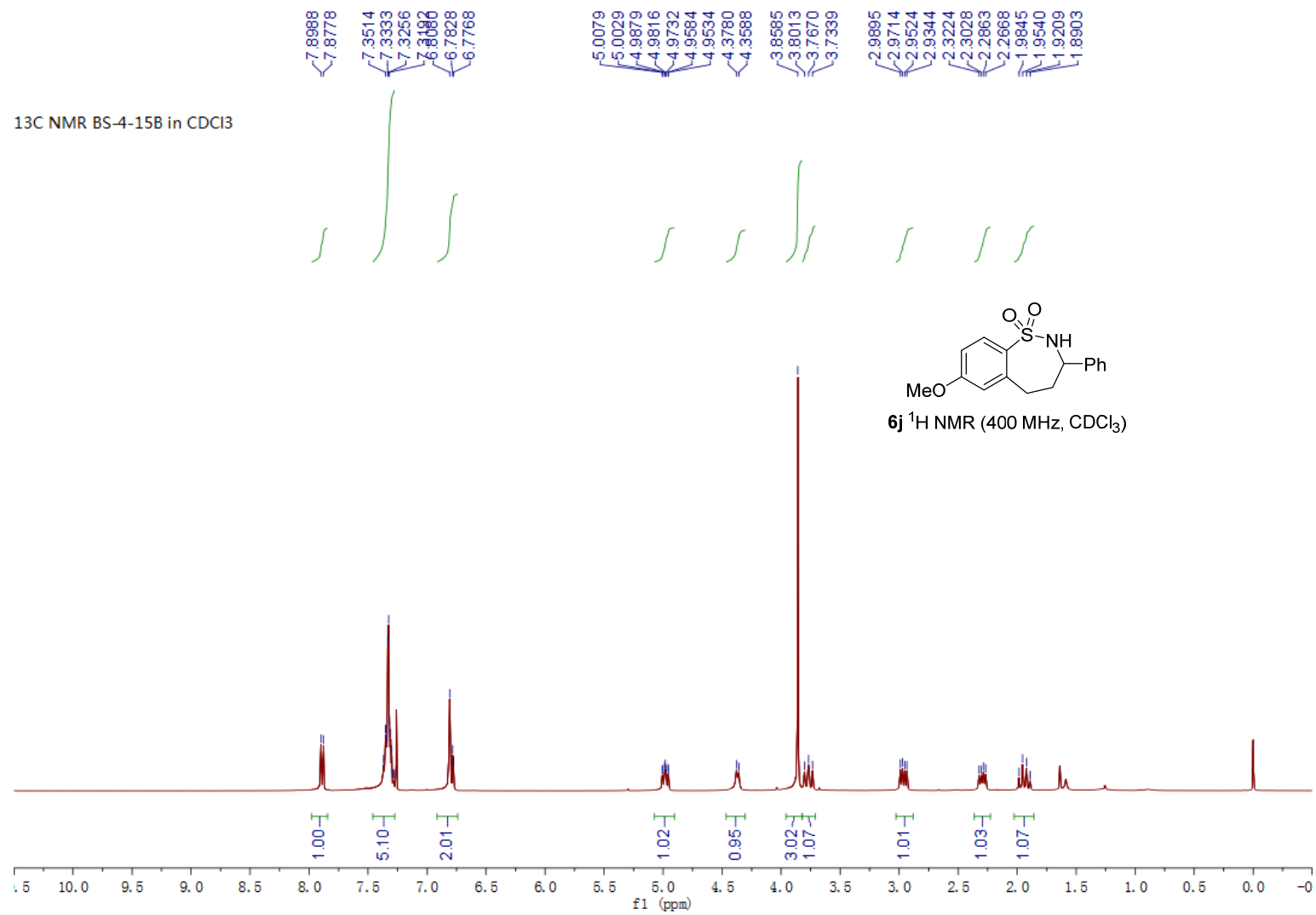


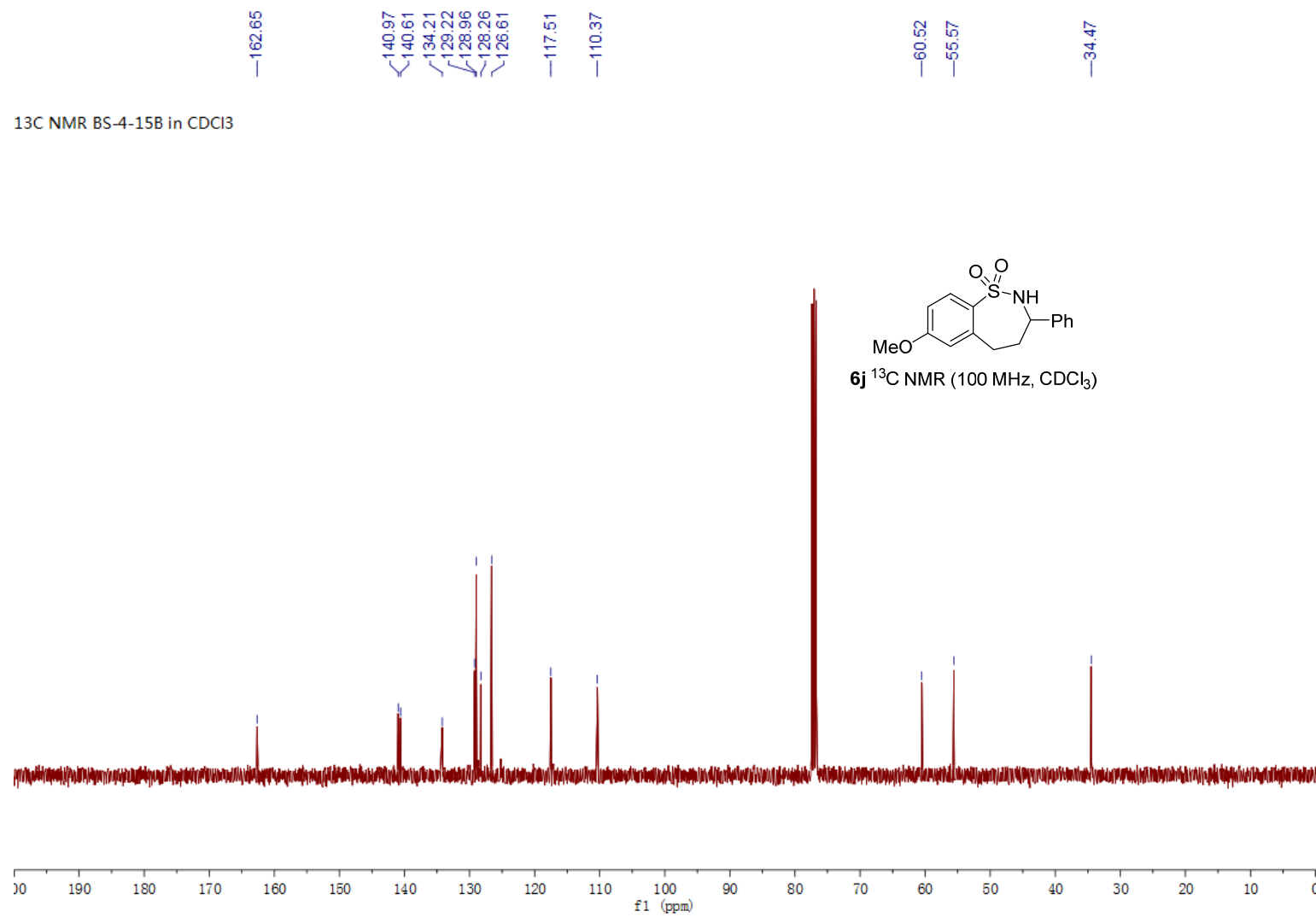
¹H NMR BS-4-12B in CDCl₃



¹³C NMR BS-4-12B in CDCl₃

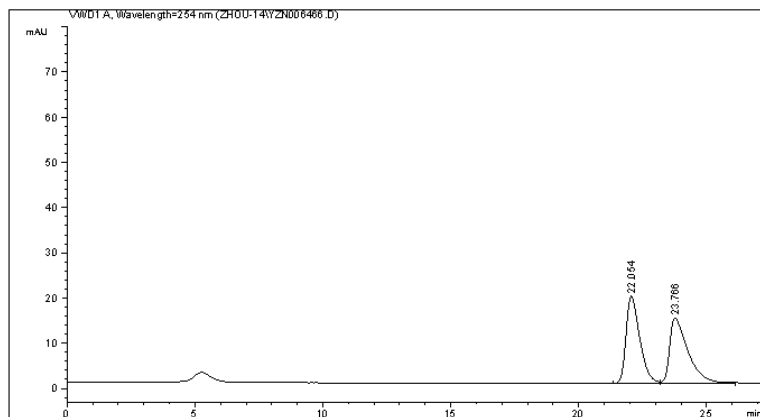






Data File C:\CHEM32\1\DATA\ZHOU-14\YZN006466.D
Sample Name: BS-3-51A(+/-)

```
=====
Acq. Operator   : Z
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 11/25/2014 10:38:33 PM
Acq. Method     : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 11/25/2014 10:38:15 PM by Z
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 7/21/2015 9:57:57 AM by
                  (modified after loading)
Sample Info     : 0J-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 254 nm
=====
```



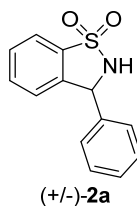
```
=====
Area Percent Report
=====
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=254 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area [mAU]	Area %
1	22.054	BV	0.5637	713.97363	19.25210	50.2596	
2	23.766	VB	0.7292	706.59686	14.41231	49.7404	

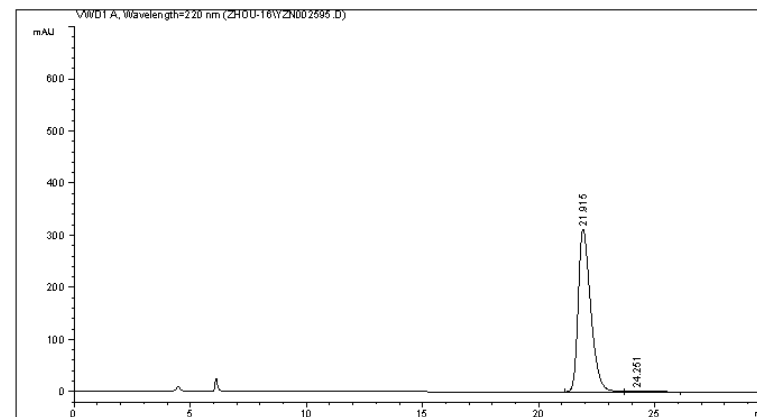
Totals : 1420.57050 33.66440

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



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN002595.D
Sample Name: BS-3-51A

```
=====
Acq. Operator   : 0
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 10/13/2016 3:51:12 PM
Acq. Method     : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 10/13/2016 3:45:28 PM by 0
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 10/19/2016 2:17:49 PM by 0
                  (modified after loading)
Sample Info     : 0J-H, Hexane/i-PrOH = 70/30, 0.7 mL/min, 30oC, 254 nm
=====
```



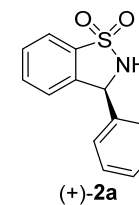
```
=====
Area Percent Report
=====
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area [mAU]	Area %
1	21.915	BV	0.5691	1.16454e4	312.24945	99.1750	
2	24.251	VB	0.9524	96.87186	1.39080	0.8250	

Totals : 1.17423e4 313.64025

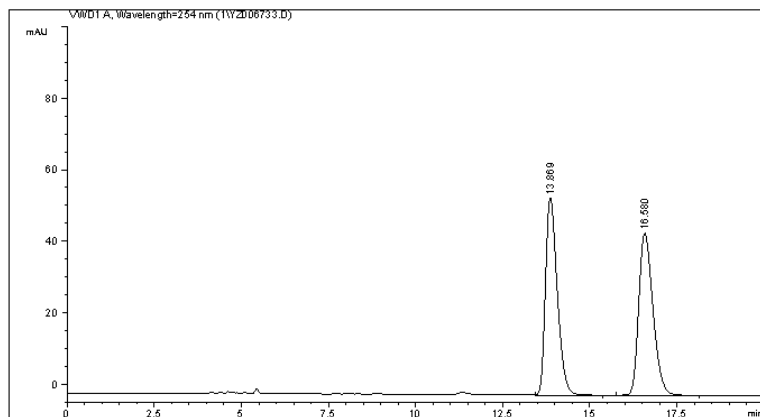
```
=====
*** End of Report ***
```



Data File C:\CHEM32\1\DATA\1\YZ006733.D
Sample Name: BS-3-55C(+/-)

=====

Acq. Operator	: ZHOU	Location	: Vial 1
Acq. Instrument	: Instrument 1		
Injection Date	: 12/3/2014 1:31:37 AM		
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC1.M		
Last changed	: 12/3/2014 1:29:16 AM by ZHOU		
	(modified after loading)		
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M		
Last changed	: 7/21/2015 10:03:06 AM by		
	(modified after loading)		
Sample Info	: OD-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 254 nm		



Area Percent Report

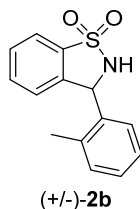
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=254 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area %
1	13.869	BB	0.3647	1300.08997	55.26591	49.9516
2	16.580	BV	0.4420	1302.60791	45.46571	50.0484

Totals : 2602.69788 100.73163

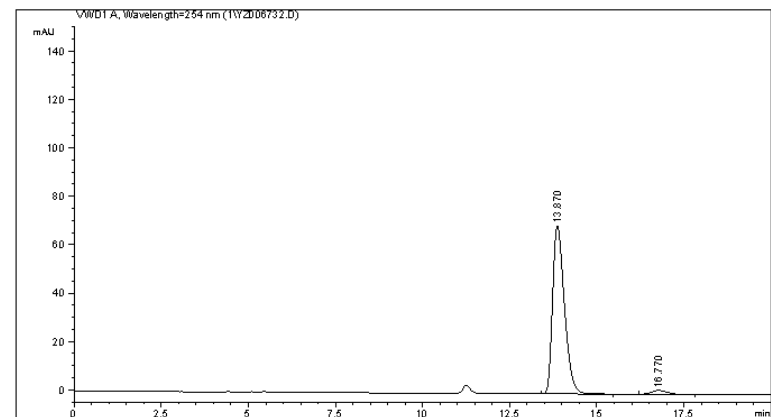
*** End of Report ***



Data File C:\CHEM32\1\DATA\1\YZ006732.D
Sample Name: BS-3-55C

=====

Acq. Operator	: ZHOU	Location	: Vial 1
Acq. Instrument	: Instrument 1		
Injection Date	: 12/3/2014 1:03:04 AM		
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC1.M		
Last changed	: 12/3/2014 12:41:15 AM by ZHOU		
	(modified after loading)		
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M		
Last changed	: 7/21/2015 10:03:28 AM by		
	(modified after loading)		
Sample Info	: OD-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 254 nm		



Area Percent Report

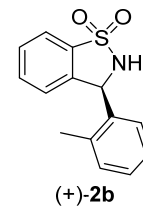
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=254 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area %
1	13.870	BB	0.3654	1638.64014	69.49314	97.0690
2	16.770	BB	0.4805	49.47792	1.58581	2.9310

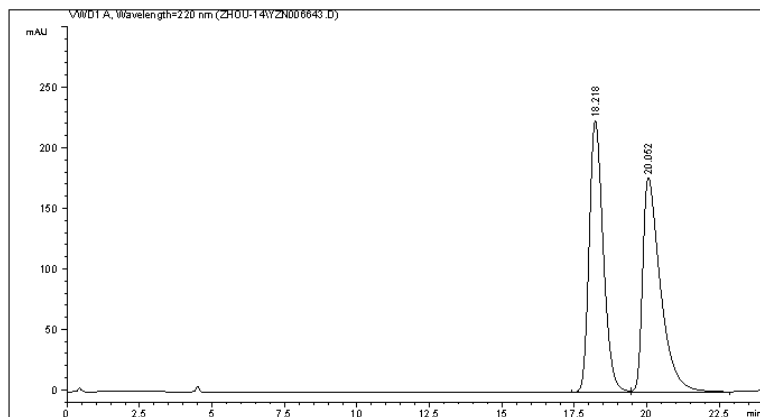
Totals : 1688.11805 71.07895

*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-14\YZN006643.D
Sample Name: BS-3-60B(+)

```
=====
Acq. Operator   : Z
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 12/14/2014 9:47:52 PM
Acq. Method     : C:\CHEM32\1\METHODS\DEF.LC.M
Last changed    : 12/14/2014 9:43:49 PM by Z
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF.LC.M
Last changed    : 7/21/2015 10:05:57 AM by
                  (modified after loading)
Sample Info     : 0J-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 220 nm
=====
```



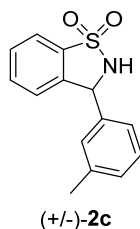
```
=====
Area Percent Report
=====
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VMD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	18.218	BV	0.5192	7586.84619	224.66182	49.8940	
2	20.052	VV	0.6355	7619.08643	177.23303	50.1060	

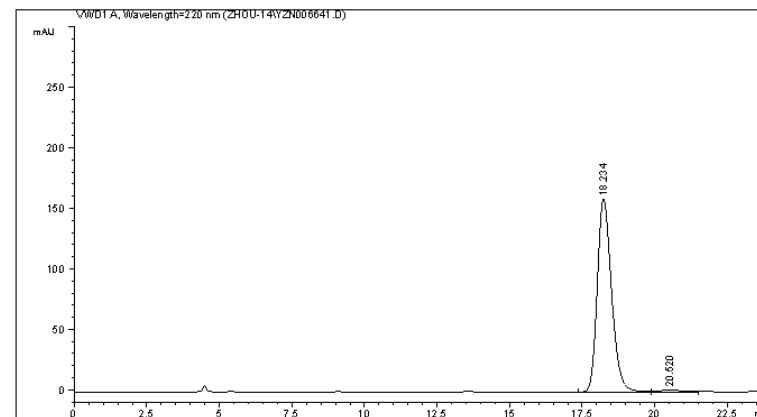
Totals : 1.52059e4 401.89485

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



Data File C:\CHEM32\1\DATA\ZHOU-14\YZN006641.D
Sample Name: BS-3-60B

```
=====
Acq. Operator   : Z
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 12/14/2014 8:25:04 PM
Acq. Method     : C:\CHEM32\1\METHODS\DEF.LC.M
Last changed    : 12/14/2014 8:24:05 PM by Z
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF.LC.M
Last changed    : 7/21/2015 10:05:57 AM by
                  (modified after loading)
Sample Info     : 0J-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 220 nm
=====
```



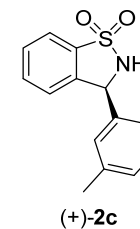
```
=====
Area Percent Report
=====
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VMD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	18.234	BV	0.5266	5451.18945	159.59685	98.4538	
2	20.520	VV	0.7196	85.60725	1.75772	1.5462	

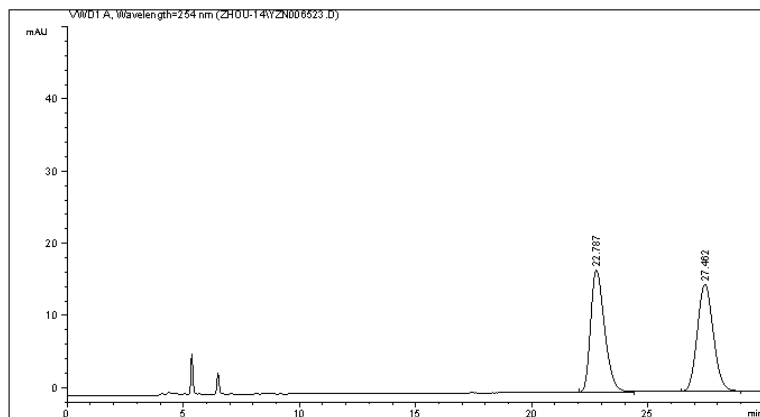
Totals : 5536.79670 161.35456

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



Data File C:\CHEM32\1\DATA\ZHOU-14\YZN006523.D
Sample Name: BS-3-55B(+/-)

```
=====
Acq. Operator   : Z
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 12/2/2014 9:37:12 AM
Acq. Method     : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 12/2/2014 9:33:22 AM by Z
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 7/21/2015 10:16:43 AM by
                  (modified after loading)
Sample Info     : OD-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 254 nm
=====
```



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

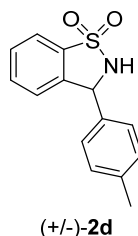
```
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=254 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area [mAU]	Area %
1	22.787	BB	0.6416	706.48566	16.91544	49.8563	
2	27.462	BB	0.7483	710.55811	14.83020	50.1437	

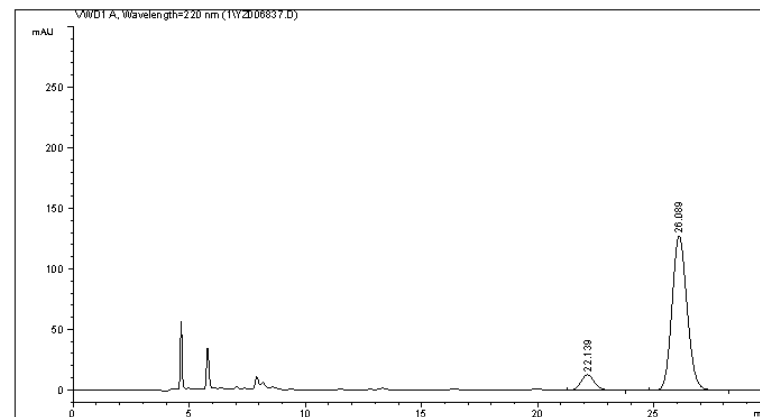
Totals : 1417.04376 31.74564

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



Data File C:\CHEM32\1\DATA\1\YZ006837.D
Sample Name: BS-3-55B

```
=====
Acq. Operator   : ZHOU
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date   : 12/16/2014 7:41:41 AM
Acq. Method      : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed     : 12/16/2014 7:16:38 AM by ZHOU
                  (modified after loading)
Analysis Method  : C:\CHEM32\1\METHODS\DEF LC.M
Last changed     : 7/21/2015 10:15:25 AM by
                  (modified after loading)
Sample Info      : OD-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 220 nm
=====
```



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

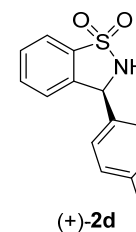
```
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area [mAU]	Area %
1	22.139	VB	0.6342	528.28247	12.92296	8.4272	
2	26.089	BB	0.7004	5740.51416	127.39679	91.5728	

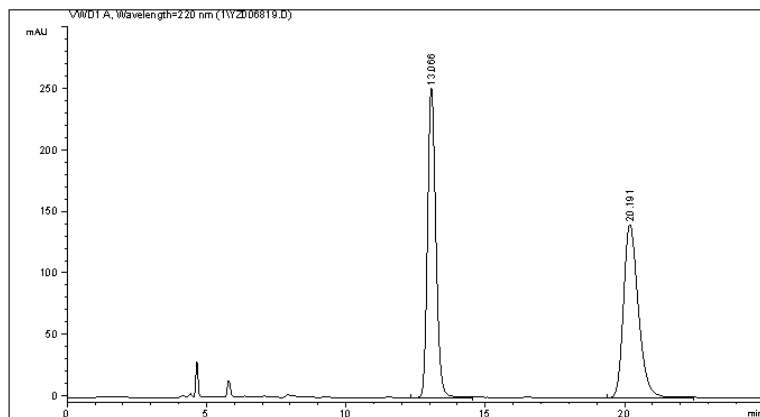
Totals : 6268.79663 140.31975

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



Data File C:\CHEM32\1\DATA\1\YZ006819.D
Sample Name: BS--3-60E(+/-)

```
=====
Acq. Operator   : ZHOU
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 12/14/2014 9:05:52 AM
Acq. Method     : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed    : 12/14/2014 9:04:39 AM by ZHOU
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 7/21/2015 10:19:47 AM by
                  (modified after loading)
Sample Info     : 0D-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 220 nm
=====
```



Area Percent Report

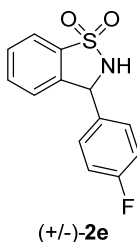
```
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area %
1	13.066	VB	0.3215	5230.54932	251.26656	50.2038
2	20.191	VB	0.5672	5188.08740	140.64362	49.7962

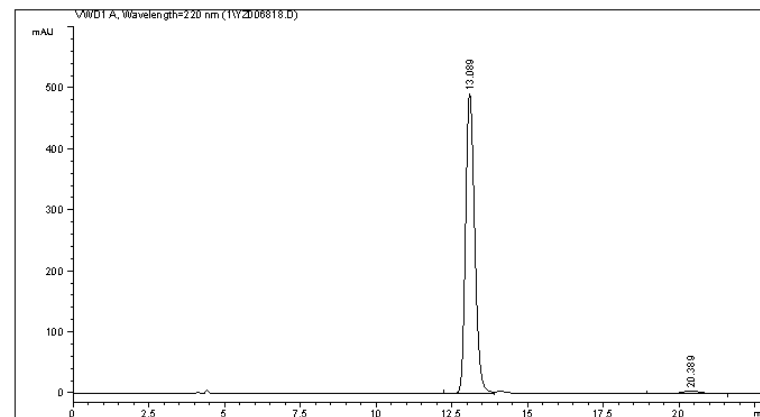
Totals : 1.04186e4 391.91017

*** End of Report ***



Data File C:\CHEM32\1\DATA\1\YZ006818.D
Sample Name: BS--3-60E

```
=====
Acq. Operator   : ZHOU
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 12/14/2014 8:41:00 AM
Acq. Method     : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed    : 12/14/2014 8:32:39 AM by ZHOU
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 7/21/2015 10:21:03 AM by
                  (modified after loading)
Sample Info     : 0D-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 220 nm
=====
```



Area Percent Report

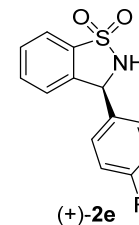
```
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area %
1	13.089	VV	0.3254	1.02738e4	491.50449	98.2261
2	20.389	VV	0.6009	185.53587	4.66423	1.7739

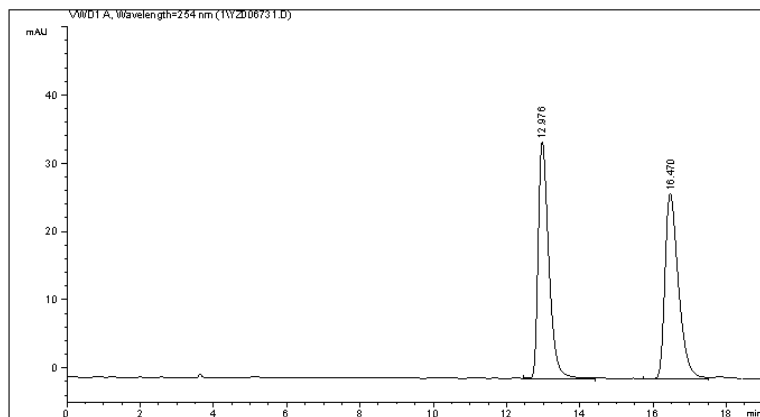
Totals : 1.04593e4 496.16872

*** End of Report ***



Data File C:\CHEM32\1\DATA\1\YZ006731.D
Sample Name: BS-3-55a

```
=====
Acq. Operator   : ZHOU
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 12/2/2014 2:35:03 PM
Acq. Method     : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed    : 12/2/2014 2:33:49 PM by ZHOU
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 7/21/2015 10:23:29 AM by
                  (modified after loading)
Sample Info     : OD-H, H/i-PrOH = 80/200, 0.8 mL/min, 30 oC, 254 nm
=====
```



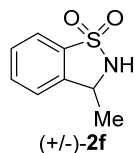
```
=====
Area Percent Report
=====
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VMD1 A, Wavelength=254 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area [mAU]	Area %
1	12.976	VB	0.3091	703.49756	34.73988	50.2870	
2	16.470	VV	0.3892	695.46686	27.14727	49.7130	

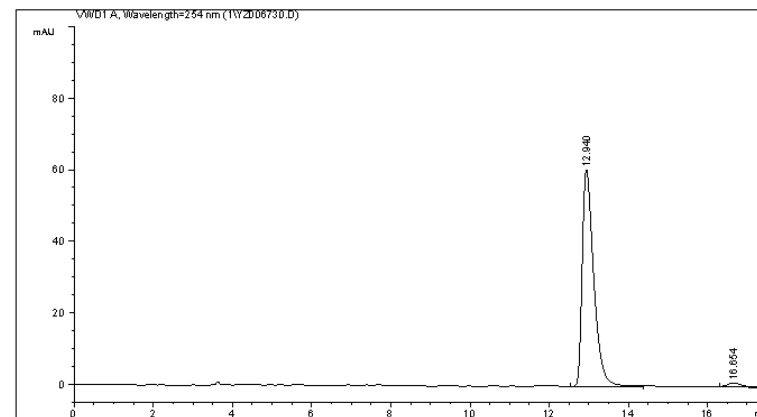
Totals : 1398.96442 61.88716

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



Data File C:\CHEM32\1\DATA\1\YZ006730.D
Sample Name: BS-3-55a

```
=====
Acq. Operator   : ZHOU
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 12/2/2014 1:54:45 PM
Acq. Method     : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed    : 12/2/2014 1:36:58 PM by ZHOU
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 7/21/2015 10:24:15 AM by
                  (modified after loading)
Sample Info     : OD-H, H/i-PrOH = 80/200, 0.8 mL/min, 30 oC, 254 nm
=====
```



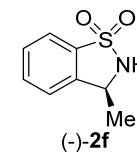
```
=====
Area Percent Report
=====
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VMD1 A, Wavelength=254 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area [mAU]	Area %
1	12.940	VV	0.3082	1239.87488	60.70688	97.4320	
2	16.654	VB	0.4318	32.67904	1.14563	2.5680	

Totals : 1272.55391 61.85251

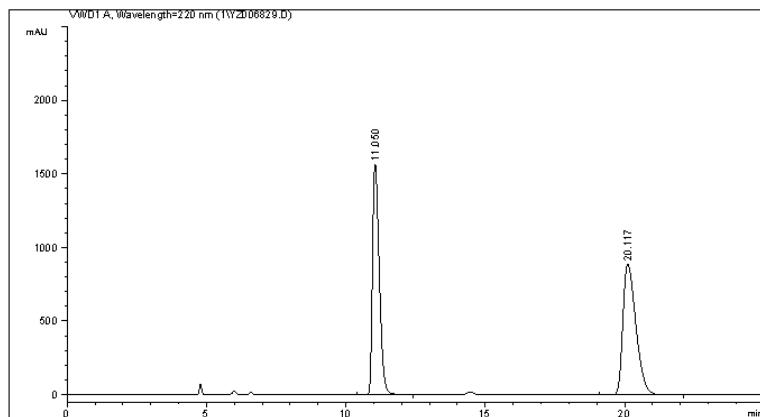
```
=====
*** End of Report ***
```



Data File C:\CHEM32\1\DATA\1\YZ006829.D
Sample Name: BS-3-56A(+)

=====

Acq. Operator	: ZHOU	Location	: Vial 1
Acq. Instrument	: Instrument 1		
Injection Date	: 12/15/2014 2:53:11 PM		
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC1.M		
Last changed	: 12/15/2014 2:49:47 PM by ZHOU		
	(modified after loading)		
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M		
Last changed	: 7/21/2015 10:26:28 AM by		
	(modified after loading)		
Sample Info	: OD-H, H ₂ O/PrOH = 80/20, 0.7 mL/min, 30 °C, 220 nm		



Area Percent Report

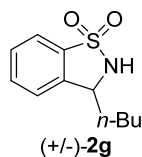
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	11.050	VV	0.2661	2.67985e4	1569.90332	48.0343	
2	20.117	VV	0.5064	2.89919e4	887.39813	51.9657	

Totals : 5.57904e4 2457.30145

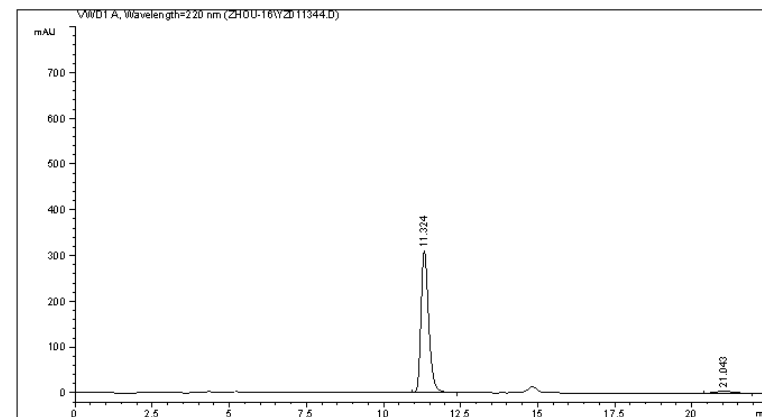
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZ011344.D
Sample Name: BS-7-5B

=====

Acq. Operator	: ZHOU	Location	: Vial 1
Acq. Instrument	: Instrument 1		
Injection Date	: 10/14/2016 5:34:15 AM		
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC1.M		
Last changed	: 10/14/2016 5:10:36 AM by		
	(modified after loading)		
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M		
Last changed	: 10/19/2016 2:04:06 PM by 0		
	(modified after loading)		
Sample Info	: OD-H, Hexane/iPrOH = 80/20, 0.7 mL/min, 30 °C, 220 nm		



Area Percent Report

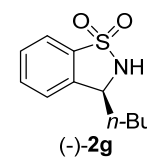
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	11.324	BB	0.2669	5386.62695	309.86606	97.2315	
2	21.043	BB	0.5082	153.37309	4.69052	2.7685	

Totals : 5540.00005 314.55658

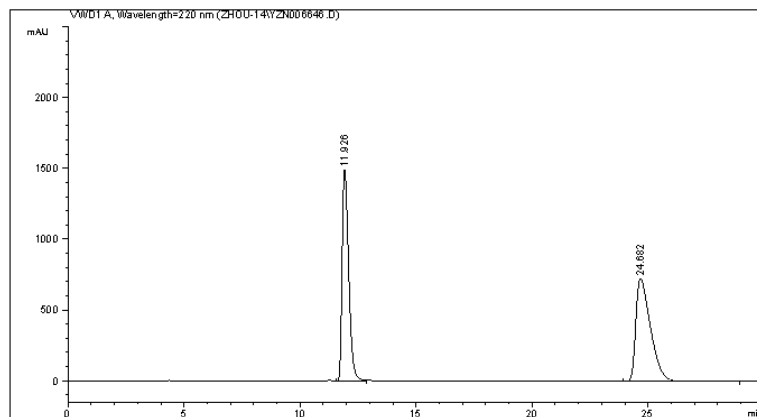
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-14\YZN006646.D
Sample Name: BS-3-56B(+)

=====

Acq. Operator	: Z	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 12/14/2014 10:57:59 PM	
Acq. Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 12/14/2014 10:55:41 PM by Z	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 10:33:54 AM by	
	(modified after loading)	
Sample Info	: OD-H, H/i-PrOH = 80/20, 0.7 mL/min, 30 oC, 220 nm	



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

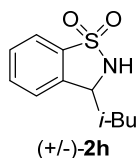
Sorted By	:	Signal
Multiplier:	:	1.0000
Dilution:	:	1.0000
Use Multiplier & Dilution Factor with ISTDs		

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.926	VV	0.3068	2.96570e4	1497.15125	48.2262
2	24.682	BB	0.6661	3.18387e4	725.80884	51.7738

Totals : 6.14957e4 2222.96008

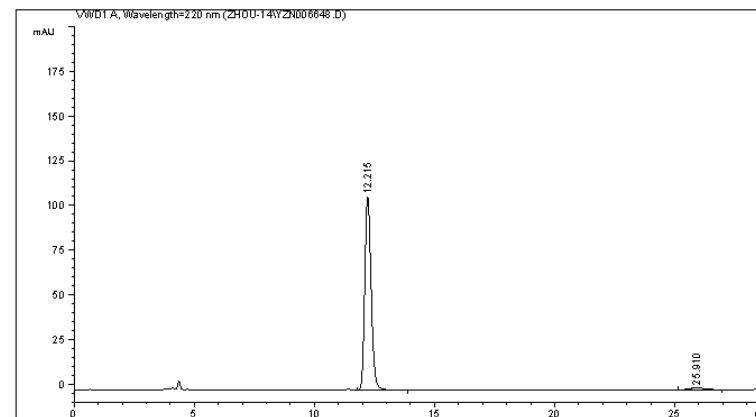
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-14\YZN006648.D
Sample Name: BS-3-56B

=====

Acq. Operator	: Z	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 12/15/2014 10:40:44 AM	
Acq. Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 12/15/2014 10:37:42 AM by Z	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 10:32:57 AM by	
	(modified after loading)	
Sample Info	: OD-H, H/i-PrOH = 80/20, 0.7 mL/min, 30 oC, 220 nm	



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

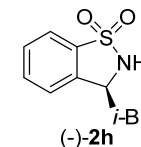
Sorted By	:	Signal
Multiplier:	:	1.0000
Dilution:	:	1.0000
Use Multiplier & Dilution Factor with ISTDs		

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.215	VB	0.2859	2004.83154	108.27026	98.0207
2	25.910	BB	0.5953	40.48392	9.98608e-1	1.9793

Totals : 2045.31547 109.26887

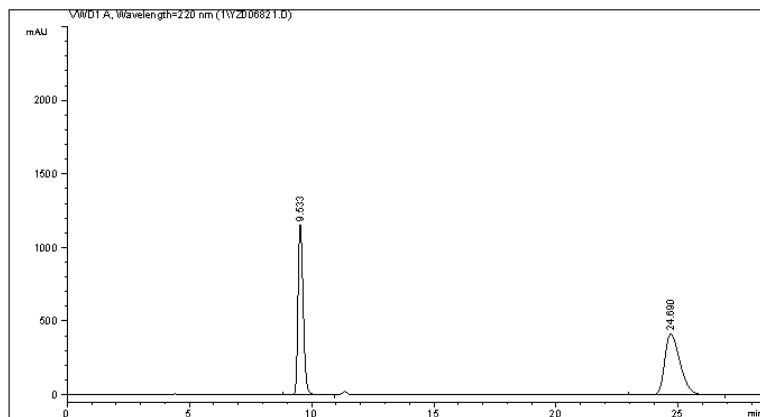
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\1\YZ006821.D
Sample Name: BS-3-60A(+/-)

=====

Acq. Operator	: ZHOU	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 12/15/2014 10:06:57 AM	
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC1.M	
Last changed	: 12/15/2014 9:43:16 AM by ZHOU	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 10:37:53 AM by	
	(modified after loading)	
Sample Info	: OD-H, H/i-PrOH = 75/25, 0.7 mL/min, 30 oC, 220 nm	



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

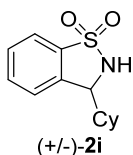
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	9.533	VV	0.2292	1.68887e4	1152.84937	48.5173	
2	24.690	VV	0.6677	1.79209e4	414.34860	51.4827	

Totals : 3.48096e4 1567.19797

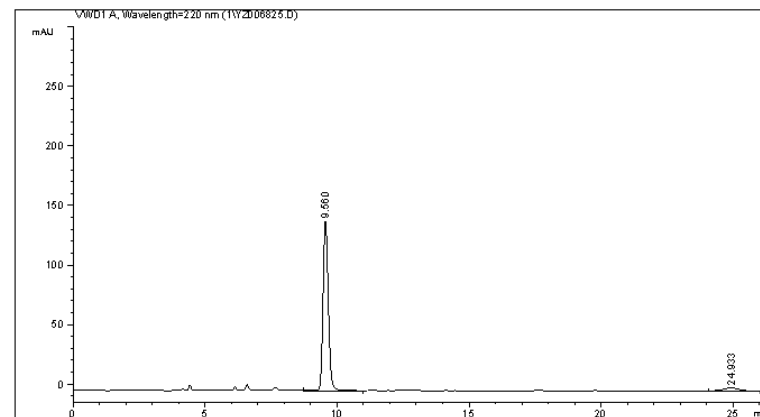
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\1\YZ006825.D
Sample Name: BS-3-60A

=====

Acq. Operator	: ZHOU	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 12/15/2014 12:19:49 PM	
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC1.M	
Last changed	: 12/15/2014 12:18:13 PM by ZHOU	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 10:39:27 AM by	
	(modified after loading)	
Sample Info	: OD-H, H/i-PrOH = 75/25, 0.7 mL/min, 30 oC, 220 nm	



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

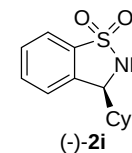
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	9.560	VB	0.2166	2004.62964	142.46884	94.6463	
2	24.933	VV	0.6897	113.39259	2.55459	5.3537	

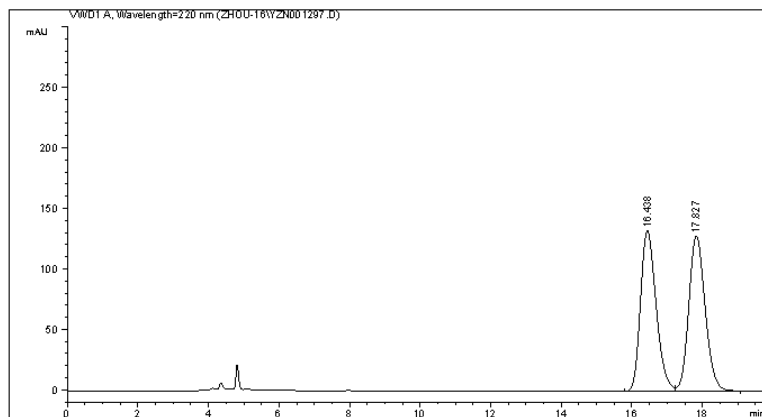
Totals : 2118.02222 145.02343

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



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001297.D
Sample Name: BS-5-97A(+)

=====
Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/9/2016 10:31:53 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/9/2016 10:29:20 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 6/29/2016 3:07:21 PM
(modified after loading)
Sample Info : OD-H, Hex/i-PrOH = 70/30, 0.7 mL/min, 30oC, 220 nm



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

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

Signal 1: VWD1 A, Wavelength=220 nm

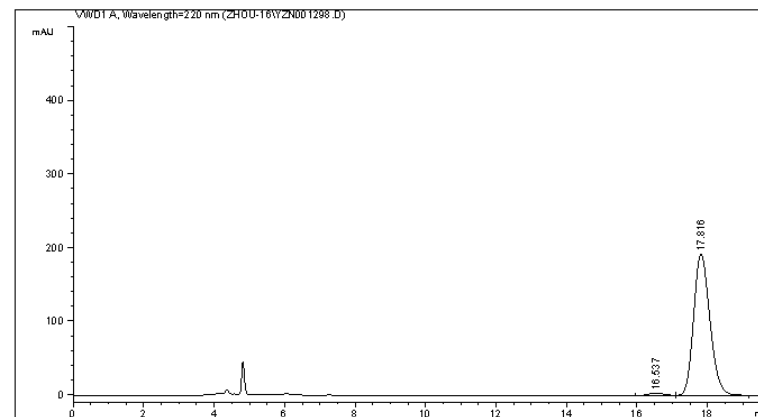
Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	16.438	EV	0.4768	4104.67383	132.89256	49.6718	
2	17.827	VB	0.5001	4158.92334	128.44221	50.3282	

Totals : 8263.59717 261.33478

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

Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001298.D
Sample Name: BS-5-97A

=====
Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/9/2016 10:54:11 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/9/2016 10:51:54 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 6/29/2016 3:08:52 PM
(modified after loading)
Sample Info : OD-H, Hex/i-PrOH = 70/30, 0.7 mL/min, 30oC, 220 nm



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

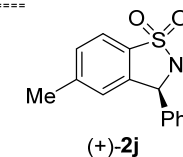
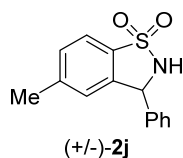
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	16.537	EV	0.4660	105.93079	3.52136	1.6849	
2	17.816	VB	0.4993	6181.00439	191.99762	98.3151	

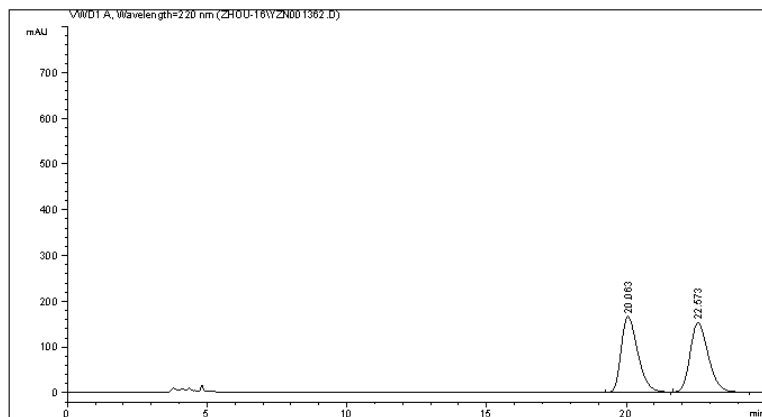
Totals : 6286.93519 195.51898

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



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001362.D
Sample Name: BS-6-4B(+)

```
=====
Acq. Operator   :
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 5/15/2016 8:51:00 PM
Acq. Method     : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed    : 5/15/2016 8:47:53 PM
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed    : 6/29/2016 3:16:36 PM
                  (modified after loading)
Sample Info     : 0D-H, Hex/i-PrOH =70/30, 0.7 mL/min, 30oC, 220 nm
=====
```



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

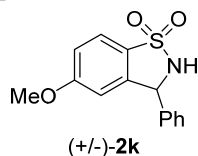
```
Sorted By      :      Signal
Multiplier:    :      1.0000
Dilution:      :      1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	20.063	BB	0.6388	6897.40186	166.11380	49.9504	
2	22.573	BB	0.7000	6911.10010	152.66460	50.0496	

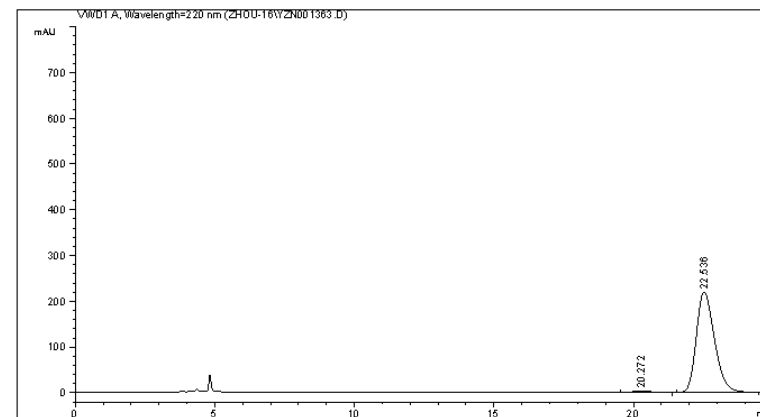
Totals : 1.38085e4 318.77840

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



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001363.D
Sample Name: BS-6-4B

```
=====
Acq. Operator   :
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 5/15/2016 9:18:14 PM
Acq. Method     : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed    : 5/15/2016 9:16:55 PM
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed    : 6/29/2016 3:16:36 PM
                  (modified after loading)
Sample Info     : 0D-H, Hex/i-PrOH =70/30, 0.7 mL/min, 30oC, 220 nm
=====
```



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

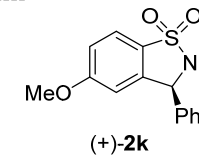
```
Sorted By      :      Signal
Multiplier:    :      1.0000
Dilution:      :      1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	20.272	BB	0.6432	159.30930	3.88305	1.5781	
2	22.536	BB	0.7005	9935.75781	219.84697	98.4219	

Totals : 1.00951e4 223.73002

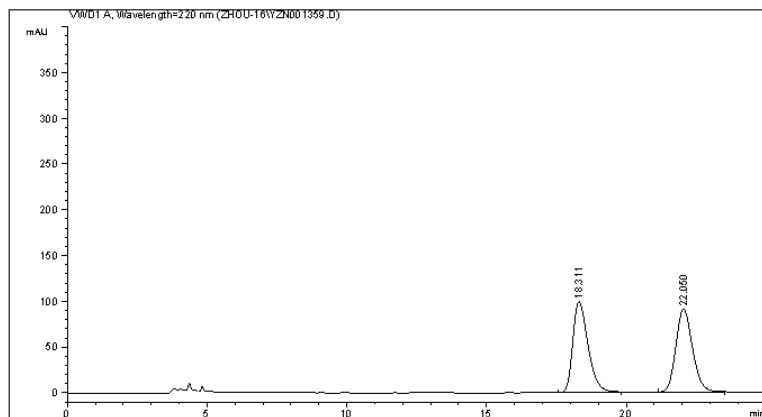
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001359.D
Sample Name: BS-6-4A(+/-)

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/15/2016 7:48:00 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/15/2016 7:30:43 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 6/29/2016 3:14:06 PM
(modified after loading)
Sample Info : OD-H, Hex/i-PrOH =70/30, 0.7 mL/min, 30oC, 220 nm



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

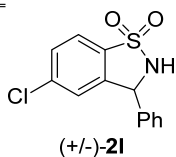
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	18.311	BB	0.5817	3756.95435	99.18011	50.1253	
2	22.050	BB	0.6376	3738.17383	90.51442	49.8747	

Totals : 7495.12817 189.69453

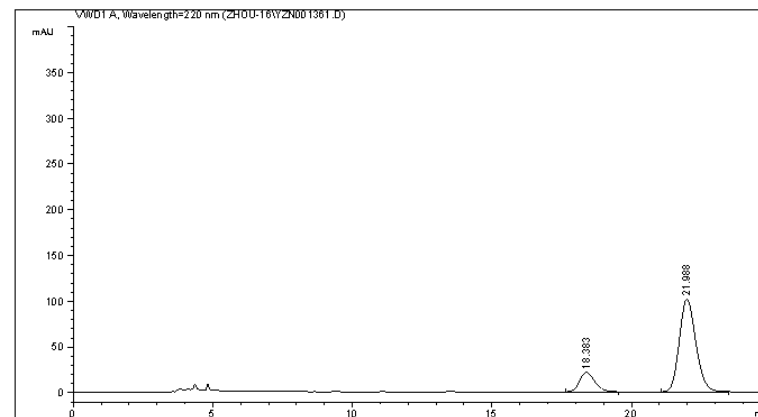
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001361.D
Sample Name: BS-6-4A

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/15/2016 8:14:50 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/15/2016 8:14:12 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 6/29/2016 3:14:06 PM
(modified after loading)
Sample Info : OD-H, Hex/i-PrOH =70/30, 0.7 mL/min, 30oC, 220 nm



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

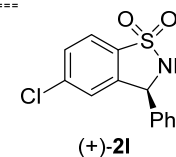
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	18.383	BB	0.5767	808.90118	21.45433	16.2255	
2	21.988	BB	0.6380	4176.46729	101.36083	83.7745	

Totals : 4985.36847 122.81516

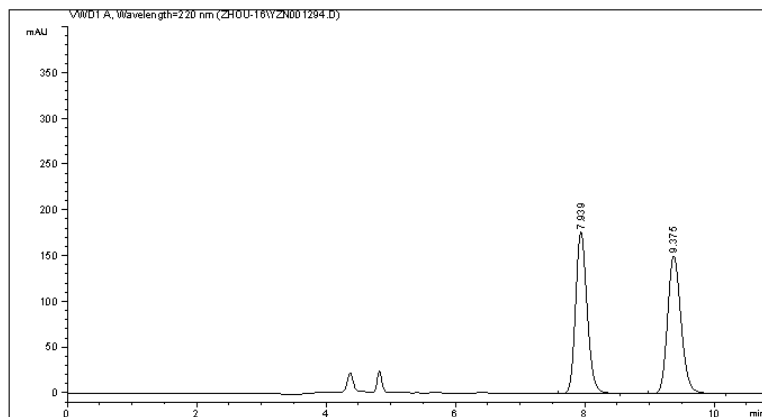
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001294.D
Sample Name: BS-5-97B(+)

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/9/2016 9:50:58 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/9/2016 9:22:38 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 6/29/2016 3:10:08 PM
(modified after loading)
Sample Info : OD-H, Hex/i-PrOH = 70/30, 0.7 mL/min, 30oC, 220 nm



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

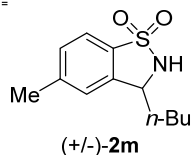
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	*s	Height [mAU]	Area %
1	7.939	VB	0.1920	2202.93799		176.70711	50.1475
2	9.375	VB	0.2249	2189.98096		150.63695	49.8525

Totals : 4392.91895 327.34406

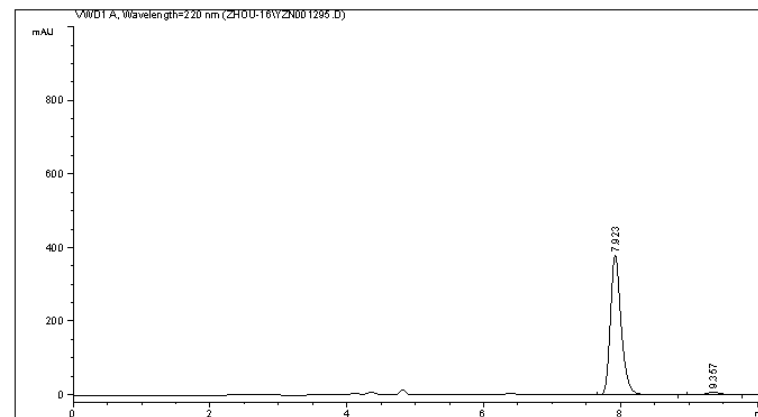
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001295.D
Sample Name: BS-5-97B

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/9/2016 10:04:55 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/9/2016 10:02:02 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 6/29/2016 3:11:04 PM
(modified after loading)
Sample Info : OD-H, Hex/i-PrOH = 70/30, 0.7 mL/min, 30oC, 220 nm



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

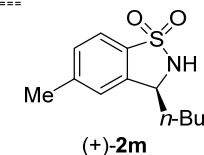
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	*s	Height [mAU]	Area %
1	7.923	EB	0.1658	4156.40332		379.07272	97.4338
2	9.357	EB	0.2305	109.46868		7.34994	2.5662

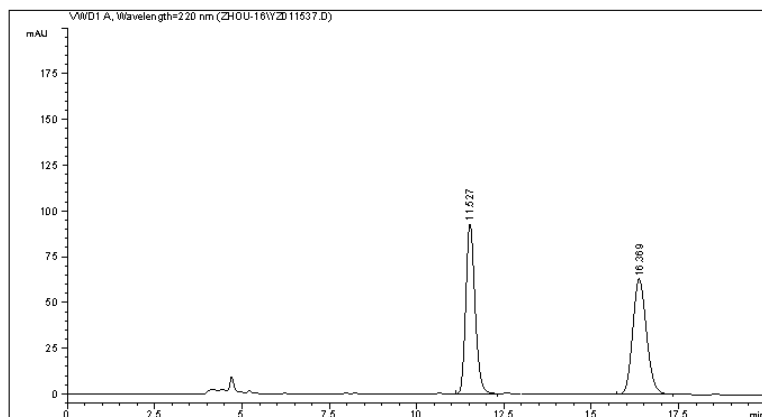
Totals : 4265.87200 386.42266

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



Data File C:\CHEM32\1\DATA\ZHOU-16\YZ011537.D
Sample Name: BS-7-54(+)

=====
Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 12/18/2016 12:47:16 PM
Acq. Method : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed : 12/18/2016 12:25:22 PM by
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 12/22/2016 10:20:50 PM by 0
(modified after loading)
Sample Info : OD-H, Hexane/iPrOH = 70/30, 0.7 mL/min, 30 oC, 220
nm



Area Percent Report

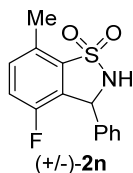
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	11.527	BB	0.2847	1725.73779	93.11014	49.8718	
2	16.369	BB	0.4258	1734.61121	63.04984	50.1282	

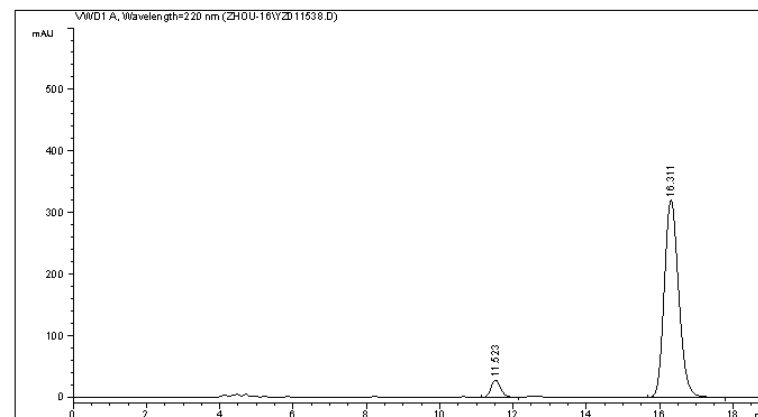
Totals : 3460.34900 156.15998

*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZ011538.D
Sample Name: BS-7-54

=====
Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 12/18/2016 1:12:23 PM
Acq. Method : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed : 12/18/2016 1:10:13 PM by
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 12/22/2016 10:22:48 PM by 0
(modified after loading)
Sample Info : OD-H, Hexane/iPrOH = 70/30, 0.7 mL/min, 30 oC, 220
nm



Area Percent Report

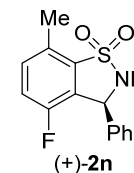
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	11.523	BB	0.2825	524.53455	28.59326	5.6453	
2	16.311	BB	0.4236	8766.94922	320.81998	94.3547	

Totals : 9291.48376 349.41324

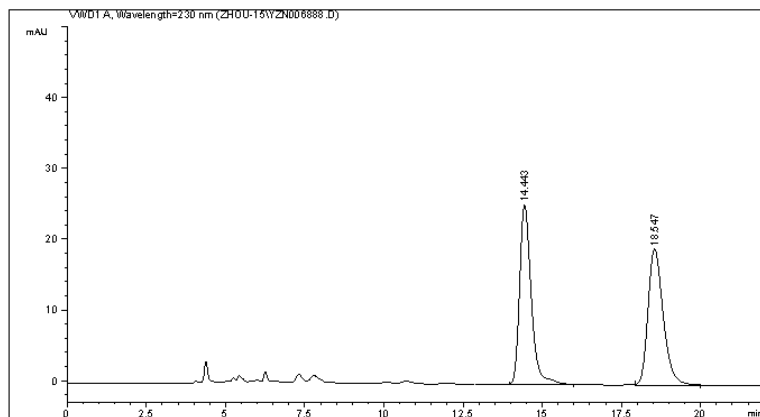
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-15\YZN006888.D
Sample Name: BS-3-70(+/-)

=====

Acq. Operator	: Z	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 1/6/2015 4:02:24 PM	
Acq. Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 1/6/2015 4:01:40 PM by Z	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 11:20:03 AM by	
	(modified after loading)	
Sample Info	: OD-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 230 nm	



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

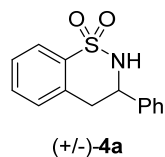
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area %
1	14.443	BB	0.3922	649.48096	25.34248	50.3105
2	18.547	BB	0.5115	641.46539	19.23088	49.6895

Totals : 1290.94635 44.57335

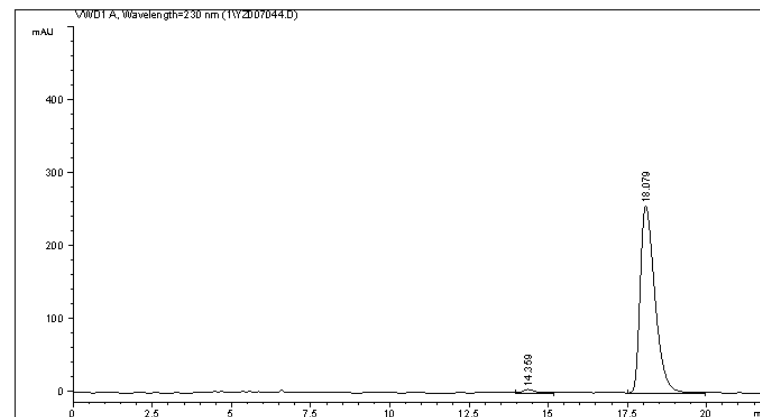
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\1\YZ007044.D
Sample Name: BS-3-70

=====

Acq. Operator	: ZHOU	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 1/18/2015 7:54:49 AM	
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC1.M	
Last changed	: 1/18/2015 7:52:24 AM by ZHOU	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 11:19:15 AM by	
	(modified after loading)	
Sample Info	: OD-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 230 nm	



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

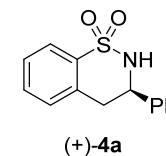
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area %
1	14.359	VB	0.3996	141.62616	5.10212	1.6764
2	18.079	VB	0.5028	8306.72266	256.67084	98.3236

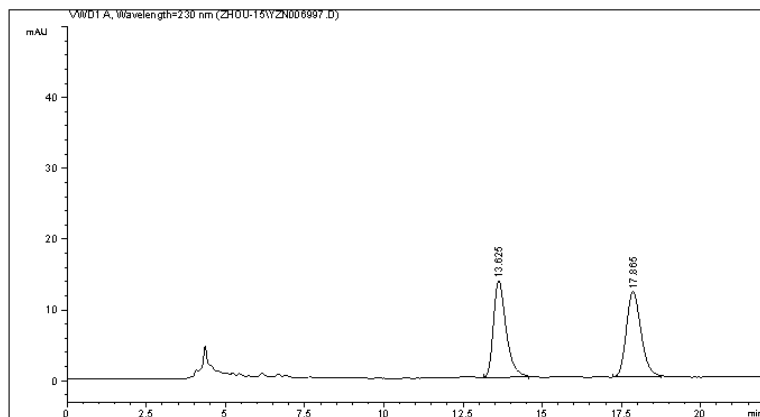
Totals : 8448.34882 261.77296

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



Data File C:\CHEM32\1\DATA\ZHOU-15\YZN006997.D
Sample Name: BS-3--73A

```
=====
Acq. Operator   : Z
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 1/16/2015 11:07:03 PM
Acq. Method     : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 1/16/2015 11:04:38 PM by Z
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 7/21/2015 11:02:52 AM by
                  (modified after loading)
Sample Info     : OD-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 230 nm
=====
```



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

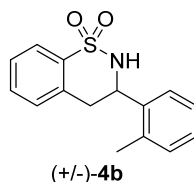
```
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Sample Amount: : 1.00000 [ng/ul] (not used in calc.)
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	13.625	BB	0.4233	378.74615	13.62572	49.4854
2	17.865	BB	0.4995	386.62296	12.00429	50.5146

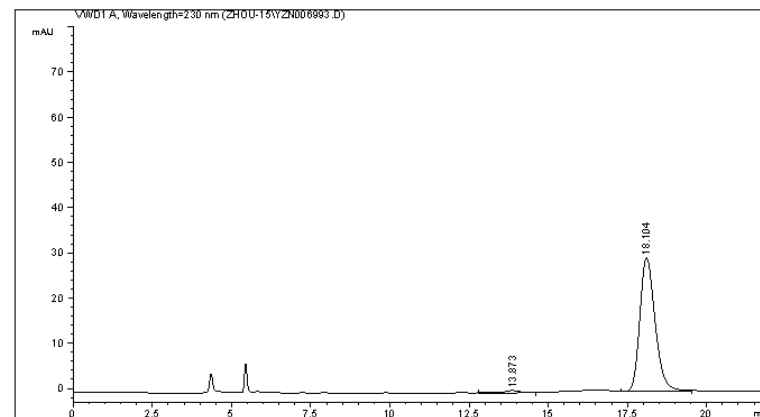
Totals : 765.36911 25.63000

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



Data File C:\CHEM32\1\DATA\ZHOU-15\YZN006993.D
Sample Name: BS-3--73A

```
=====
Acq. Operator   : Z
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date   : 1/16/2015 8:54:13 PM
Acq. Method      : C:\CHEM32\1\METHODS\DEF LC.M
Last changed     : 1/16/2015 8:35:06 PM by Z
                  (modified after loading)
Analysis Method  : C:\CHEM32\1\METHODS\DEF LC.M
Last changed     : 7/21/2015 11:03:54 AM by
                  (modified after loading)
Sample Info      : OD-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 230 nm
=====
```



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

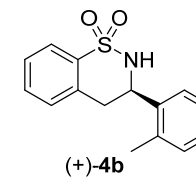
```
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Sample Amount: : 1.00000 [ng/ul] (not used in calc.)
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	13.873	BV	0.5808	22.16280	5.11960e-1	2.1823
2	18.104	VB	0.5179	993.40417	29.51423	97.8177

Totals : 1015.56698 30.02619

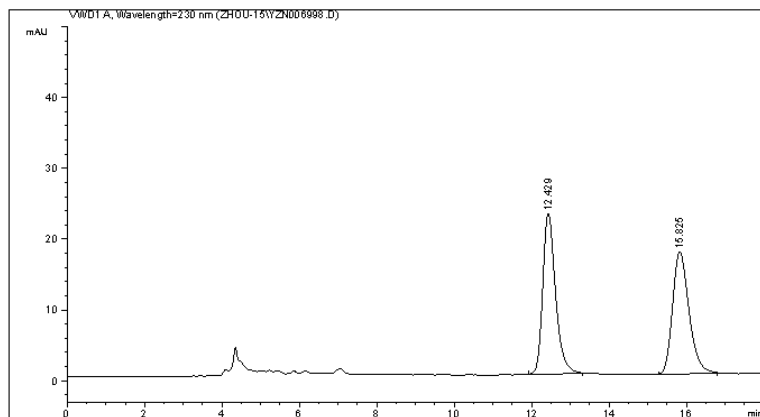
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-15\YZN006998.D
Sample Name: BS-3--73B(+)

=====

Acq. Operator	: Z	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 1/16/2015 11:33:31 PM	
Acq. Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 1/16/2015 11:31:44 PM by Z	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 11:02:08 AM by	
	(modified after loading)	
Sample Info	: 0D-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 230 nm	



Area Percent Report

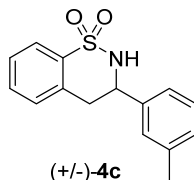
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Sample Amount: : 1.00000 [ng/ul] (not used in calc.)
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s [mAU]	Area %
1	12.429	BB	0.3480	521.40002	22.70233	50.5637
2	15.825	BB	0.4507	509.77518	17.27071	49.4363

Totals : 1031.17520 39.97304

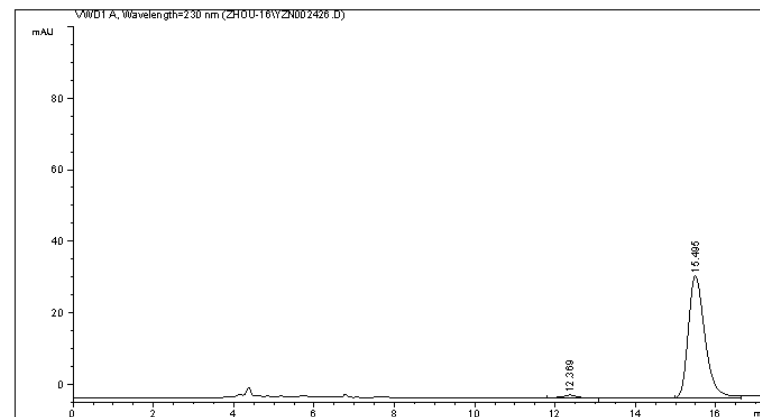
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN002426.D
Sample Name: BS-3-73B

=====

Acq. Operator	:		
Acq. Instrument	: Instrument 1	Location : Vial 1	
Injection Date	: 9/20/2016 10:34:57 PM		
Acq. Method	: C:\CHEM32\1\METHODS\DEF LC.M		
Last changed	: 9/20/2016 10:32:32 PM		
	(modified after loading)		
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M		
Last changed	: 10/19/2016 2:22:17 PM by 0		
	(modified after loading)		
Sample Info	: 0D-H, Hexane/i-PrOH = 70/30, 0.7 mL/min, 30oC, 220nm		



Area Percent Report

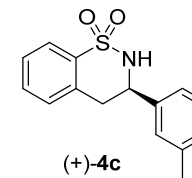
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s [mAU]	Area %
1	12.369	BB	0.4024	24.94381	9.06971e-1	2.4449
2	15.495	BB	0.4435	995.27411	34.28829	97.5551

Totals : 1020.21792 35.19526

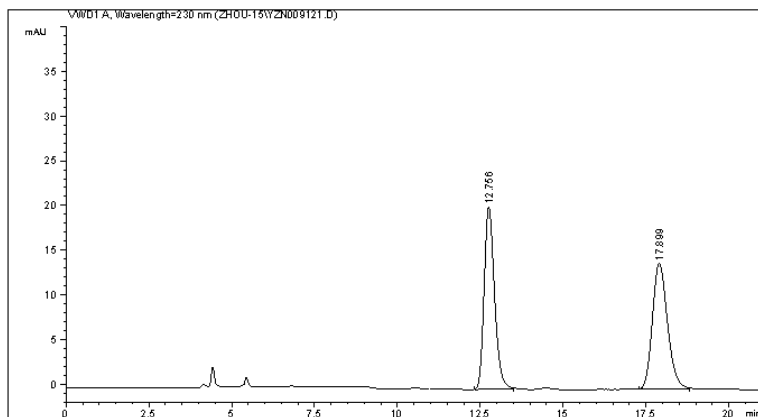
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-15\YZN009121.D
Sample Name: BS-3-73(+)

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 8/20/2015 8:53:12 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 8/20/2015 8:49:02 PM by
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 8/20/2015 9:20:09 PM by
(modified after loading)
Sample Info : 0D-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 230 nm



=====

Area Percent Report

=====

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

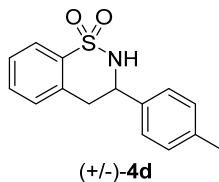
Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	12.756	BB	0.3342	445.19214	20.43899	50.1805
2	17.899	BB	0.4825	441.98898	14.03212	49.8195

Totals : 887.18112 34.47110

=====

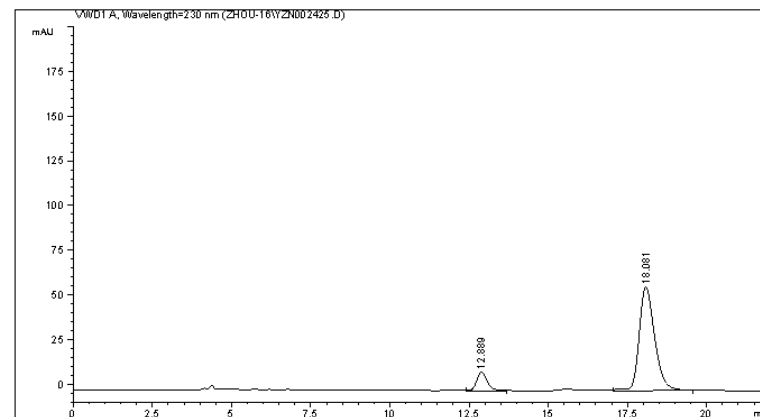
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN002425.D
Sample Name: BS-3-74C

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 9/20/2016 10:09:46 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 9/20/2016 10:06:18 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 10/19/2016 2:24:09 PM by 0
(modified after loading)
Sample Info : 0D-H, Hexane/i-PrOH = 70/30, 0.7 mL/min, 30oC, 220nm



=====

Area Percent Report

=====

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

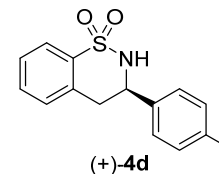
Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	12.889	VB	0.3490	235.72371	10.28322	10.6701
2	18.081	BB	0.5196	1973.47937	57.94915	89.3299

Totals : 2209.20308 68.23237

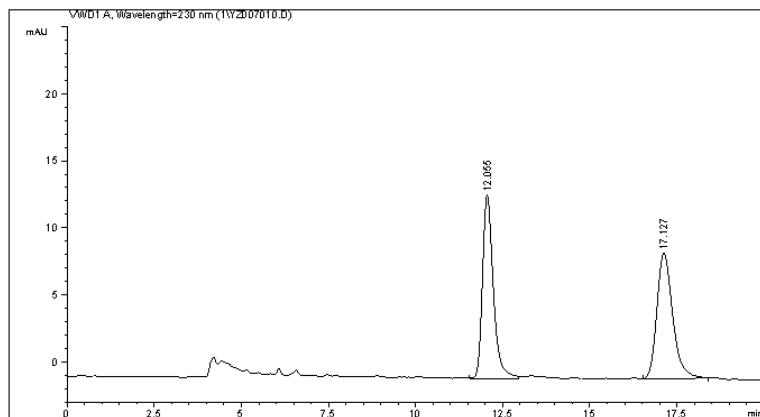
=====

*** End of Report ***



Data File C:\CHEM32\1\DATA\1\YZ007010.D
Sample Name: BS-3-74D(+)

```
=====
Acq. Operator   : ZHOU
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 1/14/2015 2:00:47 PM
Acq. Method     : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed    : 1/14/2015 1:58:23 PM by ZHOU
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 7/21/2015 11:30:40 AM by
                  (modified after loading)
Sample Info     : 0D-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 230 nm
=====
```



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

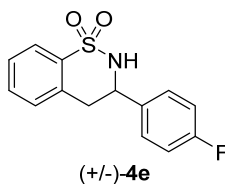
```
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area %
1	12.055	VB	0.3284	296.69174	13.69643	50.2223
2	17.127	BB	0.4805	294.06473	9.35002	49.7777

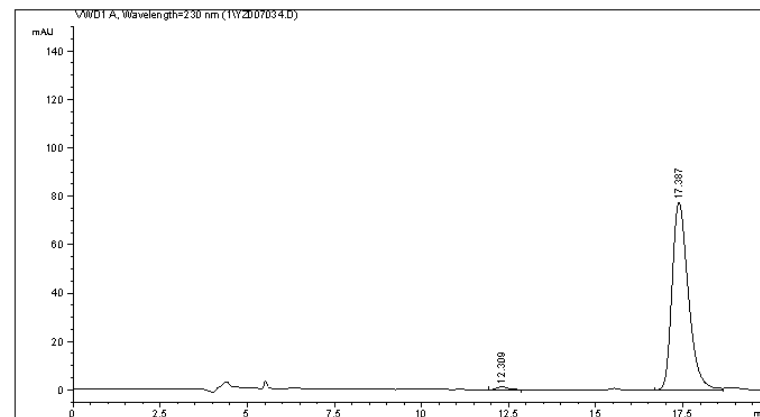
Totals : 590.75647 23.04645

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



Data File C:\CHEM32\1\DATA\1\YZ007034.D
Sample Name: BS-3-74D

```
=====
Acq. Operator   : ZHOU
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 1/17/2015 12:49:58 PM
Acq. Method     : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed    : 1/17/2015 12:32:55 PM by ZHOU
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 7/21/2015 11:33:23 AM by
                  (modified after loading)
Sample Info     : 0D-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 230 nm
=====
```



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

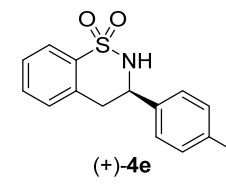
```
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area %
1	12.309	BB	0.3419	29.38605	1.28794	1.1519
2	17.387	EV	0.4988	2521.75854	77.53453	98.8481

Totals : 2551.14460 78.82247

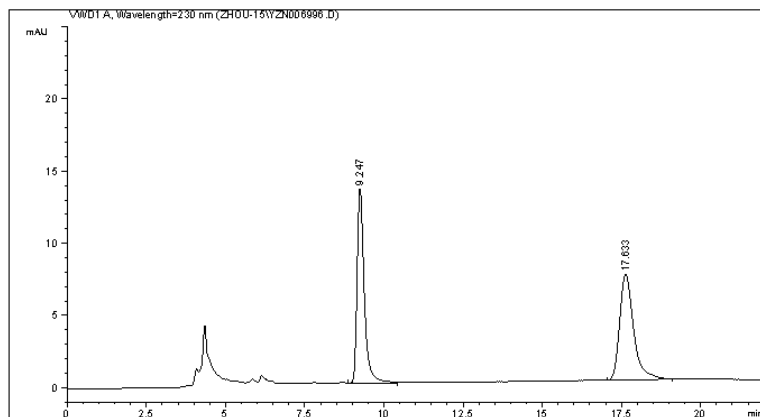
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-15\YZN006996.D
Sample Name: BS-3--78C(+)

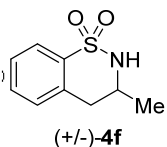
=====

Acq. Operator : Z
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 1/16/2015 10:35:30 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 1/16/2015 10:32:56 PM by Z
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 7/21/2015 10:54:42 AM by
(modified after loading)
Sample Info : OD-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 230 nm



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

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



Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	9.247	VB	0.2368	213.75076	13.52772	48.7381
2	17.633	BB	0.4668	224.81955	7.30515	51.2619

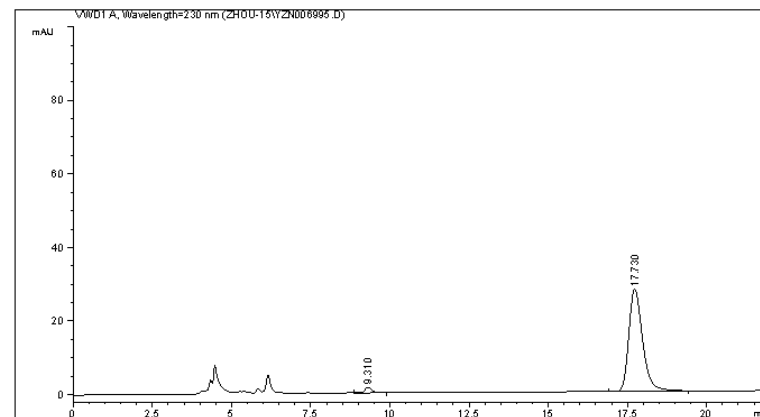
Totals : 438.57031 20.83287

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

Data File C:\CHEM32\1\DATA\ZHOU-15\YZN006995.D
Sample Name: BS-3--78C

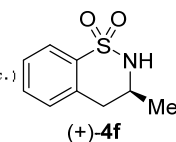
=====

Acq. Operator : Z
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 1/16/2015 10:01:11 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 1/16/2015 9:57:31 PM by Z
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 7/21/2015 10:55:15 AM by
(modified after loading)
Sample Info : OD-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 230 nm



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

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



Signal 1: VWD1 A, Wavelength=230 nm

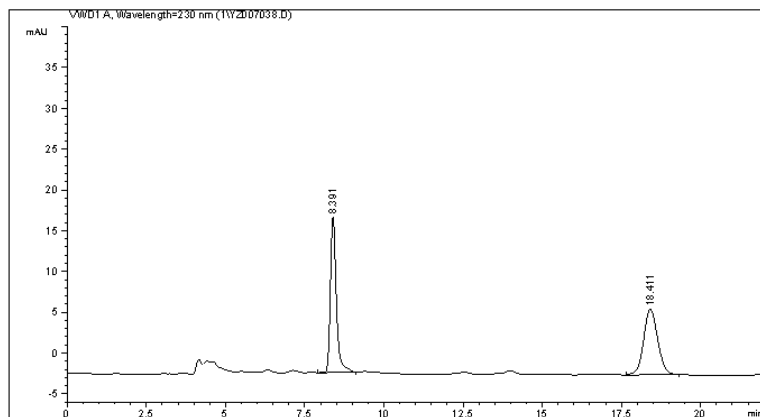
Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	9.310	VB	0.2441	24.82518	1.53405	2.8591
2	17.730	BB	0.4656	843.44727	27.72587	97.1409

Totals : 868.27244 29.25992

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

Data File C:\CHEM32\1\DATA\1\YZ007038.D
Sample Name: BS-3-78D(+)

```
=====
Acq. Operator   : ZHOU
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 1/17/2015 3:02:10 PM
Acq. Method     : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed    : 1/17/2015 3:00:26 PM by ZHOU
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 7/21/2015 2:13:12 PM by
                  (modified after loading)
Sample Info     : OD-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 230 nm
=====
```



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

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area %
1	8.391	VB	0.1987	250.57877	19.03116	50.3492
2	18.411	BB	0.4772	247.10324	8.02363	49.6508

Totals : 497.68201 27.05479

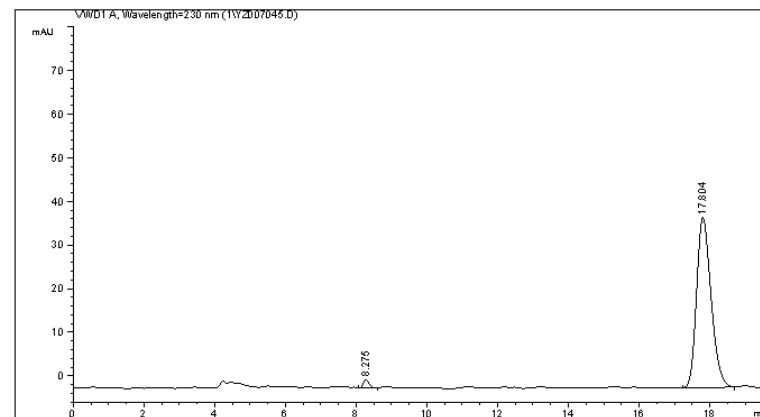
```
=====
*** End of Report ***
```

Instrument 1 7/21/2015 2:13:18 PM

Page 1 of 1

Data File C:\CHEM32\1\DATA\1\YZ007045.D
Sample Name: BS-3-78D

```
=====
Acq. Operator   : ZHOU
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date   : 1/18/2015 8:22:22 AM
Acq. Method     : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed    : 1/18/2015 8:19:41 AM by ZHOU
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 7/21/2015 2:13:57 PM by
                  (modified after loading)
Sample Info     : OD-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 230 nm
=====
```



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

Signal 1: VWD1 A, Wavelength=230 nm

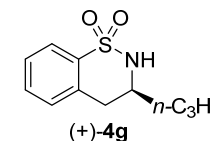
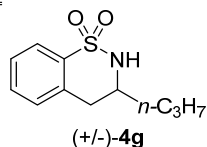
Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area %
1	8.275	BB	0.1671	19.46494	1.84034	1.7268
2	17.804	BB	0.4396	1107.76147	38.94886	98.2732

Totals : 1127.22641 40.78920

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

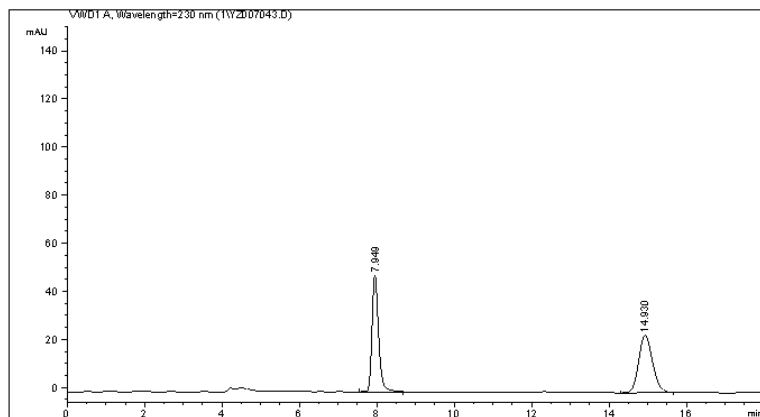
Instrument 1 7/21/2015 2:14:01 PM

Page 1 of 1



Data File C:\CHEM32\1\DATA\1\YZ007043.D
Sample Name: BS-3-78E(+)

```
=====
Acq. Operator   : ZHOU
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 1/18/2015 7:26:37 AM
Acq. Method     : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed    : 1/18/2015 7:24:39 AM by ZHOU
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 7/21/2015 2:18:23 PM by
                  (modified after loading)
Sample Info     : OD-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 230 nm
=====
```



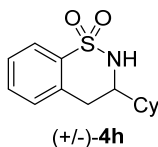
```
=====
Area Percent Report
=====
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area %
1	7.949	BB	0.1811	575.98853	48.37144	49.2671
2	14.930	BB	0.3838	593.12500	23.92973	50.7329

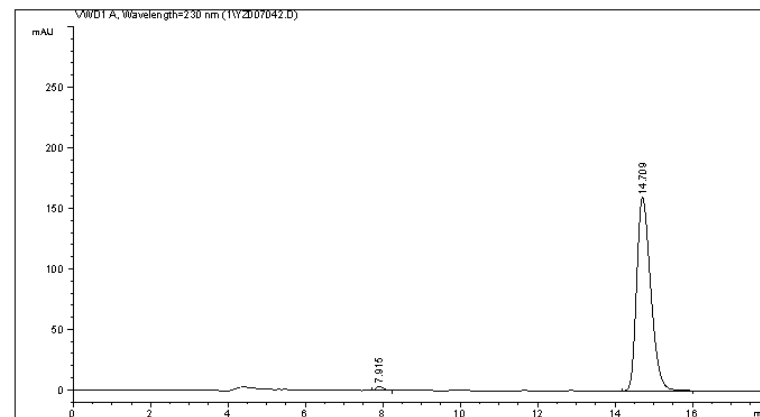
Totals : 1169.11353 72.30117

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



Data File C:\CHEM32\1\DATA\1\YZ007042.D
Sample Name: BS-3-78E

```
=====
Acq. Operator   : ZHOU
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date   : 1/18/2015 6:38:16 AM
Acq. Method      : C:\HPCHEM\1\METHODS\DEF LC1.M
Last changed     : 1/18/2015 5:56:15 AM by ZHOU
                  (modified after loading)
Analysis Method  : C:\CHEM32\1\METHODS\DEF LC.M
Last changed     : 7/21/2015 2:19:00 PM by
                  (modified after loading)
Sample Info      : OD-H, H/i-PrOH = 70/30, 0.7 mL/min, 30 oC, 230 nm
=====
```



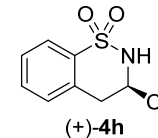
```
=====
Area Percent Report
=====
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area %
1	7.915	VB	0.1865	44.31816	3.62127	1.0906
2	14.709	BB	0.3883	4019.31396	159.67844	98.9094

Totals : 4063.63212 163.29971

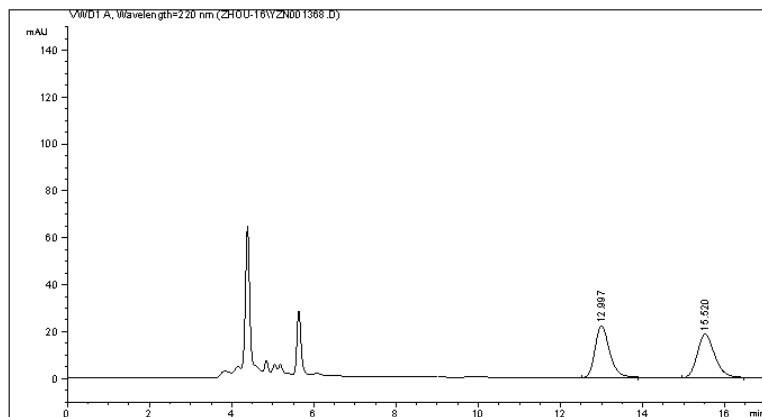
```
=====
*** End of Report ***
```



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001368.D
Sample Name: BS-6-6C(+/-)

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/15/2016 11:26:31 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/15/2016 11:23:58 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 6/29/2016 3:23:05 PM
(modified after loading)
Sample Info : 0D-H, Hex/i-PrOH =70/30, 0.7 mL/min, 30oC, 220 nm



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

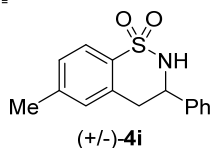
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	12.997	BB	0.3761	545.74561	22.16139	50.1036	
2	15.520	BB	0.4497	543.48840	18.54109	49.8964	

Totals : 1089.23401 40.70248

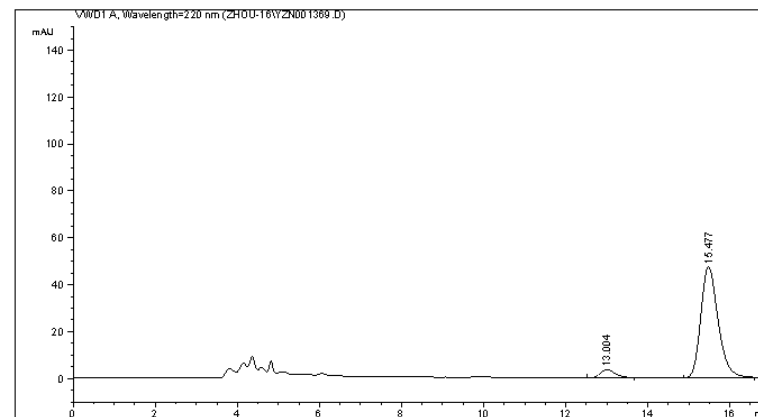
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001369.D
Sample Name: BS-6-6C

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/15/2016 11:46:01 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/15/2016 11:43:41 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 6/29/2016 3:23:05 PM
(modified after loading)
Sample Info : 0D-H, Hex/i-PrOH =70/30, 0.7 mL/min, 30oC, 220 nm



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

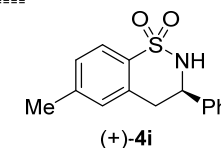
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	13.004	BB	0.3952	85.65388	3.34134	5.8268	
2	15.477	BB	0.4510	1384.35486	47.25777	94.1732	

Totals : 1470.00874 50.59911

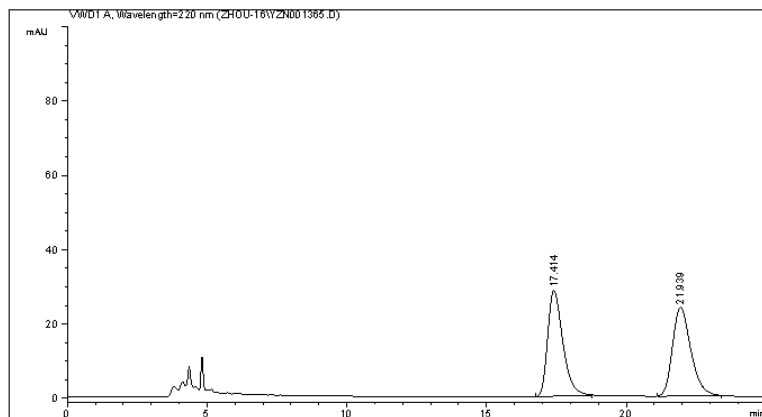
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001365.D
Sample Name: BS-6-6A(+/-)

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/15/2016 10:19:32 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/15/2016 10:16:21 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 6/29/2016 3:18:29 PM
(modified after loading)
Sample Info : OD-H, Hex/i-PrOH =70/30, 0.7 mL/min, 30oC, 220 nm



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

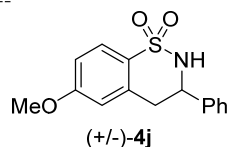
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	*s	Height [mAU]	Area %
1	17.414	BB	0.5763	1074.66589		28.53231	49.5929
2	21.939	BB	0.7008	1092.30835		23.89370	50.4071

Totals : 2166.97424 52.42601

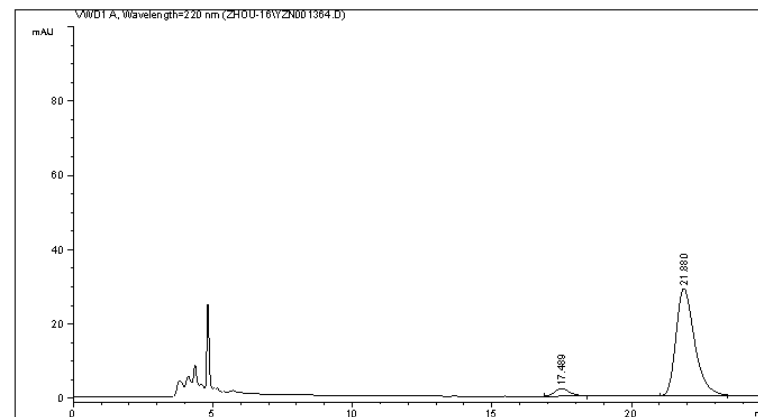
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001364.D
Sample Name: BS-6-6A

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/15/2016 9:51:06 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/15/2016 9:45:07 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 6/29/2016 3:18:29 PM
(modified after loading)
Sample Info : OD-H, Hex/i-PrOH =70/30, 0.7 mL/min, 30oC, 220 nm



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

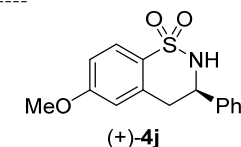
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	*s	Height [mAU]	Area %
1	17.489	BB	0.5583	76.01335		2.04100	5.3976
2	21.980	BB	0.7073	1332.27161		28.87073	94.6024

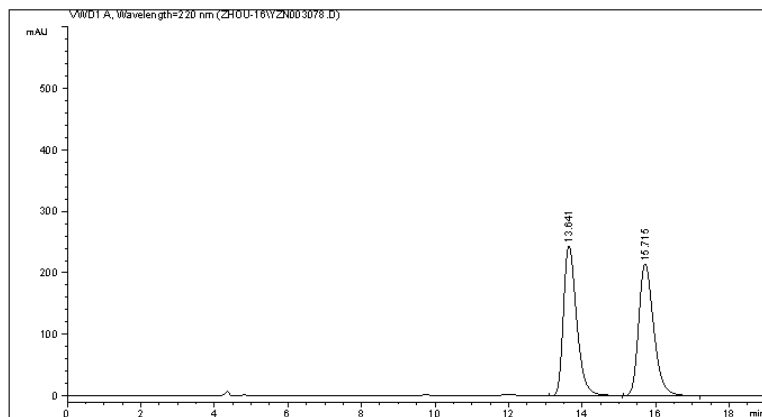
Totals : 1408.28496 30.91173

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



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN003078.D
Sample Name: BS-7-58(+)

```
=====
Acq. Operator   : 0
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 12/22/2016 11:04:06 AM
Acq. Method     : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 12/22/2016 10:49:43 AM by 0
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 12/22/2016 10:25:13 PM by 0
                  (modified after loading)
Sample Info     : 0D-H, Hexane/i-PrOH = 70/30, 0.7 mL/min, 30oC, 220 nm
=====
```



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

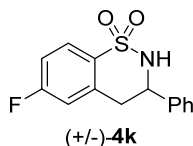
```
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	13.641	BB	0.3814	6103.41357	244.54706	50.0337	
2	15.715	BB	0.4392	6095.18262	214.51564	49.9663	

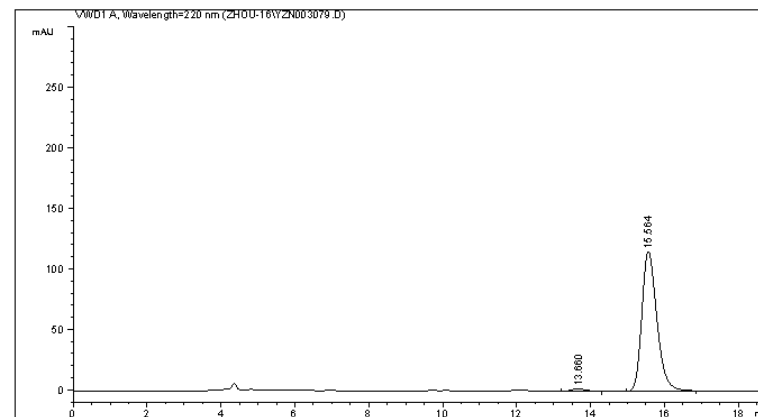
Totals : 1.21986e4 459.06270

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



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN003079.D
Sample Name: BS-7-58

```
=====
Acq. Operator   : 0
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 12/22/2016 11:24:38 AM
Acq. Method     : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 12/22/2016 11:23:46 AM by 0
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 12/22/2016 10:26:52 PM by 0
                  (modified after loading)
Sample Info     : 0D-H, Hexane/i-PrOH = 70/30, 0.7 mL/min, 30oC, 220 nm
=====
```



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

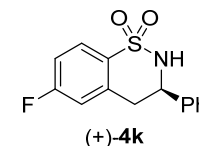
```
Sorted By      : Signal
Multiplier:    : 1.0000
Dilution:      : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	13.660	BB	0.3703	47.05296	1.98080	1.4229	
2	15.564	BB	0.4345	3259.68604	115.36879	98.5771	

Totals : 3306.73899 117.34959

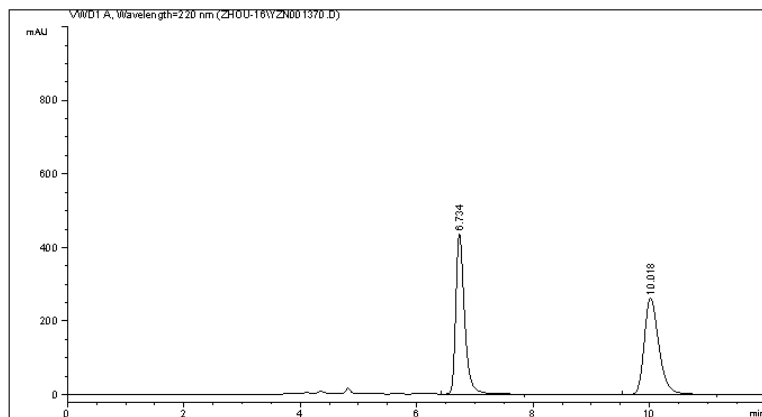
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001370.D
Sample Name: BS-6-6D(+/-)

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/16/2016 12:06:40 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/16/2016 12:04:18 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 6/29/2016 3:25:18 PM
(modified after loading)
Sample Info : 0D-H, Hex/i-PrOH =70/30, 0.7 mL/min, 30oC, 220 nm



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

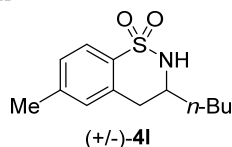
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	6.734	VB	0.1604	4637.92090	436.22708	50.4776	
2	10.018	VB	0.2671	4550.14746	261.48633	49.5224	

Totals : 9188.06836 697.71341

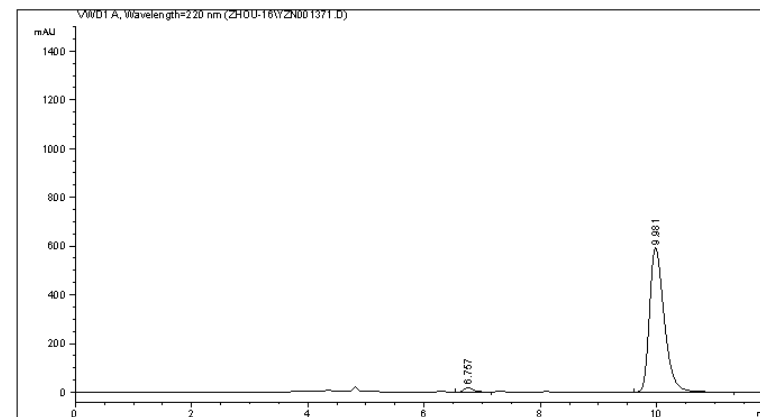
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001371.D
Sample Name: BS-6-6D

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 5/16/2016 12:21:28 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 5/16/2016 12:19:16 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 6/29/2016 3:25:42 PM
(modified after loading)
Sample Info : 0D-H, Hex/i-PrOH =70/30, 0.7 mL/min, 30oC, 220 nm



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

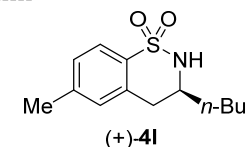
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	6.757	VV	0.1666	213.45477	19.12925	2.0143	
2	9.981	BB	0.2688	1.03836e4	591.72925	97.9857	

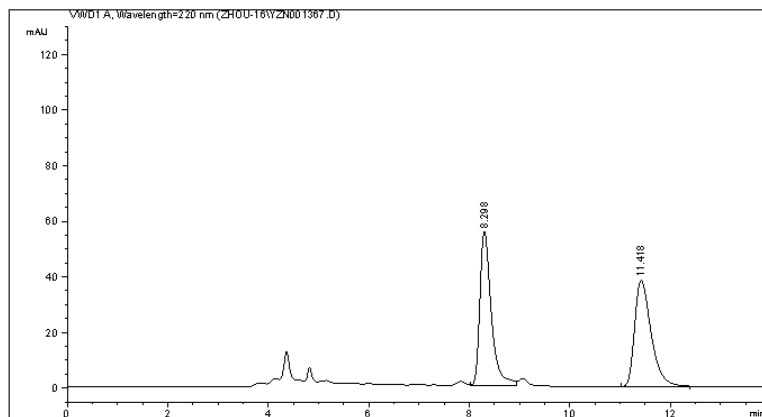
Totals : 1.05971e4 610.85850

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



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001367.D
Sample Name: BS-6-6B(+)

```
=====
Acq. Operator   :
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 5/15/2016 11:09:59 PM
Acq. Method     : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed    : 5/15/2016 11:08:07 PM
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed    : 6/29/2016 3:20:42 PM
                  (modified after loading)
Sample Info     : 0D-H, Hex/i-PrOH =70/30, 0.7 mL/min, 30oC, 220 nm
=====
```



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

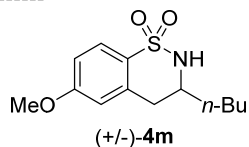
```
Sorted By      :      Signal
Multiplier:    :      1.0000
Dilution:      :      1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	*s	Height [mAU]	Area %
1	8.298	VV	0.2384	879.48865		55.61186	50.1764
2	11.418	BB	0.3472	873.30591		38.35274	49.8236

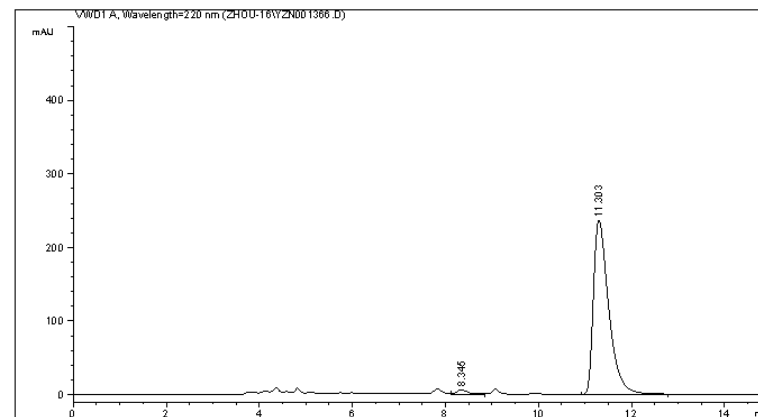
Totals : 1752.79456 93.96460

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



Data File C:\CHEM32\1\DATA\ZHOU-16\YZN001366.D
Sample Name: BS-6-6B

```
=====
Acq. Operator   :
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 5/15/2016 10:47:58 PM
Acq. Method     : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed    : 5/15/2016 10:44:50 PM
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed    : 6/29/2016 3:21:18 PM
                  (modified after loading)
Sample Info     : 0D-H, Hex/i-PrOH =70/30, 0.7 mL/min, 30oC, 220 nm
=====
```



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

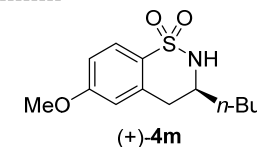
```
Sorted By      :      Signal
Multiplier:    :      1.0000
Dilution:      :      1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	*s	Height [mAU]	Area %
1	8.345	VB	0.2472	91.52423		5.56619	1.6730
2	11.303	BB	0.3438	5379.23779		236.68463	98.3270

Totals : 5470.76202 242.25082

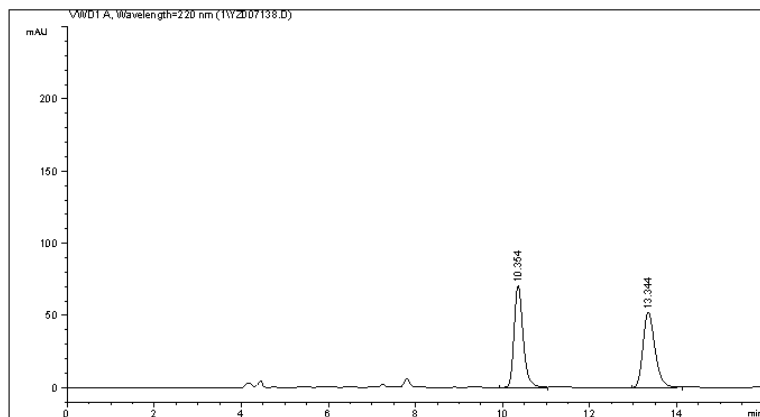
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\1\YZ007138.D
Sample Name: BS-3-82A(+)

=====

Acq. Operator	: ZHOU	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 1/30/2015 8:12:35 AM	
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC.M	
Last changed	: 1/30/2015 7:33:19 AM by ZHOU	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 2:23:17 PM by	
	(modified after loading)	
Sample Info	: OD-H, H/i-PrOH = 80/25, 0.7 mL/min, 30 oC, 220 nm	



=====

Area Percent Report

=====

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

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	10.354	BB	0.2233	1034.59033	70.64894	50.8032	
2	13.344	BB	0.2953	1001.87653	52.18114	49.1968	

Totals : 2036.46686 122.83009

=====

*** End of Report ***

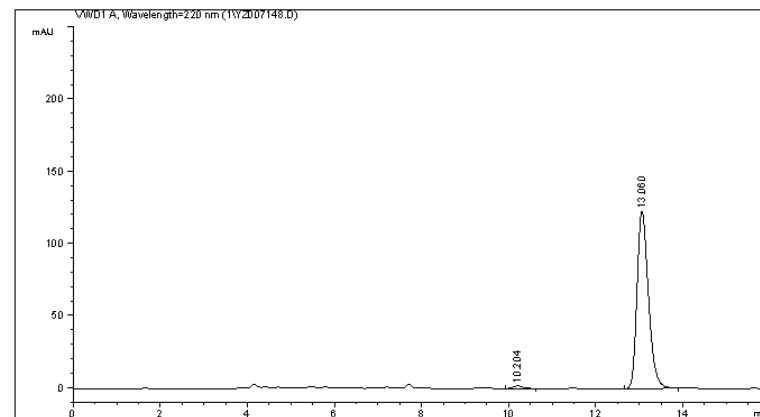
Instrument 1 7/21/2015 2:23:21 PM

Page 1 of 1

Data File C:\CHEM32\1\DATA\1\YZ007148.D
Sample Name: BS-3-86G

=====

Acq. Operator	: ZHOU	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 1/31/2015 10:53:59 AM	
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC.M	
Last changed	: 1/31/2015 9:23:15 AM by ZHOU	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 2:23:17 PM by	
	(modified after loading)	
Sample Info	: OD-H, H/i-PrOH = 80/20, 0.7 mL/min, 30 oC, 220 nm	



=====

Area Percent Report

=====

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

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s	Area [mAU]	Area %
1	10.204	BB	0.2429	28.78096	1.84744	1.2720	
2	13.060	BB	0.2791	2233.83643	122.87923	98.7280	

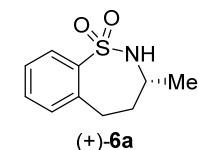
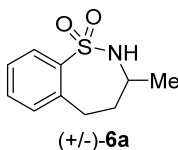
Totals : 2262.61739 124.72667

=====

*** End of Report ***

Instrument 1 7/21/2015 2:23:32 PM

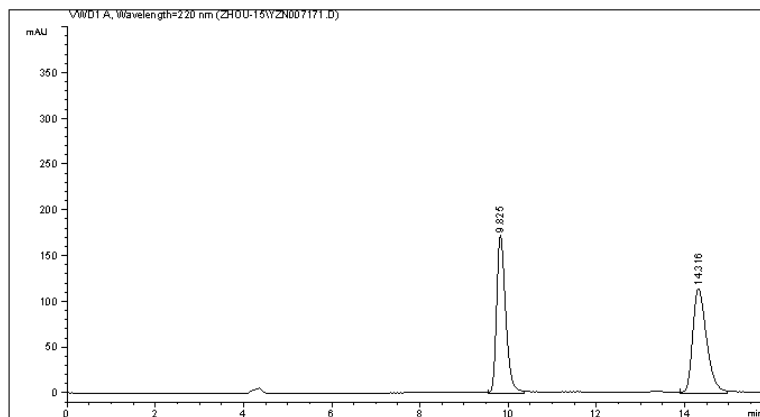
Page 1 of 1



Data File C:\CHEM32\1\DATA\ZHOU-15\YZN007171.D
Sample Name: BS-3-88A(+)

=====

Acq. Operator	: Z	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 2/1/2015 8:01:39 PM	
Acq. Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 2/1/2015 7:39:40 PM by Z	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 2:27:16 PM by	
	(modified after loading)	
Sample Info	: 0D-H, H/i-PrOH = 90/10, 0.7 mL/min, 30 oC, 220 nm	



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

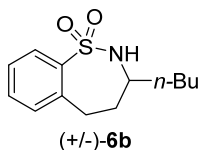
Sorted By	:	Signal	
Multiplier:	:	1.0000	
Dilution:	:	1.0000	
Sample Amount:	:	1.00000 [ng/ul]	(not used in calc.)
Use Multiplier & Dilution Factor with ISTDs			

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s [mAU]	Area %
1	9.825	VV	0.2191	2487.97339	172.65929	49.7009
2	14.316	VV	0.3362	2517.92163	114.69238	50.2991

Totals : 5005.89502 287.35167

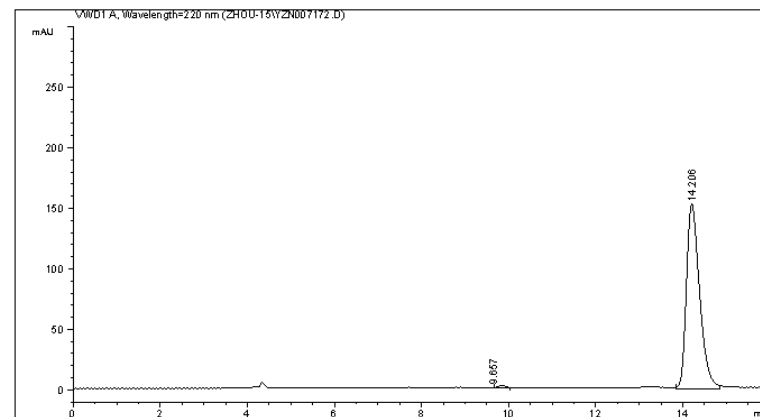
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-15\YZN007172.D
Sample Name: BS-3-88A

=====

Acq. Operator	: Z	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 2/1/2015 8:25:17 PM	
Acq. Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 2/1/2015 8:24:33 PM by Z	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 2:37:39 PM by	
	(modified after loading)	
Sample Info	: 0D-H, H/i-PrOH = 90/10, 0.7 mL/min, 30 oC, 220 nm	



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

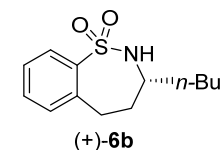
Sorted By	:	Signal	
Multiplier:	:	1.0000	
Dilution:	:	1.0000	
Sample Amount:	:	1.00000 [ng/ul]	(not used in calc.)
Use Multiplier & Dilution Factor with ISTDs			

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height *s [mAU]	Area %
1	9.657	MM R	0.2080	18.86502	2.92566e-2	0.5847
2	14.206	VV	0.3185	3207.59644	152.28621	99.4153

Totals : 3226.46146 152.31547

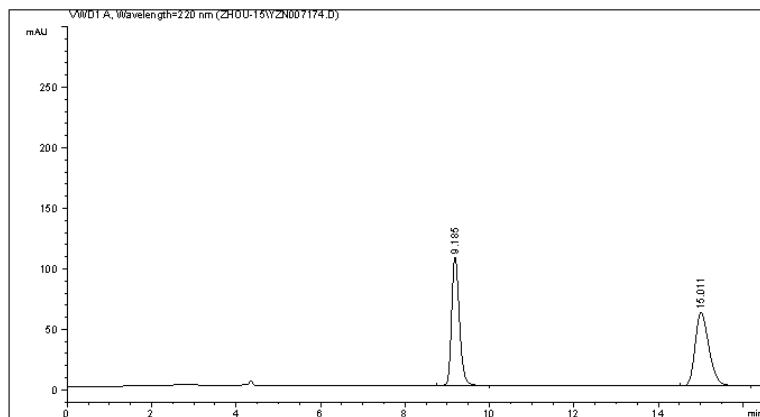
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-15\YZN007174.D
Sample Name: BS-3-88B(+)

=====

Acq. Operator : Z
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 2/1/2015 9:12:16 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 2/1/2015 9:09:53 PM by Z
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 7/21/2015 2:42:22 PM by
(modified after loading)
Sample Info : OD-H, H/i-PrOH = 90/10, 0.7 mL/min, 30 oC, 220 nm



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

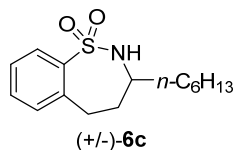
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Sample Amount: : 1.00000 [ng/ul] (not used in calc.)
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	9.185	VB	0.1960	1364.96655	106.58108	49.7854
2	15.011	VB	0.3495	1376.73608	60.60932	50.2146

Totals : 2741.70264 167.19040

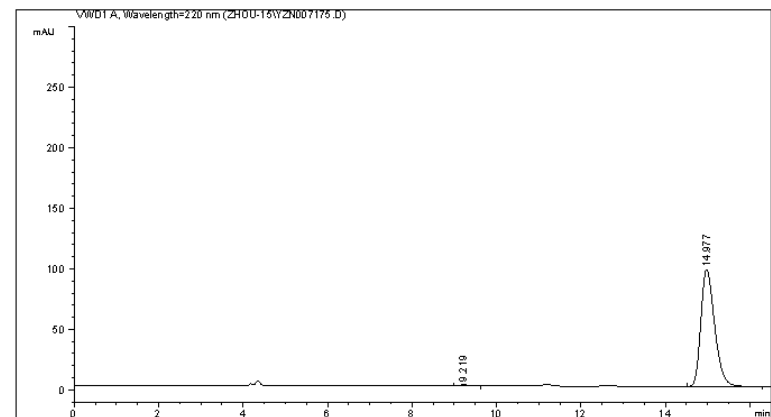
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-15\YZN007175.D
Sample Name: BS-3-88B

=====

Acq. Operator : Z
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 2/1/2015 9:34:02 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 2/1/2015 9:30:18 PM by Z
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 7/21/2015 2:42:22 PM by
(modified after loading)
Sample Info : OD-H, H/i-PrOH = 90/10, 0.7 mL/min, 30 oC, 220 nm



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

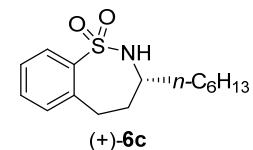
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Sample Amount: : 1.00000 [ng/ul] (not used in calc.)
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	9.219	VB	0.2109	17.20287	1.20025	0.7804
2	14.977	VB	0.3484	2187.04907	96.64989	99.2196

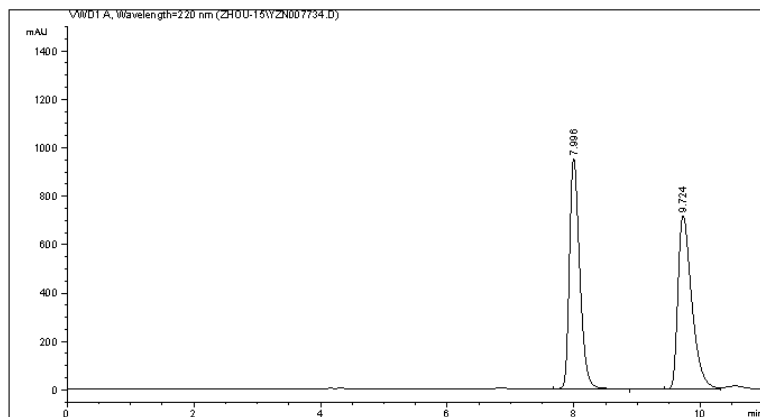
Totals : 2204.25194 97.85014

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



Data File C:\CHEM32\1\DATA\ZHOU-15\YZN007734.D
Sample Name: BS-4-12A(+)

```
=====
Acq. Operator   : Z
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 4/20/2015 7:56:09 PM
Acq. Method     : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 4/20/2015 7:53:53 PM by Z
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 7/21/2015 2:59:49 PM by
                  (modified after loading)
Sample Info     : OD-H, H/i-PrOH = 85/15, 0.7 mL/min, 30 oC, 220 nm
=====
```



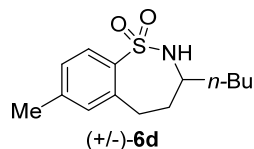
```
=====
                        Area Percent Report
=====
Sorted By      :      Signal
Multiplier:    :      1.0000
Dilution:      :      1.0000
Sample Amount: :      1.00000 [ng/ul] (not used in calc.)
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area [mAU]	Area %
1	7.996	VV	0.1747	1.08359e4	953.94543	49.8263	
2	9.724	VV	0.2313	1.09114e4	717.46490	50.1737	

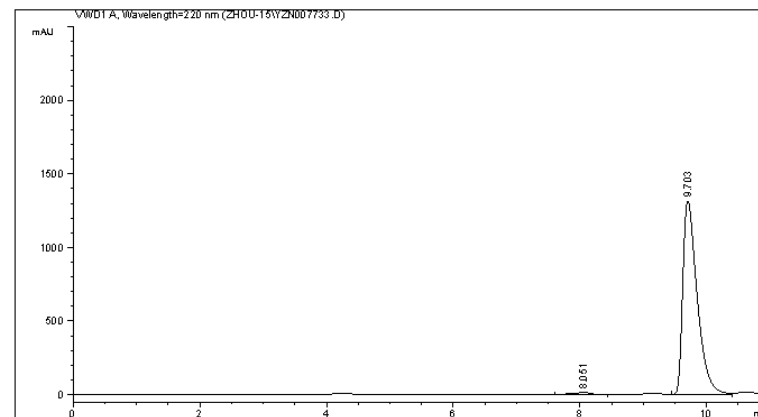
Totals : 2.17473e4 1671.41034

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



Data File C:\CHEM32\1\DATA\ZHOU-15\YZN007733.D
Sample Name: BS-4-12A

```
=====
Acq. Operator   : Z
Acq. Instrument : Instrument 1          Location : Vial 1
Injection Date  : 4/20/2015 7:38:18 PM
Acq. Method     : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 4/20/2015 7:09:16 PM by Z
                  (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed    : 7/21/2015 3:00:33 PM by
                  (modified after loading)
Sample Info     : OD-H, H/i-PrOH = 85/15, 0.7 mL/min, 30 oC, 220 nm
=====
```



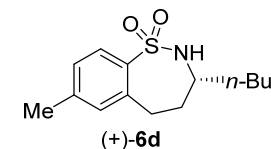
```
=====
                        Area Percent Report
=====
Sorted By      :      Signal
Multiplier:    :      1.0000
Dilution:      :      1.0000
Sample Amount: :      1.00000 [ng/ul] (not used in calc.)
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area [mAU]	Area %
1	8.051	VV	0.2357	235.93143	14.33735	1.1275	
2	9.703	VV	0.2372	2.06883e4	1316.68762	98.8725	

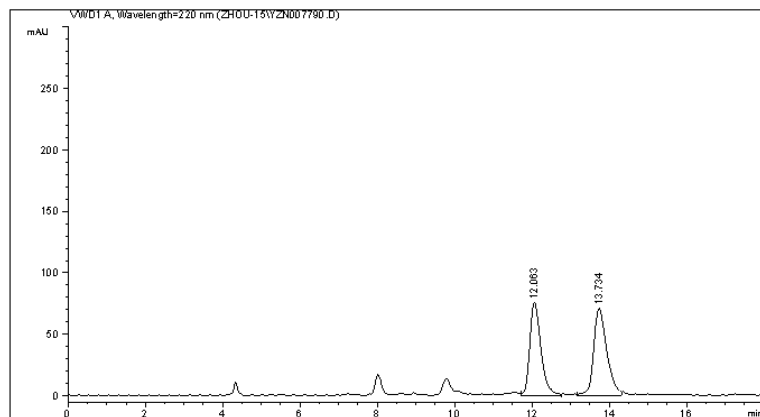
Totals : 2.09243e4 1331.02497

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



Data File C:\CHEM32\1\DATA\ZHOU-15\YZN007790.D
Sample Name: BS-4-15a(+)

=====
Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 4/23/2015 8:28:33 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 4/23/2015 7:58:59 PM by Z
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 7/21/2015 3:03:10 PM by
(modified after loading)
Sample Info : OD-H, H/i-PrOH = 85/15, 0.7 mL/min, 30 oC, 220 nm



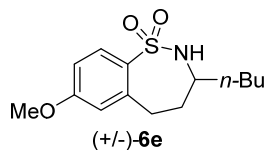
=====
Area Percent Report
=====
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Sample Amount: : 1.00000 [ng/ul] (not used in calc.)
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	12.063	VV	0.2939	1505.38696	75.46680	47.6938
2	13.734	VV	0.3571	1650.96960	70.64191	52.3062

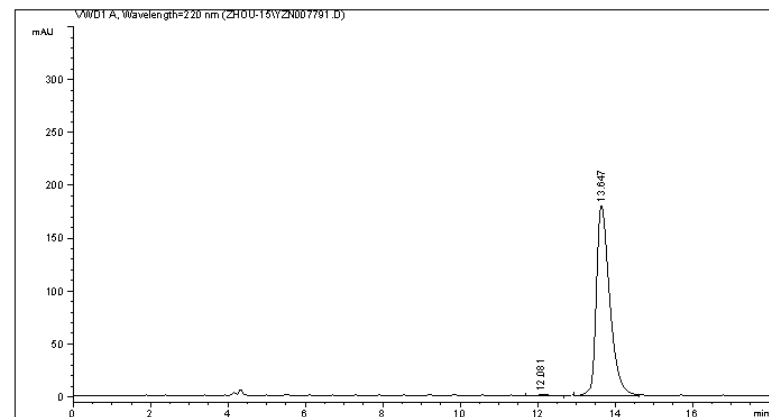
Totals : 3156.35657 146.10872

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



Data File C:\CHEM32\1\DATA\ZHOU-15\YZN007791.D
Sample Name: BS-4-15a

=====
Acq. Operator :
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 4/23/2015 9:01:11 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 4/23/2015 8:57:34 PM by
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 7/21/2015 3:04:45 PM by
(modified after loading)
Sample Info : OD-H, H/i-PrOH = 85/15, 0.7 mL/min, 30 oC, 220 nm



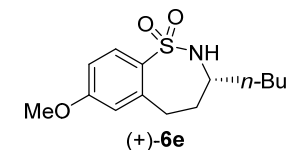
=====
Area Percent Report
=====
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Sample Amount: : 1.00000 [ng/ul] (not used in calc.)
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	12.081	BB	0.1690	22.03906	1.66512	0.5079
2	13.647	VV	0.3594	4316.92188	180.28851	99.4921

Totals : 4338.96094 181.95363

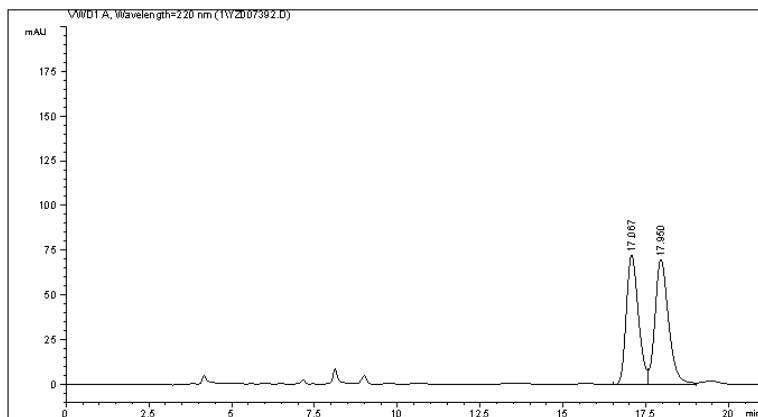
=====
*** End of Report ***



Data File C:\CHEM32\1\DATA\1\YZ007392.D
Sample Name: BS-3-93C+-)

=====

Acq. Operator	: ZHOU	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 3/27/2015 1:58:49 PM	
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC.M	
Last changed	: 3/27/2015 12:13:49 PM by ZHOU	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 2:45:14 PM by	
	(modified after loading)	
Sample Info	: 0D-H, H/i-PrOH = 85/15, 0.7 mL/min, 30 oC, 220 nm	



=====

Area Percent Report

=====

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

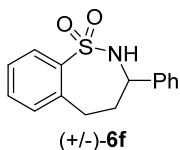
Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	17.067	VV	0.3931	1853.87573	72.82408	47.9332
2	17.950	VV	0.4368	2013.74890	70.15835	52.0668

Totals : 3867.62463 142.98243

=====

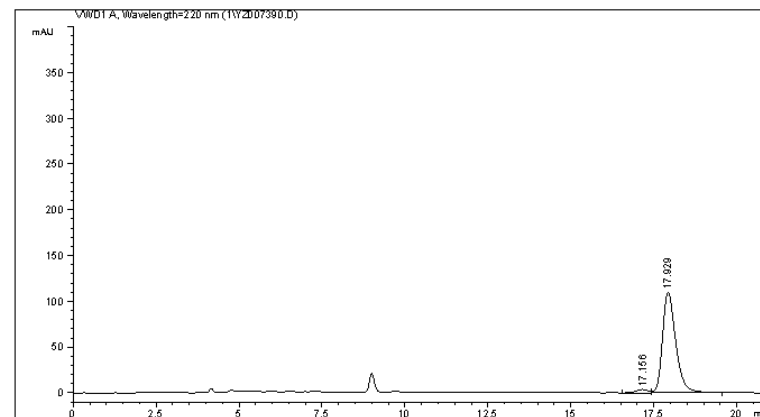
*** End of Report ***



Data File C:\CHEM32\1\DATA\1\YZ007390.D
Sample Name: BS-3-100

=====

Acq. Operator	: ZHOU	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 3/27/2015 1:00:44 PM	
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC.M	
Last changed	: 3/27/2015 12:13:49 PM by ZHOU	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 2:46:03 PM by	
	(modified after loading)	
Sample Info	: 0D-H, H/i-PrOH = 85/15, 0.7 mL/min, 30 oC, 220 nm	



=====

Area Percent Report

=====

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

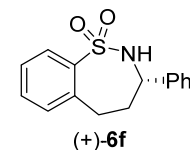
Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	17.156	VV	0.3757	90.71048	3.68929	2.8918
2	17.929	VB	0.4260	3046.06055	109.65189	97.1082

Totals : 3136.77103 113.34118

=====

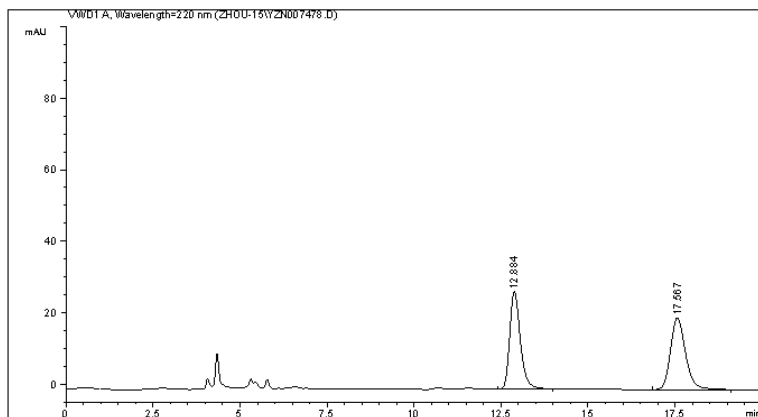
*** End of Report ***



Data File C:\CHEM32\1\DATA\ZHOU-15\YZN007478.D
Sample Name: BS-3-101A(+)

=====

Acq. Operator : Z
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 3/28/2015 4:44:02 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 3/28/2015 4:41:03 PM by Z
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 7/21/2015 2:56:35 PM by
(modified after loading)
Sample Info : OD-H, H/i-PrOH = 80/20, 0.7 mL/min, 30 oC, 220 nm



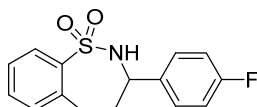
=====
Area Percent Report
=====

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

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	12.884	VB	0.3247	582.10510	27.27588	48.8346
2	17.567	BB	0.4619	609.88812	20.17571	51.1654

Totals : 1191.99323 47.45159



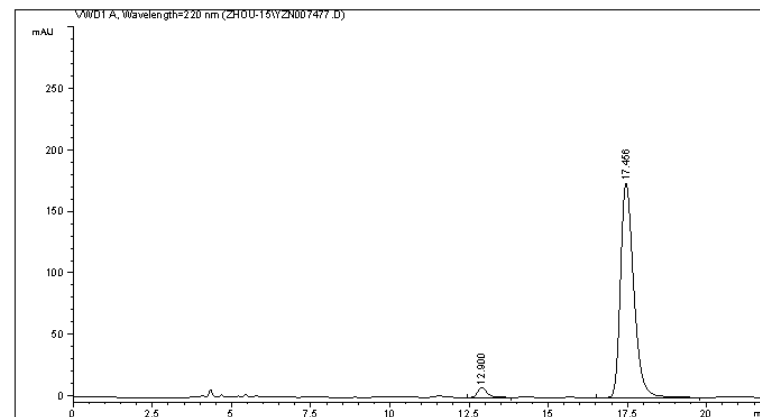
(+/-)-6g

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

Data File C:\CHEM32\1\DATA\ZHOU-15\YZN007477.D
Sample Name: BS-3-101A

=====

Acq. Operator : Z
Acq. Instrument : Instrument 1 Location : Vial 1
Injection Date : 3/28/2015 4:18:23 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 3/28/2015 4:15:04 PM by Z
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF LC.M
Last changed : 7/21/2015 2:56:02 PM by
(modified after loading)
Sample Info : OD-H, H/i-PrOH = 80/20, 0.7 mL/min, 30 oC, 220 nm



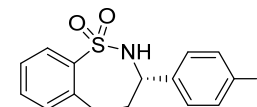
=====
Area Percent Report
=====

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

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	12.900	BB	0.3335	170.85477	7.77588	3.2646
2	17.456	VB	0.4458	5062.65332	173.97386	96.7354

Totals : 5233.50809 181.74974



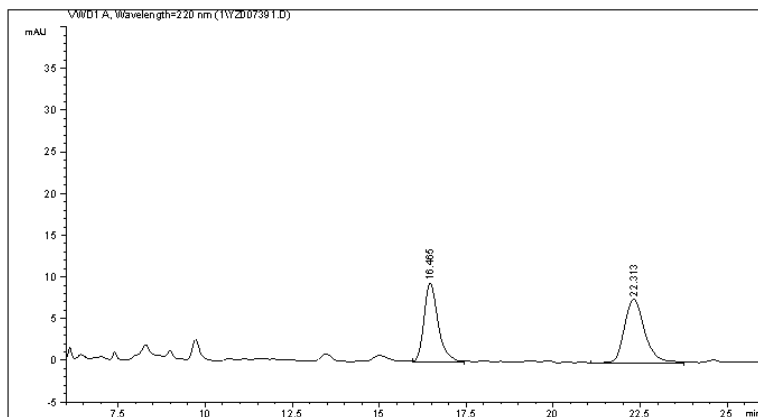
(+)-6g

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

Data File C:\CHEM32\1\DATA\1\YZ007391.D
Sample Name: BS-3-93B(+)

=====

Acq. Operator	: ZHOU	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 3/27/2015 1:30:54 PM	
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC.M	
Last changed	: 3/27/2015 12:13:49 PM by ZHOU	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 2:52:13 PM by	
	(modified after loading)	
Sample Info	: OD-H, H/i-PrOH = 85/15, 0.7 mL/min, 30 oC, 220 nm	



=====

Area Percent Report

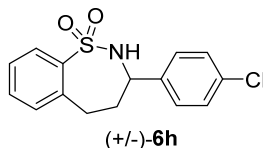
=====

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

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	16.465	VB	0.4302	263.11938	9.39260	45.8103
2	22.313	VV	0.6263	311.24750	7.64610	54.1897

Totals : 574.36688 17.03870



=====

*** End of Report ***

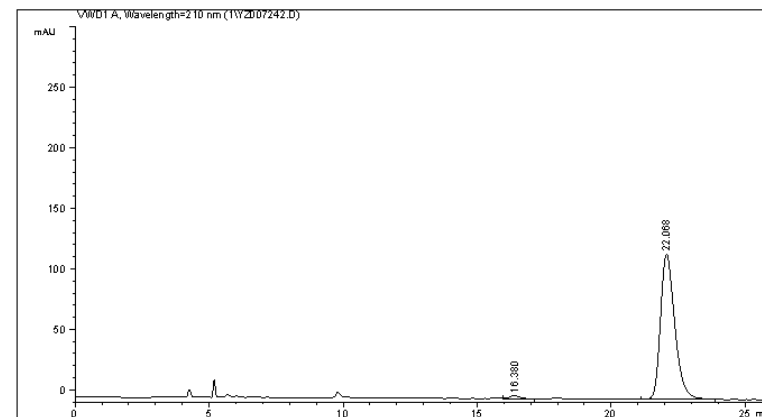
Instrument 1 7/21/2015 2:52:17 PM

Page 1 of 1

Data File C:\CHEM32\1\DATA\1\YZ007242.D
Sample Name: BS-3-93B

=====

Acq. Operator	: ZHOU	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 3/12/2015 2:59:44 AM	
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC.M	
Last changed	: 3/12/2015 1:37:05 AM by ZHOU	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 2:49:43 PM by	
	(modified after loading)	
Sample Info	: OD-H, H/i-PrOH = 85/15, 0.7 mL/min, 30 oC, 210 nm	



=====

Area Percent Report

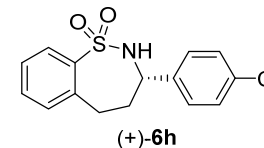
=====

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

Signal 1: VWD1 A, Wavelength=210 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	16.380	VV	0.4164	74.22246	2.65498	1.6434
2	22.068	VB	0.5675	4442.09717	119.93791	98.3566

Totals : 4516.31963 122.59289



=====

*** End of Report ***

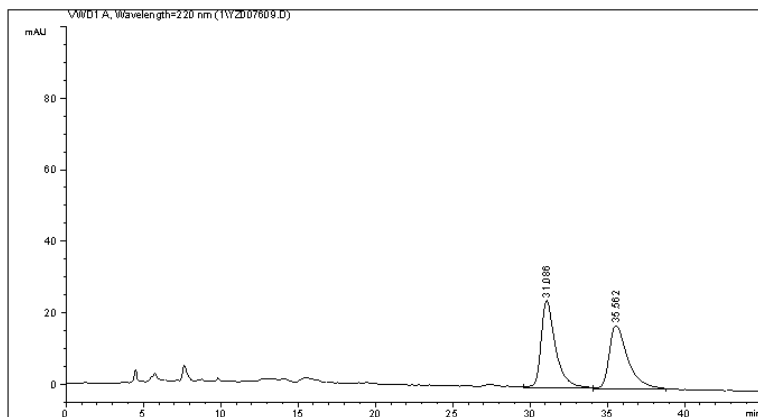
Instrument 1 7/21/2015 2:49:47 PM

Page 1 of 1

Data File C:\CHEM32\1\DATA\1\YZ007609.D
Sample Name: BS-4-12B(+)

=====

Acq. Operator	: ZHOU	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 4/21/2015 1:17:57 PM	
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC.M	
Last changed	: 4/21/2015 1:04:20 PM by ZHOU	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 3:06:53 PM by	
	(modified after loading)	
Sample Info	: 0J-H, H/i-PrOH = 85/15, 0.7 mL/min, 30 oC, 220 nm	



=====

Area Percent Report

=====

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

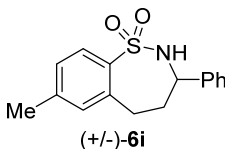
Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	31.086	VB	0.9390	1593.88074	24.42246	51.8170
2	35.562	BV	1.1891	1482.09924	17.65445	48.1830

Totals : 3075.97998 42.07691

=====

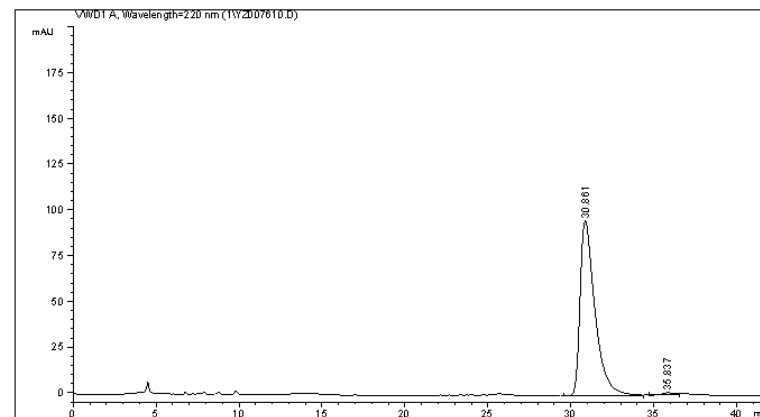
*** End of Report ***



Data File C:\CHEM32\1\DATA\1\YZ007610.D
Sample Name: BS-4-12B

=====

Acq. Operator	: ZHOU	
Acq. Instrument	: Instrument 1	Location : Vial 1
Injection Date	: 4/21/2015 2:10:05 PM	
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC.M	
Last changed	: 4/21/2015 1:04:20 PM by ZHOU	
	(modified after loading)	
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M	
Last changed	: 7/21/2015 3:07:47 PM by	
	(modified after loading)	
Sample Info	: 0J-H, H/i-PrOH = 85/15, 0.7 mL/min, 30 oC, 220 nm	



=====

Area Percent Report

=====

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

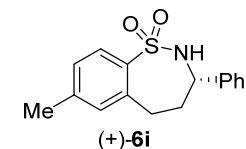
Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	30.861	VV	0.9154	5946.64941	95.36039	98.8726
2	35.837	VV	0.8024	67.80735	1.13398	1.1274

Totals : 6014.45676 96.49437

=====

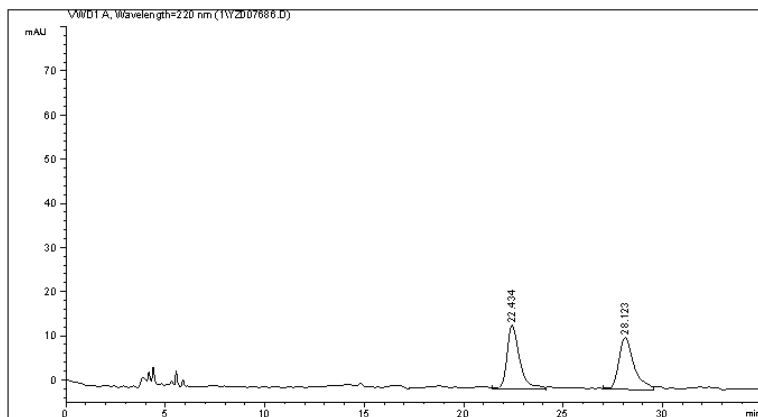
*** End of Report ***



Data File C:\CHEM32\1\DATA\1\YZ007686.D
Sample Name: BS-4-15B

=====

Acq. Operator	: ZHOU	Location	: Vial 1
Acq. Instrument	: Instrument 1		
Injection Date	: 4/30/2015 12:50:16 AM		
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC.M		
Last changed	: 4/29/2015 11:28:53 PM by ZHOU		
	(modified after loading)		
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M		
Last changed	: 7/21/2015 3:10:25 PM by		
	(modified after loading)		
Sample Info	: OD-H, H/i-PrOH = 85/15, 0.7 mL/min, 30 oC, 220 nm		



=====

Area Percent Report

=====

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

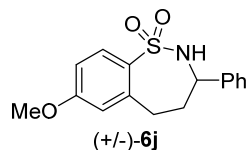
Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	22.434	VV	0.6673	640.29022	14.39853	50.4002
2	28.123	VV	0.7899	630.12213	11.78327	49.5998

Totals : 1270.41235 26.18180

=====

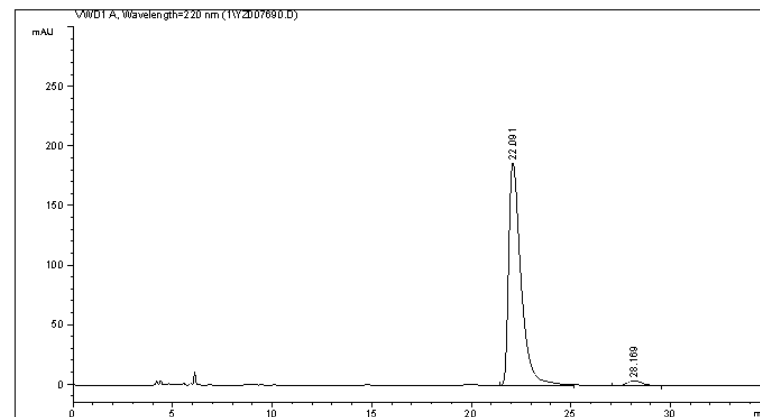
*** End of Report ***



Data File C:\CHEM32\1\DATA\1\YZ007690.D
Sample Name: BS-4-15B

=====

Acq. Operator	: ZHOU	Location	: Vial 1
Acq. Instrument	: Instrument 1		
Injection Date	: 4/30/2015 6:28:31 AM		
Acq. Method	: C:\HPCHEM\1\METHODS\DEF LC.M		
Last changed	: 4/30/2015 5:01:37 AM by ZHOU		
	(modified after loading)		
Analysis Method	: C:\CHEM32\1\METHODS\DEF LC.M		
Last changed	: 7/21/2015 3:11:05 PM by		
	(modified after loading)		
Sample Info	: OD-H, H/i-PrOH = 85/15, 0.7 mL/min, 30 oC, 220 nm		



=====

Area Percent Report

=====

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

Signal 1: VWD1 A, Wavelength=220 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height [mAU]	Area %
1	22.091	VB	0.6519	8123.09717	187.17920	96.8850
2	28.169	VB	0.8588	261.16965	4.44451	3.1150

Totals : 8384.26682 191.62371

=====

*** End of Report ***

