

Supporting Information

Transfer-Catalyst-Free Biomimetic Asymmetric Reduction of 3-Sulfonyl Coumarins with Regenerable NAD(P)H Model

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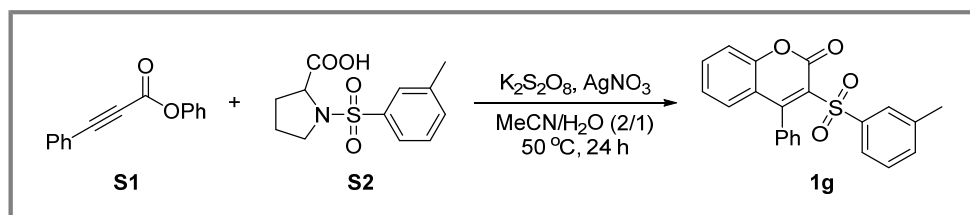
1. General and Materials

General: All reactions were carried out under an atmosphere of nitrogen using the standard Schlenk techniques, unless otherwise noted. ^1H NMR, ^{13}C NMR spectra were recorded at 400 MHz and 100 MHz with Bruker spectrometer. ^{19}F was recorded at 376 MHz with Bruker spectrometer. Chemical shifts are reported in ppm using tetramethylsilane (0) as internal standard when using CDCl_3 as solvent for ^1H NMR spectra. The following abbreviations were used to symbolize the multiplicities: s = singlet, d = doublet, t = triplet, q = quartet, m = multiplet, br = broad. Flash column chromatography was performed on silica gel (200-300 mesh). All reactions were monitored by TLC analysis. Optical rotations were measured by polarimeter. Enantiomeric excess was determined by HPLC analysis using chiral column described below in detail. High resolution mass spectra were recorded using Q-TOF time-of-flight mass spectrometer. X-ray crystallography data were collected using a Bruker D8 Venture with 3.0 ius cu and 3.0 ius mo.

Materials: Commercially available reagents and solvents were used throughout without further purification. The NAD(P)H Models CYNAMs **3a-3c** could be synthesized according to the known literature procedure.¹

2. Synthesis of Substrate 3-Sulfonyl Coumarins

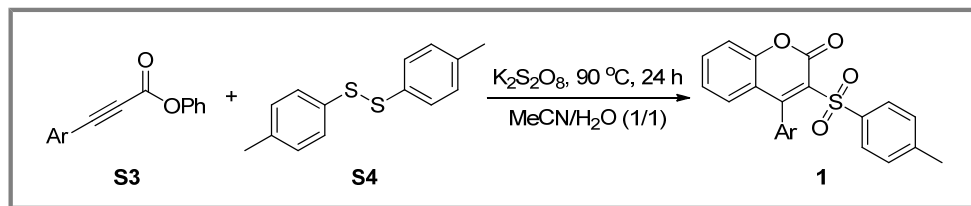
3-Sulfonyl coumarins **1** could be conveniently prepared from aryl alkynoate and *N*-tosylproline according to the known literature procedure with minor modification.²⁻³ Among them, compounds **1g**, **1l**, **1n**, **1p** and **1t** are the unknown compounds.



To a 250 mL Schlenk flask were added phenyl 3-phenylpropiolate **S1** (1.111 g, 5.0 mmol), *N*-tosylproline **S2** (1.616 g, 6.0 mmol), potassium persulfate ($\text{K}_2\text{S}_2\text{O}_8$, 2.703 g, 10.0 mmol), silver nitrate (0.043 g, 0.25 mmol, 5 mol%) and acetonitrile/water (2:1, 90 mL). Then the bottle was charged with nitrogen, and the reaction mixture was stirred at $50\text{ }^\circ\text{C}$ for 24 hours. The resulting mixture was extracted with ethyl acetate. The combined organic layers were dried over anhydrous sodium sulfate, filtered, and concentrated under the reduced pressure. The resulting residue was purified by column chromatography on silica gel using petroleum ether and ethyl acetate as eluent to afford the desired product 3-sulfonyl coumarin **1g**.

4-Phenyl-3-(*m*-tolylsulfonyl)-2H-chromen-2-one (**1g**):

0.928 g, 49% yield, white solid, new compound, mp = $166\text{--}167\text{ }^\circ\text{C}$, R_f = 0.51 (hexanes/ethyl acetate 3/1). ^1H NMR (400 MHz, CDCl_3) δ 7.81 (s, 1H), 7.79-7.70 (m, 1H), 7.63-7.56 (m, 4H), 7.41-7.31 (m, 5H), 7.22-7.16 (m, 1H), 7.05-6.99 (m, 1H), 2.41 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 159.4, 155.6, 153.9, 140.1, 138.8, 134.7, 134.5, 132.6, 130.0, 129.4, 129.3, 128.5, 128.2, 127.5, 126.2, 126.0, 124.9, 120.3, 116.8, 21.4. HRMS (ESI) m/z : $[\text{M}+\text{H}]^+$ Calcd for $\text{C}_{22}\text{H}_{17}\text{O}_4\text{S}$ 377.0842, found: 377.0846.



To a 100 mL flask were added aryl alkynoate **S3** (5.0 mmol), bis(4-methylphenyl) disulfide **S4** (7.5 mmol), potassium persulfate (4.055 g, 15.0 mmol) and acetonitrile/water (1:1, 30 mL). Then the bottle was charged with nitrogen, and the reaction mixture was stirred at 90 °C for 24 h. The resulting mixture was extracted with ethyl acetate. The combined organic layers were dried over anhydrous sodium sulfate, filtered, and concentrated under the reduced pressure. The residue was purified by column chromatography (hexanes and ethyl acetate) to afford the products **1**.

4-(4-Fluorophenyl)-3-tosyl-2H-chromen-2-one (**1l**):

0.763 g, 39% yield, white solid, new compound, mp = 165-166 °C, R_f = 0.37 (hexanes/ethyl acetate 5/1). ^1H NMR (400 MHz, CDCl_3) δ 7.87 (d, J = 8.3 Hz, 2H), 7.65-7.60 (m, 1H), 7.38-7.27 (m, 7H), 7.24-7.19 (m, 1H), 7.04-7.01 (m, 1H), 2.42 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 163.2 (d, $^1J_{\text{F-C}}$ = 247.9 Hz), 158.2, 155.5, 153.9, 145.0, 137.1, 134.8, 129.6 (d, $^3J_{\text{F-C}}$ = 13.1 Hz), 129.5, 129.4, 129.3, 128.4 (d, $^4J_{\text{F-C}}$ = 3.7 Hz), 126.6, 125.0, 120.2, 116.9, 115.6 (d, $^2J_{\text{F-C}}$ = 21.8 Hz), 21.7. ^{19}F NMR (376 MHz, CDCl_3) δ -111.27. HRMS (ESI) m/z : $[\text{M}+\text{H}]^+$ Calcd for $\text{C}_{22}\text{H}_{16}\text{FO}_4\text{S}$ 395.0748, found: 395.0747.

4-(4-Bromophenyl)-3-tosyl-2H-chromen-2-one (**1n**):

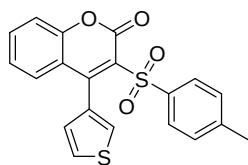
1.590 g, 40% yield, white solid, new compound, mp = 237-238 °C, R_f = 0.36 (hexanes/ethyl acetate 5/1). ^1H NMR (400 MHz, CDCl_3) δ 7.87 (d, J = 8.2 Hz, 2H), 7.72 (d, J = 8.3 Hz, 2H), 7.65-7.59 (m, 1H), 7.36-7.28 (m, 3H), 7.25-7.18 (m, 3H), 7.03-6.98 (m, 1H), 2.42 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 157.8, 155.4, 153.9, 145.0, 137.0, 134.8, 131.6, 131.6, 129.6, 129.4, 129.3, 129.1, 126.5, 125.0, 123.7, 119.9, 117.0, 21.7. HRMS (ESI) m/z : $[\text{M}+\text{H}]^+$ Calcd for $\text{C}_{22}\text{H}_{16}\text{BrO}_4\text{S}$ 454.9947 (^{79}Br) and 456.9934 (^{81}Br), found: 454.9945 (^{79}Br) and 456.9925 (^{81}Br).

4-(*o*-Tolyl)-3-tosyl-2H-chromen-2-one (**1p**):

2.956 g, 41% yield, orange solid, new compound, mp = 122-123 °C, R_f = 0.25 (hexanes/ethyl acetate 5/1). ^1H NMR (400 MHz, CDCl_3) δ 7.90 (d, J = 8.3 Hz, 2H), 7.63-7.58 (m, 1H), 7.49-7.45 (m, 1H), 7.41-7.33 (m, 3H), 7.30 (d, J = 8.2 Hz, 2H), 7.17 (t, J = 7.7 Hz, 1H), 7.11 (d, J = 7.4 Hz, 1H), 6.91 (dd, J = 8.1, 1.2 Hz, 1H), 2.41 (s, 3H), 2.21 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 159.4, 155.7, 154.0, 144.9, 137.2, 135.1, 134.7, 132.7, 130.1, 129.4, 129.4, 129.3, 129.3, 126.6, 126.4, 125.6, 125.1, 119.5, 116.9, 21.7, 20.0. HRMS (ESI) m/z : $[\text{M}+\text{H}]^+$ Calcd for $\text{C}_{23}\text{H}_{19}\text{O}_4\text{S}$ 391.0999, found: 391.1000.

4-(Thiophen-3-yl)-3-tosyl-2H-chromen-2-one (1t):

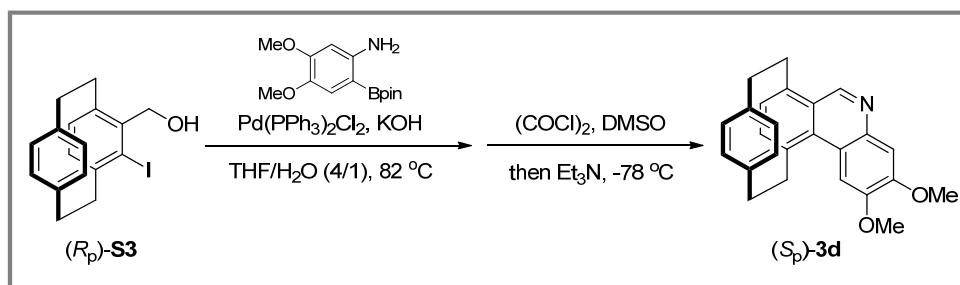
0.207 g, 11% yield, brown solid, new compound, mp = 157-158 °C, R_f = 0.35 (hexanes/ethyl acetate 5/1). ¹H NMR (400 MHz, CDCl₃) δ 7.88 (d, *J* = 8.2 Hz, 2H), 7.64-



7.55 (m, 2H), 7.41-7.37 (m, 1H), 7.34-7.27 (m, 3H), 7.25-7.14 (m, 3H), 2.41 (s, 3H). ¹³C NMR (100 MHz, CDCl₃) δ 155.5, 155.4, 153.7, 144.8, 137.3, 134.6, 131.5, 129.5, 129.3, 129.2, 128.3, 127.1, 126.2, 124.9, 120.2, 116.8, 21.7. HRMS (ESI) *m/z*: [M+H]⁺ Calcd for C₂₀H₁₅O₄S₂ 383.0406,

found: 383.0405.

3. Synthesis of the New NAD(P)H Model CYNAM



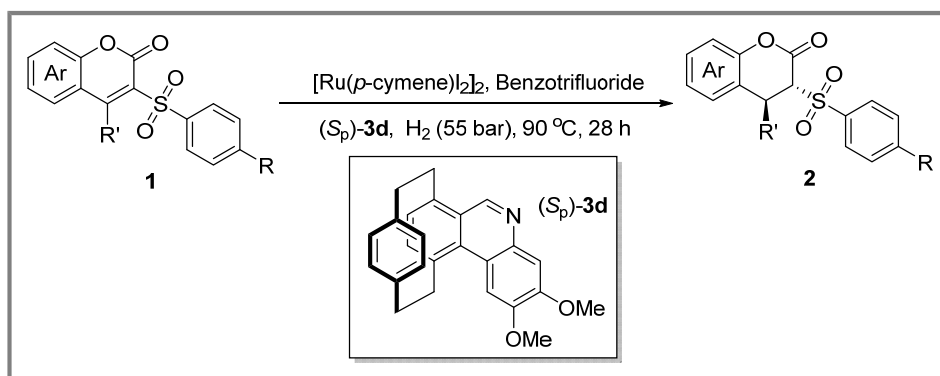
Under nitrogen, a solution of the alcohol (*R_p*)-**S3** (1.275 g, 3.5 mmol), bis(triphenylphosphine)palladium(II) chloride (491 mg, 0.7 mmol), potassium hydroxide (589 mg, 10.5 mmol) and the 2-aminoarylboronic acid pinacolate ester (1.954 g, 7.0 mmol) in tetrahydrofuran (52 mL) and water (13 mL) was heated at 82 °C for 24 h. The mixture was cooled to room temperature and extracted with ethyl acetate (20 mL×3). The combined organic layer was dried over anhydrous sodium sulfate and concentrated in *vacuo*. The residue was purified by flash column chromatography on silica gel using hexanes and ethyl acetate as eluent to achieve the intermediate (*S_p*)-**S4** (0.792 g, 58% yield).

Then, under nitrogen, to a solution of oxalyl chloride (0.434 g, 3.4 mmol, 0.29 mL) in dichloromethane (10 mL) at -78 °C was slowly added a solution of dimethyl sulfoxide (0.534 g, 6.8 mmol, 0.49 mL) in dichloromethane (5 mL). After stirring for 5 min, a solution of (*S_p*)-**S4** (0.666 g, 1.7 mmol) in dichloromethane (5 mL) was added dropwise. After the mixture was stirred for 15 min, triethylamine (1.730 g, 17.1 mmol, 2.4 mL) was then added dropwise. 10 min later, the mixture was warmed to room temperature and stirred for 3 h. The reaction mixture was quenched with water (10 mL). After being extracted with dichloromethane and washed with water twice, the combined organic layer was dried over anhydrous sodium sulfate and concentrated in *vacuo*. The residue was purified by flash column chromatography using hexanes and ethyl acetate as eluent to achieve the product NAD(P)H model (*S_p*)-**3d** (0.224 g, 39% yield).

(*S_p*)-2,3-Dimethoxy[2]paracyclo[2](7,10)phenanthridinophane (**3d**):

0.244 g, 39% yield, yellow solid, new compound, mp = 86-87 °C, *R_f* = 0.15 (hexanes/ethyl acetate 1/1), >99% ee, $[\alpha]_D^{20} = +254.23$ (c 0.80, CHCl₃). ¹H NMR (400 MHz, CDCl₃) δ 9.07 (s, 1H), 7.89 (s, 1H), 7.60 (s, 1H), 7.06 (d, *J* = 7.4 Hz, 1H), 6.87 (d, *J* = 7.4 Hz, 1H), 6.56-6.47 (m, 2H), 5.92-5.71 (m, 1H), 5.37-5.29 (m, 1H), 4.35-4.28 (m, 1H), 4.12 (s, 3H), 4.10 (s, 3H), 4.05-3.97 (m, 1H), 3.37-3.25 (m, 2H), 3.20-3.03 (m, 3H), 2.85-2.76 (m, 1H). ¹³C NMR (100 MHz, CDCl₃) δ 150.2, 148.7, 148.3, 141.0, 138.4, 138.3, 138.0, 137.8, 134.6, 134.4, 132.2, 132.1, 131.8, 130.0, 128.1, 128.0, 119.3, 109.6, 106.1, 56.2, 56.1, 38.7, 34.7, 34.1, 32.5. HRMS (ESI) *m/z*: [M+H]⁺ Calcd for C₂₅H₂₄NO₂ 370.1802, found: 370.1801.

4. Biomimetic Asymmetric Reduction of 3-Sulfonyl Coumarins

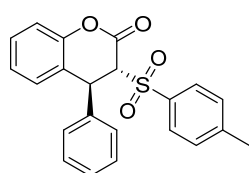


A mixture of $[\text{Ru}(p\text{-cymene})\text{I}_2]_2$ (2.2 mg, 0.00225 mmol), chiral NAD(P)H model (S_p)-**3d** (5.5 mg, 0.015 mmol) and 3-sulfonyl coumarins **1** (0.15 mmol) in benzotrifluoride (3.0 mL) was stirred at room temperature for 5 min in glove box and then the mixture was transferred to an autoclave. The hydrogenation was performed at 90 °C under hydrogen gas (55 bar) for 28 h. After carefully release of the hydrogen gas, the autoclave was opened. The mixture was concentrated under the reduced pressure, and then purified by column chromatography on silica gel using hexanes and ethyl acetate as eluent to give the biomimetic reductive products **2**.

The enantiomeric excesses of the reductive products **2** were determined by chiral HPLC. The racemates could be synthesized under the above condition with readily available phenanthridine instead of chiral NAD(P)H model.

(+)-4-Phenyl-3-tosylchroman-2-one (**2a**):

56 mg, 99% yield, colorless oil, new compound, $R_f = 0.51$ (hexanes/ethyl acetate 5/1), d.r. > 20:1,

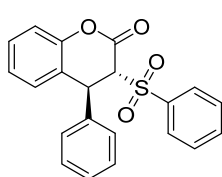


98% ee, $[\alpha]_D^{20} = +160.08$ (c 1.12, CHCl_3), ^1H NMR (400 MHz, CDCl_3) δ 7.71-7.64 (m, 2H), 7.35-7.26 (m, 7H), 7.21-7.16 (m, 1H), 7.09-7.04 (m, 2H), 7.00-6.94 (m, 1H), 5.17 (s, 1H), 4.41 (s, 1H), 2.43 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 159.9, 150.9, 146.0, 139.4, 133.9, 129.9, 129.6, 129.4, 129.2, 129.1, 128.2, 127.1, 125.7, 120.9, 116.9, 72.1, 42.3, 21.8.

HPLC (Chiralpak IA column, $\lambda = 230$ nm, 30 °C, $n\text{-Hexane}/i\text{-PrOH} = 70/30$, flow rate = 1.0 mL/min) retention time = 7.4 min (minor) and 10.5 min (major). HRMS (ESI) m/z : $[\text{M}+\text{NH}_4]^+$ Calcd for $\text{C}_{22}\text{H}_{22}\text{NO}_4\text{S}$ 396.1264, found: 396.1265.

(+)-4-Phenyl-3-(phenylsulfonyl)chroman-2-one (**2b**):

54 mg, 99% yield, colorless oil, new compound, $R_f = 0.51$ (hexanes/ethyl acetate 5/1), d.r. = 32:1,



98% ee, $[\alpha]_D^{20} = +145.18$ (c 1.08, CHCl_3), ^1H NMR (400 MHz, CDCl_3) δ 7.96-7.92 (m, 0.06H), 7.83-7.76 (m, 1.94H), 7.67-7.61 (m, 1H), 7.53-7.46 (m, 2H), 7.34-7.24 (m, 5H), 7.23-7.16 (m, 1H), 7.16-7.13 (m, 0.06H), 7.09-7.05 (m, 1.94H), 7.03-6.88 (m, 1H), 5.19 (s, 0.97H), 5.15 (d, $J = 6.0$ Hz, 0.03H), 4.66 (d, $J = 5.9$ Hz, 0.03H), 4.45 (d, $J = 0.6$ Hz, 0.97H). ^{13}C NMR (100 MHz, CDCl_3) δ 159.8, 150.9, 139.3, 136.9, 134.7, 129.7, 129.6, 129.3, 129.2, 129.1, 128.2,

127.1, 125.8, 120.8, 117.0, 72.0, 42.2. HPLC (Chiralpak IA column, $\lambda = 230$ nm, 30 °C, $n\text{-Hexane}$

i-PrOH = 70/30, flow rate = 1.0 mL/min) retention time = 6.9 min (minor) and 11.4 min (major). HRMS (ESI) *m/z*: [M+NH₄]⁺ Calcd for C₂₁H₂₀NO₄S 382.1108, found: 382.1108.

(+)-3-((4-Methoxyphenyl)sulfonyl)-4-phenylchroman-2-one (2c):

58 mg, 98% yield, colorless oil, new compound, *R_f* = 0.37 (hexanes/ethyl acetate 5/1), d.r. > 20:1, 98% ee, [α]_D²⁰ = +161.71 (*c* 1.16, CHCl₃), ¹H NMR (400 MHz, CDCl₃) δ 7.75-7.62 (m, 2H), 7.33-7.24 (m, 5H), 7.20-7.15 (m, 1H), 7.09-7.04 (m, 2H), 6.99-6.94 (m, 1H), 6.93-6.88 (m, 2H), 5.16 (s, 1H), 4.42 (d, *J* = 0.5 Hz, 1H), 3.87 (s, 3H). ¹³C NMR (100 MHz, CDCl₃) δ 164.5, 160.1, 150.8, 139.5, 131.3, 129.6, 129.5, 129.2, 128.1, 128.1, 127.1, 125.7, 121.0, 116.9, 114.5, 72.2, 55.8, 42.4. HPLC (Chiralpak IA column, λ = 230 nm, 30 °C, *n*-Hexane/*i*-PrOH = 70/30, flow rate = 1.0 mL/min) retention time = 9.0 min (minor) and 13.7 min (major). HRMS (ESI) *m/z*: [M+NH₄]⁺ Calcd for C₂₂H₂₂NO₅S 412.1213, found: 412.1213.

(+)-3-((4-Fluorophenyl)sulfonyl)-4-phenylchroman-2-one (2d):

57 mg, 99% yield, pale yellow solid, new compound, mp = 118-119 °C, *R_f* = 0.52 (hexanes/ethyl acetate 5/1), d.r. = 32:1, 98% ee, [α]_D²⁰ = +131.83 (*c* 1.14, CHCl₃), ¹H NMR (400 MHz, CDCl₃) δ 7.91-7.86 (m, 0.06H), 7.84-7.73 (m, 1.94H), 7.35-7.25 (m, 5H), 7.23-7.18 (m, 1H), 7.14 (t, *J* = 8.5 Hz, 2H), 7.07 (d, *J* = 7.1 Hz, 2H), 6.97 (d, *J* = 8.2 Hz, 1H), 5.19 (s, 0.97H), 5.11 (d, *J* = 6.0 Hz, 0.03H), 4.67 (d, *J* = 6.0 Hz, 0.03H), 4.44 (s, 0.97H). ¹³C NMR (100 MHz, CDCl₃) δ 166.4 (d, ¹*J*_{F-C} = 257.0 Hz), 159.8, 150.8, 139.1, 132.8 (d, ⁴*J*_{F-C} = 3.1 Hz), 132.1 (d, ³*J*_{F-C} = 9.8 Hz), 129.7, 129.6, 129.2, 128.3, 127.0, 125.9, 120.8, 116.8 (d, ²*J*_{F-C} = 19.5 Hz), 116.5, 72.2, 42.3. ¹⁹F NMR (376 MHz, CDCl₃) δ -101.16, -102.69. HPLC (Chiralpak IA, λ = 230 nm, 30 °C, *n*-Hexane/*i*-PrOH = 70/30, flow rate = 1.0 mL/min) retention time = 6.8 min (minor) and 9.7 min (major). HRMS (ESI) *m/z*: [M+NH₄]⁺ Calcd for C₂₁H₁₉FNO₄S 400.1013, found: 400.1018.

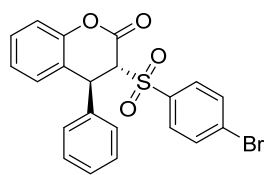
(+)-3-((4-Chlorophenyl)sulfonyl)-4-phenylchroman-2-one (2e):

59 mg, 99% yield, colorless oil, new compound, *R_f* = 0.42 (hexanes/ethyl acetate 5/1), d.r. = 32:1, 98% ee, [α]_D²⁰ = +149.90 (*c* 1.18, CHCl₃), ¹H NMR (400 MHz, CDCl₃) δ 7.77 (d, *J* = 8.6 Hz, 0.06H), 7.68 (d, *J* = 8.6 Hz, 1.94H), 7.41 (d, *J* = 8.6 Hz, 2H), 7.35-7.23 (m, 5H), 7.21-7.15 (m, 1H), 7.13 (d, *J* = 7.9 Hz, 0.06H), 7.04 (d, *J* = 7.0 Hz, 1.94H), 6.98 (d, *J* = 8.0 Hz, 0.03H), 6.94 (d, *J* = 8.1 Hz, 0.97H), 5.15 (s, 0.97H), 5.08 (d, *J* = 6.0 Hz, 0.03H), 4.64 (d, *J* = 6.0 Hz, 0.03H), 4.40 (s, 0.97H). ¹³C NMR (100 MHz, CDCl₃) δ 159.7, 150.7, 141.7, 139.1, 135.2, 130.5, 129.7, 129.7, 129.6, 129.2, 128.3, 127.0, 125.9, 120.7, 117.0, 72.1, 42.2. HPLC (Chiralpak IA column, λ = 230 nm, 30 °C, *n*-Hexane/*i*-PrOH = 70/30, flow rate = 1.0 mL/min) retention time = 7.3 min (minor) and 10.2 min (major). HRMS (ESI) *m/z*: [M+NH₄]⁺ Calcd for C₂₁H₁₉ClNO₄S 416.0718 (³⁵Cl) and 418.0700 (³⁷Cl), found: 416.0715 (³⁵Cl) and 418.0689 (³⁷Cl).

(+)-3-((4-Bromophenyl)sulfonyl)-4-phenylchroman-2-one (2f):

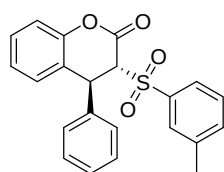
65 mg, 98% yield, colorless solid, new compound, mp = 127-128 °C, *R_f* = 0.43 (hexanes/ethyl acetate 5/1), d.r. > 20:1, 98% ee, [α]_D²⁰ = +145.22 (*c* 1.30, CHCl₃), ¹H NMR (400 MHz, CDCl₃) δ 7.66-7.56 (m, 4H), 7.36-7.26 (m, 5H), 7.23-7.18 (m, 1H), 7.07 (d, *J* = 7.0 Hz, 2H), 6.96 (d, *J* = 8.2

Hz, 1H), 5.17 (s, 1H), 4.43 (s, 1H). ^{13}C NMR (100 MHz, CDCl_3) δ 159.7, 150.7, 139.0, 135.7, 132.6, 130.5, 129.7, 129.7, 129.2, 128.3, 127.0, 125.9, 120.7, 117.0, 72.1, 42.2. HPLC (Chiralpak IA column, λ = 230 nm, 30 °C, *n*-Hexane/*i*-PrOH = 70/30, flow rate = 1.0 mL/min) retention time = 7.7 min (minor) and 10.8 min (major). HRMS (ESI) m/z : $[\text{M}+\text{NH}_4]^+$ Calcd for $\text{C}_{21}\text{H}_{19}\text{BrNO}_4\text{S}$ 460.0213 (^{79}Br) and 462.0200 (^{81}Br), found: 460.0214 (^{79}Br) and 462.0196 (^{81}Br).



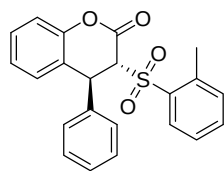
(+)-4-Phenyl-3-(*m*-tolylsulfonyl)chroman-2-one (2g):

56 mg, 99% yield, colorless oil, new compound, R_f = 0.51 (hexanes/ethyl acetate 5/1), d.r. > 20:1, 98% ee, $[\alpha]_D^{20}$ = +147.85 (*c* 1.12, CHCl_3), ^1H NMR (400 MHz, CDCl_3) δ 7.58 (d, *J* = 7.6 Hz, 1H), 7.52 (s, 1H), 7.42-7.32 (m, 2H), 7.32-7.27 (m, 2H), 7.26-7.20 (m, 3H), 7.17-7.13 (m, 1H), 7.04 (d, *J* = 7.1 Hz, 2H), 6.92 (d, *J* = 8.1 Hz, 1H), 5.15 (s, 1H), 4.41 (s, 1H), 2.34 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 159.8, 150.9, 139.7, 139.4, 136.7, 135.5, 129.6, 129.6, 129.3, 129.2, 129.1, 128.2, 127.1, 126.3, 125.6, 120.9, 116.9, 72.1, 42.4, 21.3. HPLC (Chiralpak IA column, λ = 230 nm, 30 °C, *n*-Hexane/*i*-PrOH = 80/20, flow rate = 0.8 mL/min) retention time = 9.2 min (minor) and 12.8 min (major). HRMS (ESI) m/z : $[\text{M}+\text{NH}_4]^+$ Calcd for $\text{C}_{22}\text{H}_{22}\text{NO}_4\text{S}$ 396.1264, found: 396.1263.



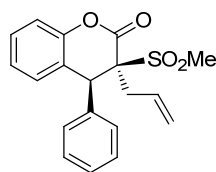
(+)-4-Phenyl-3-(*o*-tolylsulfonyl)chroman-2-one (2h):

40 mg, 70% yield, colorless oil, new compound, R_f = 0.51 (hexanes/ethyl acetate 5/1), d.r. > 20:1, 98% ee, $[\alpha]_D^{20}$ = +178.71 (*c* 0.78, CHCl_3), ^1H NMR (400 MHz, CDCl_3) δ 7.90-7.82 (m, 1H), 7.54-7.49 (m, 1H), 7.36-7.22 (m, 7H), 7.20-7.15 (m, 1H), 7.14-7.09 (m, 1H), 7.06-7.01 (m, 2H), 5.17 (s, 1H), 4.40 (d, *J* = 0.9 Hz, 1H), 2.69 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 159.8, 151.1, 139.6, 138.6, 135.5, 134.8, 133.1, 131.5, 129.7, 129.7, 129.2, 128.2, 127.0, 126.8, 125.8, 121.0, 117.0, 71.2, 41.8, 20.8. HPLC (Chiralpak IA column, λ = 230 nm, 30 °C, *n*-Hexane/*i*-PrOH = 80/20, flow rate = 0.8 mL/min) retention time = 8.8 min (minor) and 18.6 min (major). HRMS (ESI) m/z : $[\text{M}+\text{NH}_4]^+$ Calcd for $\text{C}_{22}\text{H}_{22}\text{NO}_4\text{S}$ 396.1264, found: 396.1266.



(+)-3-Allyl-3-(methylsulfonyl)-4-phenylchroman-2-one (2i):

Procedure for allyl alkylation: Sodium carbonate (63.6 mg, 0.60 mmol), allyl iodide (100.8 mg, 54.8 μL , 0.60 mmol) and 18-crown-6 (11.8 mg, 10.2 μL , 0.045 mmol) were added to the stirred solution of the reductive product of **1i** (45 mg, 0.15 mmol) in tetrahydrofuran (3.0 mL). The mixture was stirred at ambient temperature for 24 h. The mixture was concentrated under the reduced pressure, and then purified by column chromatography on silica gel using hexanes and ethyl acetate as eluent to afford the product (+)-**2i**, 47 mg, 91% yield, white solid, new compound, mp = 63-64 °C, R_f = 0.61 (hexanes/ethyl acetate 10/1), d.r. > 20:1, 98% ee, $[\alpha]_D^{20}$ = +250.09 (*c* 0.94, CHCl_3), ^1H NMR (400 MHz, CDCl_3) δ 7.79-6.87 (m, 9H), 6.16-5.97 (m, 1H), 5.19 (d, *J* = 10.0 Hz, 1H), 4.80 (d, *J* = 17.0 Hz, 1H), 4.53 (s, 1H), 3.06-2.97 (m, 1H), 2.86-2.77 (m, 1H), 2.16 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 164.3, 149.6, 137.1, 130.6, 129.5, 129.2, 128.8, 126.0, 124.8, 121.3, 116.9, 73.6, 49.2, 41.4, 35.8. HPLC (Chiralpak AD-H column, λ = 230 nm, 30 °C,



n-Hexane/*i*-PrOH = 80/20, flow rate = 1.0 mL/min) retention time = 9.0 min (major) and 10.2 min (minor). HRMS (ESI) *m/z*: [M+NH₄]⁺ Calcd for C₁₉H₂₂NO₄S [M+NH₄]⁺ 360.1264, found: 360.1262.

(+)-4-(*p*-Tolyl)-3-tosylchroman-2-one (2j):

56 mg, 95% yield, colorless oil, new compound, *R*_f = 0.37 (hexanes/ethyl acetate 5/1), d.r. > 20:1, 98% ee, [α]²⁰_D = +158.56 (*c* 1.05, CHCl₃), ¹H NMR (400 MHz, CDCl₃) δ 7.63 (d, *J* = 8.3 Hz, 2H), 7.28-7.21 (m, 4H), 7.16-7.12 (m, 1H), 7.08 (d, *J* = 8.0 Hz, 2H), 6.95-6.88 (m, 3H), 5.10 (s, 1H), 4.37 (s, 1H), 2.40 (s, 3H), 2.27 (s, 3H). ¹³C NMR (100 MHz, CDCl₃) δ 160.0, 150.8, 145.9, 138.0, 136.5, 133.9, 130.2, 129.9, 129.3, 129.2, 129.1, 126.9, 125.6, 121.2, 116.9, 72.2, 41.9, 21.7, 21.0. HPLC (Chiralpak ID column, λ = 230 nm, 30 °C, *n*-Hexane/*i*-PrOH = 70/30, flow rate = 1.0 mL/min) retention time = 14.9 min (major) and 18.0 min (minor). HRMS (ESI) *m/z*: [M+NH₄]⁺ Calcd for C₂₃H₂₄NO₄S 410.1421, found: 410.1423.

(+)-4-(4-Methoxyphenyl)-3-tosylchroman-2-one (2k):

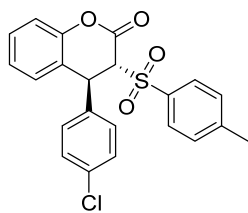
55 mg, 90% yield, pale yellow solid, new compound, mp = 132-133 °C, *R*_f = 0.30 (hexanes/ethyl acetate 5/1), d.r. > 20:1, 98% ee, [α]²⁰_D = +156.08 (*c* 1.10, CHCl₃), ¹H NMR (400 MHz, CDCl₃) δ 7.62 (d, *J* = 8.1 Hz, 2H), 7.28-7.20 (m, 4H), 7.17-7.12 (m, 1H), 7.00-6.89 (m, 3H), 6.79 (d, *J* = 8.6 Hz, 2H), 5.08 (s, 1H), 4.36 (s, 1H), 3.72 (s, 3H), 2.39 (s, 3H). ¹³C NMR (100 MHz, CDCl₃) δ 160.0, 159.3, 150.8, 145.9, 133.9, 131.4, 129.9, 129.3, 129.1, 129.0, 128.2, 125.7, 121.3, 116.9, 114.9, 72.3, 55.3, 41.6, 21.7. HPLC (Chiralpak IA column, λ = 230 nm, 30 °C, *n*-Hexane/*i*-PrOH = 70/30, flow rate = 1.0 mL/min) retention time = 11.1 min (minor) and 13.5 min (major). HRMS (ESI) *m/z*: [M+NH₄]⁺ Calcd for C₂₃H₂₄NO₅S 426.1370, found: 426.1370.

(+)-(3*R*,4*S*)-4-(4-Fluorophenyl)-3-tosylchroman-2-one (2l):

59 mg, 99% yield, pale yellow solid, new compound, mp = 117-118 °C, *R*_f = 0.37 (hexanes/ethyl acetate 5/1), d.r. > 20:1, 98% ee, [α]²⁰_D = +151.60 (*c* 1.18, CHCl₃), ¹H NMR (400 MHz, CDCl₃) δ 7.63 (d, *J* = 8.3 Hz, 2H), 7.31-7.22 (m, 4H), 7.19-7.14 (m, 1H), 7.03-6.93 (m, 5H), 5.13 (s, 1H), 4.34 (s, 1H), 2.40 (s, 3H). ¹³C NMR (100 MHz, CDCl₃) δ 162.3 (d, ¹*J*_{F-C} = 246.3 Hz), 159.8, 150.8, 146.1, 135.2 (d, ⁴*J*_{F-C} = 3.5 Hz), 133.8, 129.9, 129.6, 129.1, 129.1, 128.8 (d, ³*J*_{F-C} = 8.1 Hz), 125.8, 120.8, 117.0, 116.5 (d, ²*J*_{F-C} = 21.6 Hz), 72.1, 41.6, 21.7. ¹⁹F NMR (376 MHz, CDCl₃) δ -113.53. HPLC (Chiralpak IA column, λ = 230 nm, 30 °C, *n*-Hexane/*i*-PrOH = 70/30, flow rate = 1.0 mL/min) retention time = 7.8 min (minor) and 10.2 min (major). HRMS (ESI) *m/z*: [M+NH₄]⁺ Calcd for C₂₂H₂₁FNO₄S 414.1170, found: 414.1165.

(+)-4-(4-Chlorophenyl)-3-tosylchroman-2-one (2m):

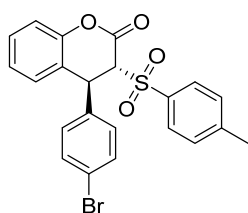
61 mg, 99% yield, colorless oil, new compound, *R*_f = 0.50 (hexanes/ethyl acetate 5/1), d.r. > 20:1, 98% ee, [α]²⁰_D = +147.37 (*c* 1.22, CHCl₃), ¹H NMR (400 MHz, CDCl₃) δ 7.63 (d, *J* = 8.2 Hz, 2H),



7.32-7.21 (m, 6H), 7.19-7.14 (m, 1H), 7.00-6.93 (m, 3H), 5.12 (s, 1H), 4.33 (s, 1H), 2.41 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 159.7, 150.8, 146.1, 137.9, 134.2, 133.7, 129.9, 129.7, 129.1, 129.1, 128.5, 125.8, 120.5, 117.0, 71.8, 41.6, 21.8. HPLC (Chiralpak IA column, λ = 230 nm, 30 °C, *n*-Hexane/*i*-PrOH = 75/25, flow rate = 1.0 mL/min) retention time = 10.1 min (minor) and 13.0 min (major). HRMS (ESI) m/z : $[\text{M}+\text{H}]^+$ Calcd for $\text{C}_{22}\text{H}_{18}\text{ClO}_4\text{S}$ 413.0609 (^{35}Cl) and 415.0592 (^{37}Cl), found: 413.0609 (^{35}Cl) and 415.0576 (^{37}Cl).

(+)-4-(4-Bromophenyl)-3-tosylchroman-2-one (2n):

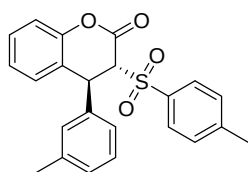
68 mg, 99% yield, colorless oil, new compound, R_f = 0.54 (hexanes/ethyl acetate 5/1), d.r. > 20:1,



98% ee, $[\alpha]_D^{20}$ = +135.43 (c 1.36, CHCl_3), ^1H NMR (400 MHz, CDCl_3) δ 7.63 (d, J = 7.9 Hz, 2H), 7.41 (d, J = 8.0 Hz, 2H), 7.31-7.20 (m, 4H), 7.19-7.14 (m, 1H), 6.97-6.86 (m, 3H), 5.11 (s, 1H), 4.34 (s, 1H), 2.40 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 159.7, 150.8, 146.1, 138.4, 133.7, 132.7, 129.9, 129.7, 129.1, 129.1, 128.8, 125.8, 122.3, 120.4, 117.0, 71.7, 41.7, 21.8. HPLC (Chiralpak IA column, λ = 230 nm, 30 °C, *n*-Hexane/*i*-PrOH = 75/25, flow rate = 1.0 mL/min) retention time = 10.6 min (minor) and 14.0 min (major). HRMS (ESI) m/z : $[\text{M}+\text{Na}]^+$ Calcd for $\text{C}_{22}\text{H}_{17}\text{BrNaO}_4\text{S}$ 478.9923 (^{79}Br) and 480.9910 (^{81}Br), found: 478.9923 (^{79}Br) and 480.9894 (^{81}Br).

(+)-4-(*m*-Tolyl)-3-tosylchroman-2-one (2o):

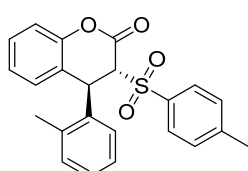
58 mg, 99% yield, pale yellow oil, new compound, R_f = 0.54 (hexanes/ethyl acetate 5/1), d.r. >



20:1, 98% ee, $[\alpha]_D^{20}$ = +160.25 (c 1.16, CHCl_3), ^1H NMR (400 MHz, CDCl_3) δ 7.64 (d, J = 8.2 Hz, 2H), 7.29-7.21 (m, 4H), 7.17-7.12 (m, 2H), 7.03 (d, J = 7.5 Hz, 1H), 6.93 (d, J = 8.1 Hz, 1H), 6.85 (s, 1H), 6.80 (d, J = 7.7 Hz, 1H), 5.10 (s, 1H), 4.38 (s, 1H), 2.39 (s, 3H), 2.26 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 160.0, 150.8, 146.0, 139.4, 139.4, 133.9, 129.9, 129.4, 129.4, 129.2, 129.1, 128.9, 127.7, 125.7, 124.1, 121.0, 116.9, 72.1, 42.2, 21.8, 21.5. HPLC (Chiralpak IA column, λ = 230 nm, 30 °C, *n*-Hexane/*i*-PrOH = 75/25, flow rate = 1.0 mL/min) retention time = 7.0 min (minor) and 8.4 min (major). HRMS (ESI) m/z : $[\text{M}+\text{H}]^+$ Calcd for $\text{C}_{23}\text{H}_{21}\text{O}_4\text{S}$ 393.1155, found: 393.1154.

(+)-4-(*o*-Tolyl)-3-tosylchroman-2-one (2p):

14 mg, 24% yield, colorless oil, new compound, R_f = 0.52 (hexanes/ethyl acetate 5/1), d.r. > 20:1,



98% ee, $[\alpha]_D^{20}$ = +113.92 (c 0.28, CHCl_3), ^1H NMR (400 MHz, CDCl_3) δ 7.66 (d, J = 8.2 Hz, 2H), 7.33-7.26 (m, 2H), 7.25-7.21 (m, 2H), 7.20-7.11 (m, 3H), 7.04-6.97 (m, 2H), 6.55 (d, J = 7.7 Hz, 1H), 5.39 (s, 1H), 4.21 (s, 1H), 2.54 (s, 3H), 2.41 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 159.8, 151.5, 146.0, 137.1, 134.9, 133.9, 131.5, 129.9, 129.4, 129.3, 129.0, 128.1, 127.3, 127.2, 125.8, 121.1, 116.8, 70.7, 38.7, 21.8, 19.4. HPLC (Chiralpak IA column, λ = 230 nm, 30 °C, *n*-Hexane/*i*-PrOH = 75/25, flow rate = 1.0 mL/min) retention time = 6.7 min (minor) and 8.4 min (major). HRMS (ESI) m/z : $[\text{M}+\text{H}]^+$ Calcd for $\text{C}_{23}\text{H}_{21}\text{O}_4\text{S}$ 393.1155, found: 393.1152.

(+)-4-(Naphthalen-2-yl)-3-tosylchroman-2-one (2q):

64 mg, 99% yield, pale yellow oil, new compound, $R_f = 0.55$ (hexanes/ethyl acetate 5/1), d.r. > 20:1, 97% ee, $[\alpha]_D^{20} = +156.55$ (c 1.28, CHCl_3), ^1H NMR (400 MHz, CDCl_3) δ 7.80-7.73 (m, 2H), 7.70-7.63 (m, 3H), 7.46-7.41 (m, 2H), 7.37 (s, 1H), 7.33-7.25 (m, 2H), 7.22 (d, $J = 8.5$ Hz, 3H), 7.18-7.13 (m, 1H), 6.98 (d, $J = 8.0$ Hz, 1H), 5.31 (s, 1H), 4.50 (s, 1H), 2.37 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 160.0, 151.0, 146.1, 136.7, 133.8, 133.4, 132.8, 129.9, 129.8, 129.6, 129.4, 129.1, 128.0, 127.7, 126.8, 126.7, 126.1, 125.8, 124.7, 120.9, 117.0, 71.9, 42.4, 21.8. HPLC: (Chiralpak IA column, $\lambda = 230$ nm, 30 $^\circ\text{C}$, n -Hexane/ i -PrOH = 75/25, flow rate = 1.0 mL/min) retention time = 12.2 min (minor) and 14.7 min (major). HRMS (ESI) m/z : $[\text{M}+\text{H}]^+$ Calcd for $\text{C}_{26}\text{H}_{21}\text{O}_4\text{S}$ 429.1155, found: 429.1151.

(+)-7-Methyl-4-phenyl-3-tosylchroman-2-one (2r):

49 mg, 83% yield, colorless oil, new compound, $R_f = 0.50$ (hexanes/ethyl acetate 5/1), d.r. > 20:1, 98% ee, $[\alpha]_D^{20} = +156.62$ (c 0.98, CHCl_3), ^1H NMR (400 MHz, CDCl_3) δ 7.63 (d, $J = 7.9$ Hz, 2H), 7.31-7.18 (m, 5H), 7.14-6.94 (m, 4H), 6.73 (s, 1H), 5.09 (s, 1H), 4.36 (s, 1H), 2.40 (s, 3H), 2.32 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 160.2, 150.8, 145.9, 140.0, 139.6, 134.0, 129.8, 129.5, 129.1, 128.9, 128.1, 127.1, 126.5, 117.8, 117.3, 72.2, 42.0, 21.8, 21.2. HPLC (Chiralpak IA column, $\lambda = 230$ nm, 30 $^\circ\text{C}$, n -Hexane/ i -PrOH = 70/30, flow rate = 1.0 mL/min) retention time = 7.5 min (minor) and 9.9 min (major). HRMS (ESI) m/z : $[\text{M}+\text{H}]^+$ Calcd for $\text{C}_{23}\text{H}_{21}\text{O}_4\text{S}$ 393.1155, found: 393.1152.

(+)-6,8-Dimethyl-4-phenyl-3-tosylchroman-2-one (2s):

58 mg, 95% yield, colorless oil, new compound, $R_f = 0.53$ (hexanes/ethyl acetate 5/1), d.r. > 20:1, 99% ee, $[\alpha]_D^{20} = +128.61$ (c 1.16, CHCl_3), ^1H NMR (400 MHz, CDCl_3) δ 7.58 (d, $J = 8.2$ Hz, 2H), 7.30-7.16 (m, 5H), 7.03 (d, $J = 7.3$ Hz, 2H), 6.86 (d, $J = 4.3$ Hz, 2H), 5.03 (s, 1H), 4.39 (s, 1H), 2.38 (s, 3H), 2.24 (s, 3H), 2.09 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 160.3, 147.1, 145.8, 139.5, 134.9, 133.9, 131.6, 129.6, 129.5, 129.0, 128.1, 127.1, 127.0, 125.9, 120.1, 72.5, 42.8, 21.7, 20.7, 15.4. HPLC (Chiralpak IA column, $\lambda = 230$ nm, 30 $^\circ\text{C}$, n -Hexane/ i -PrOH = 70/30, flow rate = 1.0 mL/min) retention time = 5.9 min (minor) and 7.4 min (major). HRMS (ESI) m/z : $[\text{M}+\text{H}]^+$ Calcd for $\text{C}_{24}\text{H}_{23}\text{O}_4\text{S}$ 407.1312, found: 407.1305.

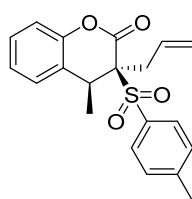
(+)-4-(Thiophen-3-yl)-3-tosylchroman-2-one (2t):

50 mg, 87% yield, colorless oil, new compound, $R_f = 0.53$ (hexanes/ethyl acetate 5/1), d.r. > 20:1, 98% ee, $[\alpha]_D^{20} = +128.47$ (c 0.92, CHCl_3), ^1H NMR (400 MHz, CDCl_3) δ 7.64-7.59 (m, 2H), 7.30-7.21 (m, 5H), 7.18-7.13 (m, 1H), 6.91-6.86 (m, 2H), 6.84-6.76 (m, 1H), 5.17 (s, 1H), 4.46 (s, 1H), 2.40 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 160.1, 150.5, 146.0, 139.6, 134.0, 129.9, 129.4, 129.0, 128.9, 127.9, 125.9, 125.6, 122.4, 121.4, 117.0, 71.2, 38.1, 21.7. HPLC (Chiralpak AD-H column, $\lambda = 230$ nm, 30 $^\circ\text{C}$, n -Hexane/ i -PrOH = 70/30, flow rate = 0.8 mL/min) retention time =

13.5 min (minor) and 19.1 min (major). HRMS (ESI) m/z : $[M+NH_4]^+$ Calcd for $C_{20}H_{20}NO_4S_2$ 402.0828, found: 402.0829.

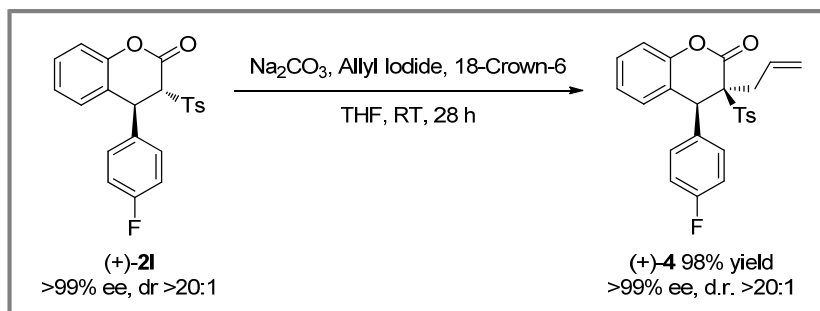
(-)-3-Allyl-4-methyl-3-tosylchroman-2-one (2u):

Procedure for allyl alkylation: Potassium carbonate (82.8 mg, 0.60 mmol), allyl iodide (100.8 mg, 54.8 μ L, 0.60 mmol), 18-crown-6 (11.8 mg, 10.2 μ L, 0.045 mmol) and tetrahydrofuran (3.0 mL) were added to the mixture of the reduction system. The mixture was stirred at ambient temperature for 48 h. The mixture was concentrated under the reduced pressure, and then purified by column chromatography on silica gel using hexanes and ethyl acetate as eluent to afford the product (-)-**2u**, 30 mg, 56% yield, colorless oil, new compound, R_f = 0.60 (hexanes/ethyl acetate



5/1), d.r. > 20:1, 93% ee, $[\alpha]_D^{20}$ = -22.25 (c 0.40, $CHCl_3$), 1H NMR (400 MHz, $CDCl_3$) δ 7.69 (d, J = 8.3 Hz, 2H), 7.30-7.26 (m, 2H), 7.25-7.20 (m, 2H), 7.19-7.14 (m, 1H), 6.87 (dd, J = 7.9, 1.1 Hz, 1H), 5.65-5.49 (m, 1H), 5.19-5.10 (m, 2H), 3.65 (q, J = 7.1 Hz, 1H), 3.02 (dd, J = 14.4, 5.7 Hz, 1H), 2.78 (dd, J = 14.4, 8.8 Hz, 1H), 2.41 (s, 3H), 1.76 (d, J = 7.1 Hz, 3H). ^{13}C NMR (100 MHz, $CDCl_3$) δ 163.6, 150.1, 145.6, 134.0, 130.2, 130.1, 129.5, 128.4, 126.1, 125.5, 125.3, 121.6, 115.8, 73.2, 35.3, 34.6, 21.7, 13.4. HPLC (Chiralpak AD-H column, λ = 230 nm, 30 $^{\circ}C$, n -Hexane/ i -PrOH = 90/10, flow rate = 1.0 mL/min) retention time = 12.4 min (minor) and 13.2 min (major). HRMS (ESI) m/z : $[M+NH_4]^+$ Calcd for $C_{20}H_{24}NO_4S$ 374.1421, found: 374.1413.

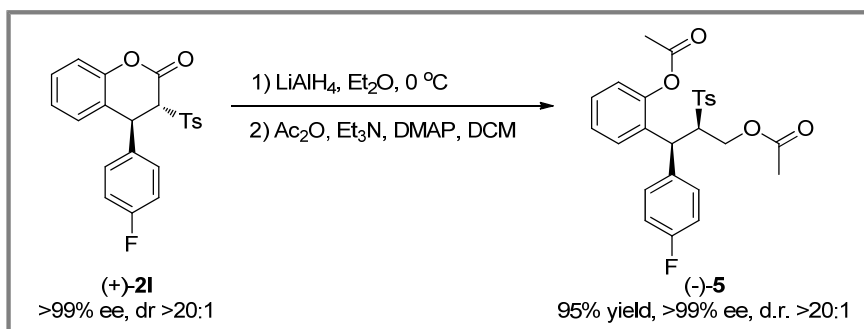
5. Transformation of the Biomimetic Reductive Products



Sodium carbonate (42.4 mg, 0.40 mmol), allyl iodide (67.2 mg, 36.5 μL , 0.40 mmol) and 18-crown-6 (7.9 mg, 6.8 μL , 0.03 mmol) were added to the stirred solution of (+)-**2I** (39.6 mg, 0.10 mmol, 99% ee) in tetrahydrofuran (2.0 mL). The mixture was stirred at ambient temperature for 28 hours. The mixture was concentrated under the reduced pressure, and then purified by column chromatography on silica gel using hexanes and ethyl acetate as eluent to afford the desirable alkylation product (+)-**4**.

(+)-(3*S*,4*S*)-3-Allyl-4-(4-fluorophenyl)-3-tosylchroman-2-one (**4**):

43 mg, 98% yield, colorless oil, new compound, R_f = 0.41 (hexanes/ethyl acetate 5/1), >99% ee, $[\alpha]_D^{20}$ = +176.03 (c 0.86, CHCl_3), ^1H NMR (400 MHz, CDCl_3) δ 7.74 (d, J = 8.1 Hz, 2H), 7.30 (t, J = 7.2 Hz, 3H), 7.20 (d, J = 8.1 Hz, 2H), 7.15-7.05 (m, 3H), 6.91 (t, J = 8.3 Hz, 2H), 5.83-5.68 (m, 7.3 Hz, 1H), 5.15 (d, J = 10.0 Hz, 1H), 4.79 (d, J = 16.9 Hz, 1H), 4.70 (s, 1H), 2.84-2.72 (m, 2H), 2.39 (s, 3H). ^{13}C NMR (100 MHz, CDCl_3) δ 163.5, 162.5 (d, $^1J_{\text{F-C}}$ = 246.2 Hz), 149.4, 145.2, 134.8, 133.1 (d, $^4J_{\text{F-C}}$ = 3.1 Hz), 131.3 (d, $^3J_{\text{F-C}}$ = 6.7 Hz), 131.1, 130.0, 129.2, 129.2, 128.7, 125.6, 125.0, 121.7, 116.7, 115.7 (d, $^2J_{\text{F-C}}$ = 21.3 Hz), 74.6, 48.0, 37.5, 21.7. ^{19}F NMR (376 MHz, CDCl_3) δ -113.68. HPLC (Chiralpak IC column, 230 nm, 30 $^\circ\text{C}$, n -Hexane/ i -PrOH = 80/20, flow = 1.0 mL/min) retention time = 14.9 min (minor) and 16.2 min (major). HRMS (ESI) m/z : $[\text{M}+\text{NH}_4]^+$ Calcd for $\text{C}_{25}\text{H}_{25}\text{FNO}_4\text{S}$ 454.1483, found: 454.1486.



To a solution of (+)-**2I** (39.6 mg, 0.10 mmol, 99% ee) in ethyl ether (3.0 mL) was added lithium aluminium tetrahydride (5.7 mg, 0.15 mmol) at 0 $^\circ\text{C}$. The reaction mixture was stirred for 3 h. The solution was quenched with water (3.0 mL) and 1N HCl (3.0 mL). The aqueous layer was extracted with ethyl acetate (10 mL $\times$ 3), washed with brine, dried over anhydrous sodium sulfate, and concentrated under the reduced pressure. Then, to the solution of the concentrated mixture above in dichloromethane (3.0 mL) was added triethylamine (101 mg, 0.14 mL, 1.0 mmol), 4-

dimethylaminopyridine (6.1 mg, 0.05 mmol) and acetic anhydride (51.1 mg, 47.3 μ L, 0.5 mmol) at ambient temperature. The reaction mixture was stirred overnight. The solution was quenched with saturated ammonium chloride aqueous solution (10 mL). The aqueous layer was extracted with dichloromethane, washed with brine, dried over anhydrous sodium sulfate, and concentrated under the reduced pressure. The residue was purified by silica gel column chromatography using hexanes and ethyl acetate as eluent to afford the desirable product (-)-5.

(-)-2-((1S,2R)-3-Acetoxy-1-(4-fluorophenyl)-2-tosylpropyl)phenyl acetate (5):

46 mg, 95% yield, colorless oil, new compound, R_f = 0.31 (hexanes/ethyl acetate 2/1), >99% ee, $[\alpha]_D^{20}$ = -44.45 (*c* 0.92, CHCl₃), ¹H NMR (400 MHz, CDCl₃) δ 7.45 (d, *J* = 8.2 Hz, 2H), 7.26-7.12 (m, 7H), 7.03 (d, *J* = 8.0 Hz, 1H), 6.81 (t, *J* = 8.6 Hz, 2H), 4.94 (d, *J* = 9.8 Hz, 1H), 4.55-4.47 (m, 1H), 4.30-4.21 (m, 2H), 2.39 (s, 3H), 2.38 (s, 3H), 1.84 (s, 3H). ¹³C NMR (100 MHz, CDCl₃) δ 170.1, 169.1, 161.7 (d, ¹*J*_{F-C} = 244.8 Hz), 148.1, 144.5, 137.0, 135.4 (d, ⁴*J*_{F-C} = 2.9 Hz), 132.1, 130.0 (d, ³*J*_{F-C} = 8.1 Hz), 129.5, 128.7, 128.4, 128.3, 126.5, 123.4, 115.3 (d, ²*J*_{F-C} = 21.5 Hz), 66.2, 60.9, 41.0, 21.6, 21.1, 20.4. ¹⁹F NMR (376 MHz, CDCl₃) δ -115.37. HPLC (Chiralpak AD-H column, 230 nm, 30 °C, *n*-Hexane/*i*-PrOH = 60/40, flow = 1.0 mL/min) retention time = 9.7 min (minor) and 15.2 min (major). HRMS (ESI) *m/z*: [M+NH₄]⁺ Calcd for C₂₆H₂₉FNO₆S 502.1694, found: 502.1696.

6. Determination of the Absolute Configuration

To determine the absolute configuration of (+)-4-(4-fluorophenyl)-3-tosylchroman-2-one (**21**) (98% ee): firstly, (+)-**21** was upgraded to >99% ee by recrystallization with *n*-hexane/dichloromethane and completely dissolved in dichloromethane (2 mL). Then part of dichloromethane was evaporated and *n*-hexane (3 mL) was added slowly at ambient temperature. The solvent was slowly evaporated and the single crystal of was obtained after one day. The structure in **Figure S1** showed the absolute configuration of (+)-**21** is (3*R*,4*S*). The CCDC number is 2069554. These details can be obtained free of charge via www.ccdc.com.ac.uk/data_request/cif from the Cambridge Crystallographic Data Centre.

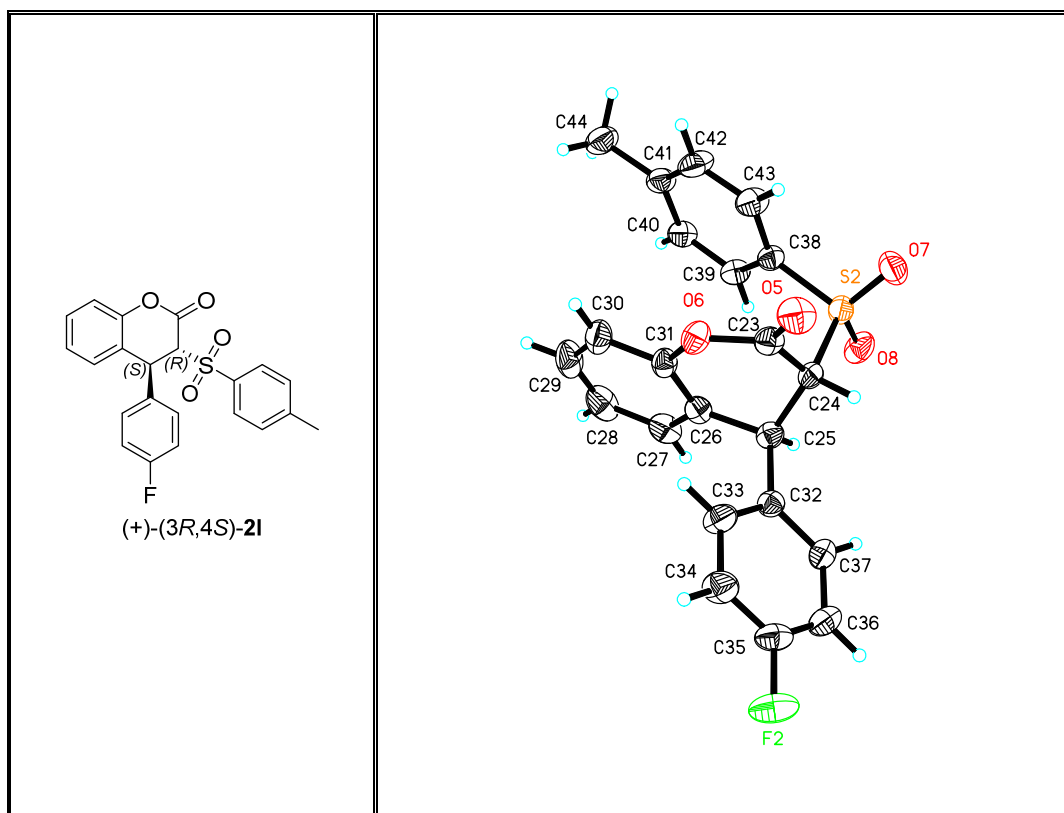


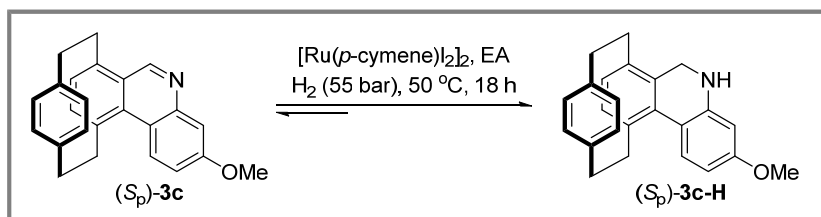
Figure S1. X-ray Crystallographic Analysis of (+)-(3*R*,4*S*)-**21**

Crystal Data and Structure Refinement for mo_d8v21165_0m for (+)-(3*R*,4*S*)-21

Identification code	mo_d8v21165_0m
Empirical formula	C ₂₂ H ₁₇ F O ₄ S
Formula weight	396.41
Temperature	293(2) K
Wavelength	0.71073 Å
Crystal system	Monoclinic
Space group	P 21
Unit cell dimensions	a = 8.3425(2) Å α = 90° b = 24.9386(7) Å β = 99.1170(10)° c = 9.4740(3) Å γ = 90°
Volume	1946.17(9) Å ³
Z	4
Density (calculated)	1.353 Mg/m ³
Absorption coefficient	0.201 mm ⁻¹
F(000)	824
Crystal size	0.200 x 0.150 x 0.120 mm ³
Theta range for data collection	2.473 to 25.999°
Index ranges	-10 ≤ h ≤ 10, -30 ≤ k ≤ 30, -9 ≤ l ≤ 11
Reflections collected	20019
Independent reflections	7567 [R(int) = 0.0238]
Completeness to theta = 25.242°	99.1 %
Absorption correction	Semi-empirical from equivalents
Max. and min. transmission	0.7456 and 0.6411
Refinement method	Full-matrix least-squares on F ²
Data / restraints / parameters	7567 / 1 / 508
Goodness-of-fit on F ²	1.020
Final R indices [I > 2σ(I)]	R ₁ = 0.0333, wR ₂ = 0.0780
R indices (all data)	R ₁ = 0.0404, wR ₂ = 0.0838
Absolute structure parameter	0.020(18)
Extinction coefficient	0.018(4)
Largest diff. peak and hole	0.157 and -0.165 e.Å ⁻³

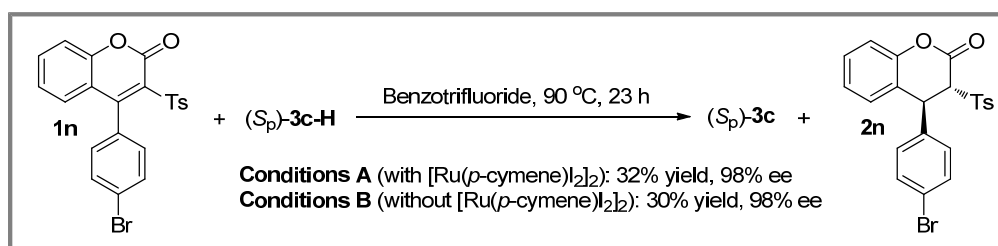
7. Mechanistic Investigation

7.1 The Reduction of NAD(P)H Model (*S_p*)-3c with Hydrogen Gas



A mixture of $[\text{Ru}(p\text{-cymene})\text{I}_2]_2$ (4.9 mg, 0.005 mmol), (*S_p*)-**3c** (33.9 mg, 0.10 mmol) in ethyl acetate (2 mL) was stirred at room temperature for 5 min in glove box and then the mixture was transferred to an autoclave. The hydrogenation was performed at 50 °C under hydrogen gas (55 bar) for 18 h. After careful release of the gas, the autoclave was opened, and it gave the reductive (*S_p*)-**3c-H** (about 95% yield). Because (*S_p*)-**3c-H** was unstable (easy rearomatization to **3c**), it directly participated in the next reaction without further separation.

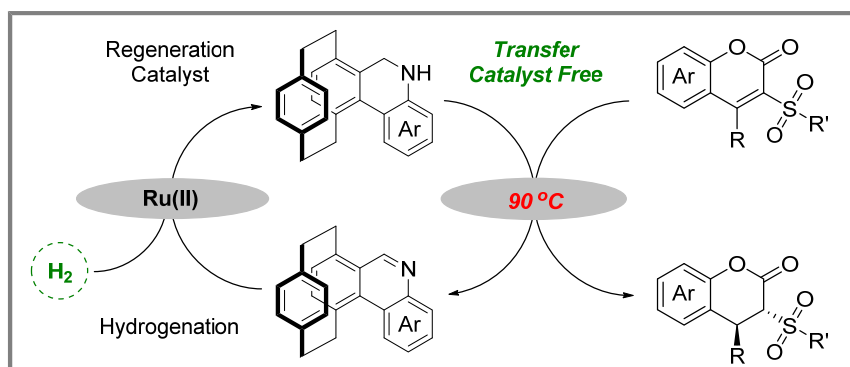
7.2 Control Experiments of Ruthenium Complex



Conditions A: in a 10 mL sealed tube, a mixture of (*S_p*)-**3c-H** (34.2 mg, 0.10 mmol), $[\text{Ru}(p\text{-cymene})\text{I}_2]_2$ (1.5 mg, 0.0015 mmol) and substrate **1n** (45.5, 0.10 mmol) in benzotrifluoride (3.0 mL) was stirred at 90 °C for 23 h. The final reaction mixture was concentrated in vacuo and gave product **2n** with 32% yield and 98% ee.

Conditions B: in a 10 mL sealed tube, a mixture of (*S_p*)-**3c-H** (34.2 mg, 0.10 mmol) and substrate **1n** (45.5, 0.10 mmol) in benzotrifluoride (3.0 mL) was stirred at 90 °C for 23 h. The final reaction mixture was concentrated in vacuo and gave product **2n** with 30% yield and 98% ee.

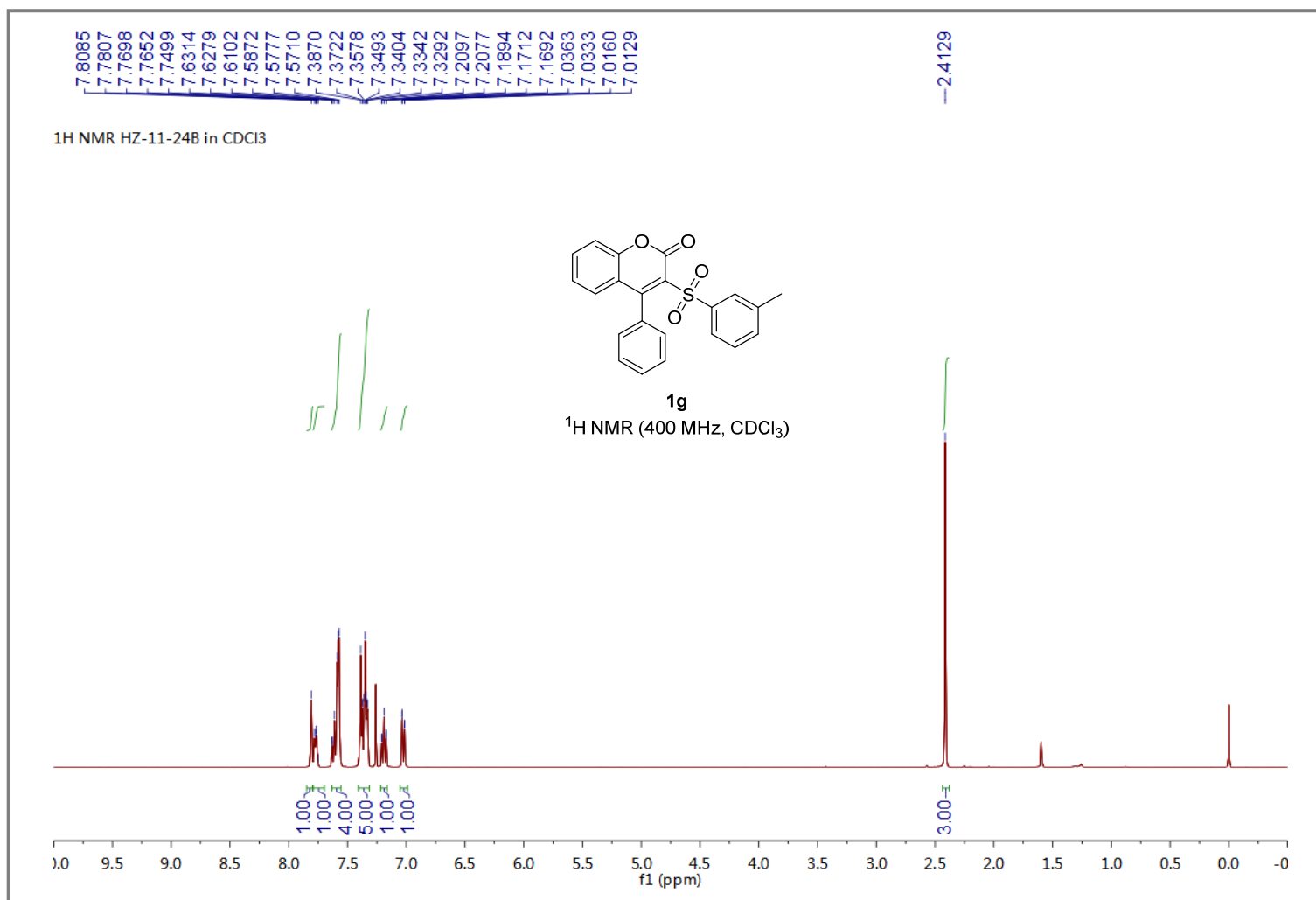
Similar results could be achieved for conditions A and B, which indicated that reduced (*S_p*)-**3c-H** reagent participated in the hydrogenation process without the ruthenium complex. Based on the experimental results and our previous work^{1,4-5}, a plausible mechanism for the biomimetic asymmetric reduction of 3-sulfonyl coumarins was illustrated in the scheme below.

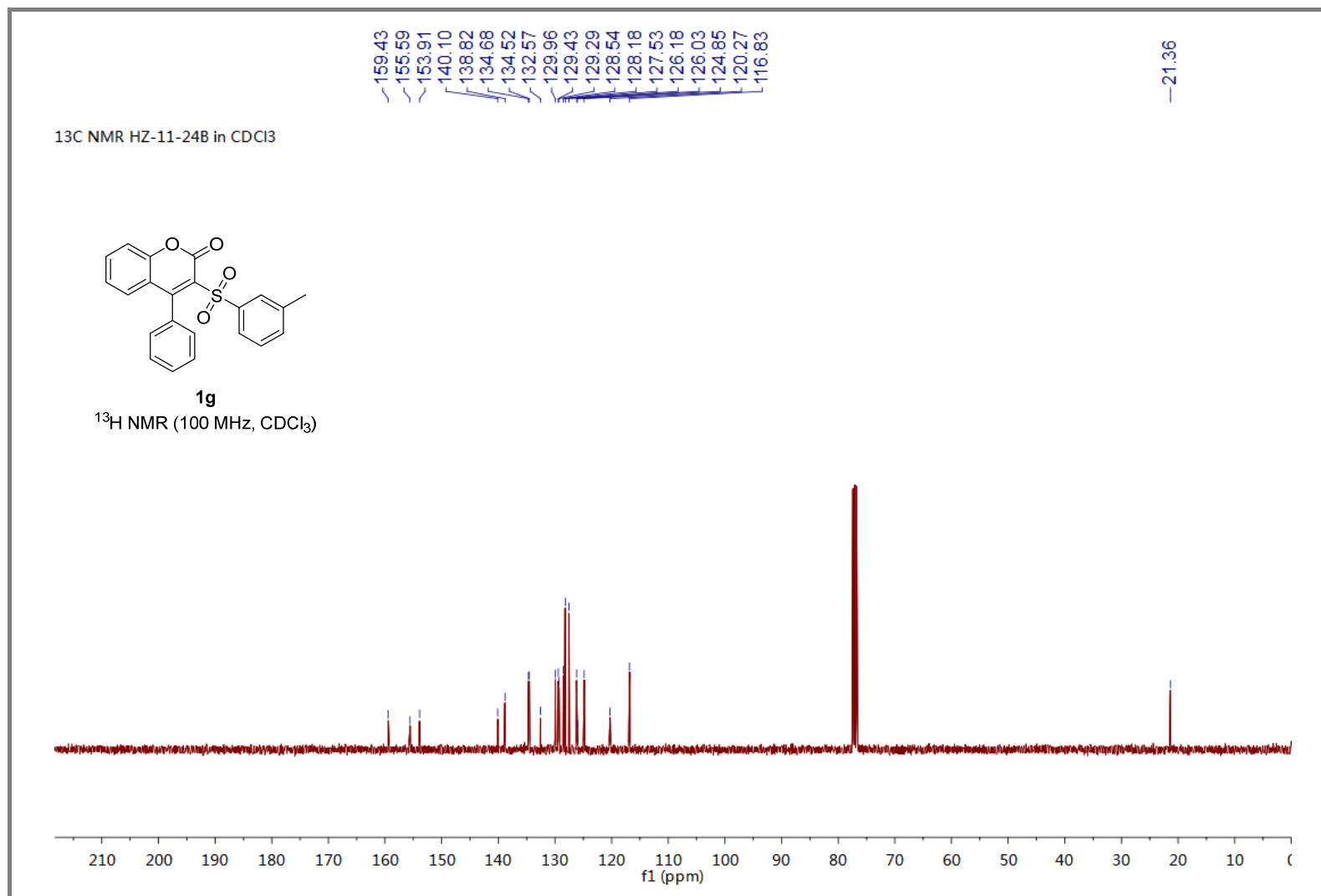


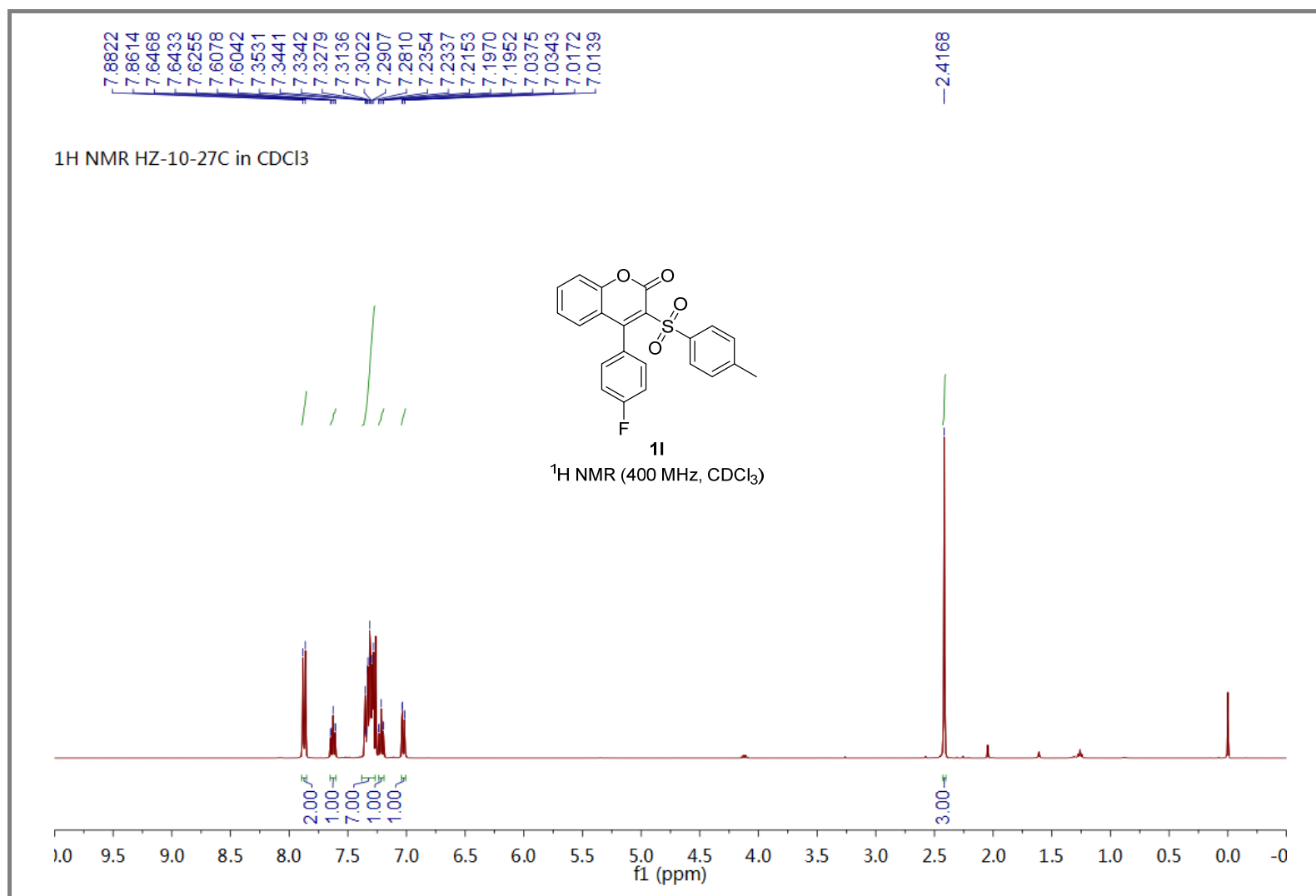
8. References

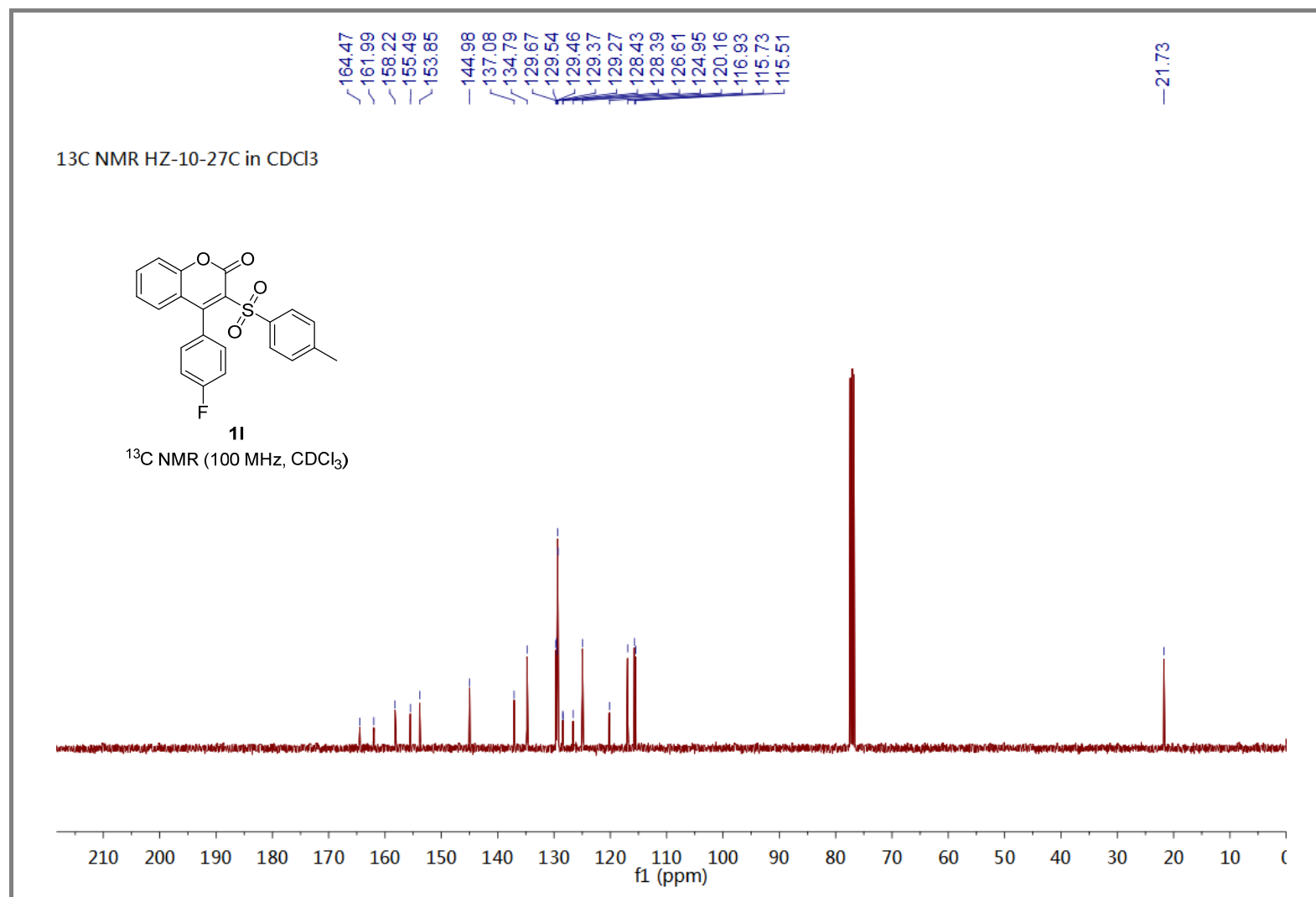
- 1) Zhu, Z.-H.; Ding, Y.-X.; Wu, B.; Zhou, Y.-G. Design and Synthesis of Chiral and Regenerable [2.2]Paracyclophane-Based NAD(P)H Models and Application in Biomimetic Reduction of Flavonoids. *Chem. Sci.* **2020**, *11*, 10220-10224.
- 2) Ren, H.; Zhang, M.; Zhang, A. Q. Synthesis of 3-Sulfonyl Coumarins through Radical Sulfonylation with Disulfides under Catalyst-Free Conditions. *Tetrahedron* **2018**, *74*, 4435-4444.
- 3) Kanyiva, K. S.; Hamada, D.; Makino, S.; Takano, H.; Shibata, T. α -Amino Acid Sulfonamides as Versatile Sulfonylation Reagents: Silver-Catalyzed Synthesis of Coumarins and Oxindoles by Radical Cyclization. *Eur. J. Org. Chem.* **2018**, *43*, 5905-5909.
- 4) Chen, Q.-A.; Gao, K.; Duan, Y.; Ye, Z.-S.; Shi, L.; Yang, Y.; Zhou, Y.-G. Dihydrophenanthridine: A New and Easily Regenerable NAD(P)H Model for Biomimetic Asymmetric Hydrogenation. *J. Am. Chem. Soc.* **2012**, *134*, 2442-2448.
- 5) Wang, J.; Zhao, Z.-B.; Zhao, Y. Luo, G.; Zhu, Z.-H.; Luo, Y.; Zhou, Y.-G. Chiral and Regenerable NAD(P)H Models Enabled Biomimetic Asymmetric Reduction: Design, Synthesis, Scope, and Mechanistic Studies. *J. Org. Chem.* **2020**, *85*, 2355-2368.

9. Copy of NMR and HPLC Spectra

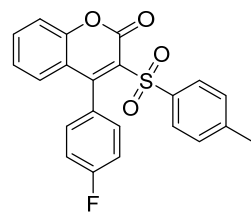






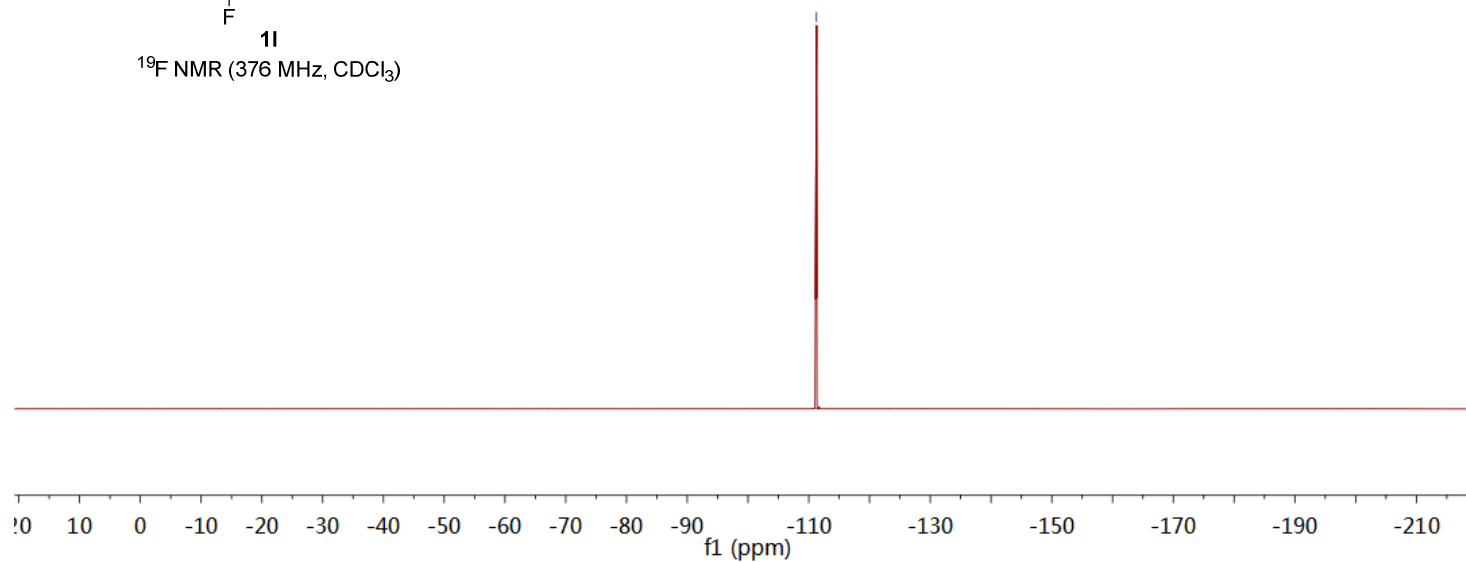


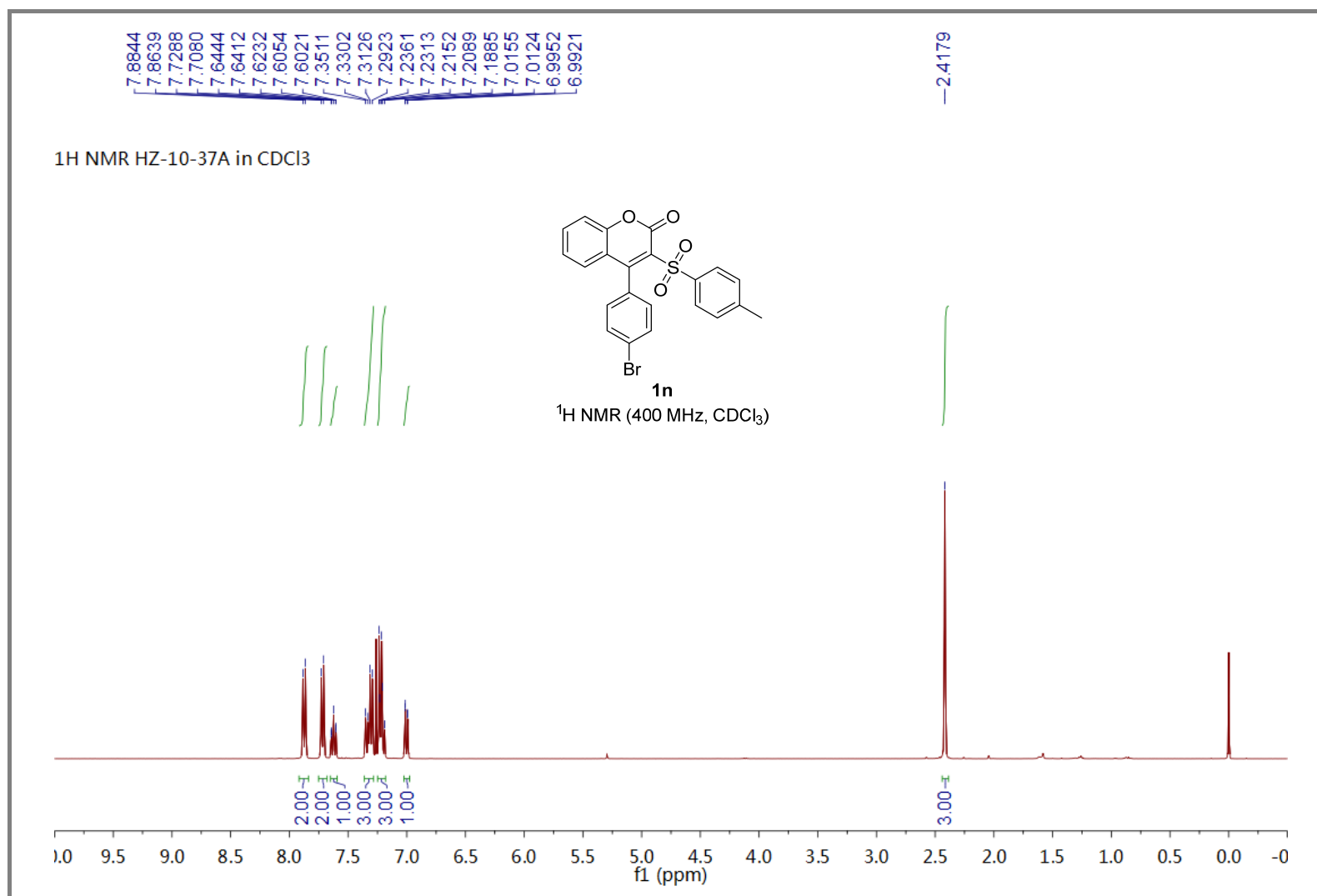
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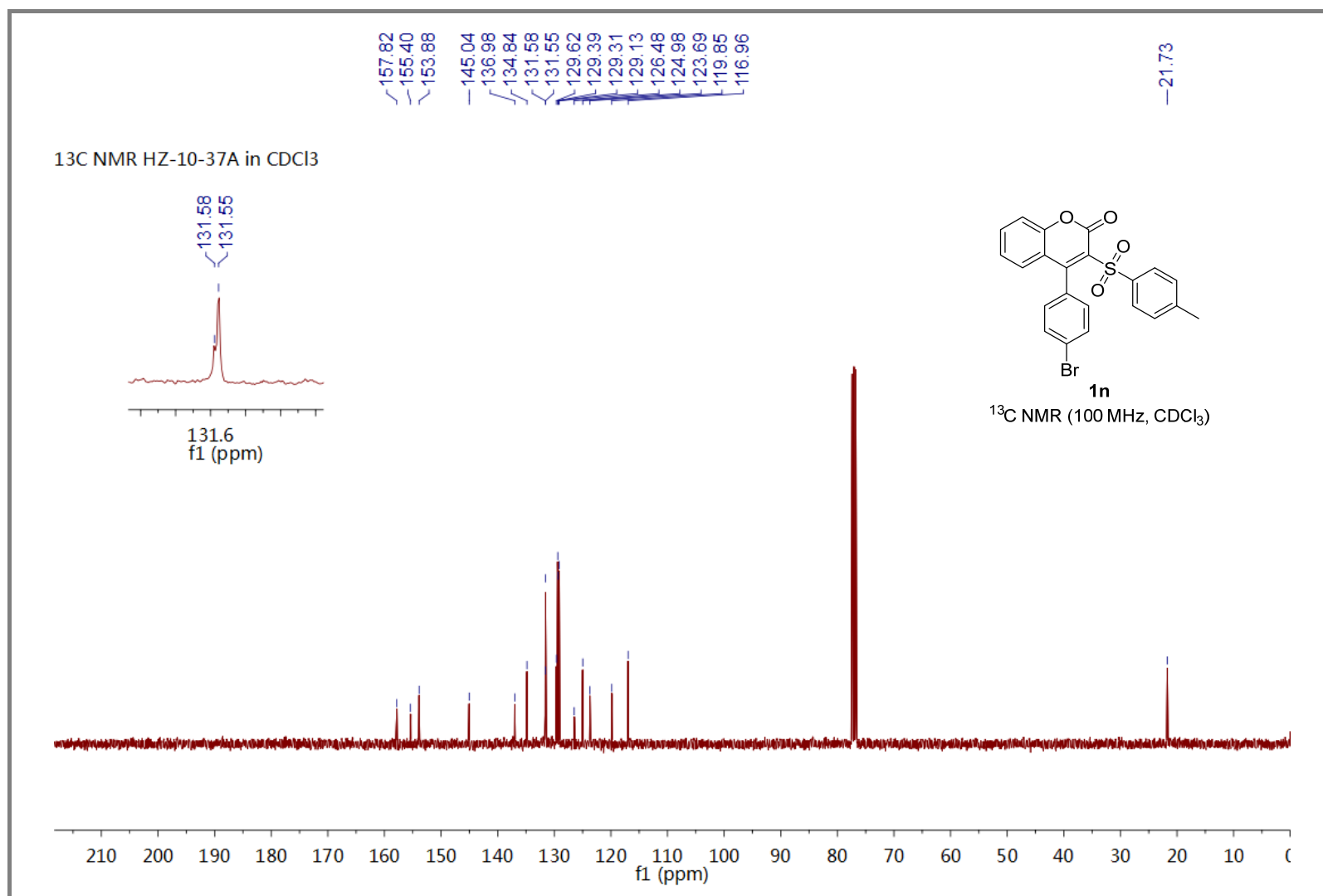


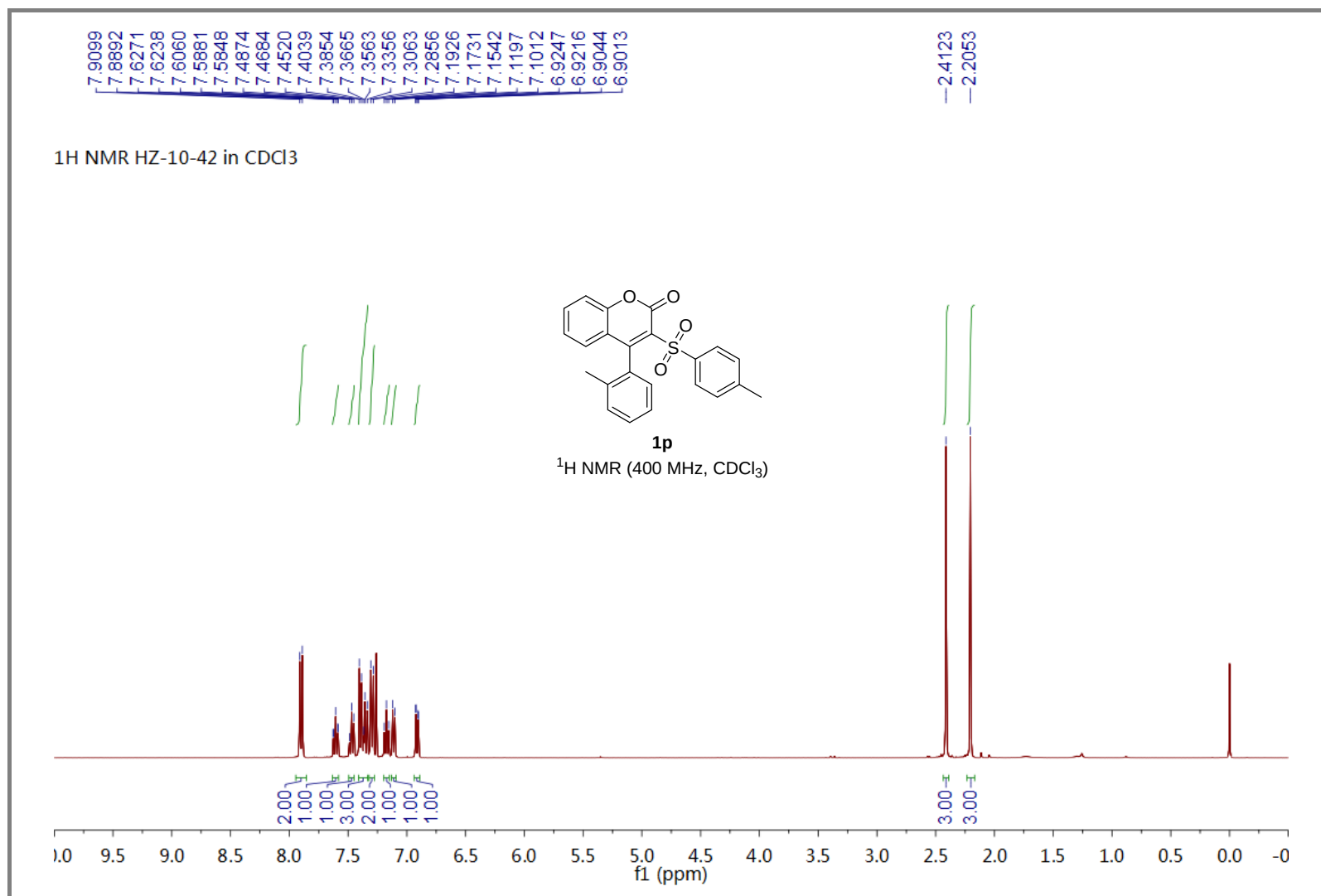
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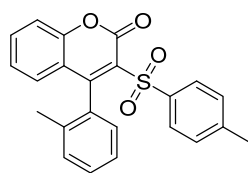






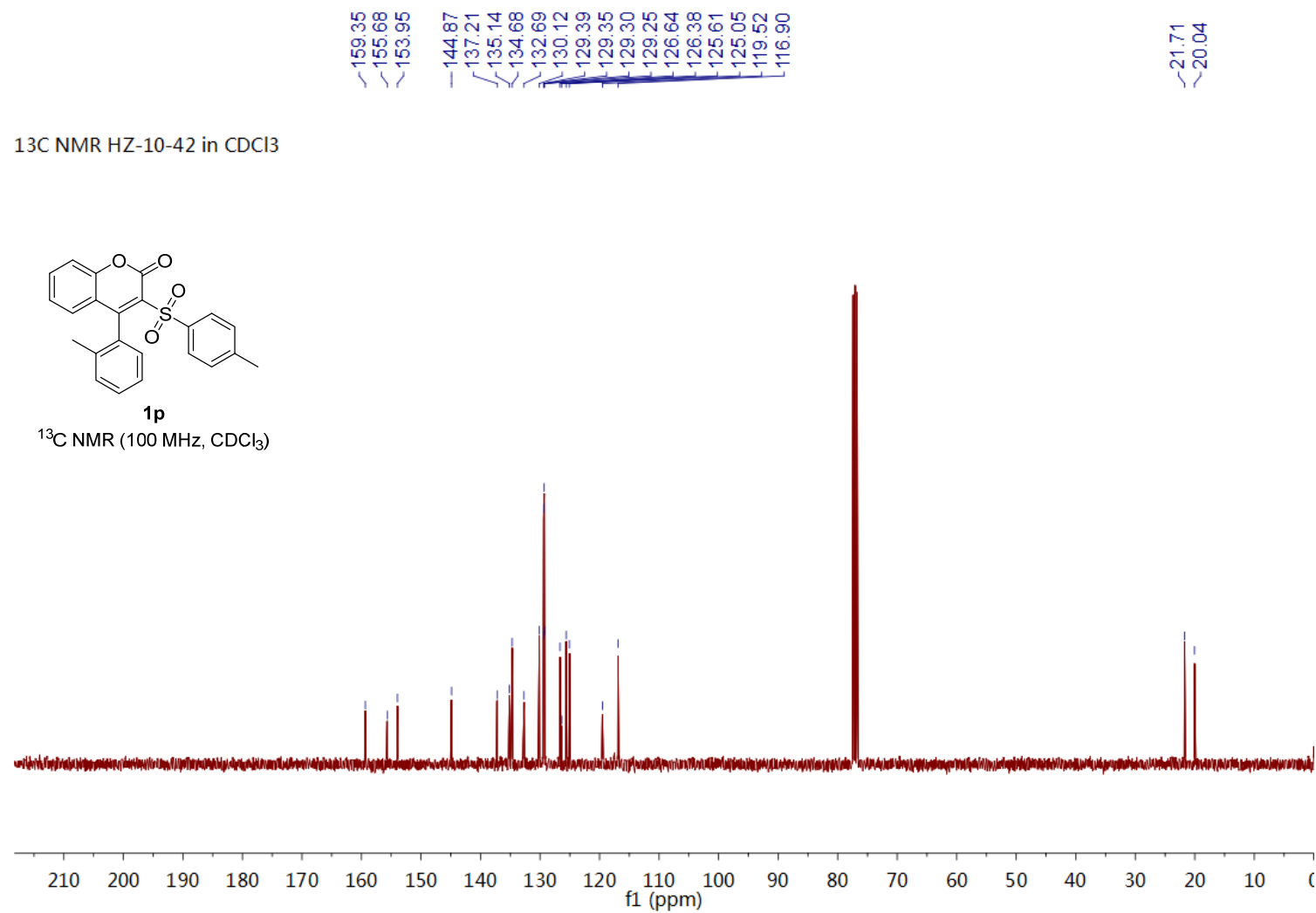


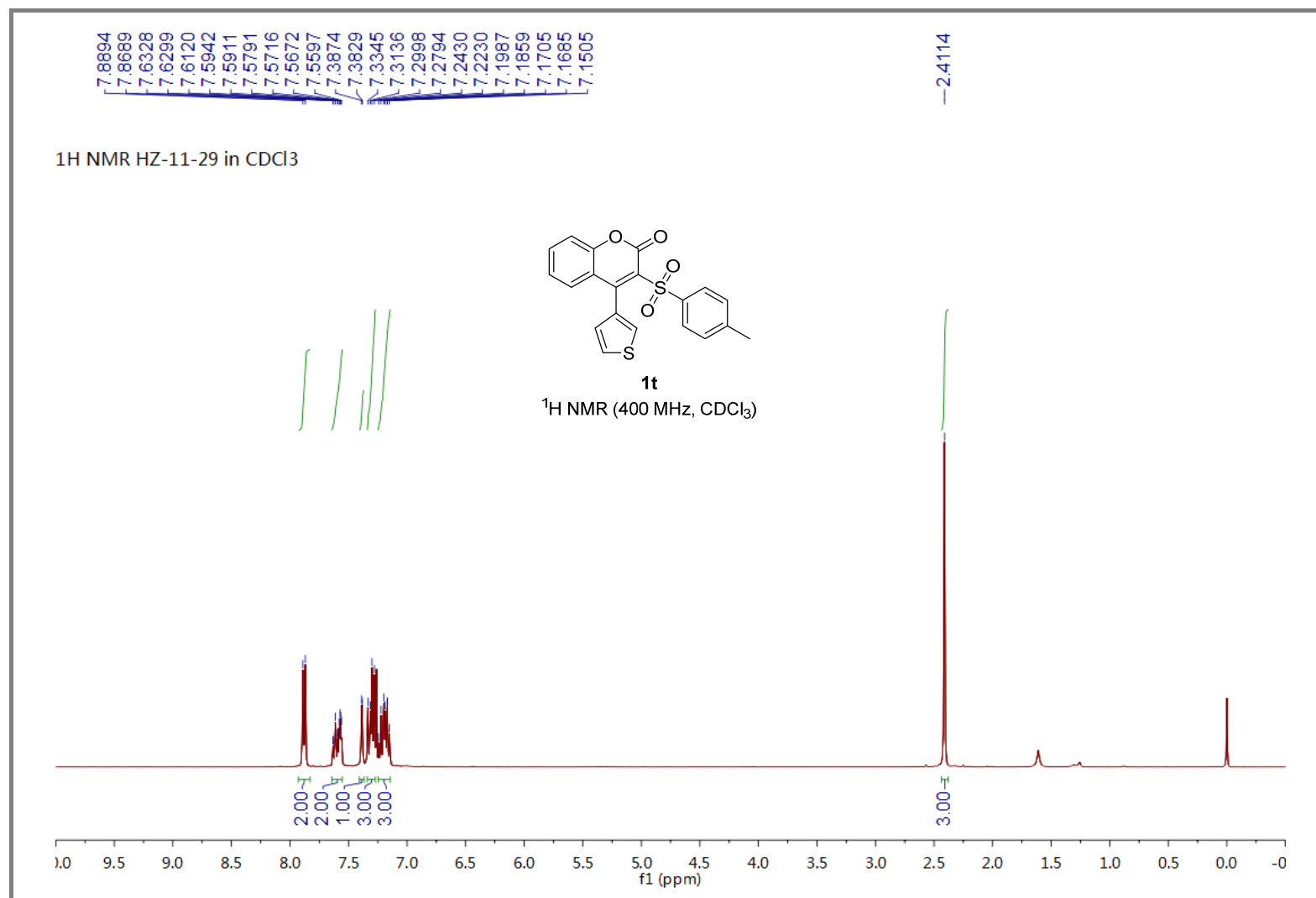
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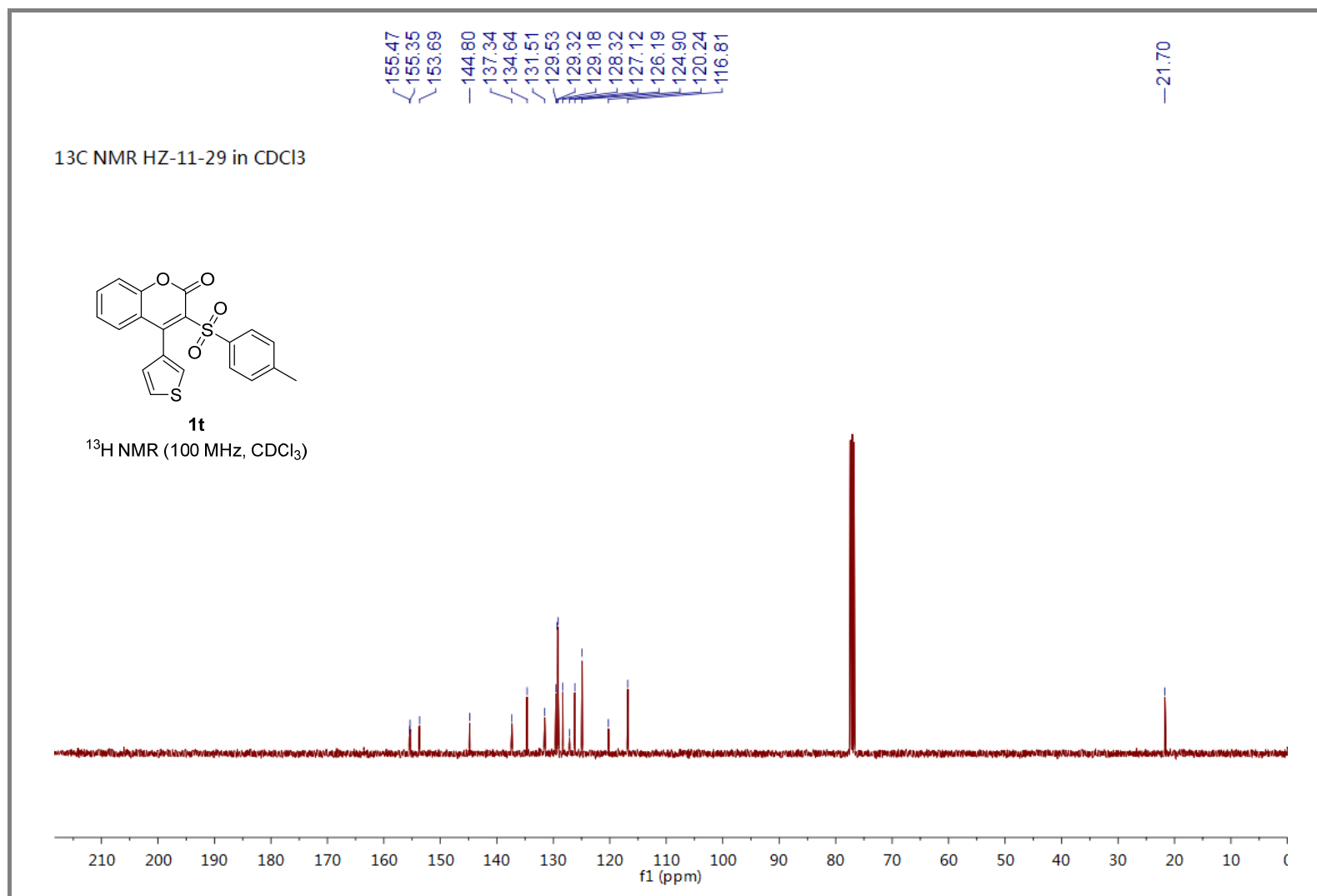


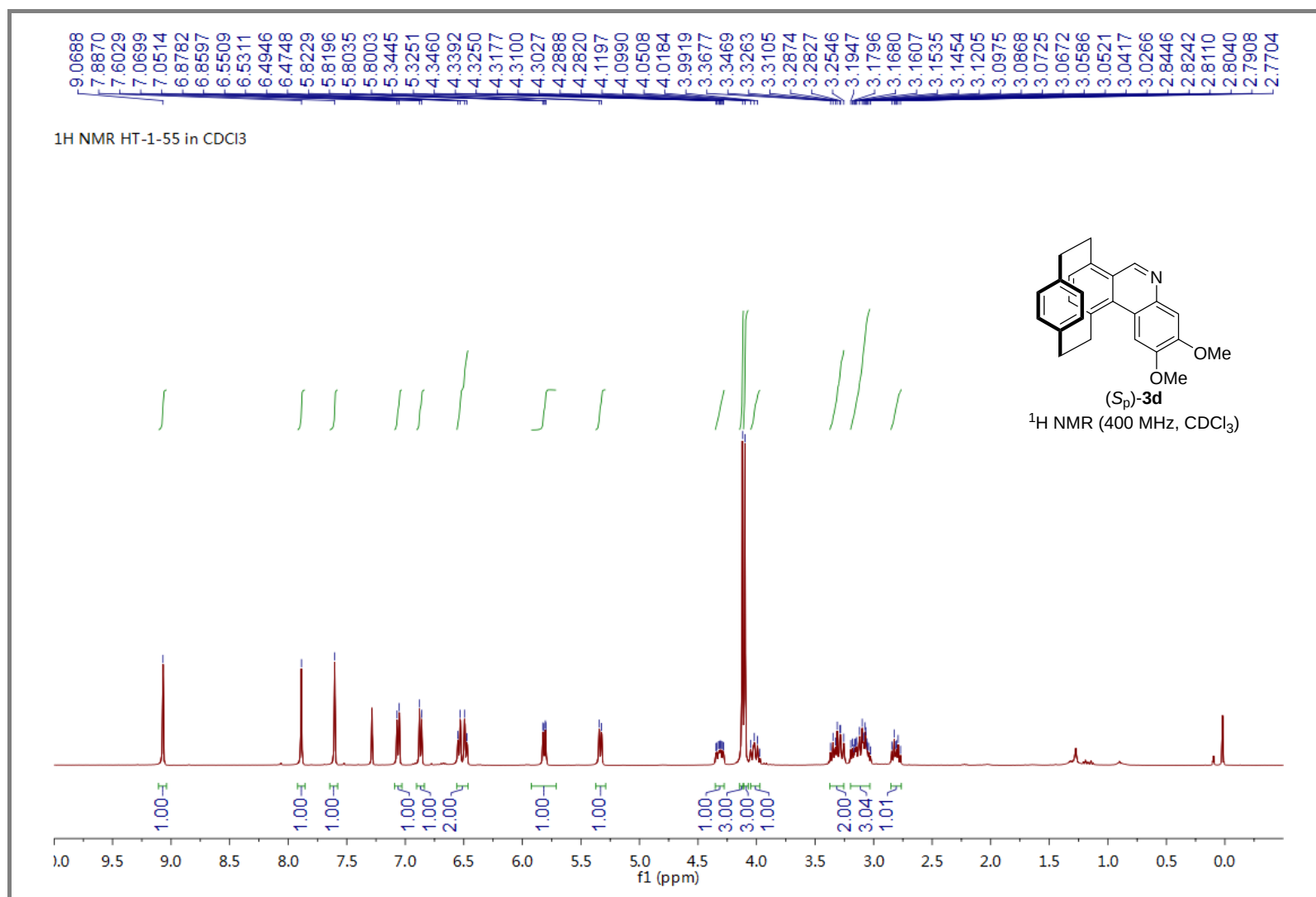
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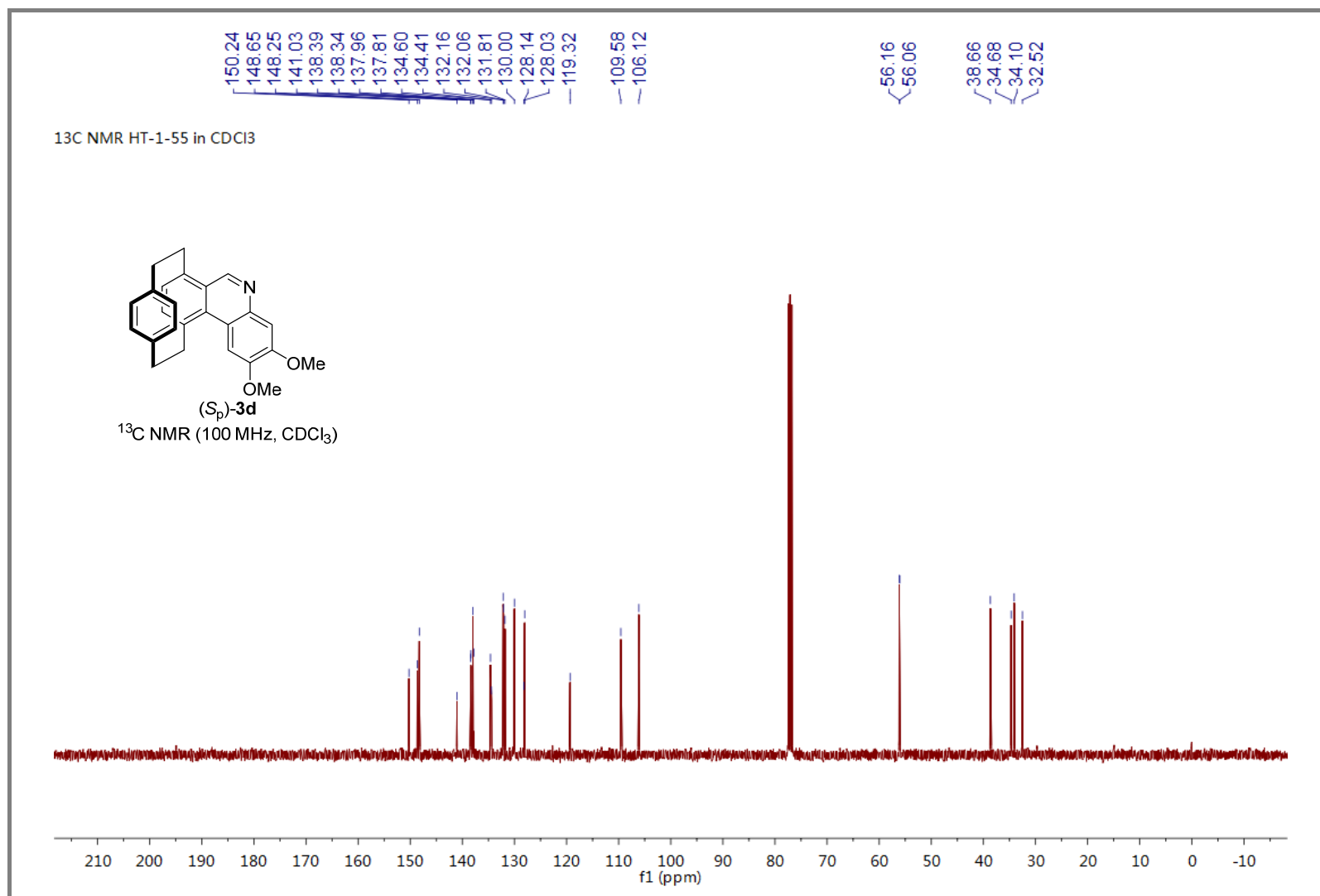
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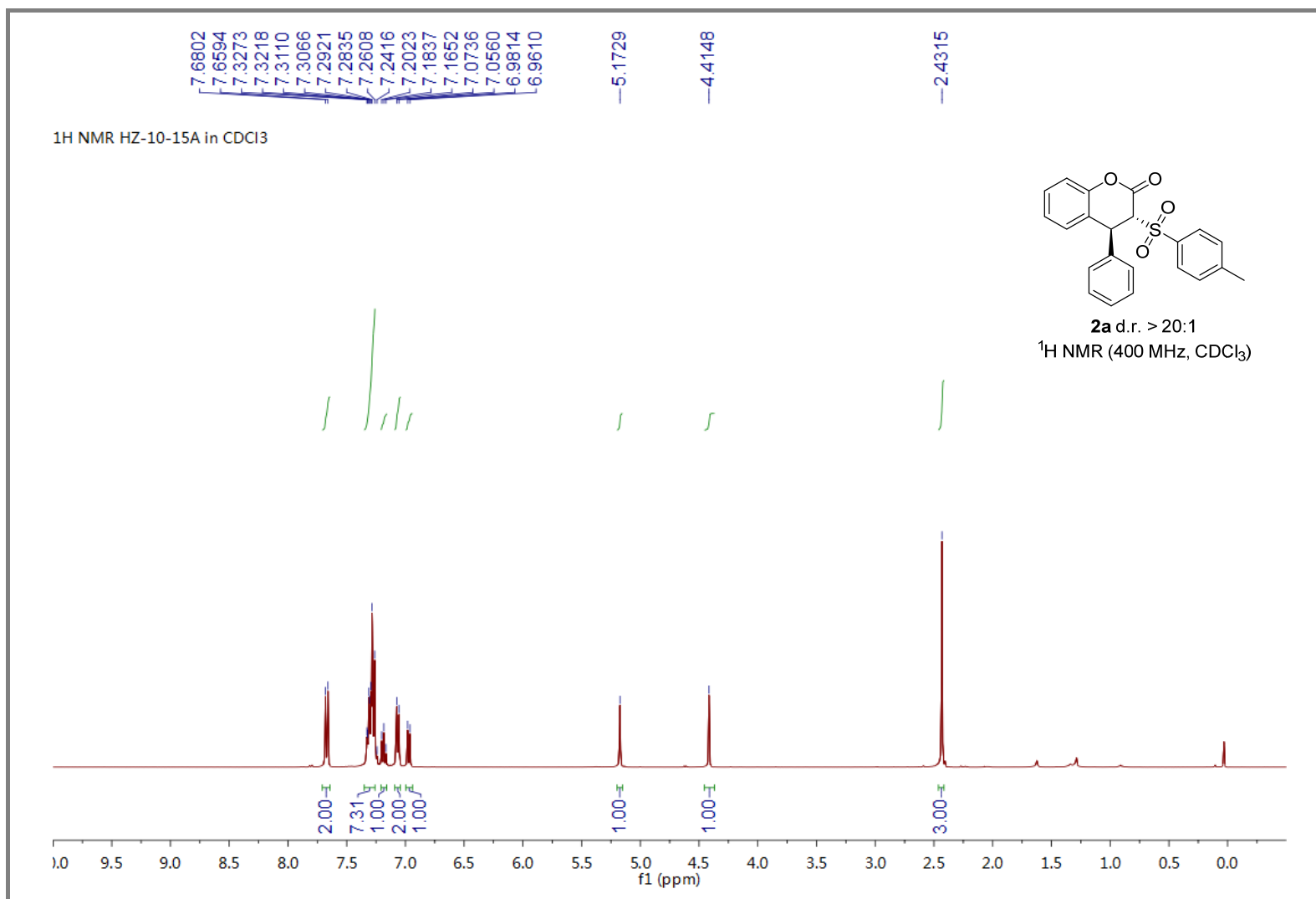


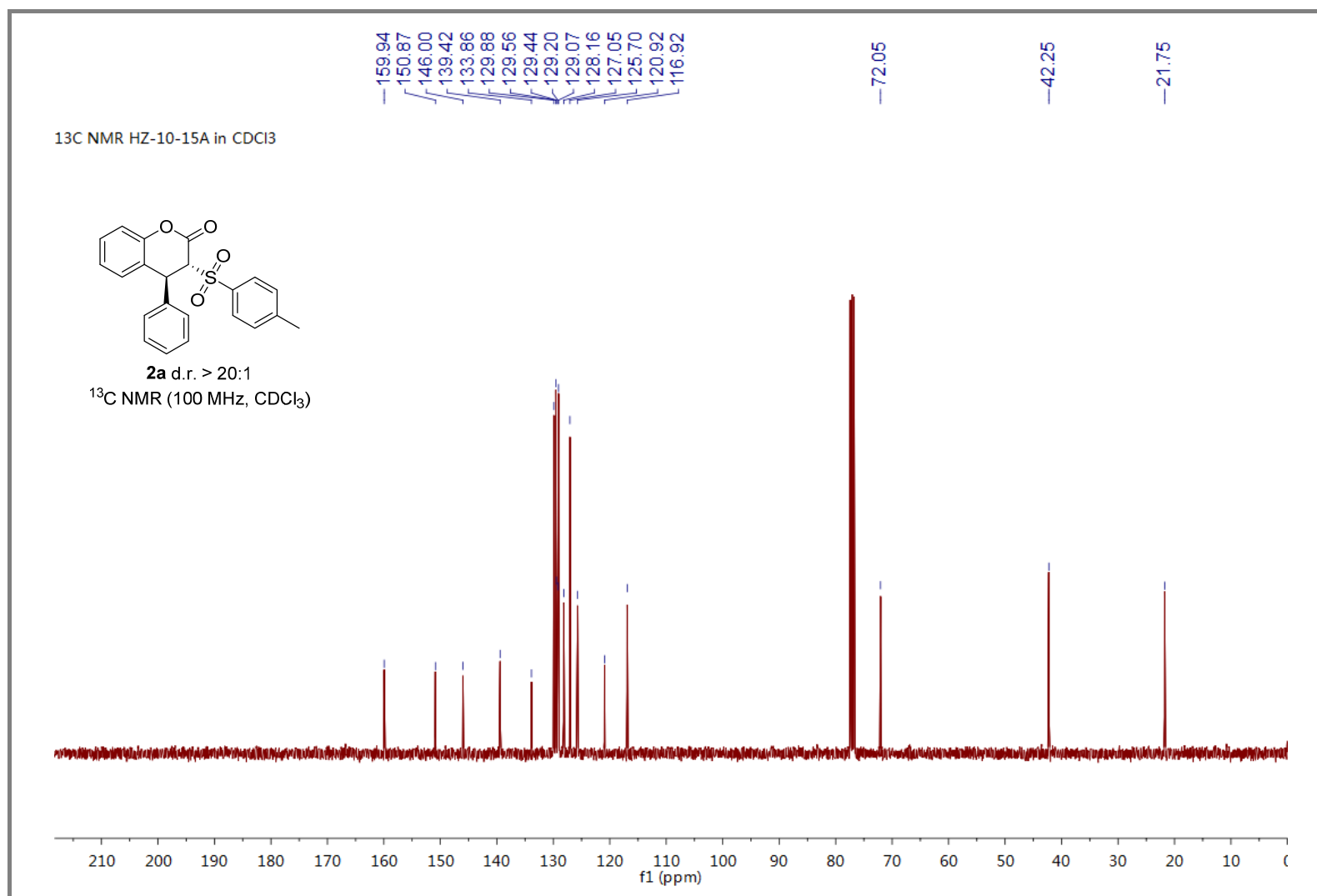


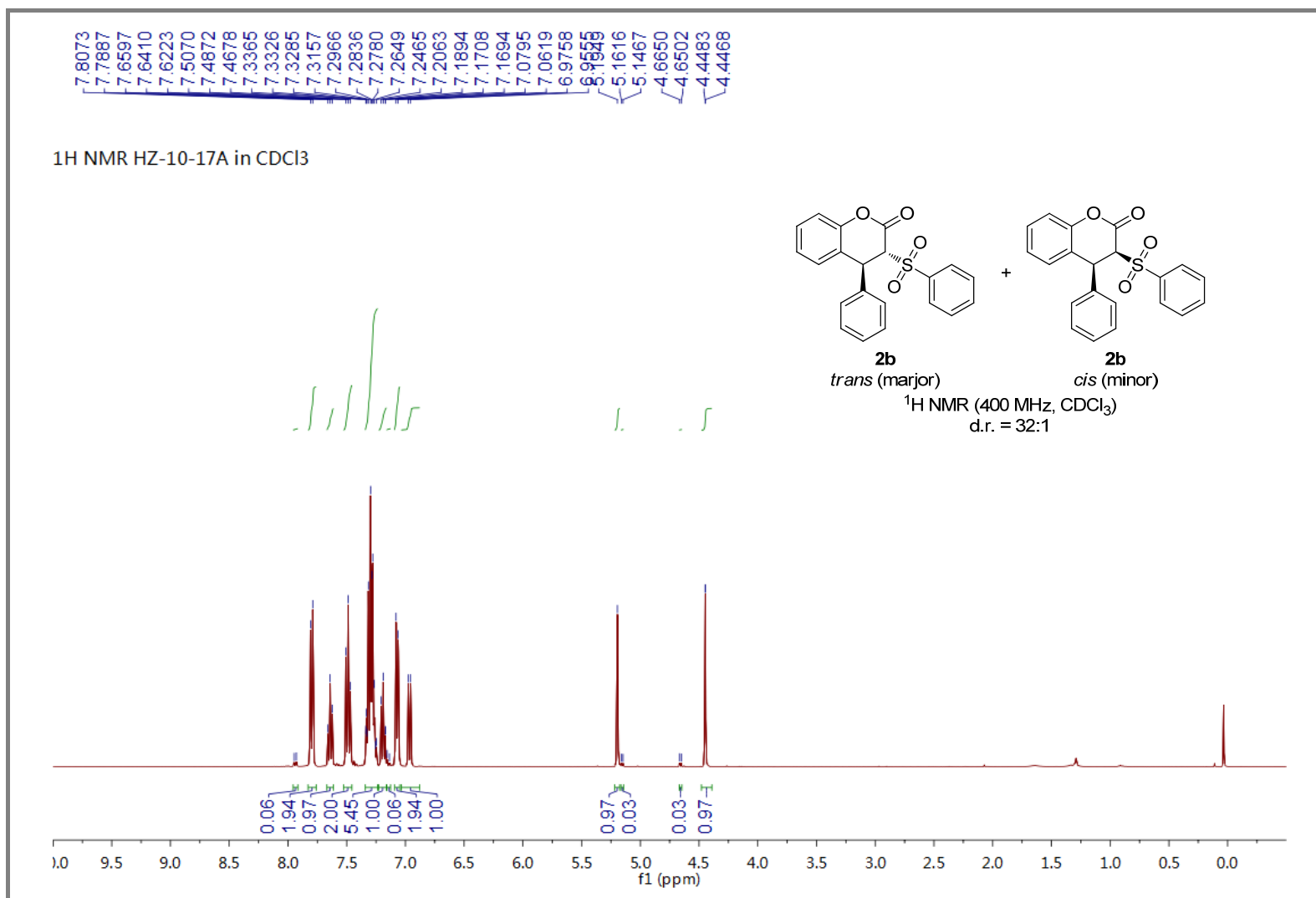


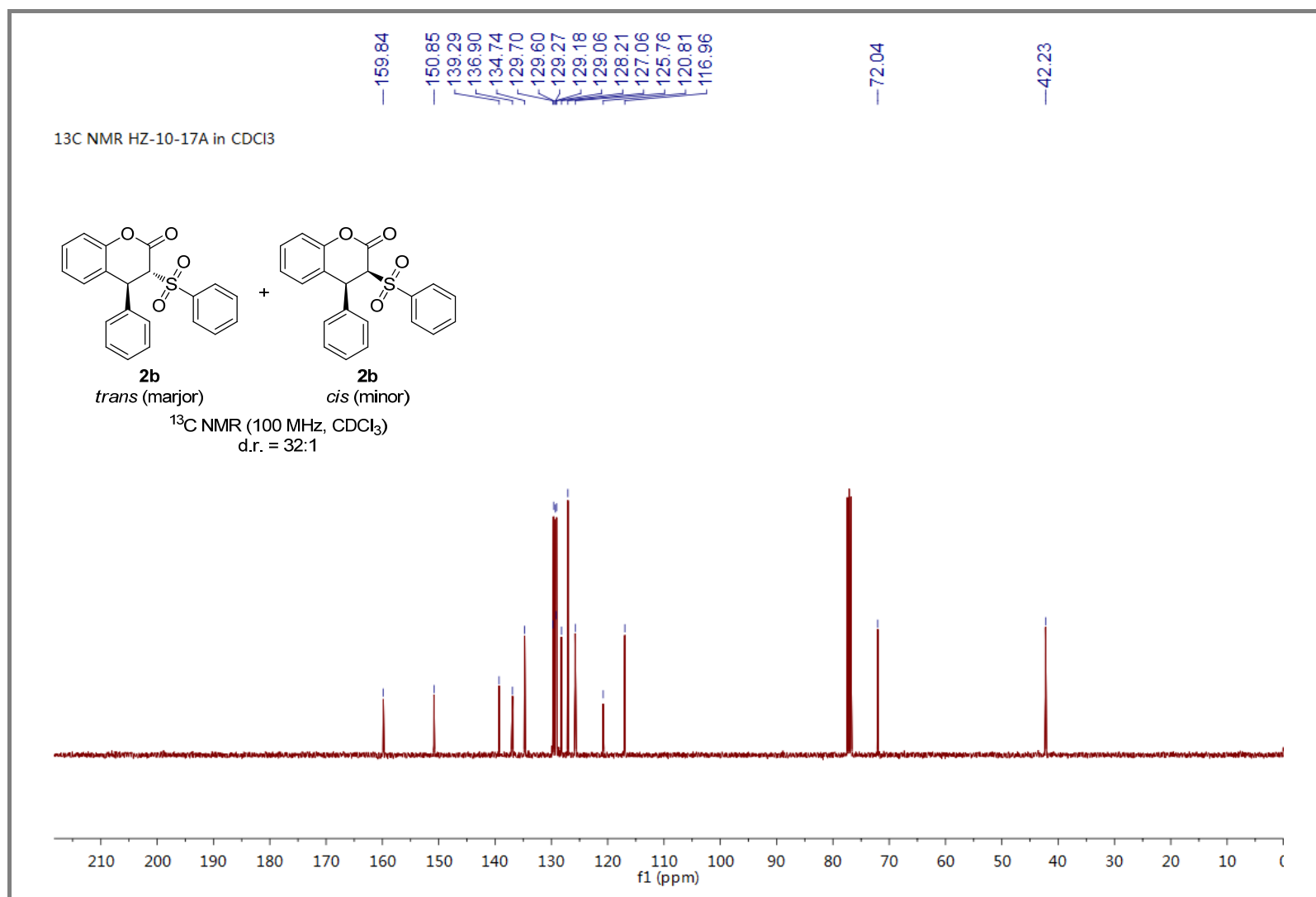


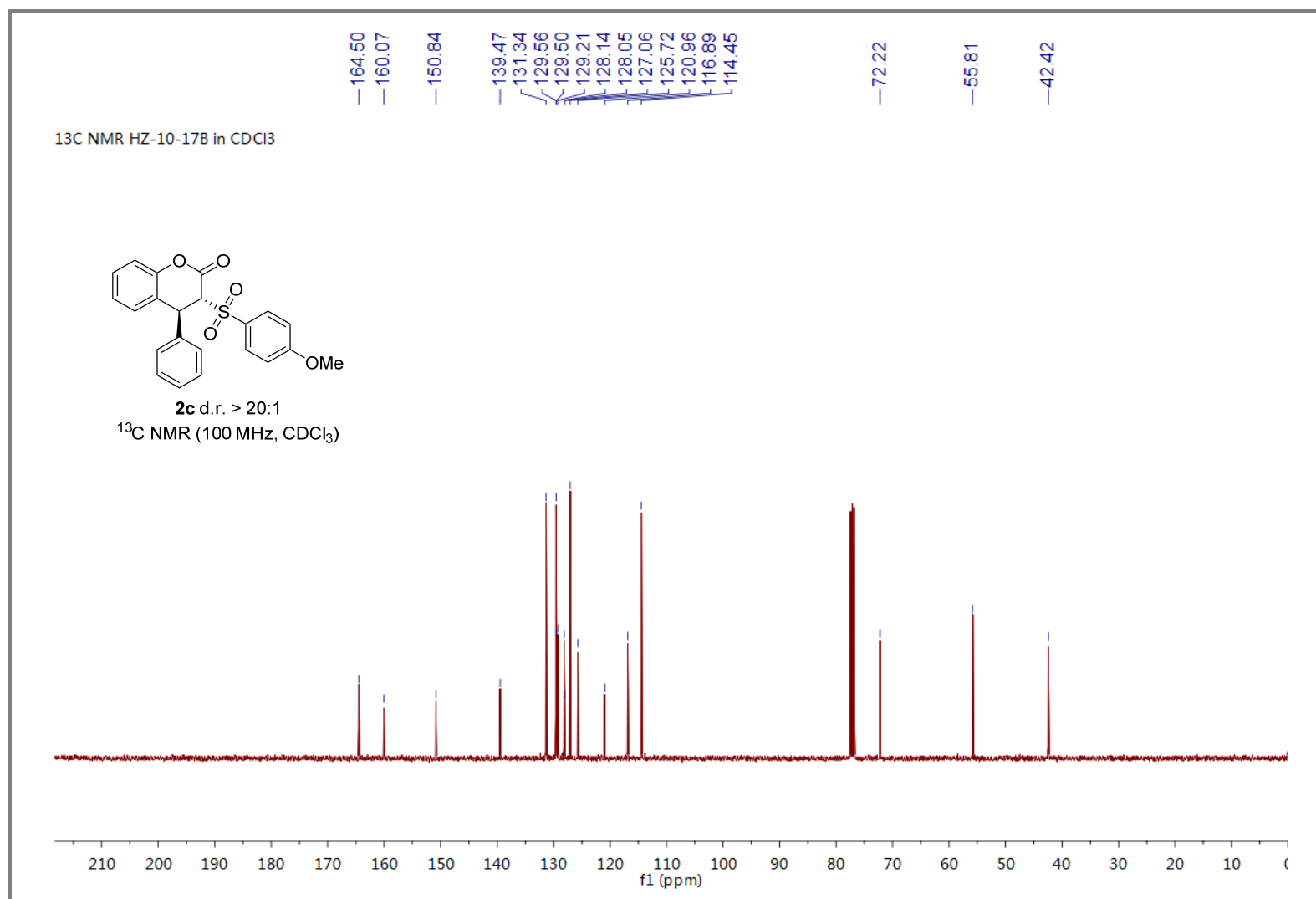


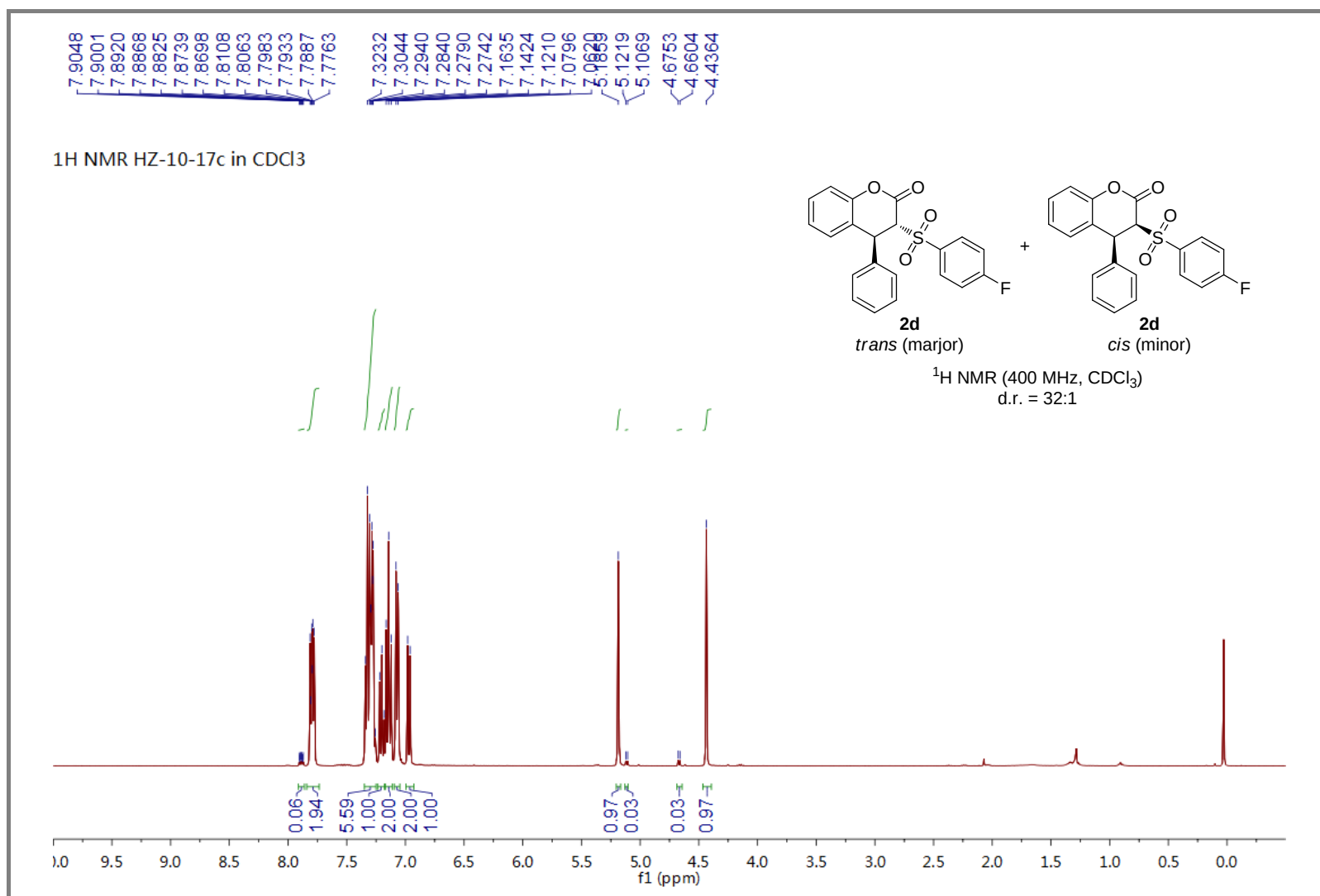


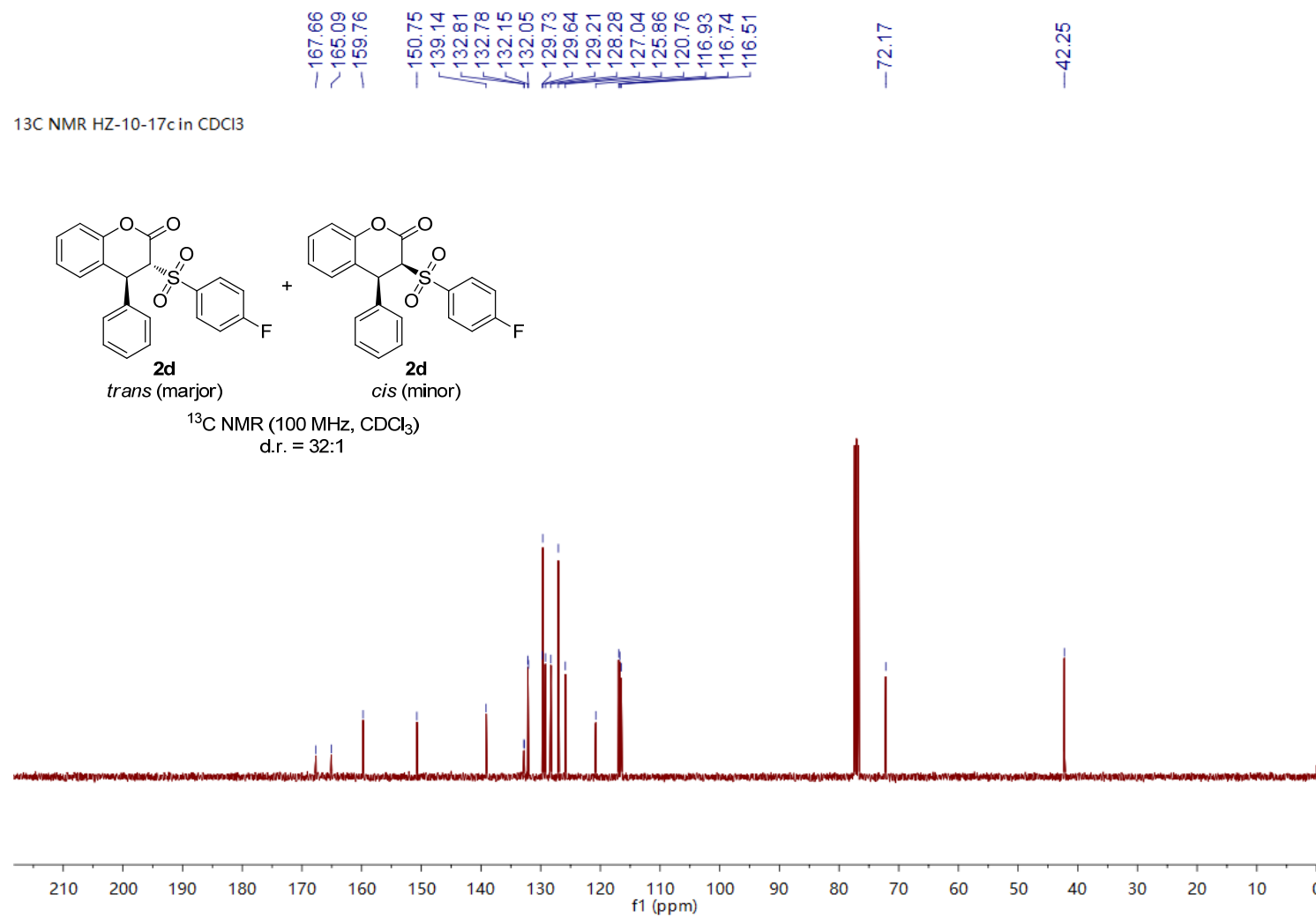




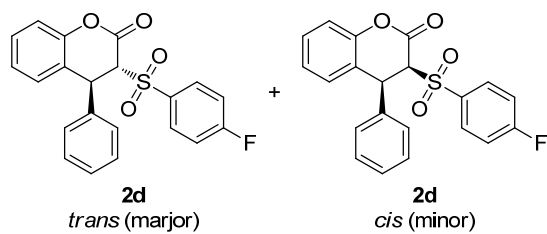




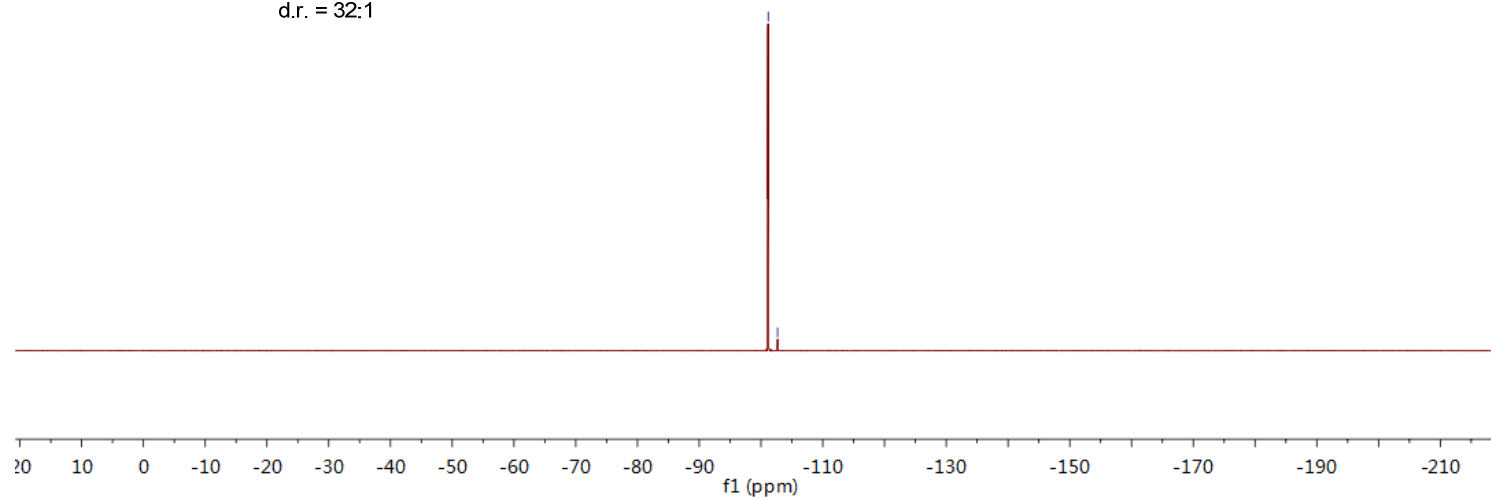


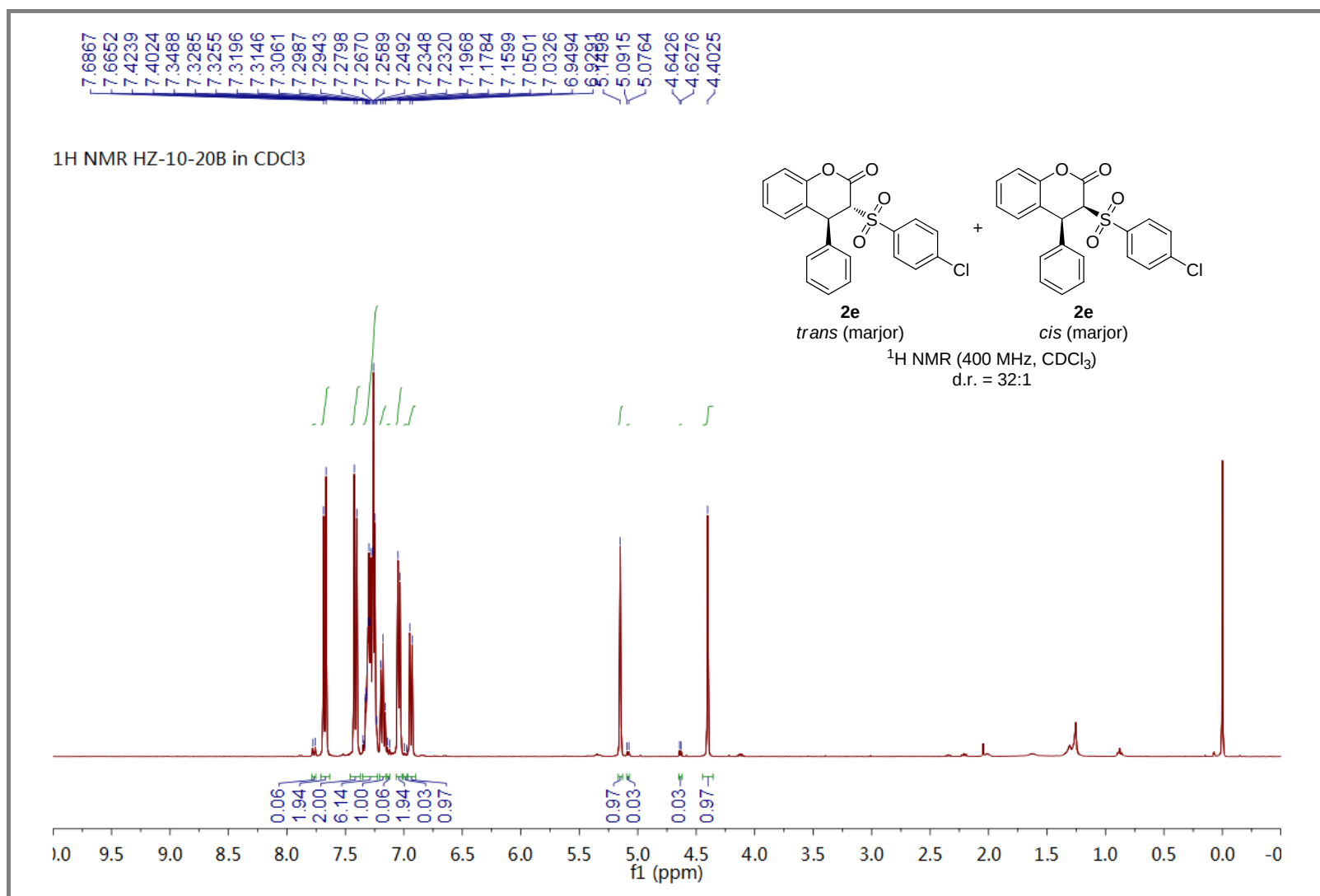


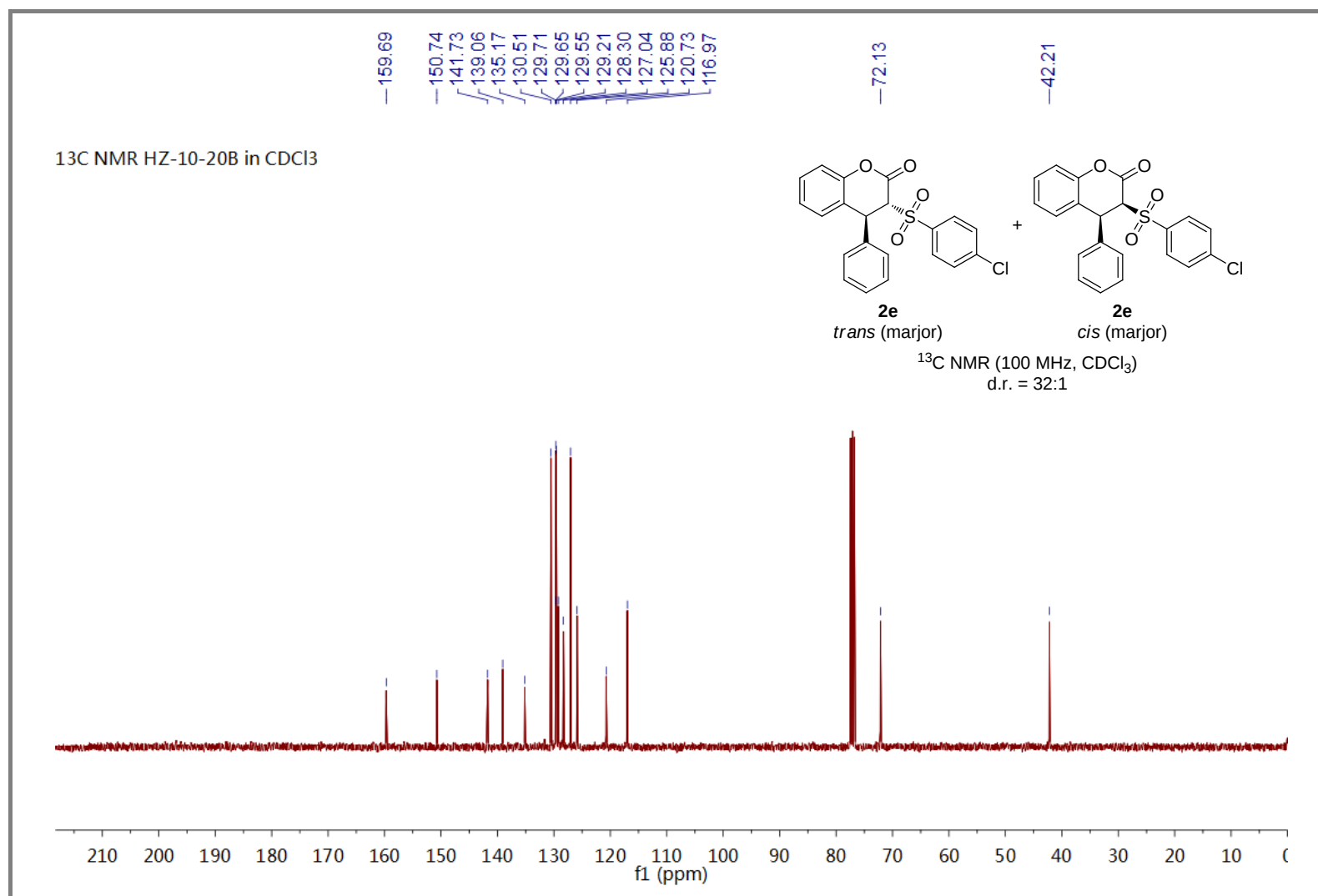
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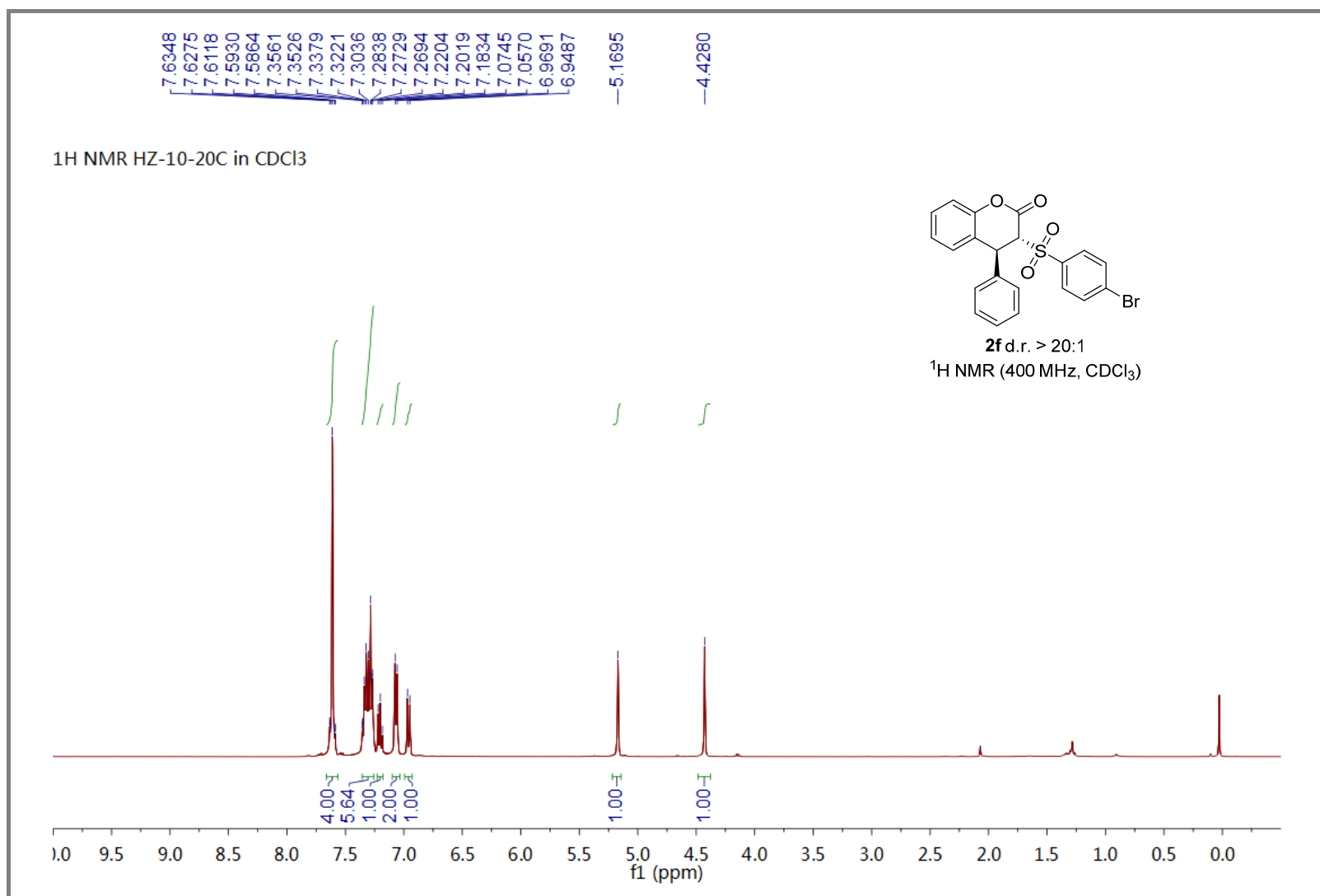


¹⁹F NMR (376 MHz, CDCl₃)
d.r. = 32:1

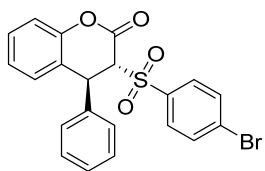






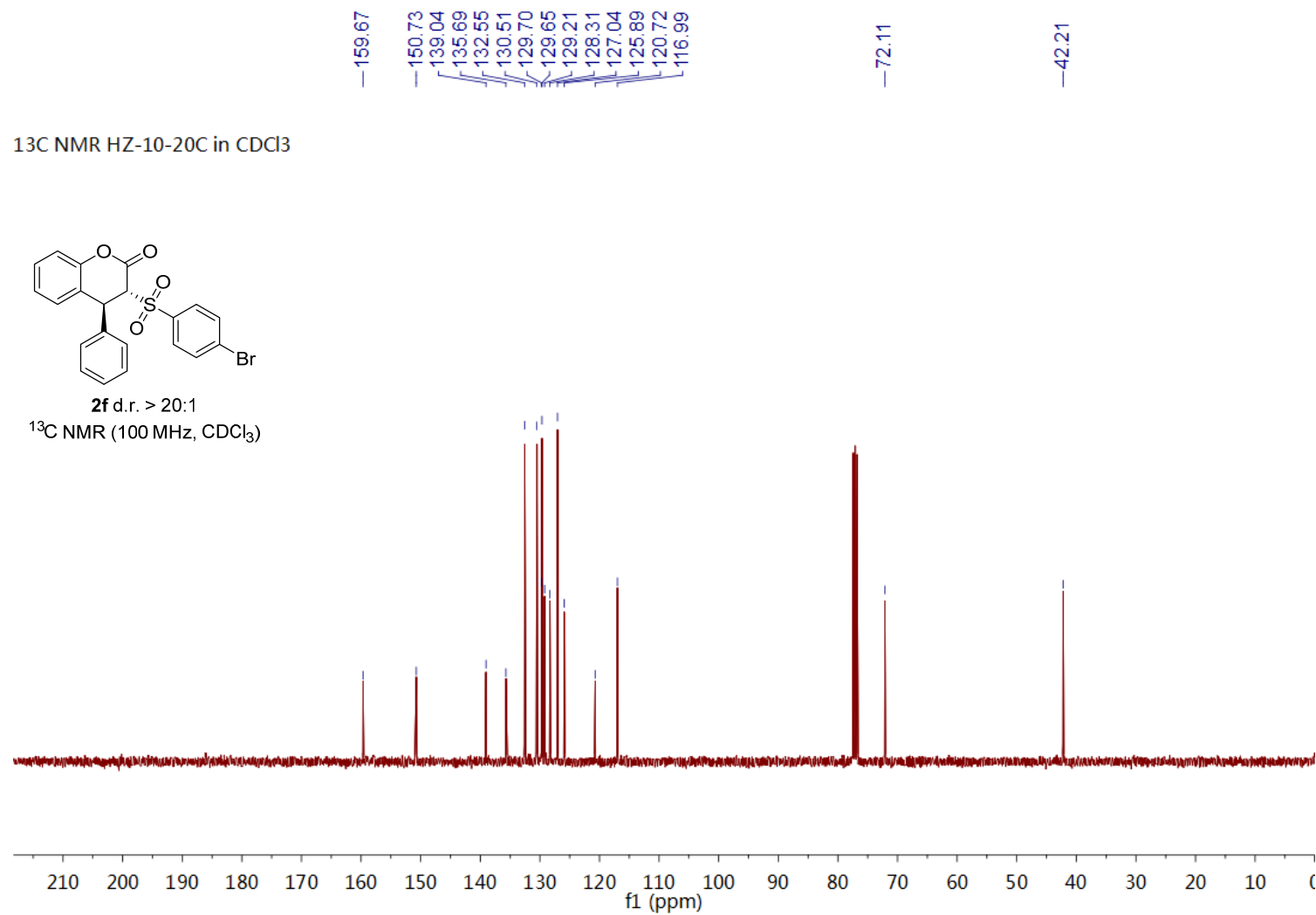


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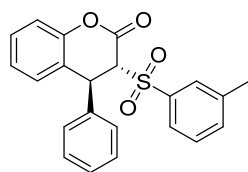


2f d.r. > 20:1

¹³C NMR (100 MHz, CDCl₃)

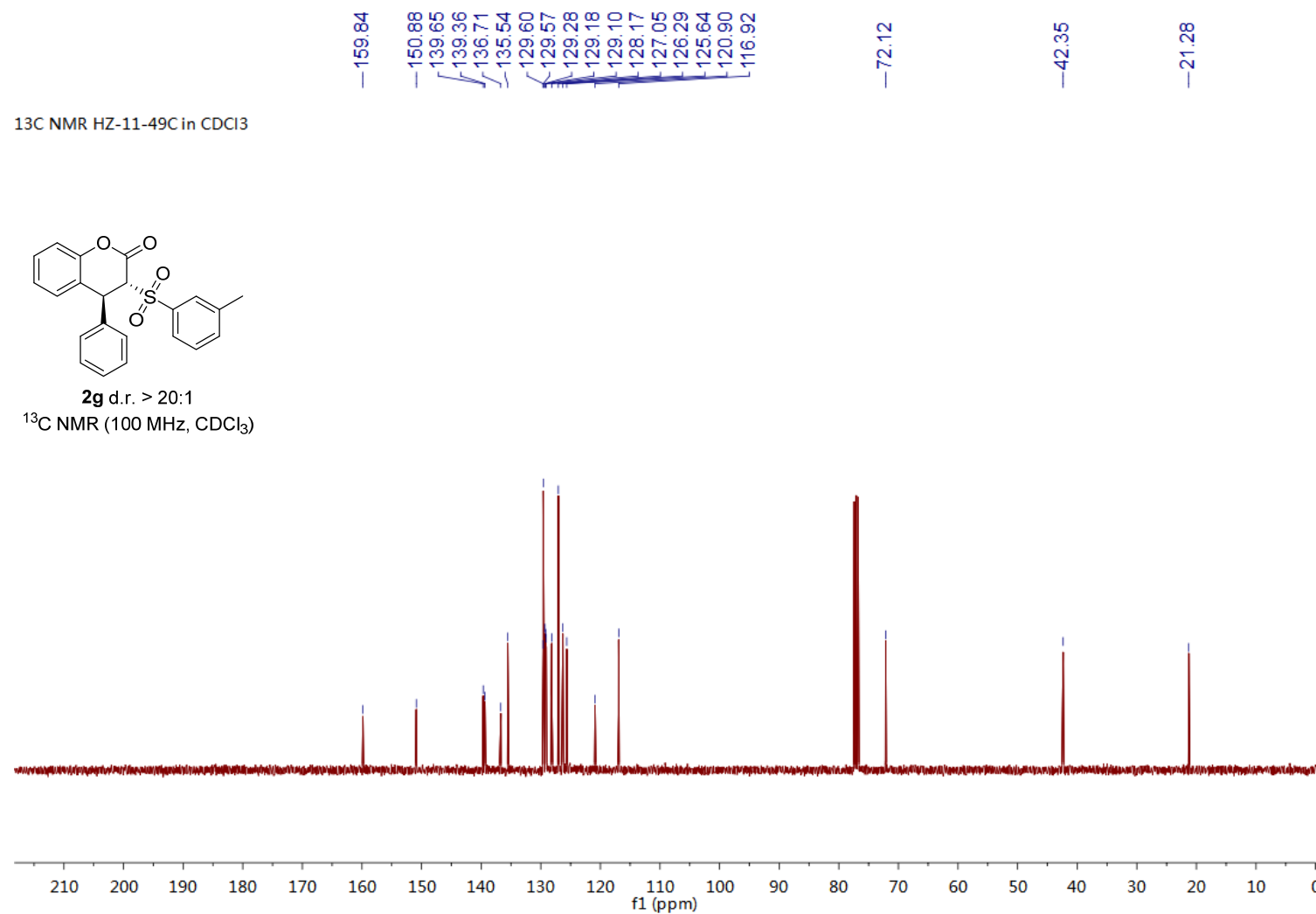


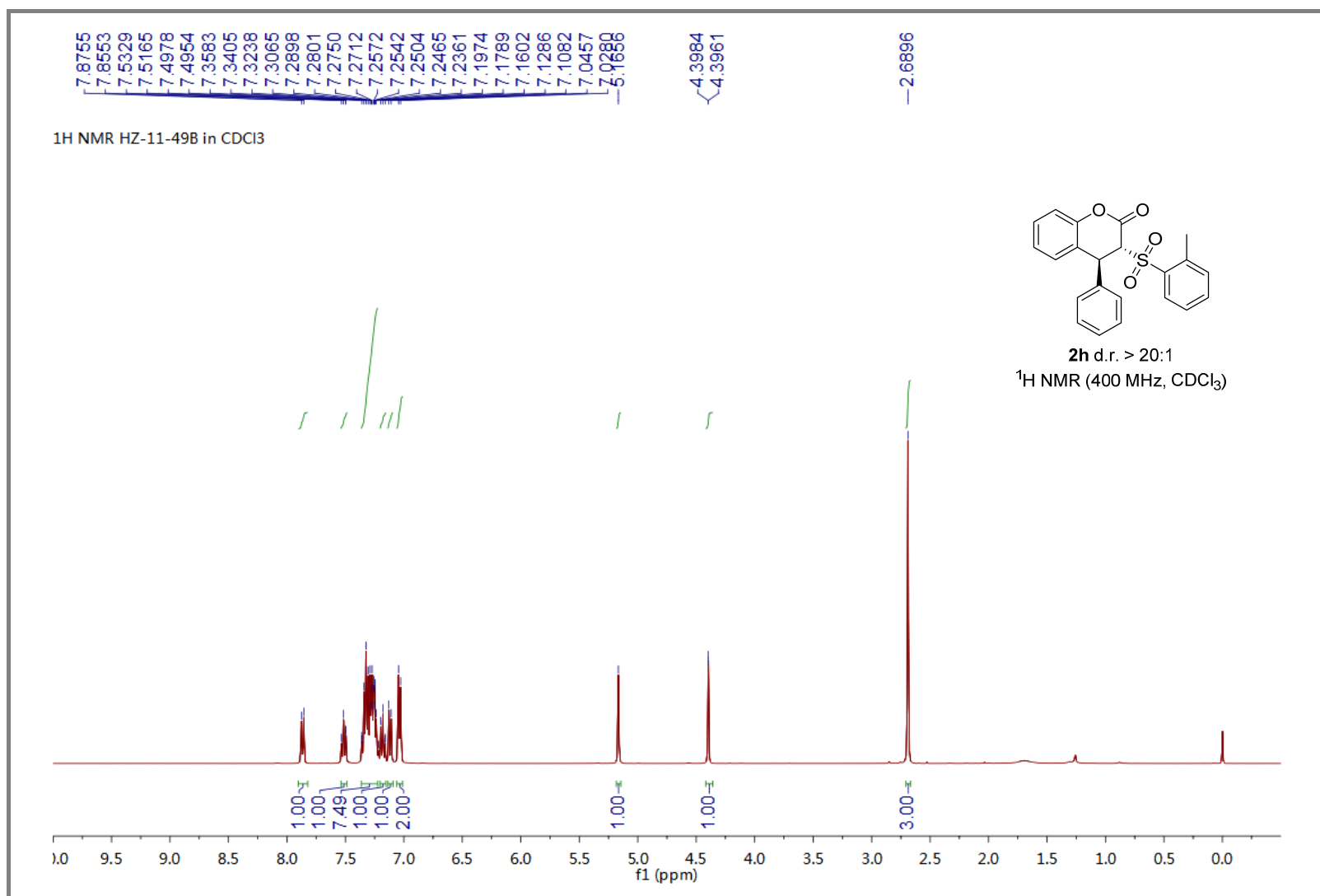
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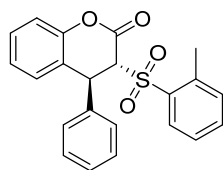
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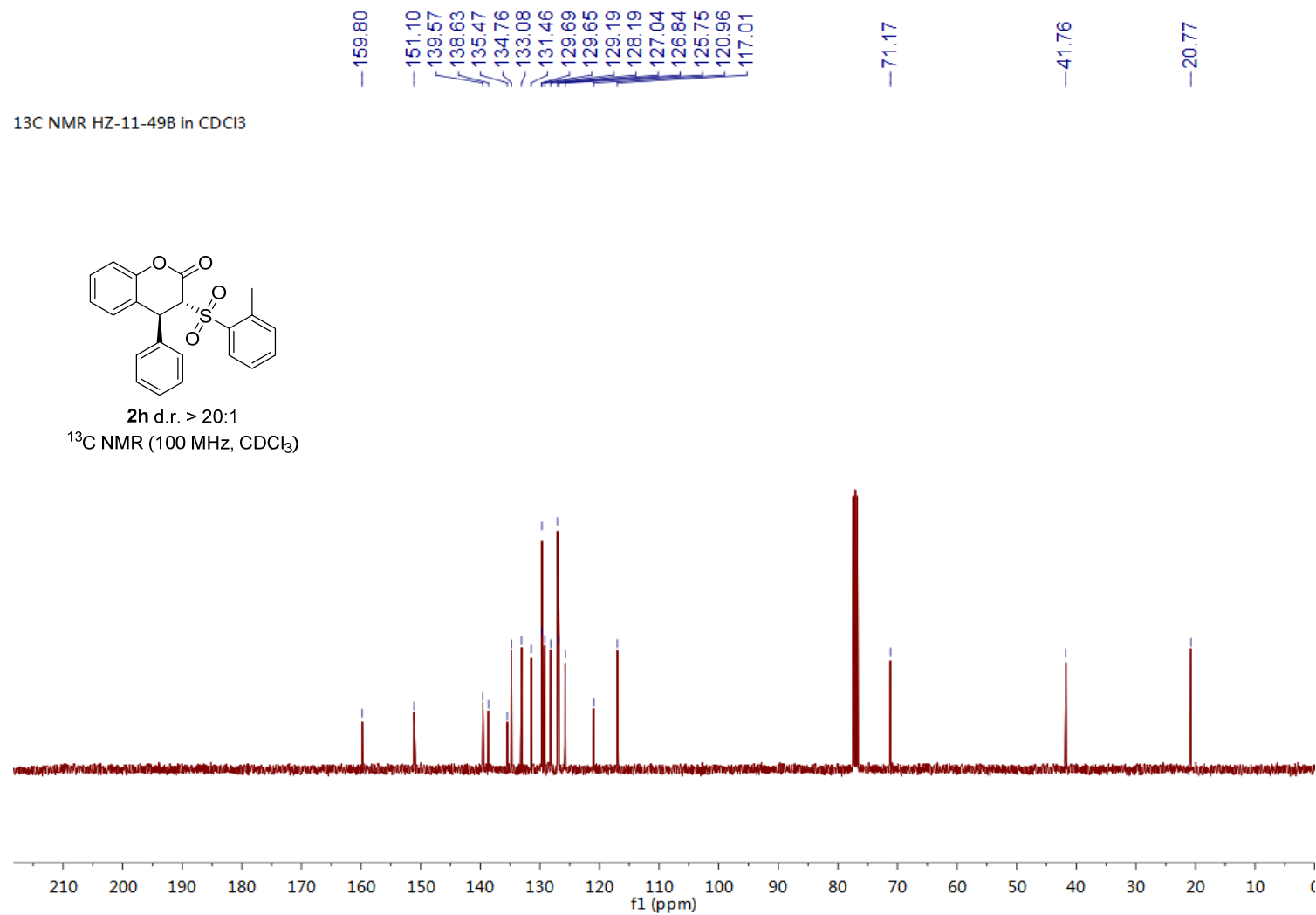


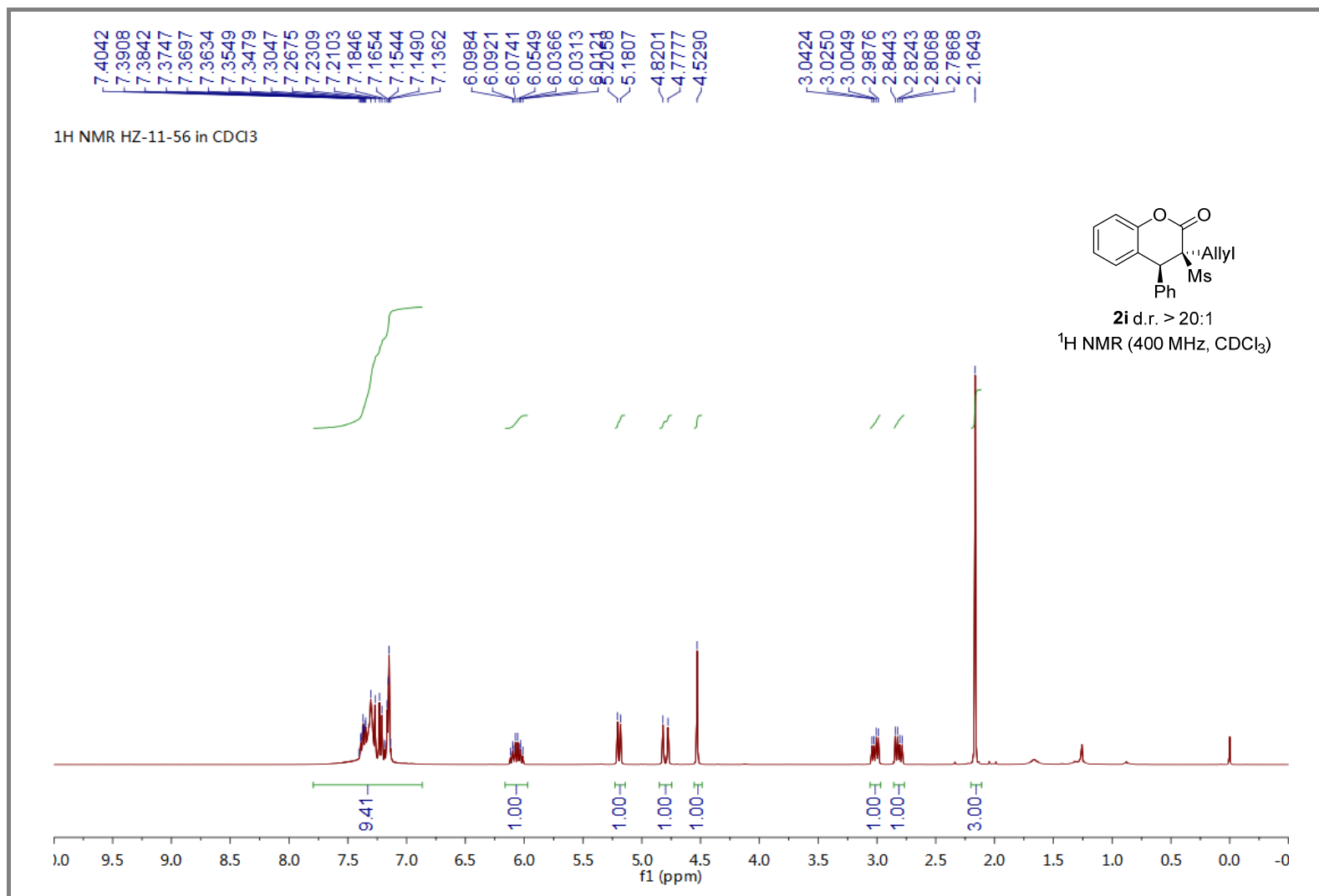
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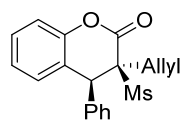
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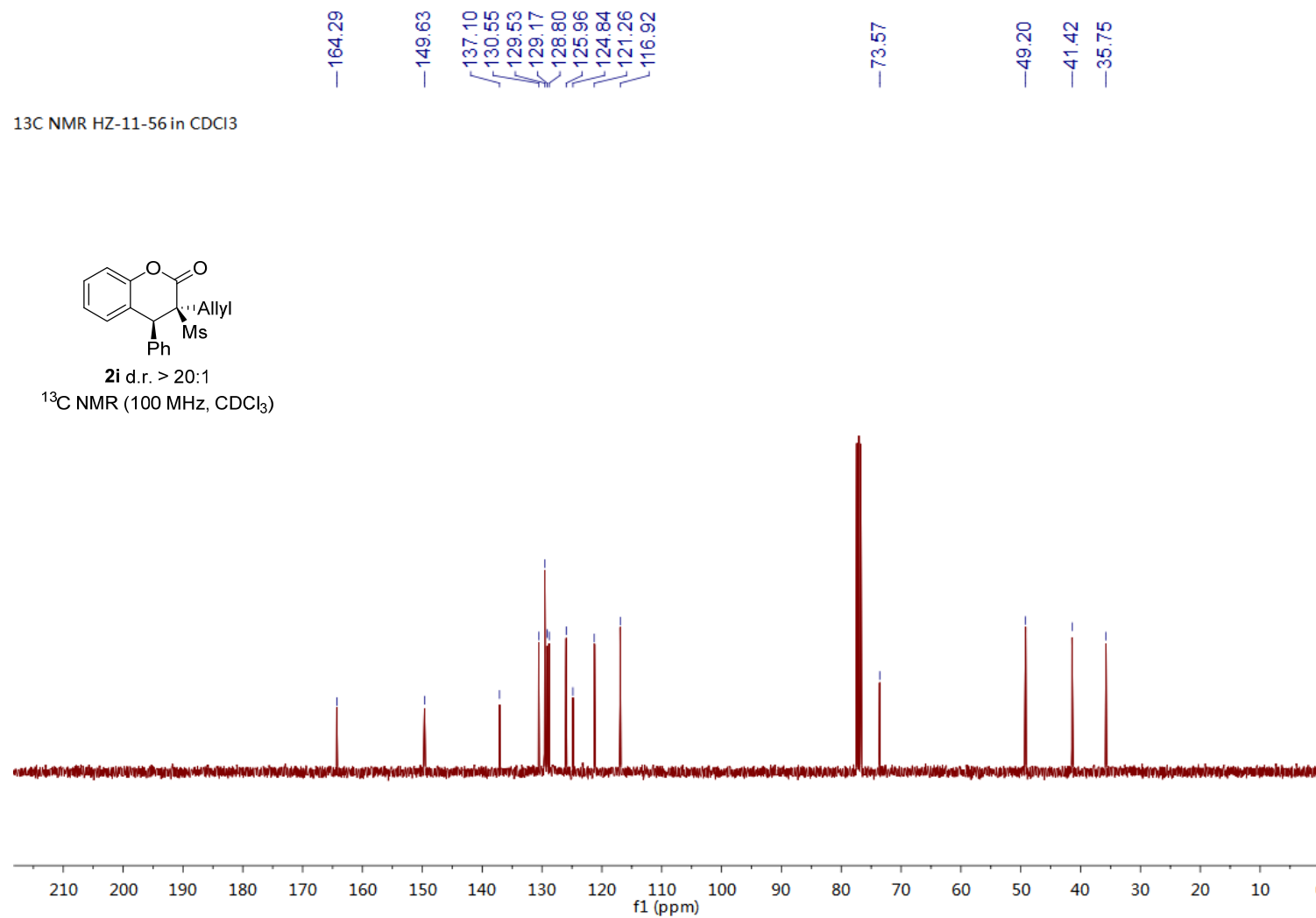


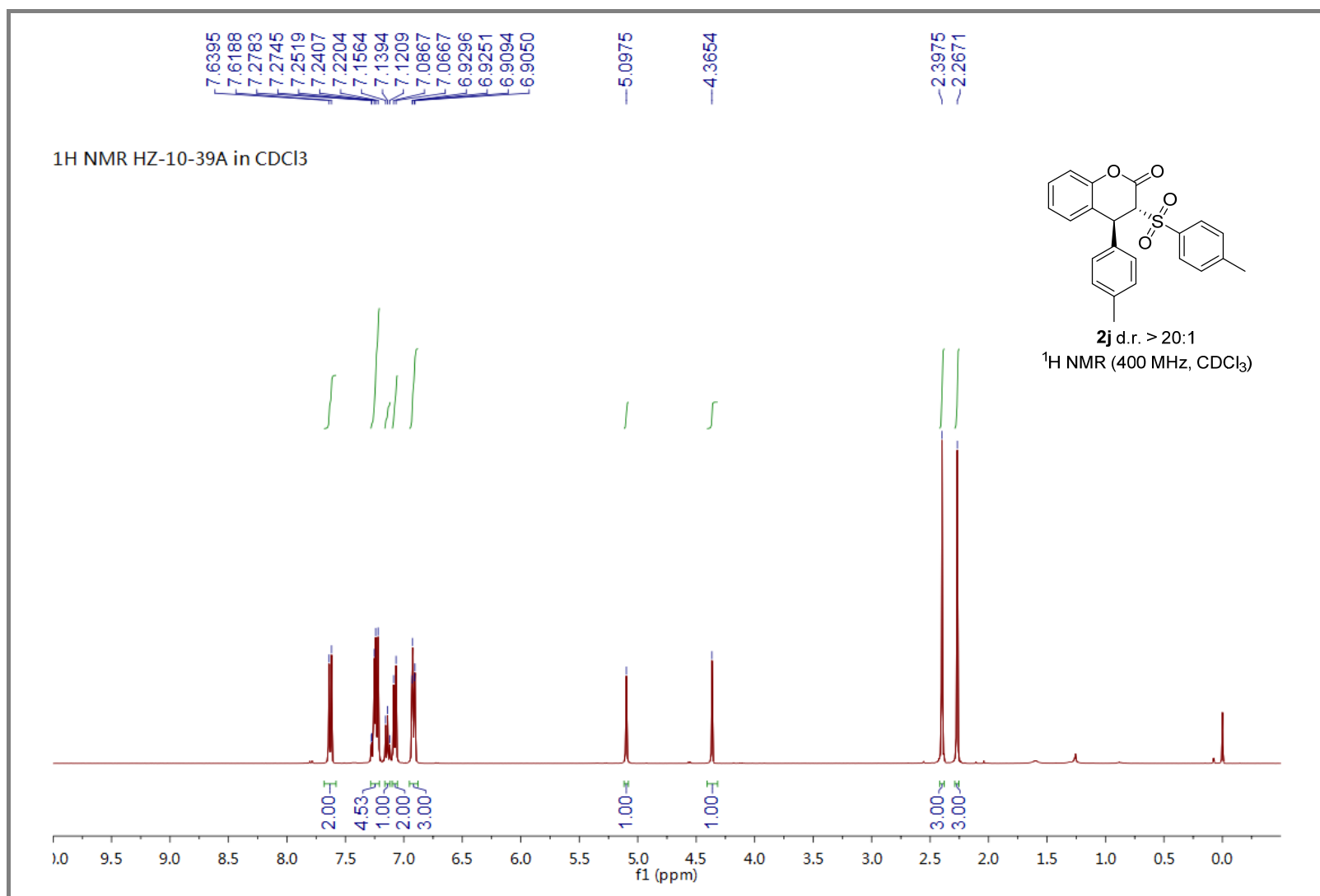
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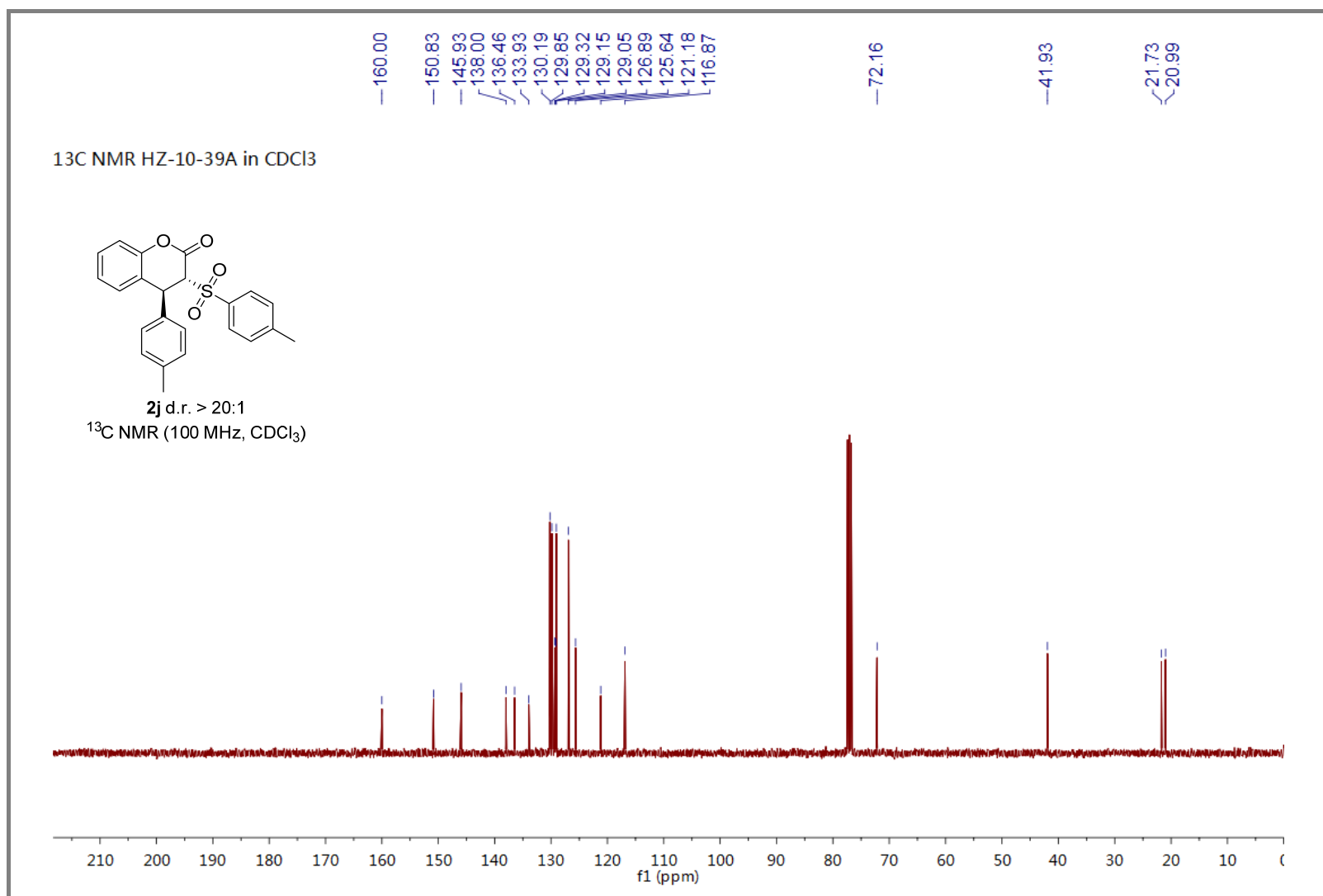


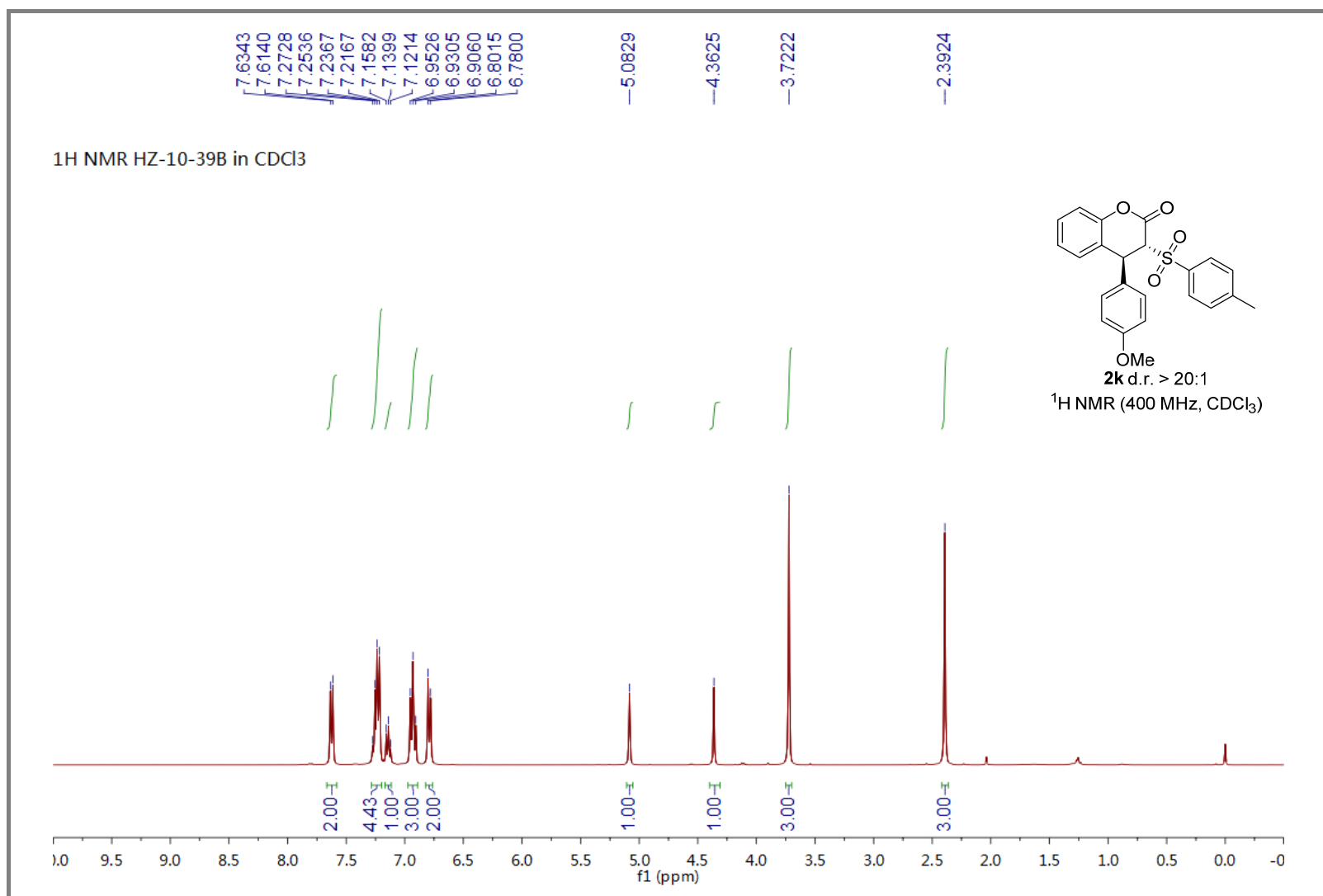
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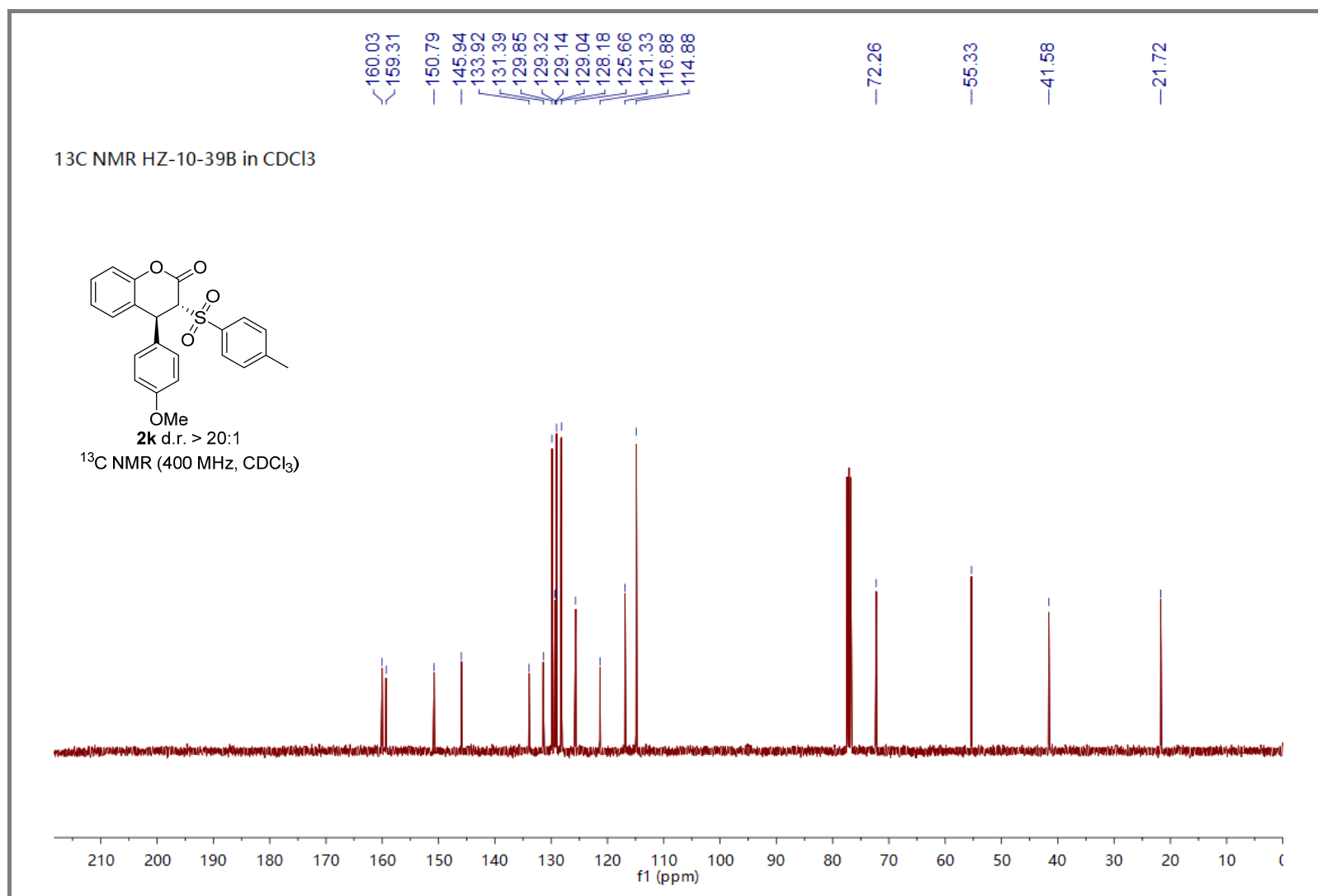
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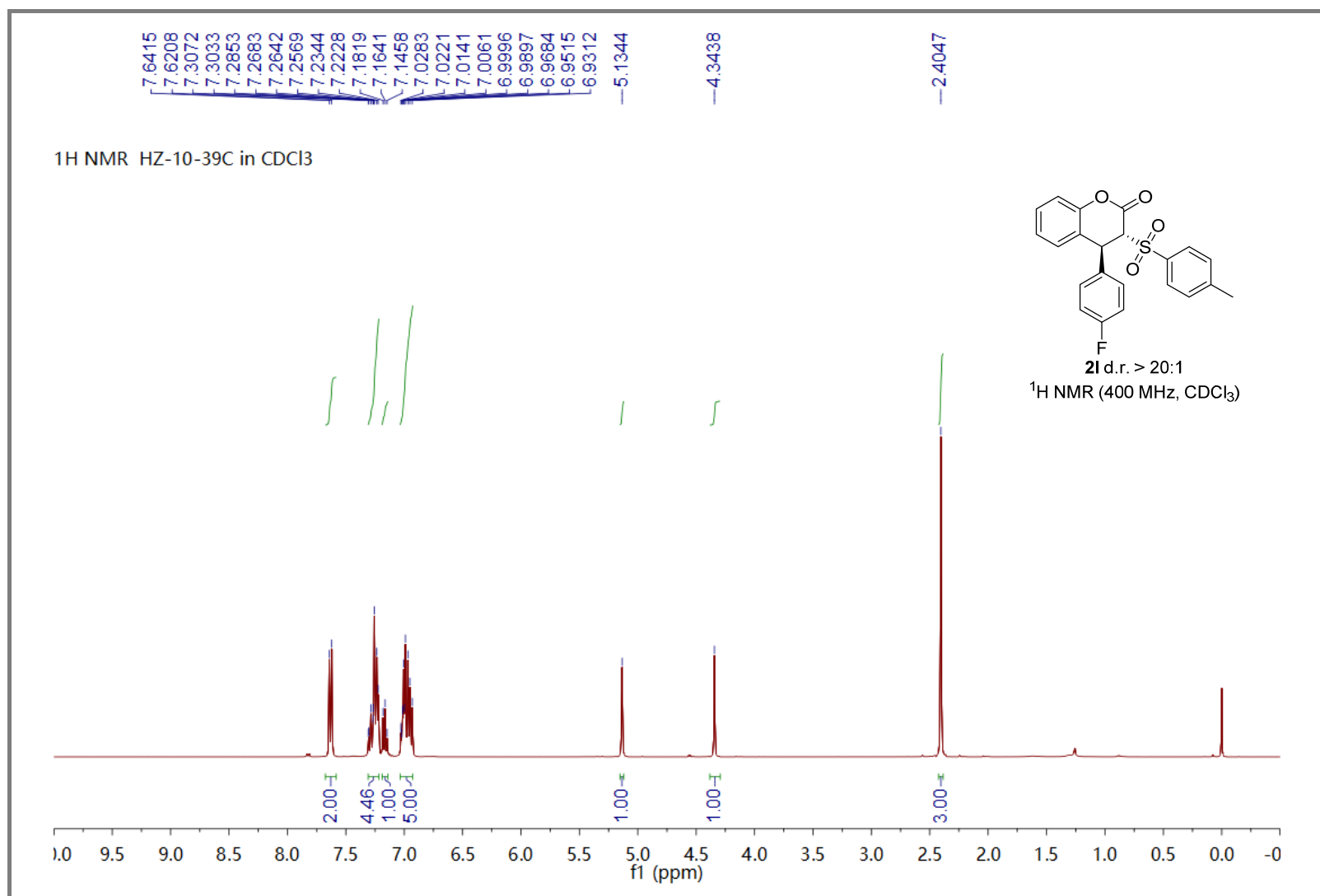




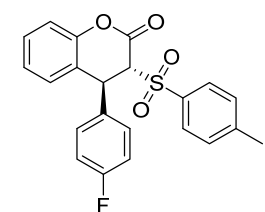




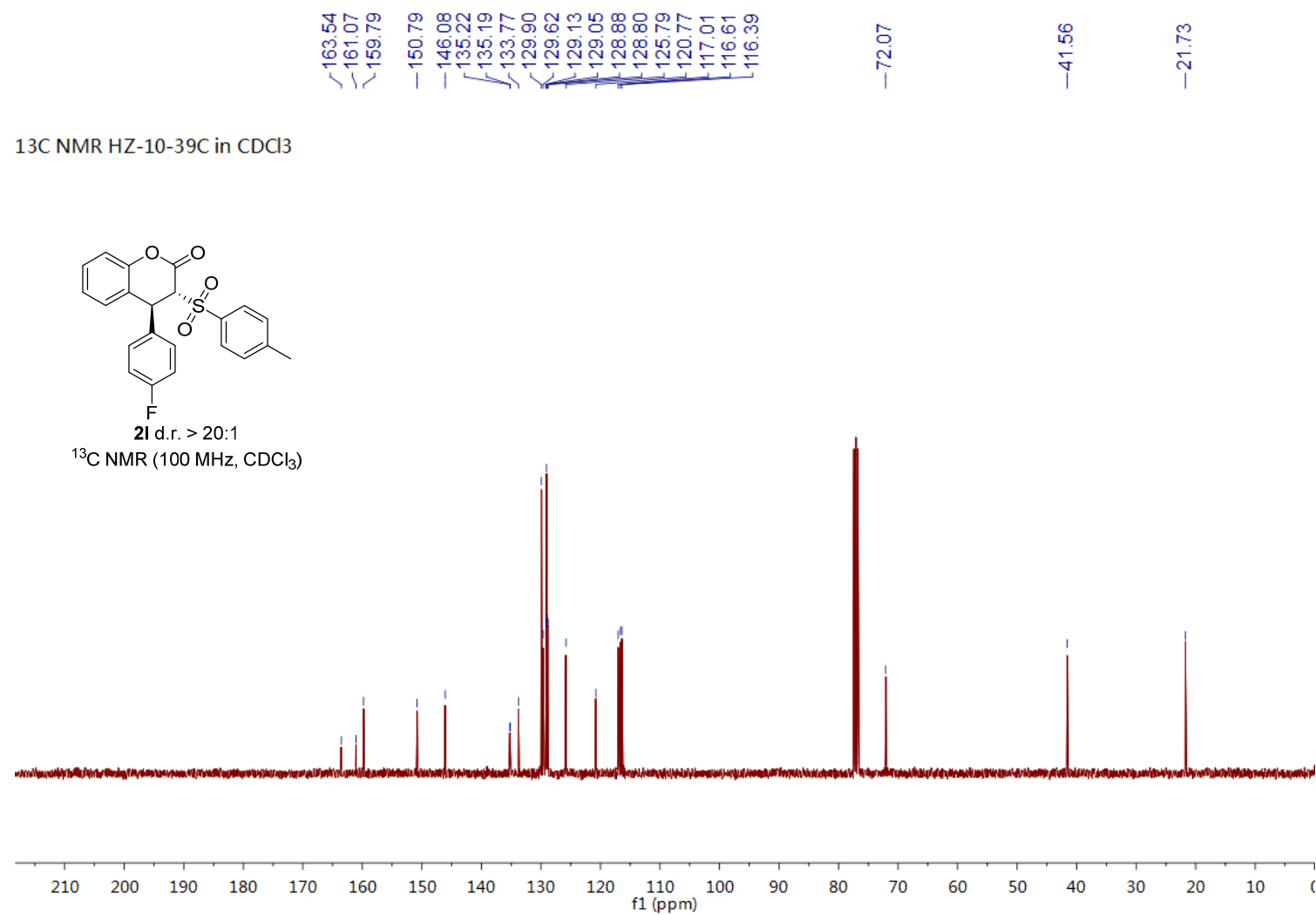




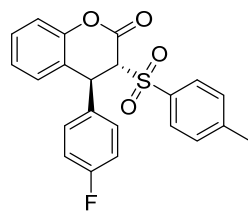
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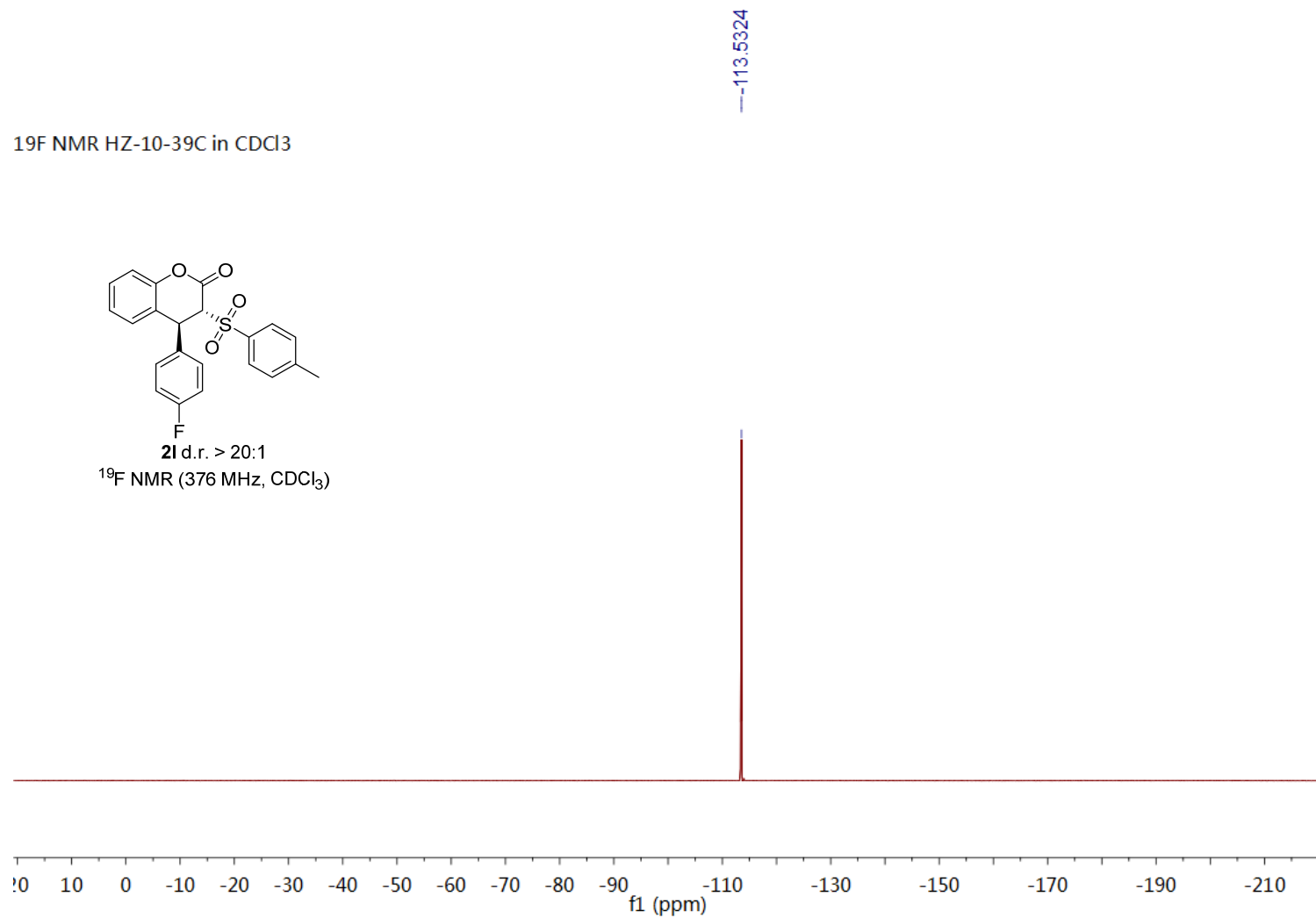


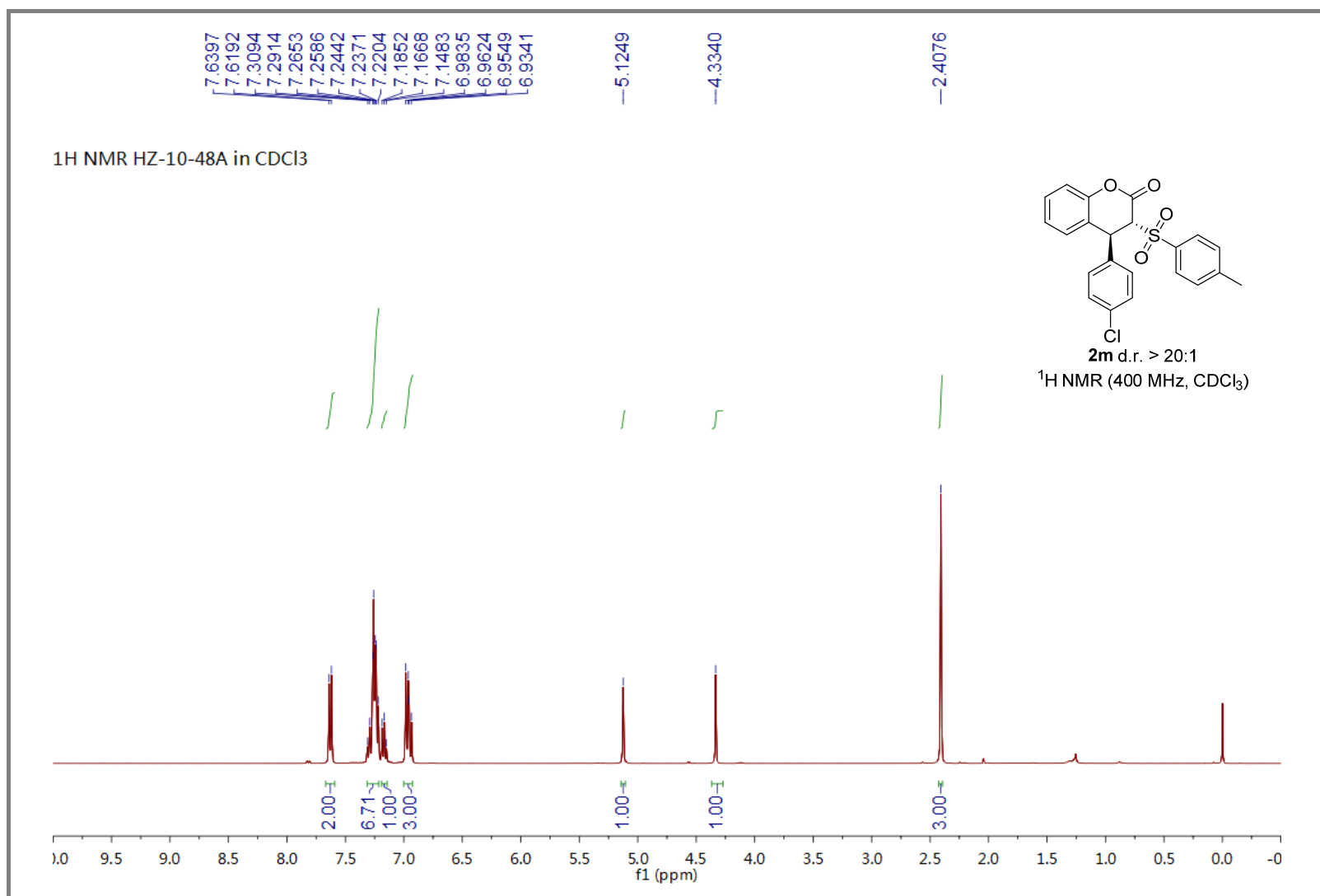
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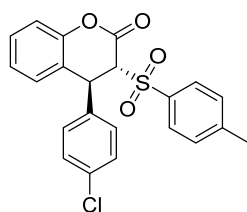
21 d.r. > 20:1

¹⁹F NMR (376 MHz, CDCl₃)

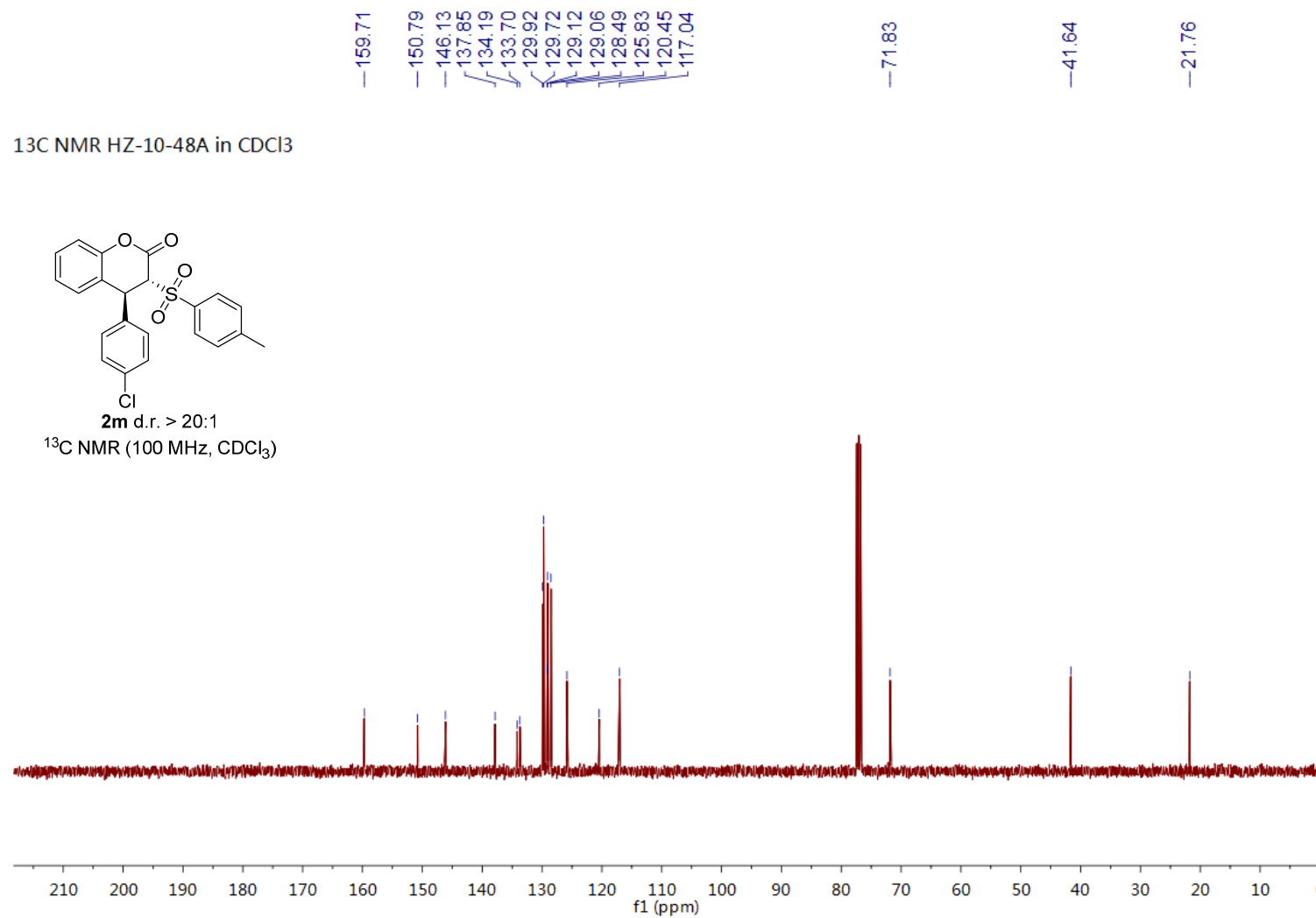


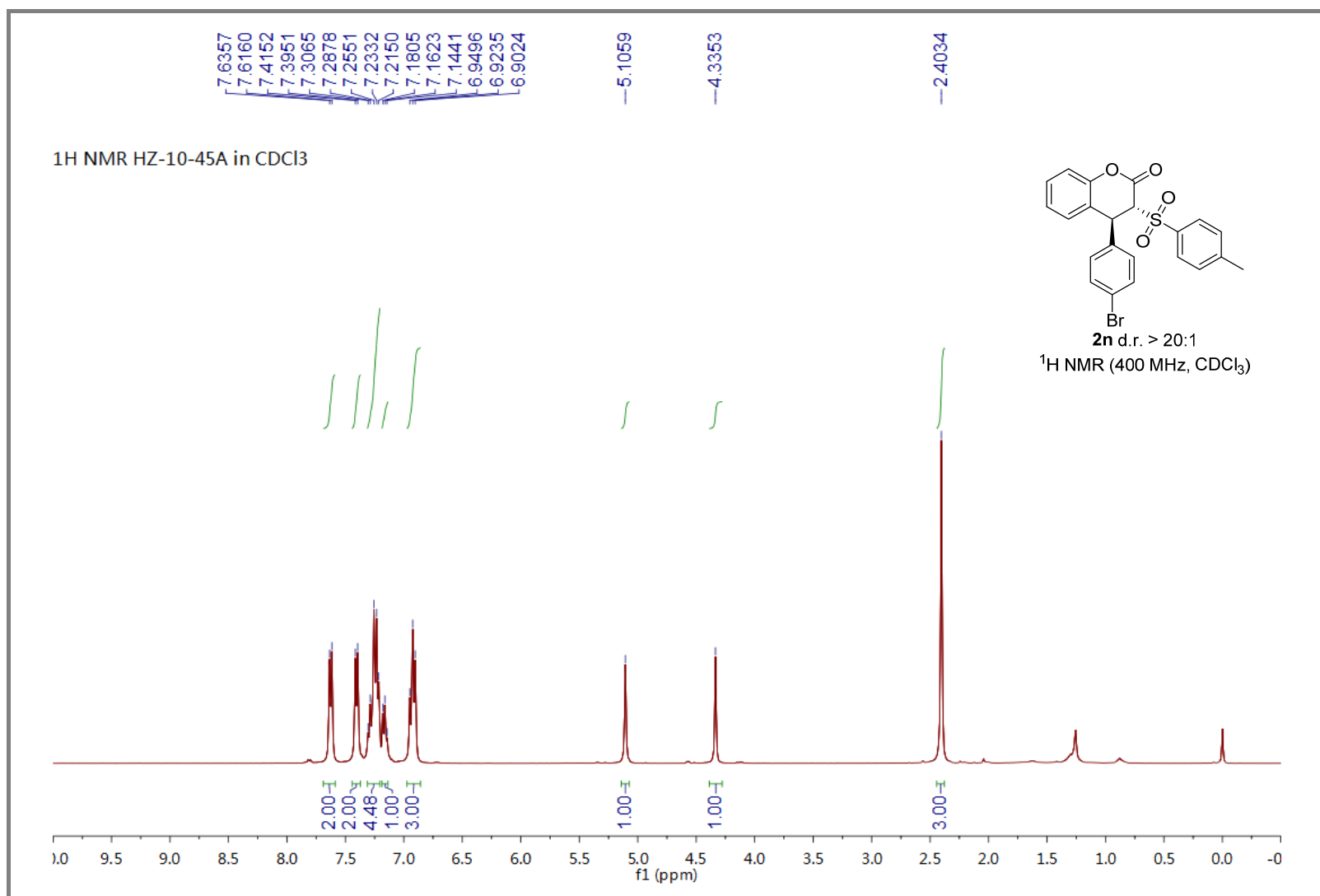


¹³C NMR HZ-10-48A in CDCl₃

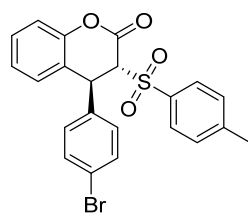


¹³C NMR (100 MHz, CDCl₃)

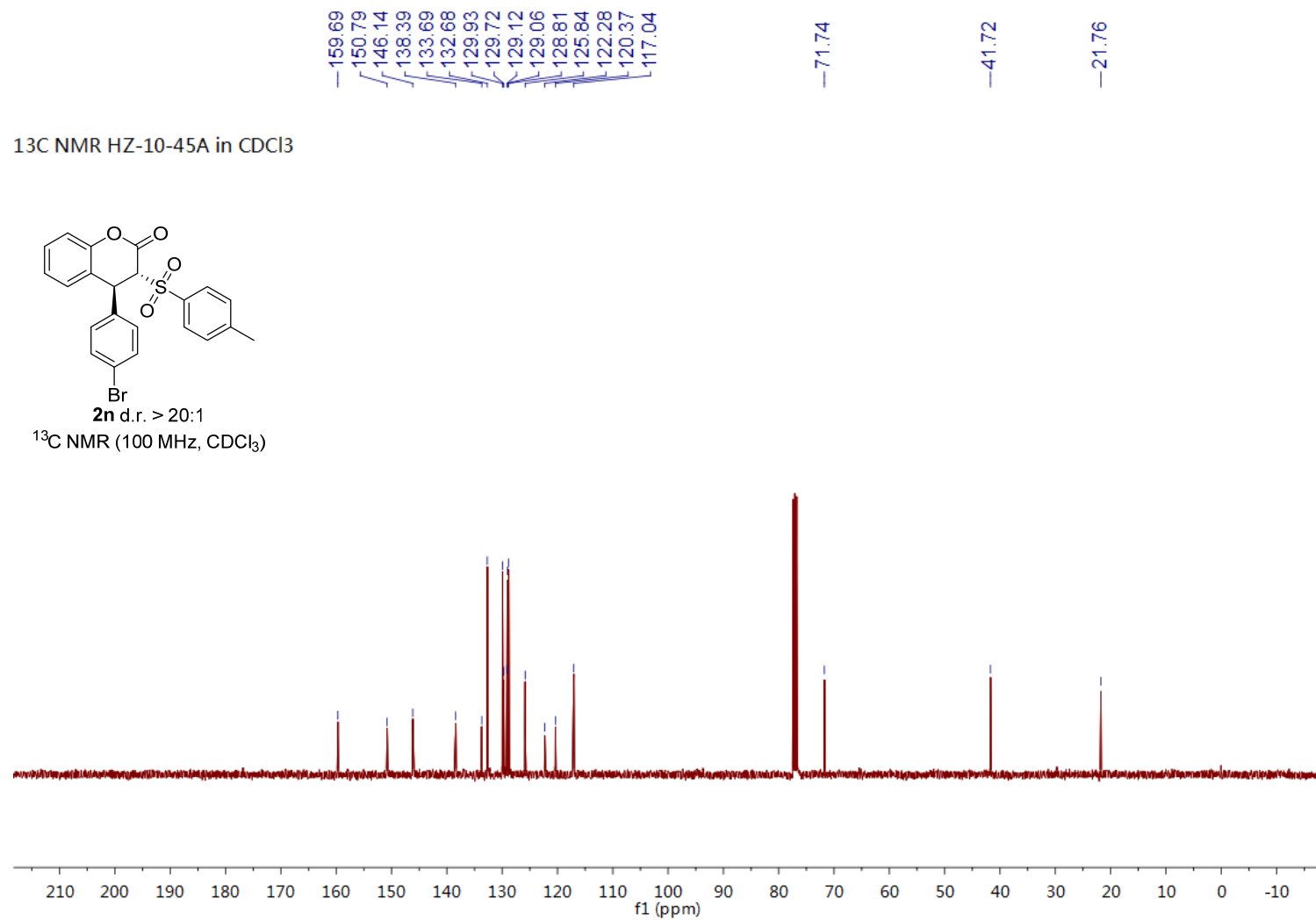




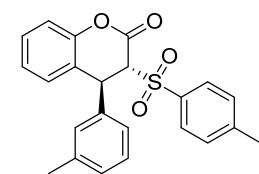
¹³C NMR HZ-10-45A in CDCl₃



¹³C NMR (100 MHz, CDCl₃)

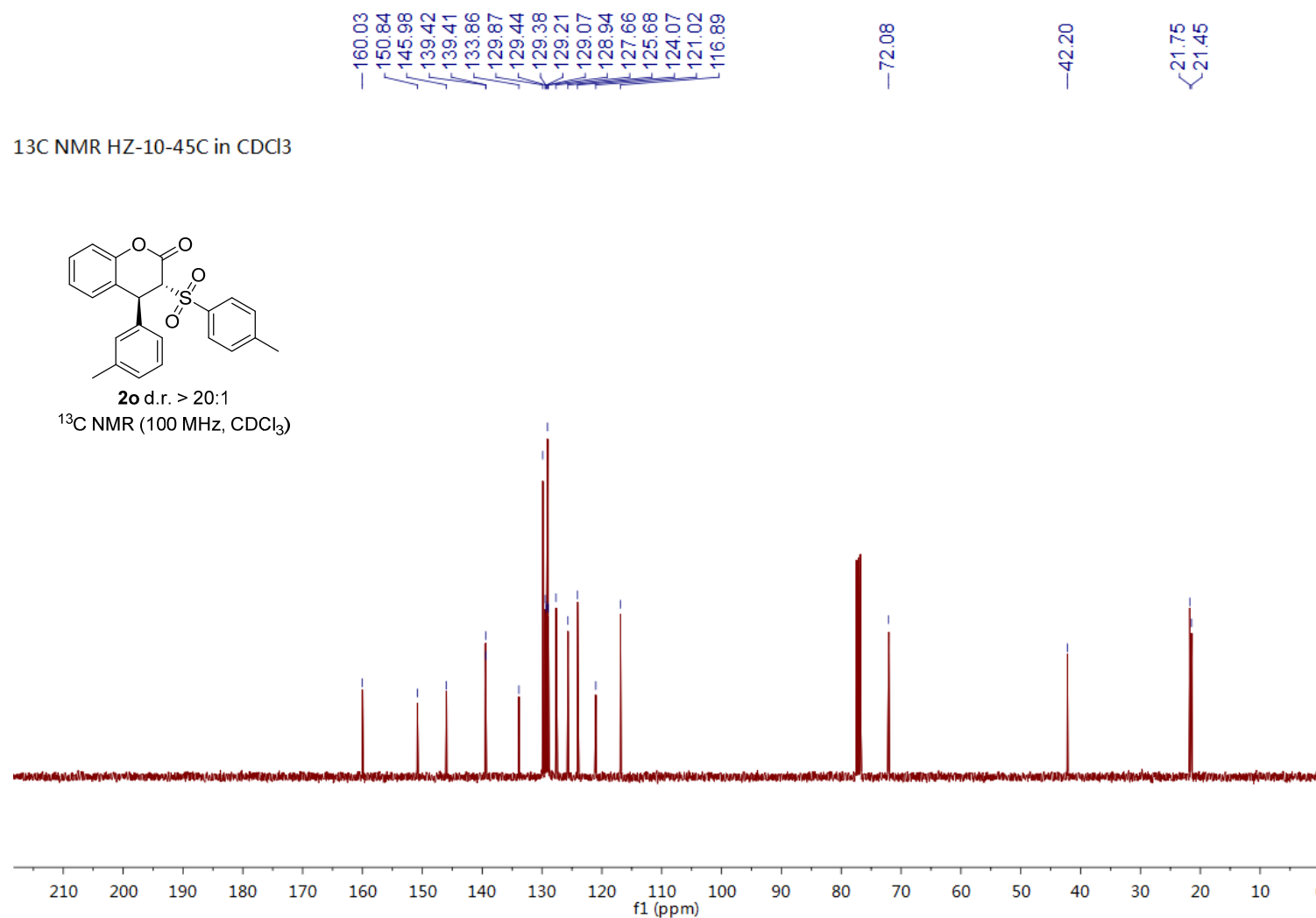


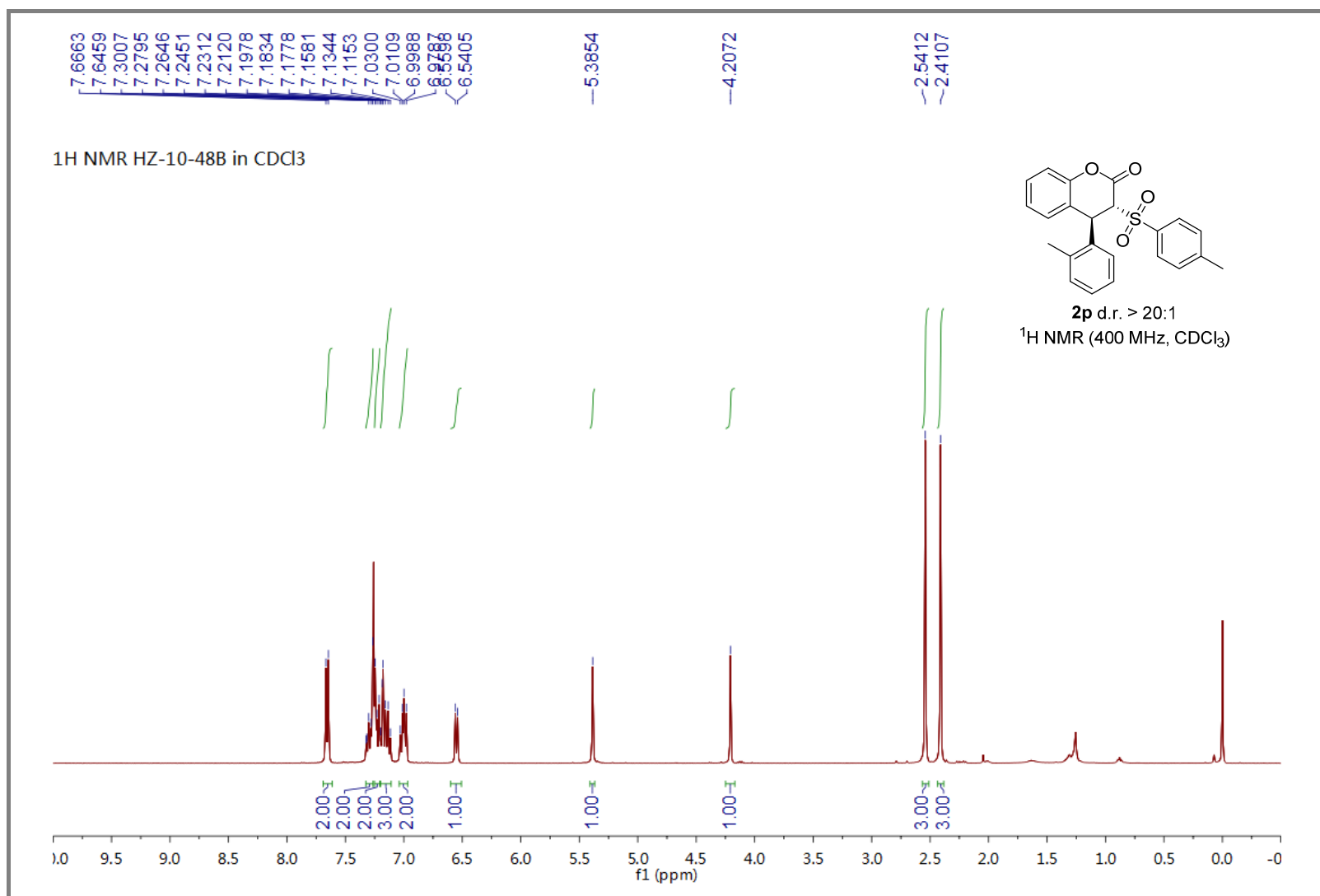
¹³C NMR HZ-10-45C in CDCl₃

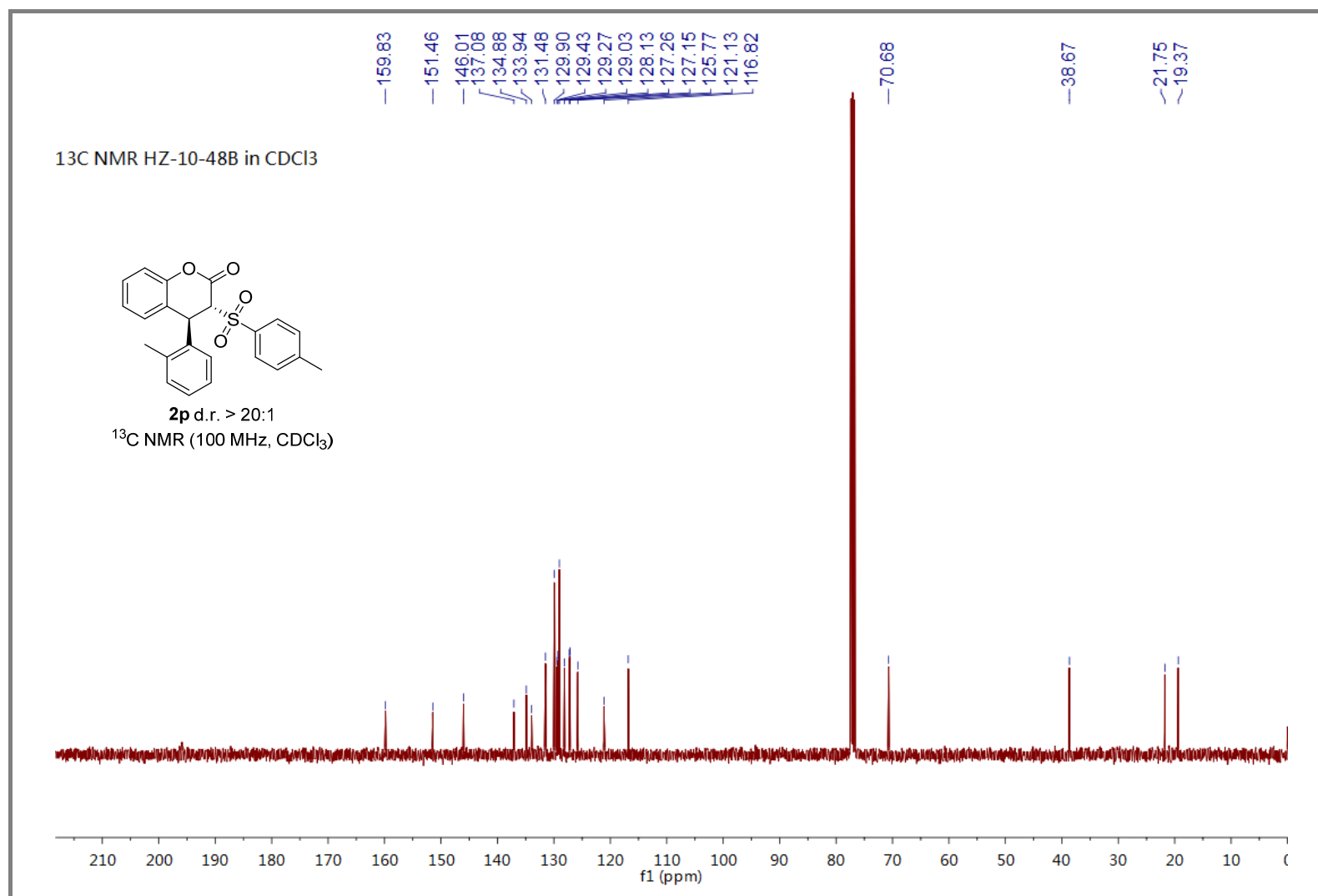


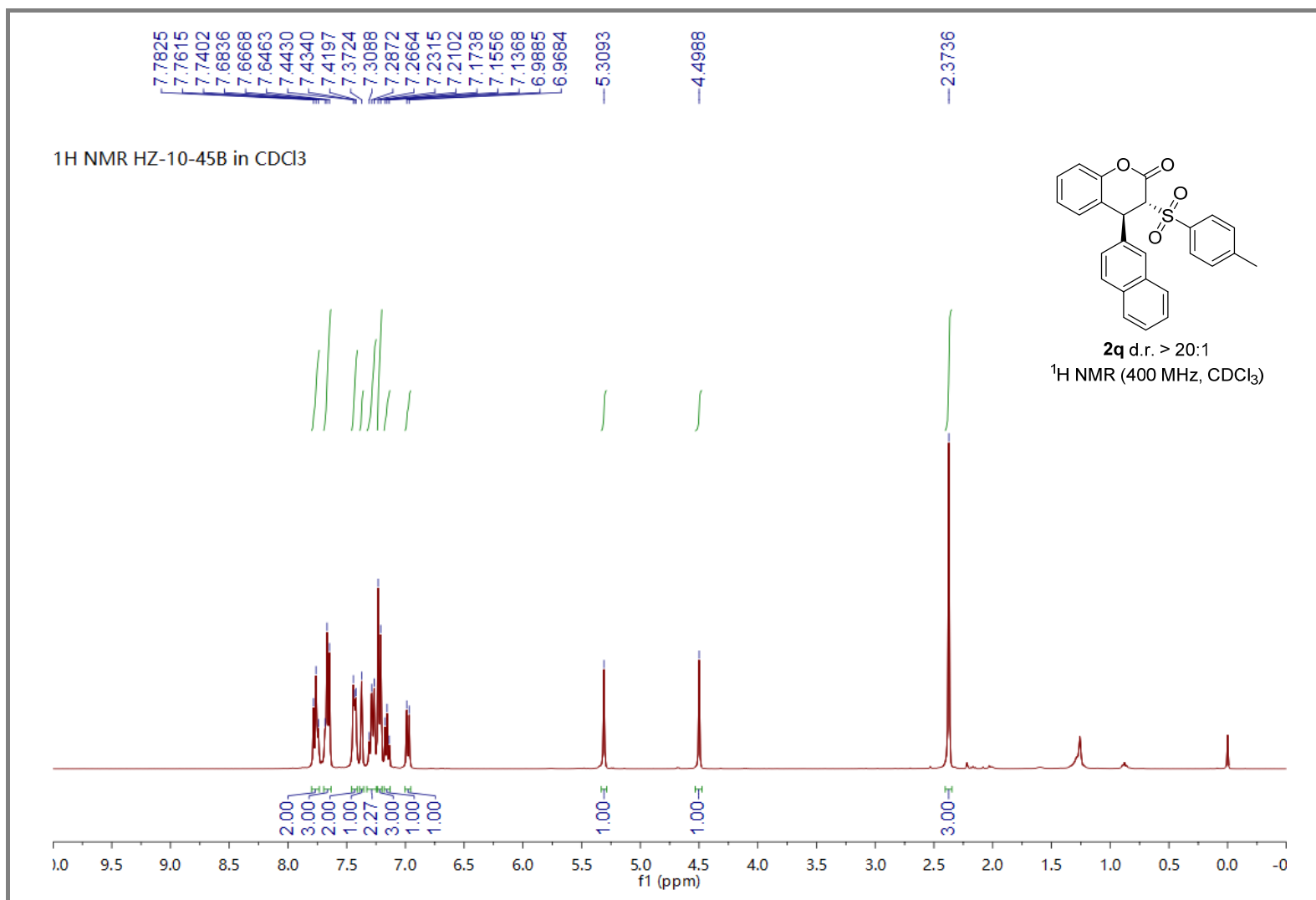
2o d.r. > 20:1

¹³C NMR (100 MHz, CDCl₃)

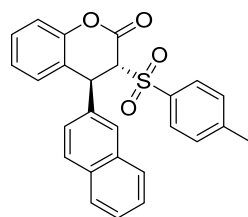






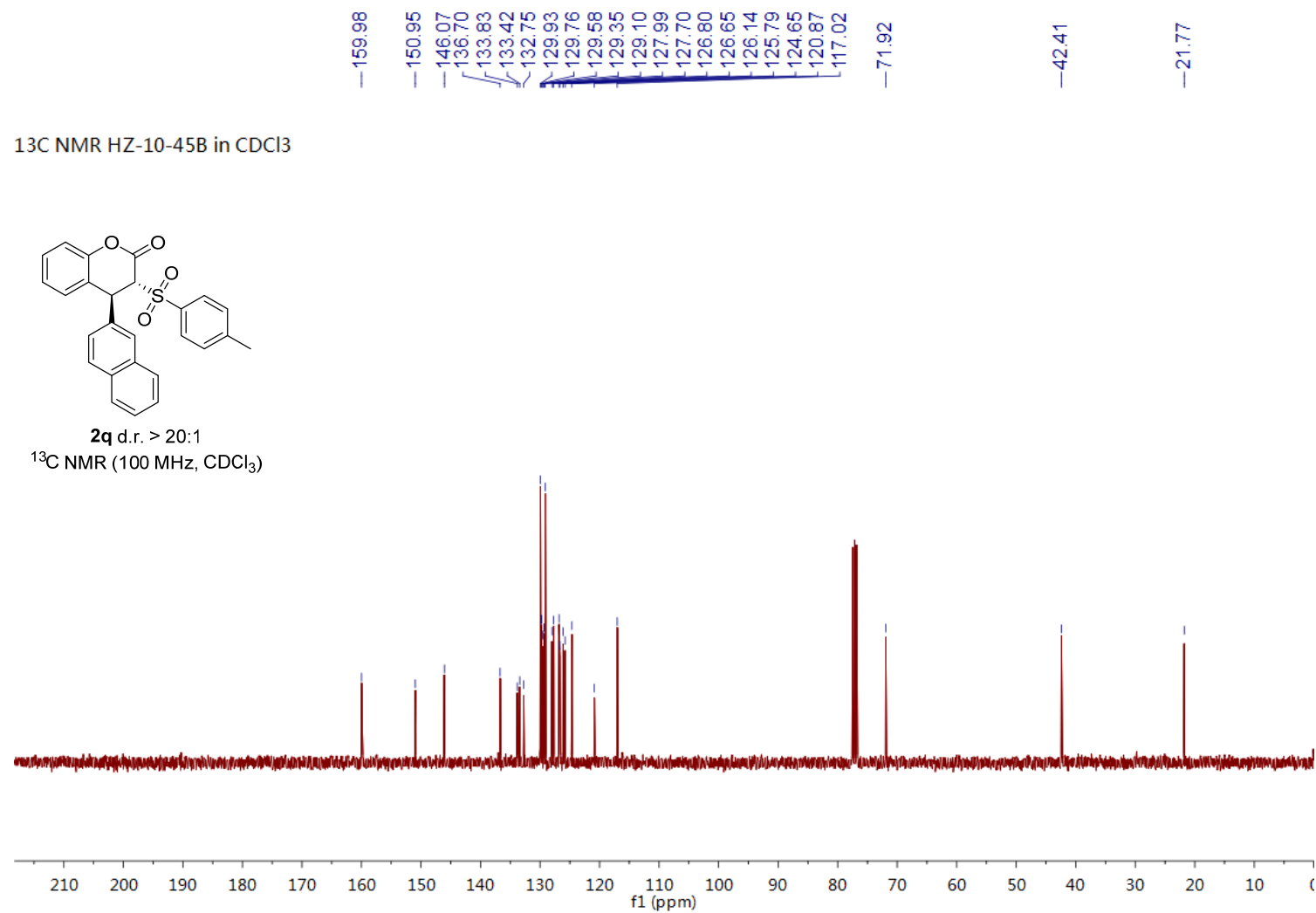


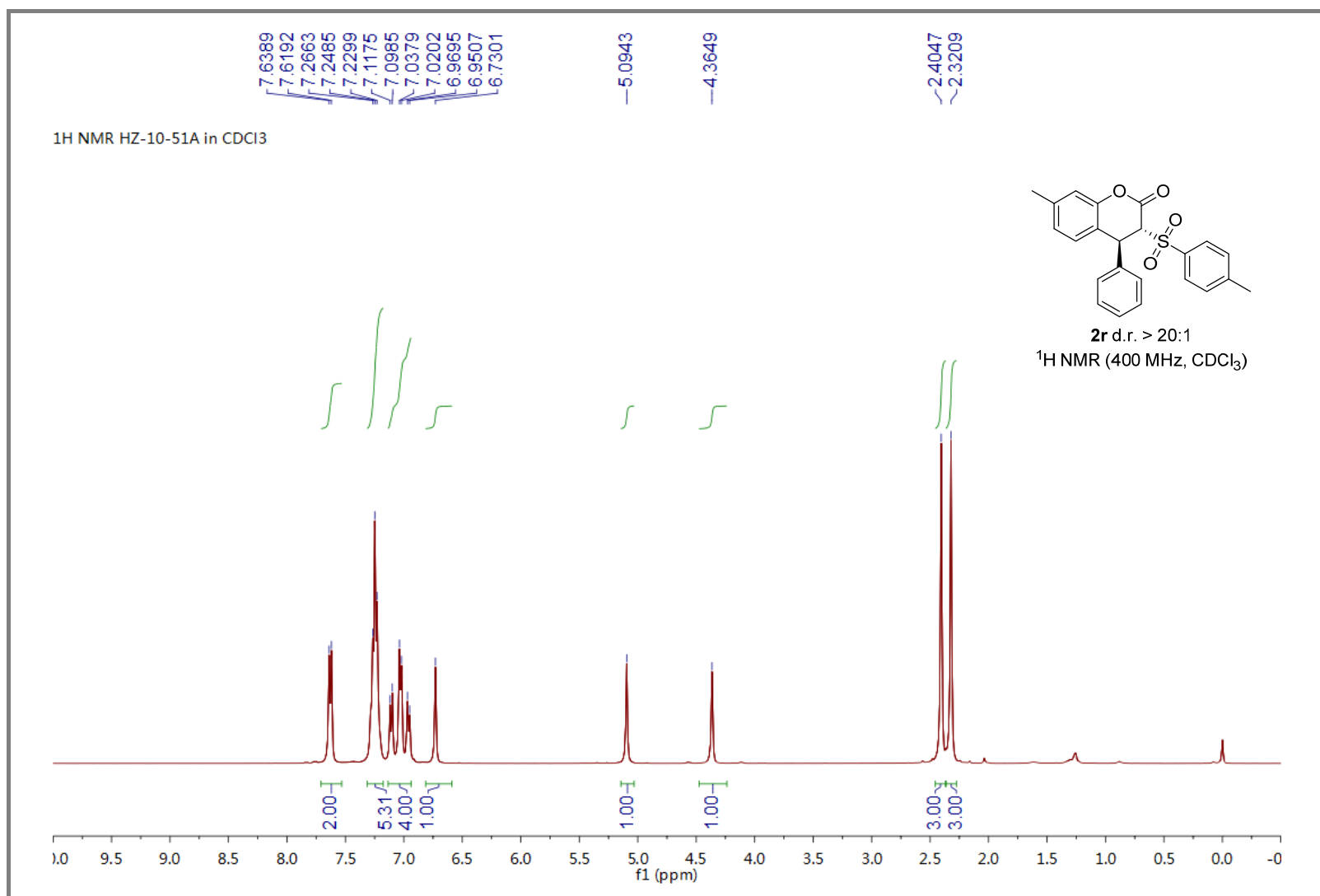
¹³C NMR HZ-10-45B in CDCl₃



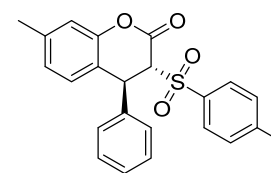
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¹³C NMR (100 MHz, CDCl₃)

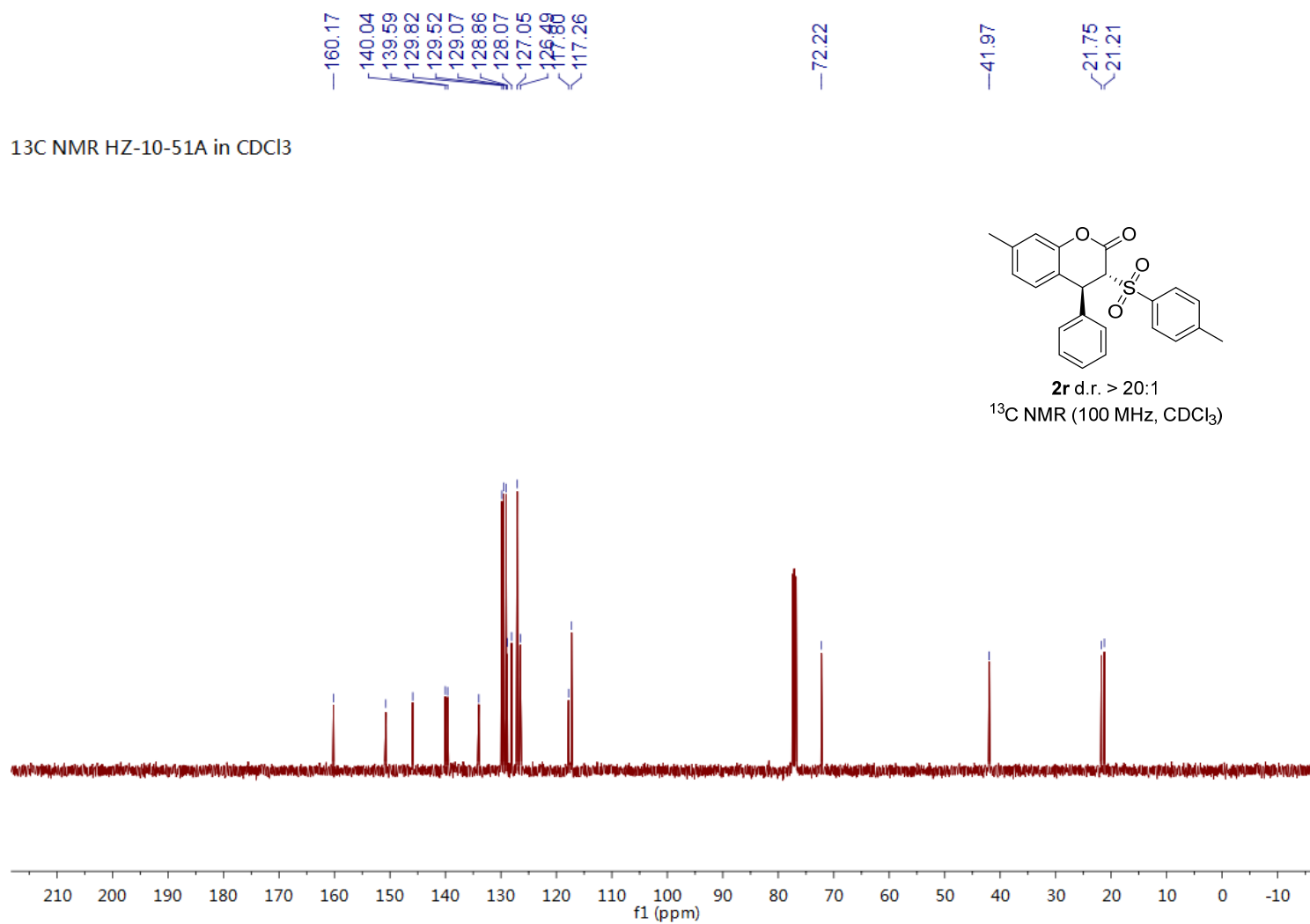


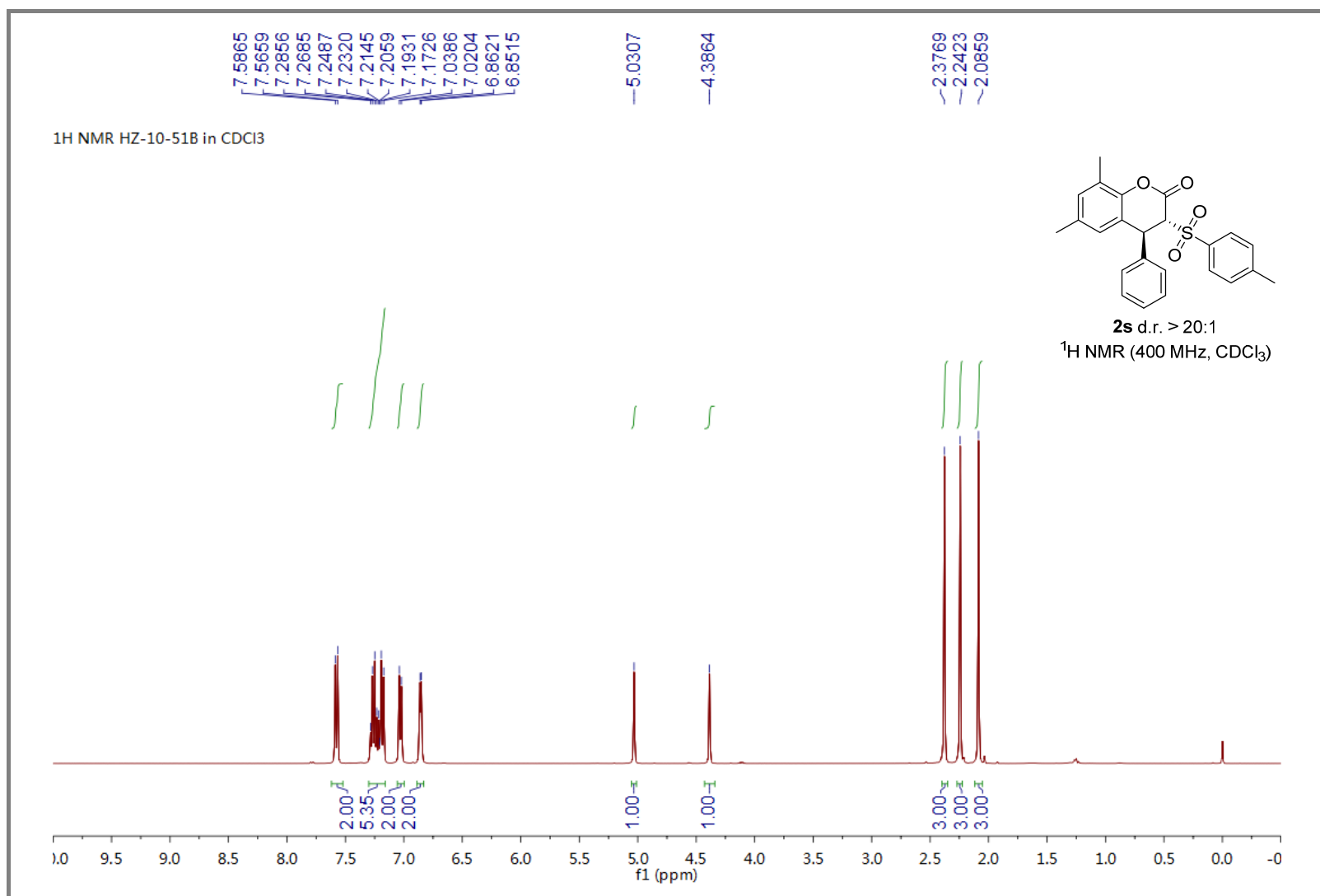


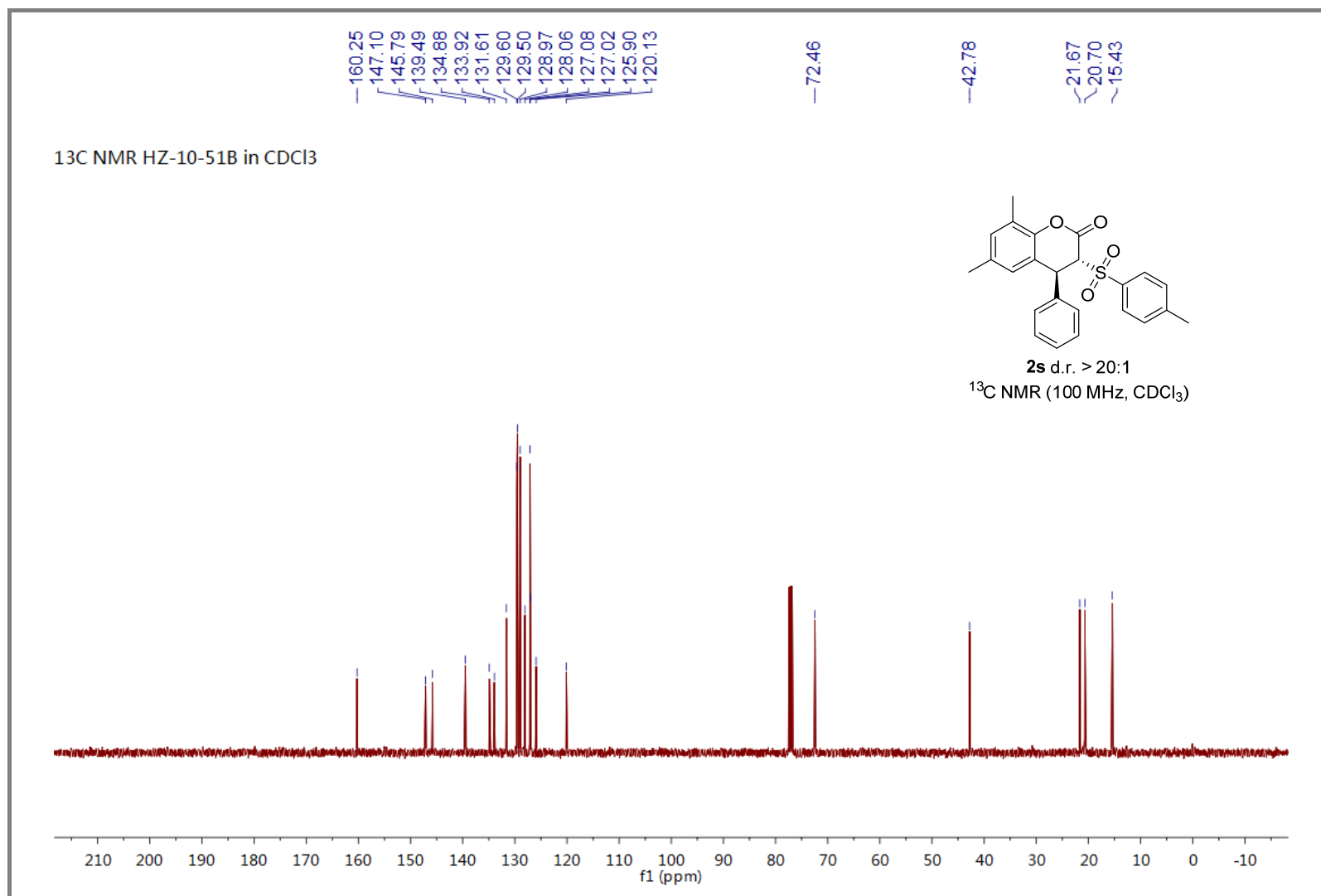
¹³C NMR HZ-10-51A in CDCl₃

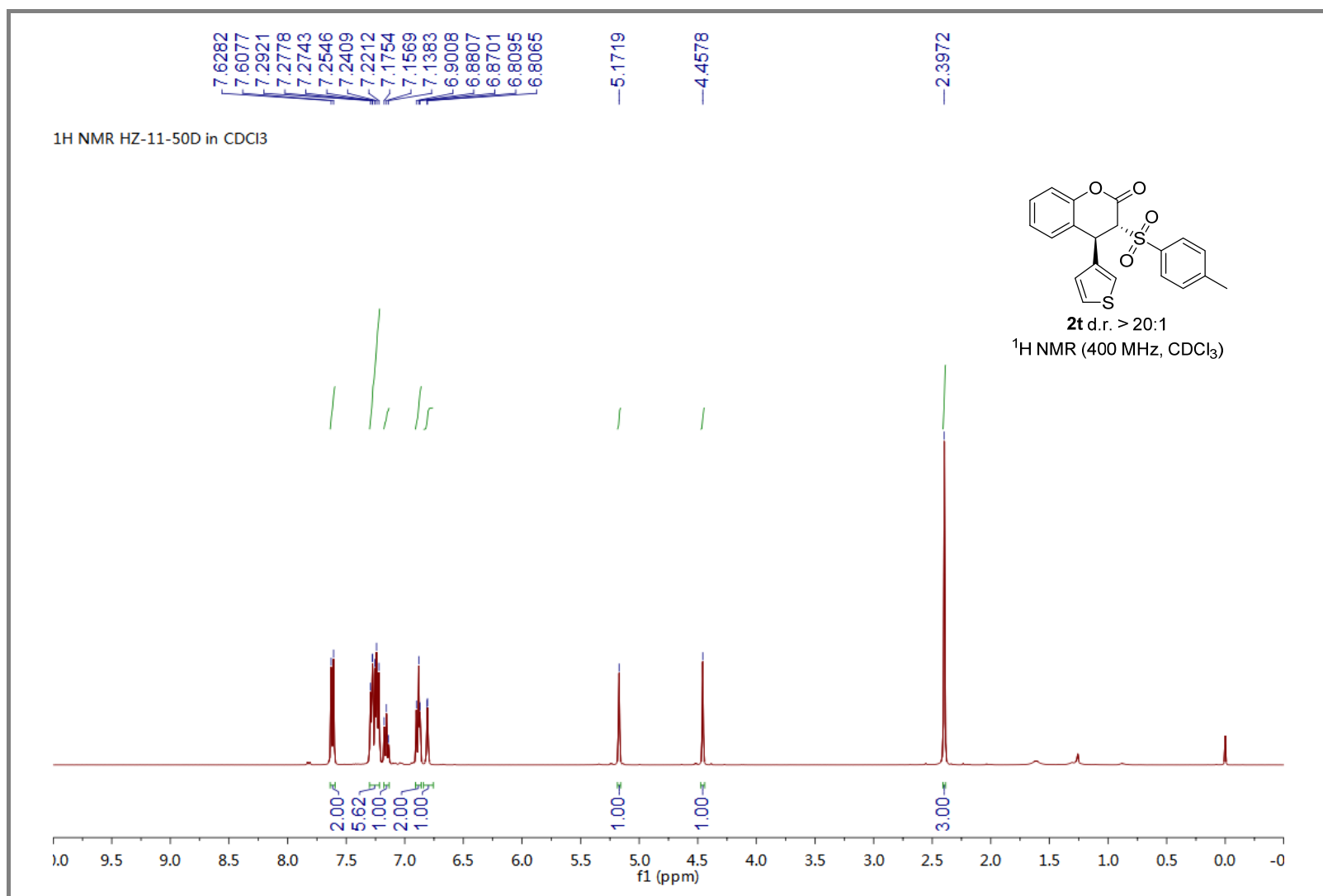


2r d.r. > 20:1
¹³C NMR (100 MHz, CDCl₃)

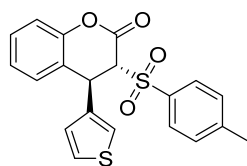




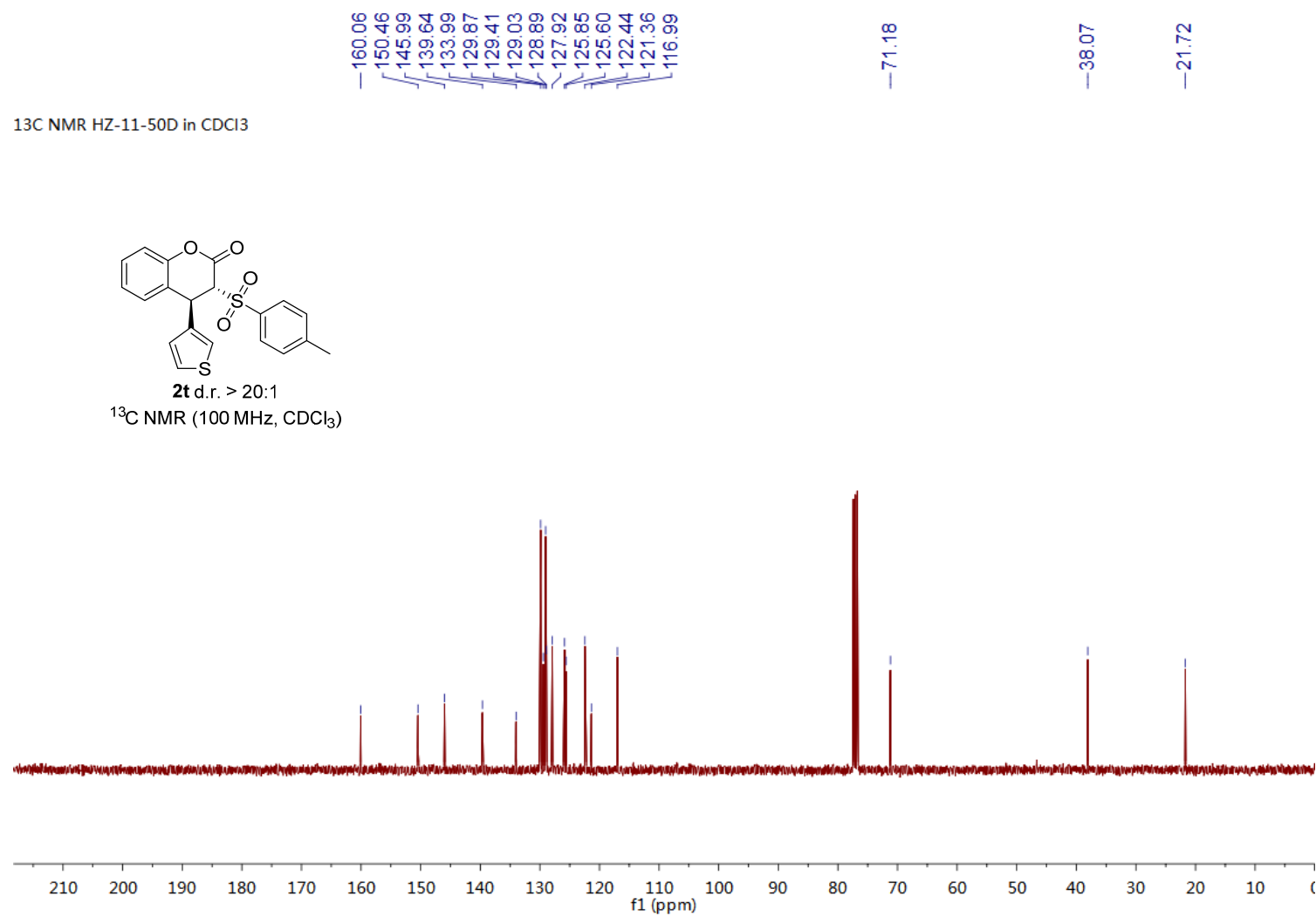


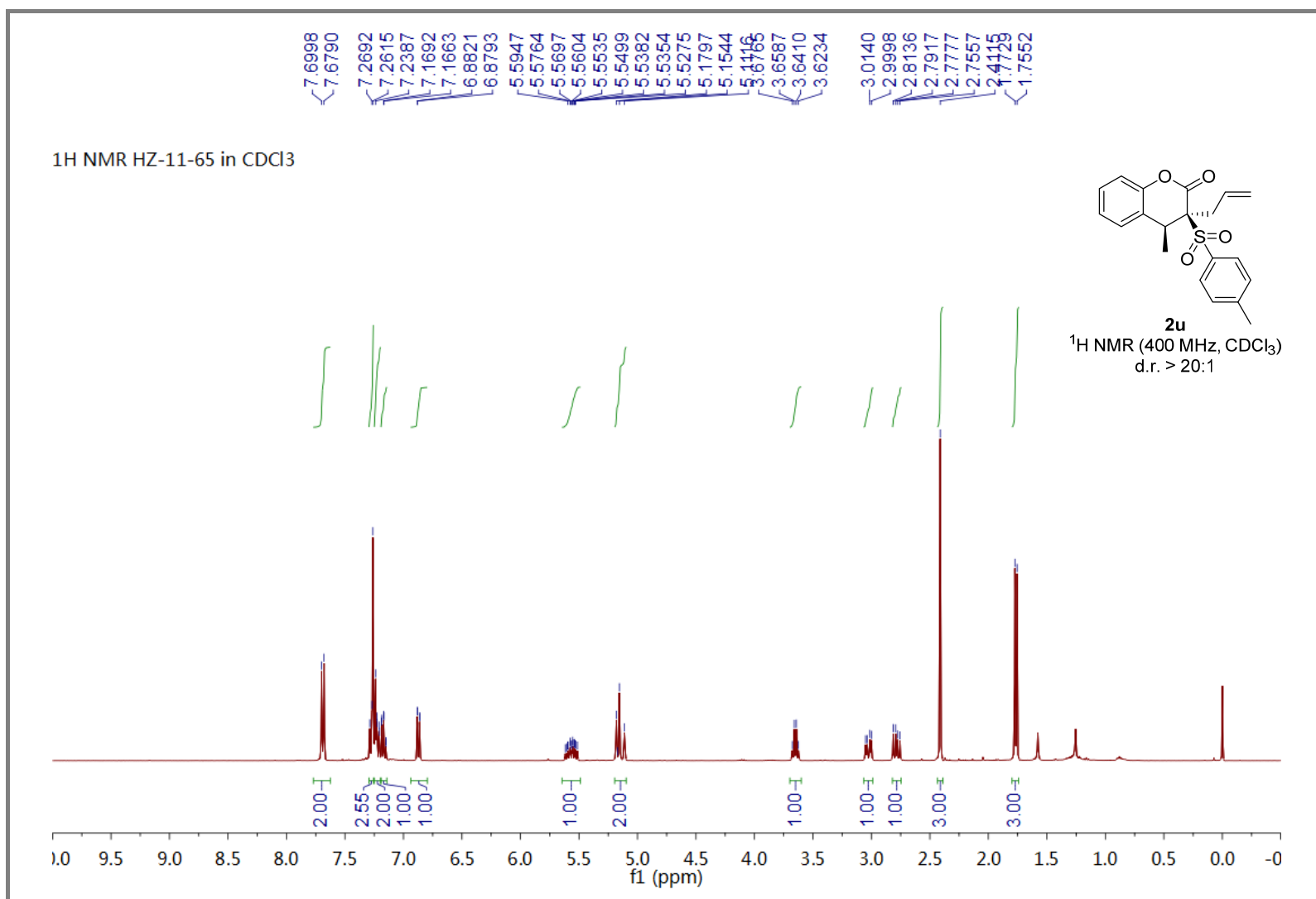


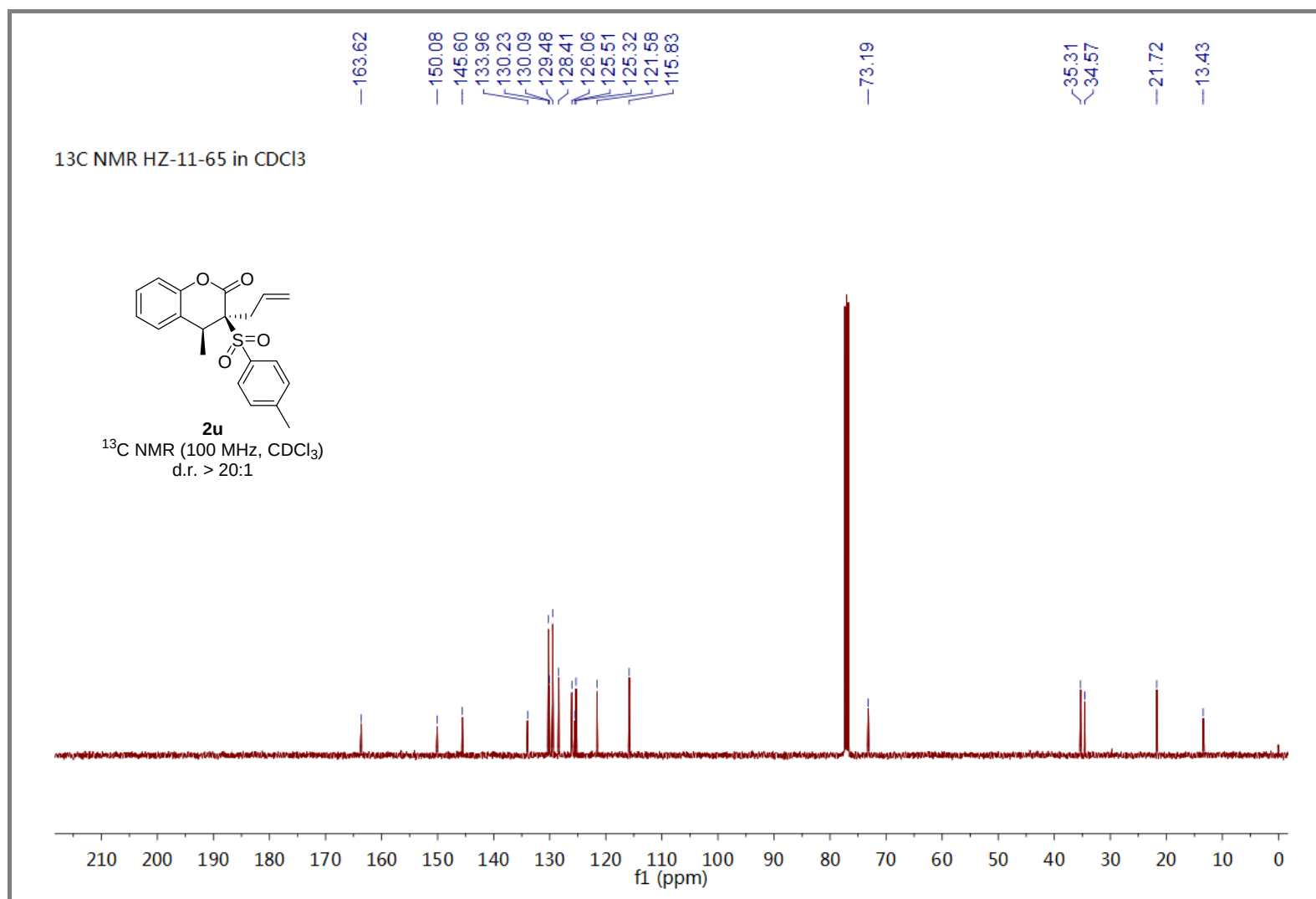
¹³C NMR HZ-11-50D in CDCl₃

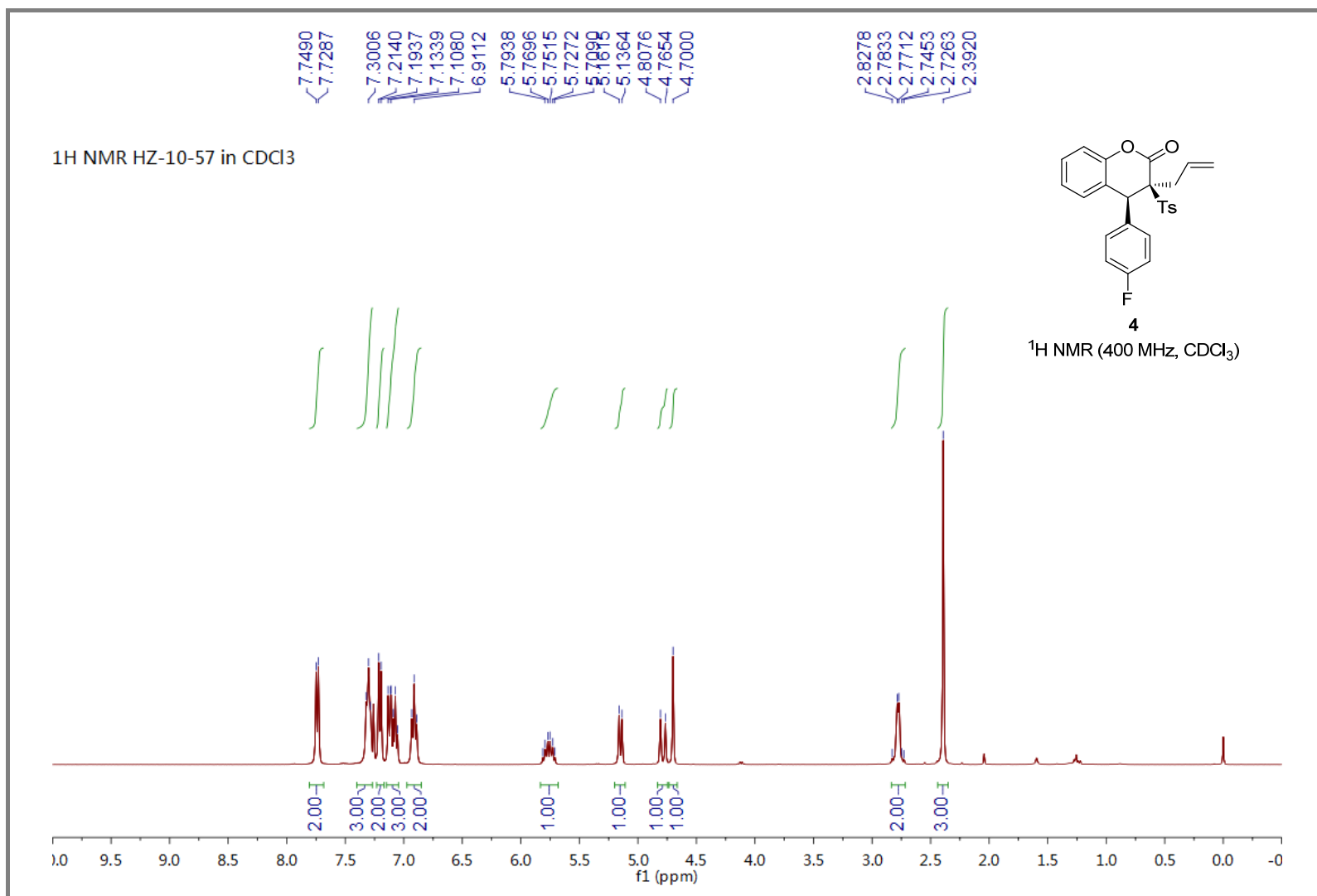


2t d.r. > 20:1
¹³C NMR (100 MHz, CDCl₃)

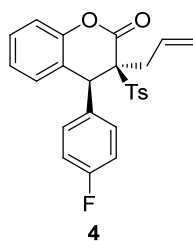




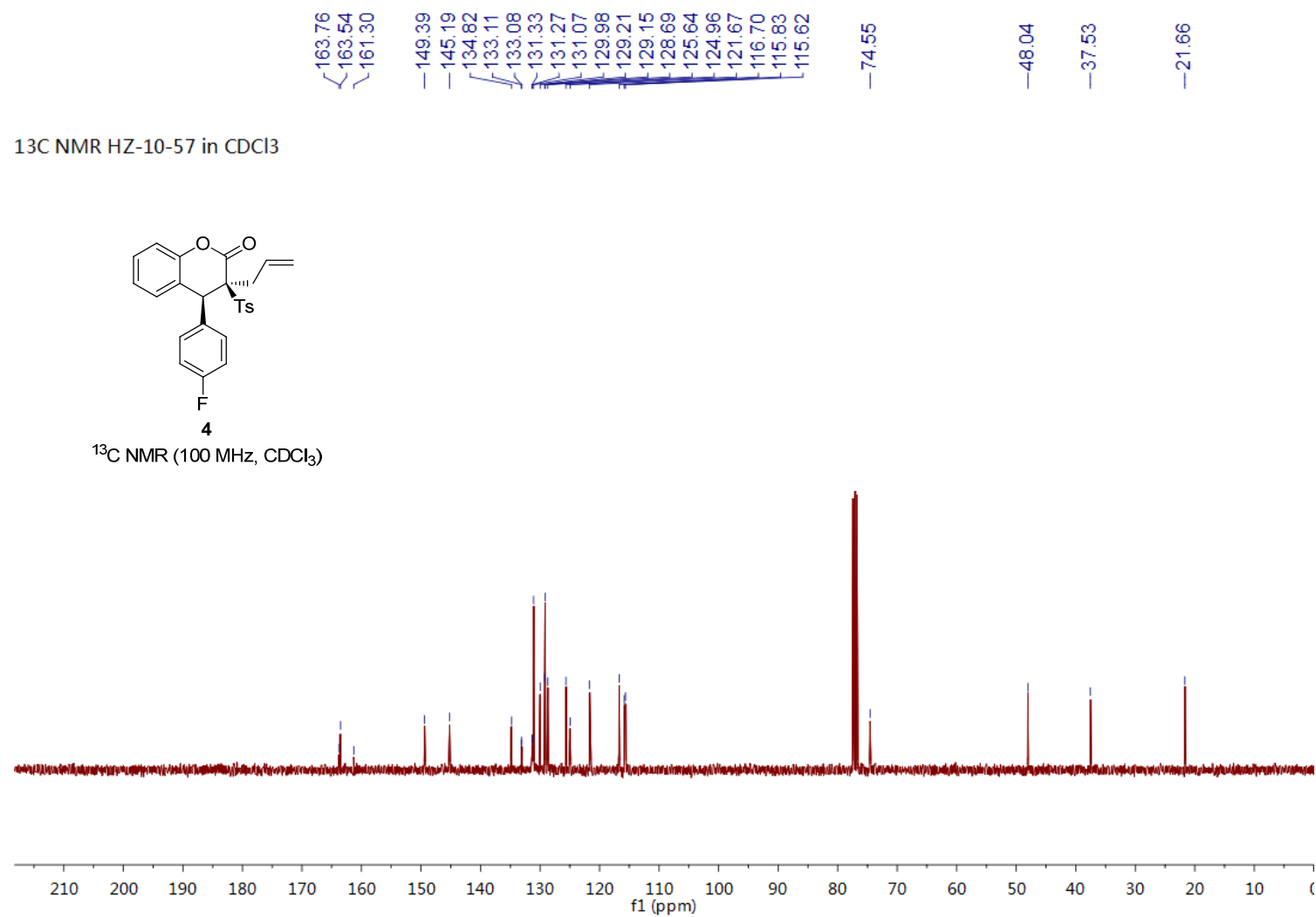




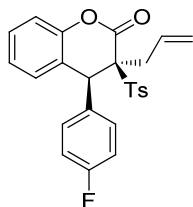
¹³C NMR HZ-10-57 in CDCl₃



¹³C NMR (100 MHz, CDCl₃)

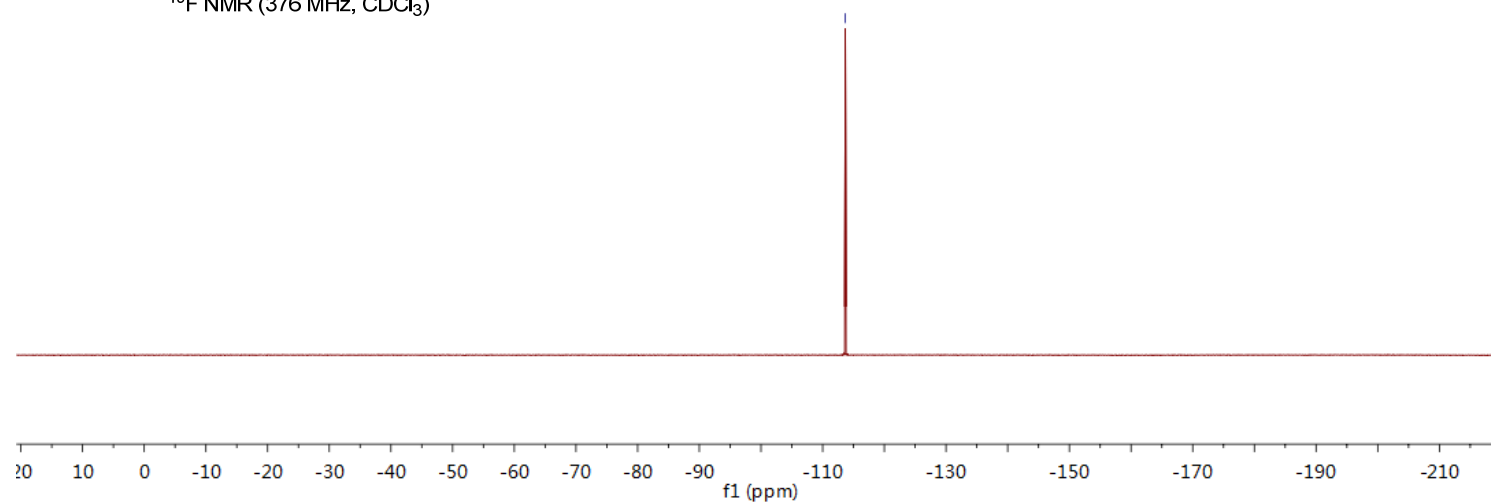


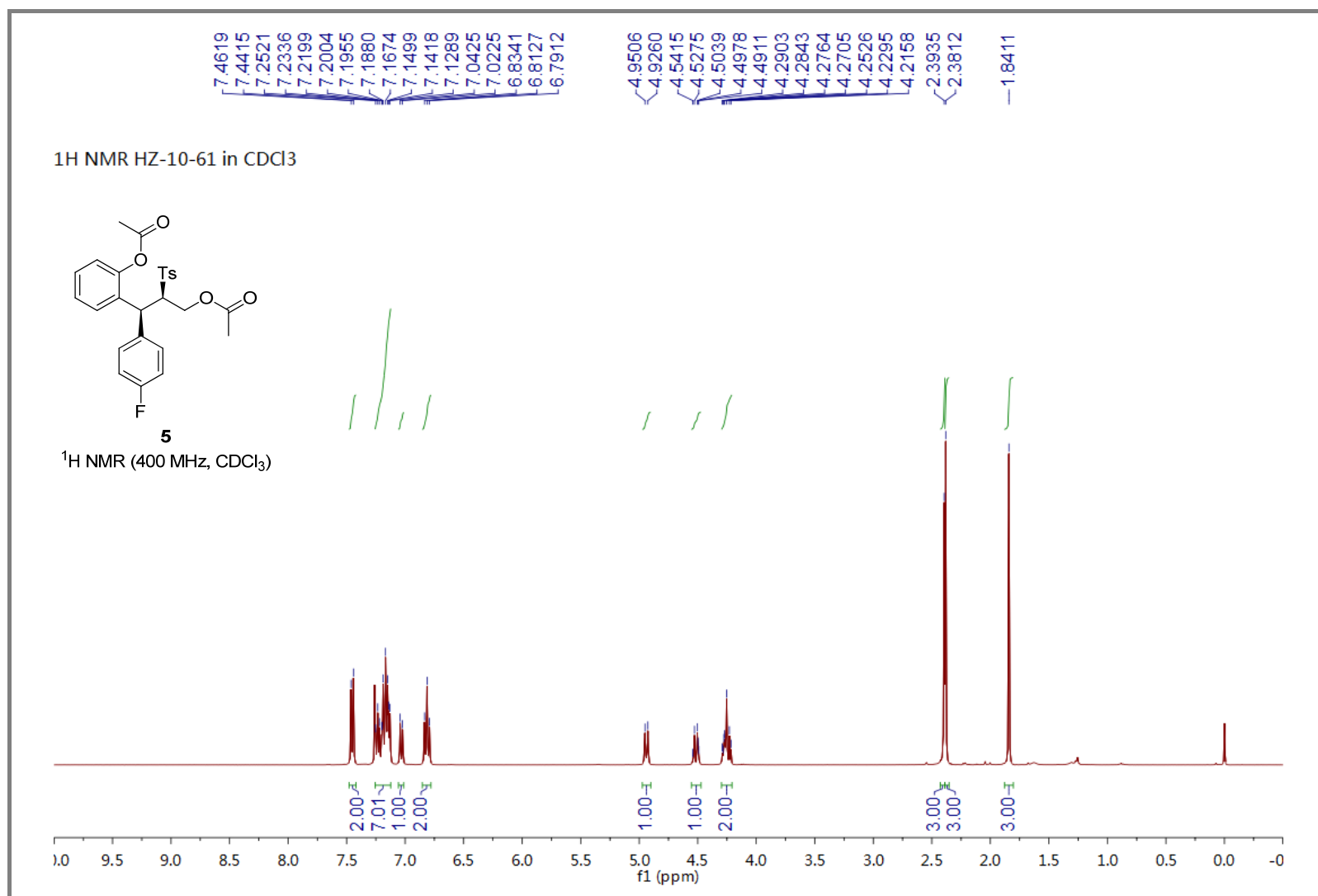
¹⁹F NMR HZ-10-57 in CDCl₃



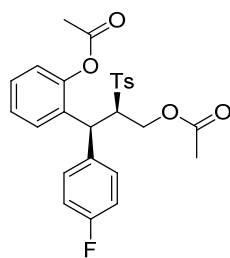
4

¹⁹F NMR (376 MHz, CDCl₃)



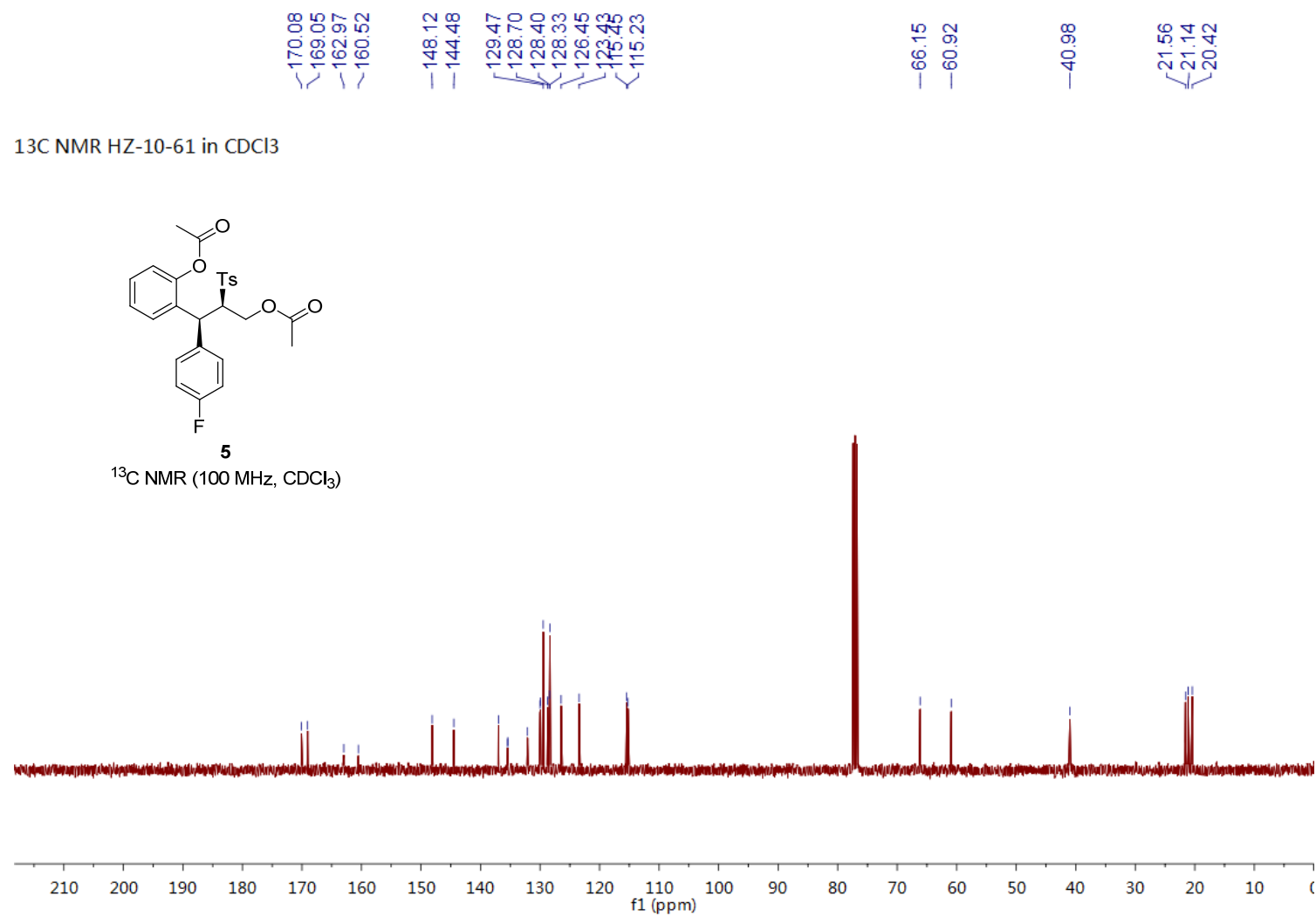


¹³C NMR HZ-10-61 in CDCl₃

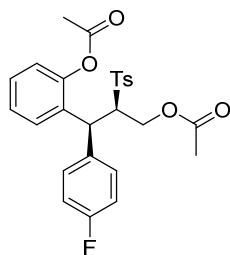


5

¹³C NMR (100 MHz, CDCl₃)

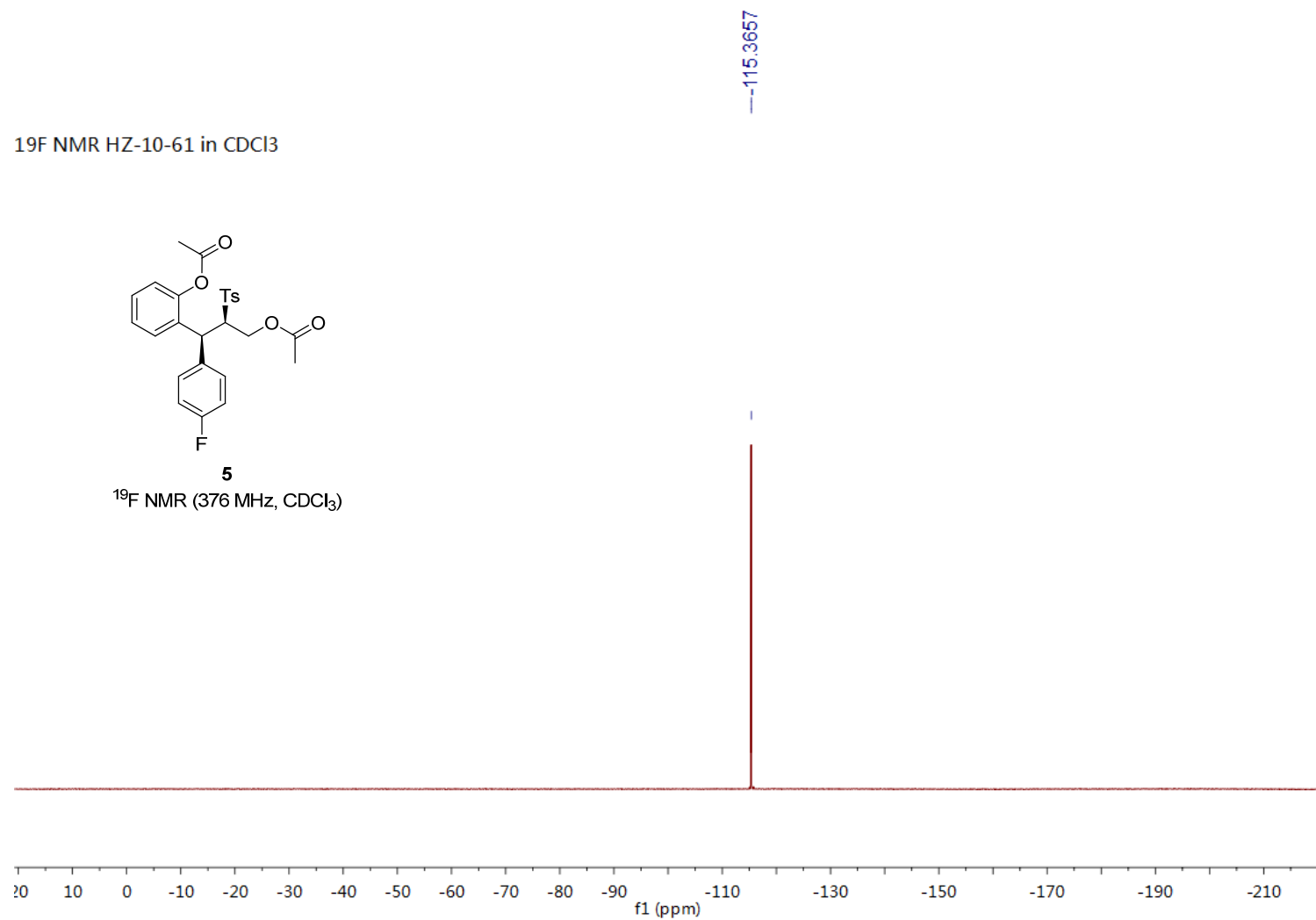


¹⁹F NMR HZ-10-61 in CDCl₃



5

¹⁹F NMR (376 MHz, CDCl₃)



Data File C:\CHEM32\1\DATA\ZHOU2019\SIG1006178.D
Sample Name: HZ-10-15B+-

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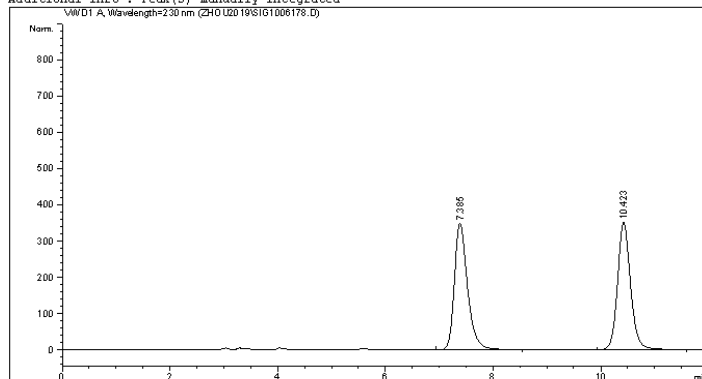
Acq. Operator :
Acq. Instrument : 仪器 1 Location : Vial 1
Injection Date : 12/25/20 20:03:41 Inj Volume : 5.000 µl

Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/25/20 19:56:53
(modified after loading)

Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/25/20 20:45:15
(modified after loading)

Sample Info : IA, n-hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

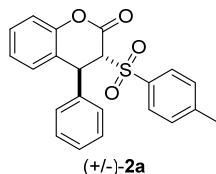


Area Percent Report

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

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.385	BB	0.2555	5882.63135	349.26910	49.9647
2	10.423	BB	0.2544	5890.93848	351.73358	50.0353



仪器 1 12/25/20 20:45:18

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Data File C:\CHEM32\1\DATA\ZHOU2019\SIG1006177.D
Sample Name: HZ-10-15A

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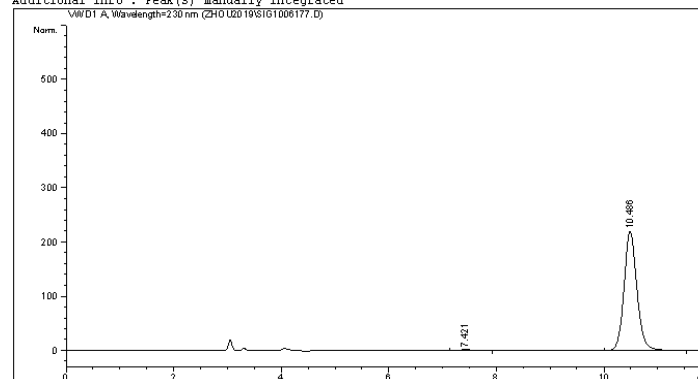
Acq. Operator :
Acq. Instrument : 仪器 1 Location : Vial 1
Injection Date : 12/25/20 19:42:47 Inj Volume : 5.000 µl

Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/25/20 19:17:08
(modified after loading)

Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/25/20 20:47:41
(modified after loading)

Sample Info : IA, n-hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

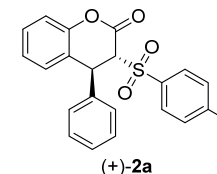


Area Percent Report

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

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.421	BB	0.2578	32.00377	1.86918	0.8641
2	10.486	BB	0.2538	3671.84814	219.95047	99.1359



仪器 1 12/25/20 20:47:44

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Data File C:\CHEM32\1\DATA\ZHOU2019\SIG1006223.D
Sample Name: HZ-10-18A+-

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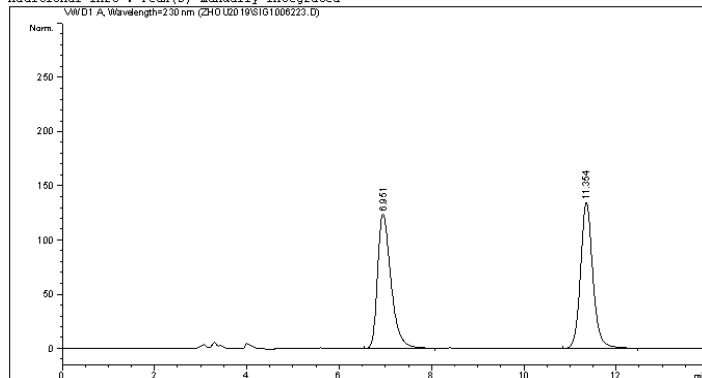
Acq. Operator :
Acq. Instrument : 仪器 1 Location : Vial 1
Injection Date : 12/29/20 22:04:14 Inj Volume : 5.000 µl

Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/29/20 22:03:16
(modified after loading)

Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/29/20 22:23:04
(modified after loading)

Sample Info : IA, n-hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

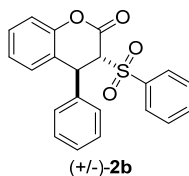


Area Percent Report

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

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.951	BB	0.3062	2492.81250	124.18739	49.9379
2	11.354	BB	0.2813	2499.00806	134.76805	50.0621



仪器 1 12/29/20 22:23:09

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Data File C:\CHEM32\1\DATA\ZHOU2019\SIG1006222.D
Sample Name: HZ-10-17A

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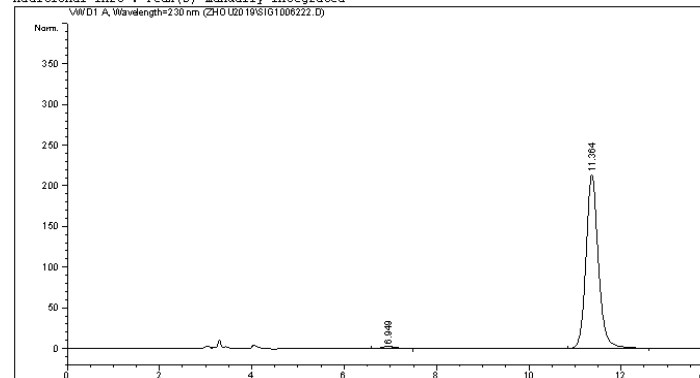
Acq. Operator :
Acq. Instrument : 仪器 1 Location : Vial 1
Injection Date : 12/29/20 21:44:49 Inj Volume : 5.000 µl

Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/29/20 21:42:31
(modified after loading)

Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/29/20 22:06:03
(modified after loading)

Sample Info : IA, n-hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

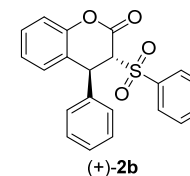


Area Percent Report

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

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.949	BB	0.3208	47.95851	2.30435	1.1954
2	11.364	BB	0.2805	3963.91797	213.60262	98.8046



仪器 1 12/29/20 22:06:07

Page 1 of 2

Data File C:\CHEM32\1\DATA\ZHOU2019\SIG1006225.D
Sample Name: HZ-10-18B+-

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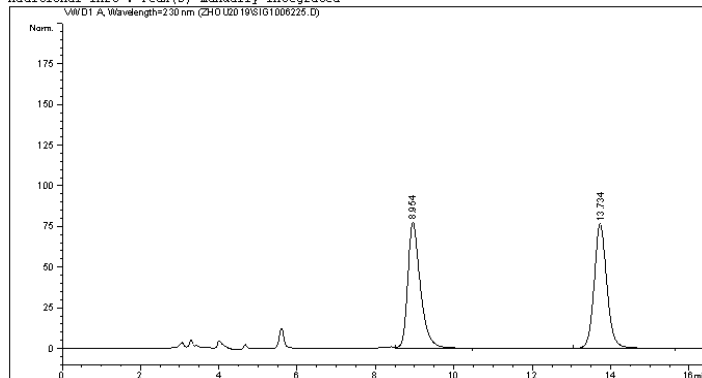
Acq. Operator :
Acq. Instrument : 仪器 1 Location : Vial 1
Injection Date : 12/29/20 22:41:28 Inj Volume : 5.000 µl

Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/29/20 22:40:12
(modified after loading)

Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/29/20 22:58:57
(modified after loading)

Sample Info : IA, n-hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

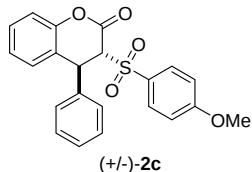


Area Percent Report

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

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.954	VB	0.3343	1729.60388	77.51499	50.1036
2	13.734	BB	0.3400	1722.45288	76.69793	49.8964



仪器 1 12/29/20 22:59:03

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Data File C:\CHEM32\1\DATA\ZHOU2019\SIG1006224.D
Sample Name: HZ-10-17B

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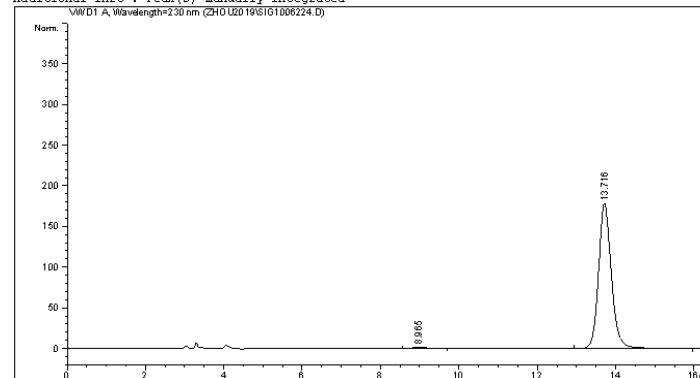
Acq. Operator :
Acq. Instrument : 仪器 1 Location : Vial 1
Injection Date : 12/29/20 22:22:37 Inj Volume : 5.000 µl

Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/29/20 22:20:48
(modified after loading)

Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/29/20 22:44:27
(modified after loading)

Sample Info : IA, n-hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

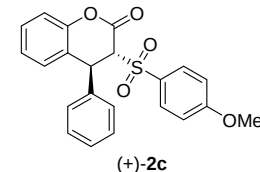


Area Percent Report

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

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.965	BB	0.3329	34.26607	1.55608	0.8440
2	13.716	BB	0.3411	4025.63892	178.51363	99.1560



仪器 1 12/29/20 22:44:31

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Data File C:\CHEM32\1\DATA\ZHOU2019\SIG1006227.D
Sample Name: HZ-10-18C+-

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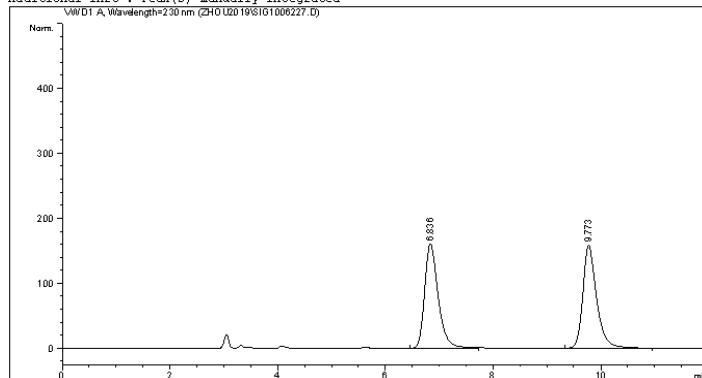
Acq. Operator :
Acq. Instrument : 仪器 1 Location : Vial 1
Injection Date : 12/29/20 23:16:15 Inj Volume : 5.000 µl

Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/29/20 23:15:21
(modified after loading)

Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/29/20 23:17:59
(modified after loading)

Sample Info : IA, n-hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

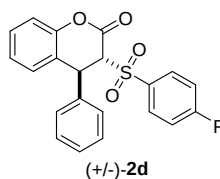


Area Percent Report

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

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.836	BB	0.2653	2793.39331	160.29990	49.8917
2	9.773	BB	0.2635	2805.52563	158.52328	50.1083



仪器 1 12/29/20 23:29:30

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Data File C:\CHEM32\1\DATA\ZHOU2019\SIG1006226.D
Sample Name: HZ-10-17C

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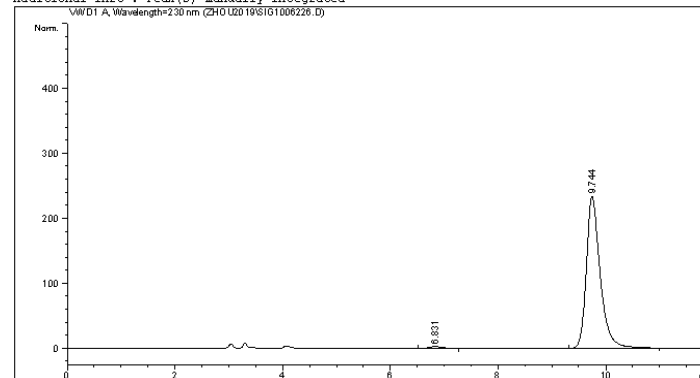
Acq. Operator :
Acq. Instrument : 仪器 1 Location : Vial 1
Injection Date : 12/29/20 23:00:53 Inj Volume : 5.000 µl

Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/29/20 22:58:13
(modified after loading)

Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/29/20 23:17:59
(modified after loading)

Sample Info : IA, n-hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

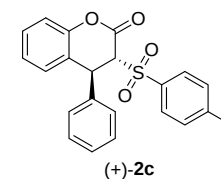


Area Percent Report

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

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.831	BB	0.2616	40.42218	2.37487	0.9646
2	9.744	BB	0.2643	4150.32227	233.60210	99.0354



仪器 1 12/29/20 23:18:02

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Data File C:\CHEM32\1\DATA\ZHOU2019\SIG1006246.D
Sample Name: HZ-10-21B+-

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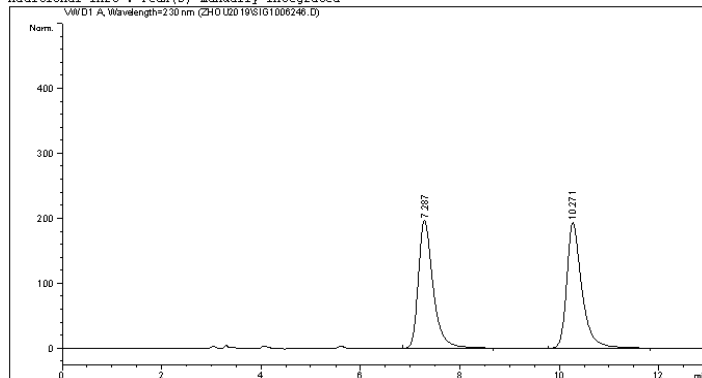
Acq. Operator :
Acq. Instrument : 仪器 1 Location : Vial 1
Injection Date : 12/30/20 20:20:06 Inj Volume : 5.000 µl

Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/30/20 20:16:27
(modified after loading)

Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/30/20 20:37:24
(modified after loading)

Sample Info : IA, n-hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

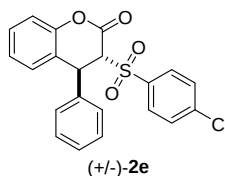


Area Percent Report

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

Signal 1: WVD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.287	BB	0.3106	4033.83936	196.38701	49.7976
2	10.271	BB	0.3128	4066.63696	193.04802	50.2024



仪器 1 12/30/20 20:37:26

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Data File C:\CHEM32\1\DATA\ZHOU2019\SIG1006245.D
Sample Name: HZ-10-20B

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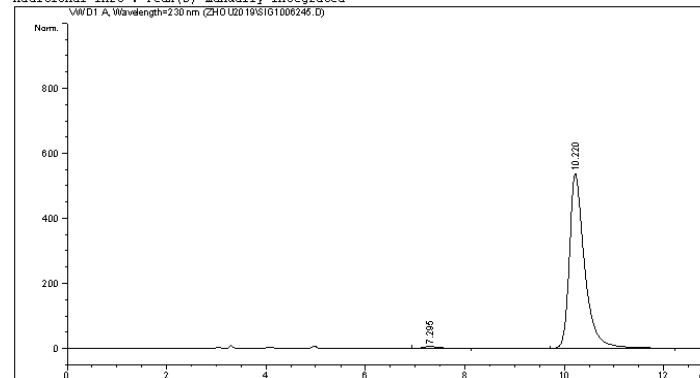
Acq. Operator :
Acq. Instrument : 仪器 1 Location : Vial 1
Injection Date : 12/30/20 20:00:59 Inj Volume : 5.000 µl

Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/30/20 19:59:11
(modified after loading)

Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/30/20 20:26:50
(modified after loading)

Sample Info : IA, n-hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

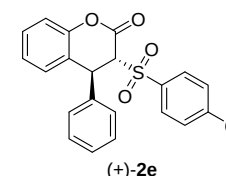


Area Percent Report

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

Signal 1: WVD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.295	BB	0.3394	123.07564	5.32936	1.0742
2	10.220	BB	0.3114	1.13343e4	536.60602	98.9258



仪器 1 12/30/20 20:26:53

Page 1 of 2

Data File C:\CHEM32\1\DATA\ZHOU2019\SIG1006248.D
Sample Name: HZ-10-21C+-

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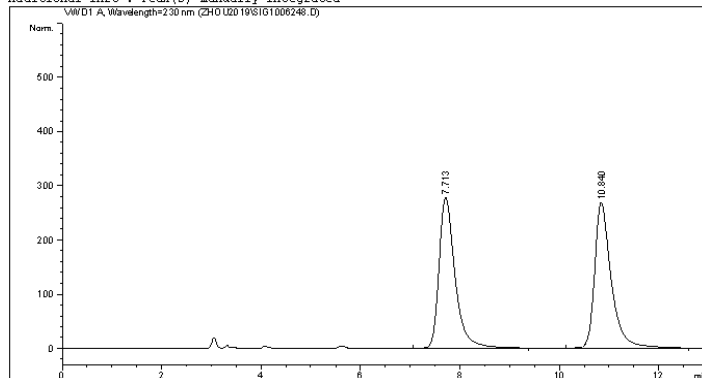
Acq. Operator :
Acq. Instrument : 仪器 1 Location : Vial 1
Injection Date : 12/30/20 20:58:43 Inj Volume : 5.000 µl

Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/30/20 20:57:21
(modified after loading)

Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/30/20 21:13:55
(modified after loading)

Sample Info : IA, n-hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

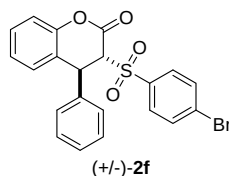


Area Percent Report

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

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.713	BB	0.3360	6209.98340	278.64737	50.0357
2	10.840	BB	0.3414	6201.12549	268.58310	49.9643



仪器 1 12/30/20 21:13:59

Page 1 of 2

Data File C:\CHEM32\1\DATA\ZHOU2019\SIG1006247.D
Sample Name: HZ-10-20C

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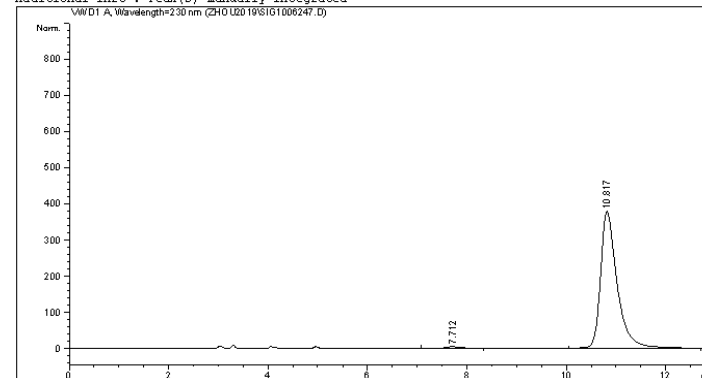
Acq. Operator :
Acq. Instrument : 仪器 1 Location : Vial 1
Injection Date : 12/30/20 20:40:09 Inj Volume : 5.000 µl

Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/30/20 20:36:51
(modified after loading)

Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 12/30/20 21:01:32
(modified after loading)

Sample Info : IA, n-hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

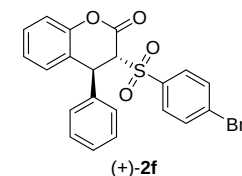


Area Percent Report

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

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.712	BB	0.3526	95.74295	4.05339	1.0826
2	10.817	BB	0.3413	8747.65625	379.03619	98.9174



仪器 1 12/30/20 21:01:36

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Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3002780.D
Sample Name: HZ-11-49C+-

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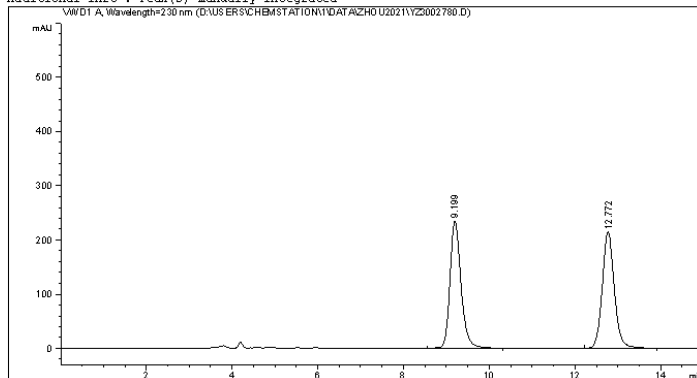
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 7/10/2021 4:05:38 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 7/10/2021 2:44:54 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 7/10/2021 4:26:06 PM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 80/20, 0.8 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

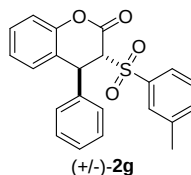


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.199	VB	0.2713	4202.14648	235.38873	50.1100
2	12.772	BB	0.2974	4183.70264	214.61815	49.8900



Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3002781.D
Sample Name: HZ-11-49C

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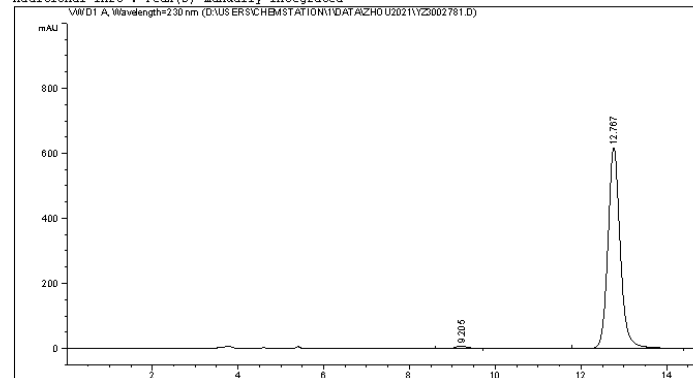
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 7/10/2021 4:25:38 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 7/10/2021 2:44:54 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 7/10/2021 4:44:46 PM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 80/20, 0.8 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

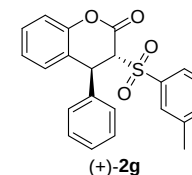


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.205	BB	0.2738	124.24998	6.84344	1.0057
2	12.767	VB R	0.3013	1.22303e4	616.87054	98.9943



Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3002779.D
Sample Name: HZ-11-49B+-

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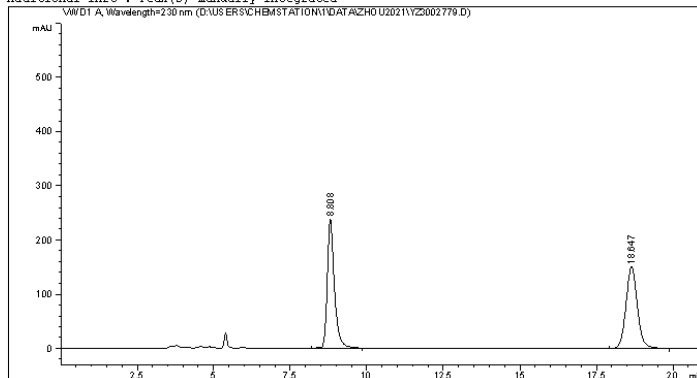
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 7/10/2021 3:41:22 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 7/10/2021 2:44:54 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 7/10/2021 4:04:02 PM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 80/20, 0.8 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

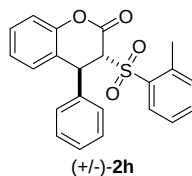


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.808	VB R	0.2371	3811.44092	237.67677	49.7203
2	18.647	BB	0.3927	3854.32495	150.54021	50.2797



1260II 7/10/2021 4:04:08 PM SYSTEM

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Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3002778.D
Sample Name: HZ-11-49B

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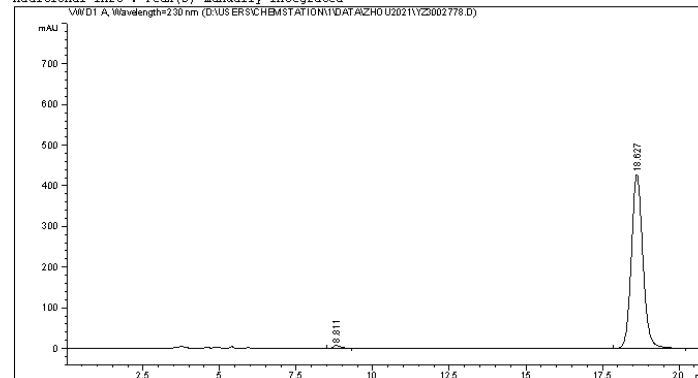
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 7/10/2021 3:17:09 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 7/10/2021 2:44:54 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 7/10/2021 3:44:38 PM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 80/20, 0.8 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

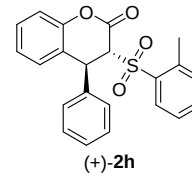


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.811	VB	0.2342	110.96770	7.02949	0.9892
2	18.627	BB	0.3992	1.11067e4	427.37448	99.0108



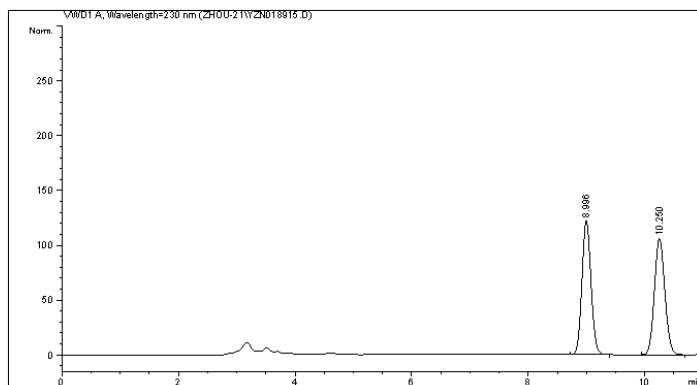
1260II 7/10/2021 3:44:40 PM SYSTEM

Page 1 of 2

Data File C:\CHEM32\1\DATA\ZHOU-21\YZN018915.D
Sample Name: HZ-11-56+-

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : -
Injection Date : 7/15/2021 3:34:41 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/15/2021 3:16:11 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/15/2021 5:05:12 AM
(modified after loading)
Sample Info : AD-H, Hexane/i-PrOH = 80/20, 1.0 mL/min, 30 oC, 230 nm



Area Percent Report

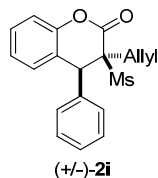
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area [mAU]	Area %
1	8.996	BB	0.1669	1323.21594	122.38219	50.0224	
2	10.250	BB	0.1936	1322.02893	105.98039	49.9776	

Totals : 2645.24487 228.36259

*** End of Report ***



(+/-)-2i

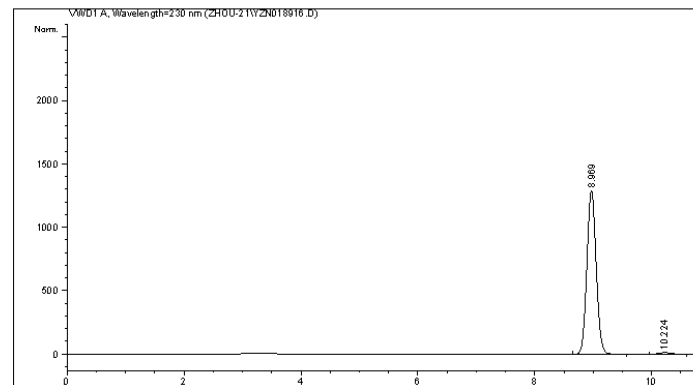
Instrument 1 7/15/2021 5:05:14 AM

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-21\YZN018916.D
Sample Name: HZ-11-56

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : -
Injection Date : 7/15/2021 4:01:58 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/15/2021 4:01:01 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 7/15/2021 5:04:47 AM
(modified after loading)
Sample Info : AD-H, Hexane/i-PrOH = 80/20, 1.0 mL/min, 30 oC, 230 nm



Area Percent Report

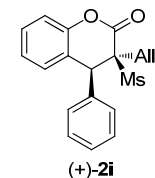
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area [mAU]	Area %
1	8.969	BB	0.1705	1.41857e4	1290.67273	98.9766	
2	10.224	BB	0.1945	146.67892	11.80422	1.0234	

Totals : 1.43324e4 1302.47695

*** End of Report ***



(+)-2i

Instrument 1 7/15/2021 5:04:53 AM

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU2021\SIG1006473.D
Sample Name: HZ-10-40A+-

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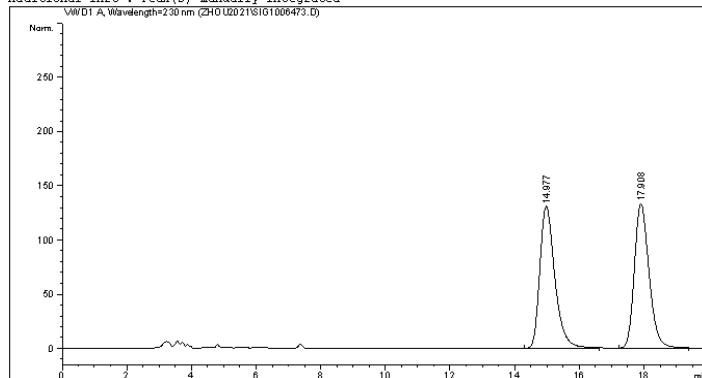
Acq. Operator :
Acq. Instrument : 仪器 1 Location : Vial 1
Injection Date : 1/20/21 17:11:39 Inj Volume : 5.000 µl

Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 1/20/21 17:09:32
(modified after loading)

Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 1/20/21 17:33:49
(modified after loading)

Sample Info : ID, n-hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

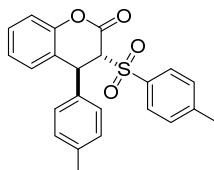


Area Percent Report

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

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	14.977	BB	0.4917	4211.36475	130.79126	50.1334
2	17.908	BB	0.4853	4188.95752	133.05103	49.8666



(+/-)-2j

仪器 1 1/20/21 17:33:53

Page 1 of 2

Data File C:\CHEM32\1\DATA\ZHOU2021\SIG1006471.D
Sample Name: HZ-10-39A

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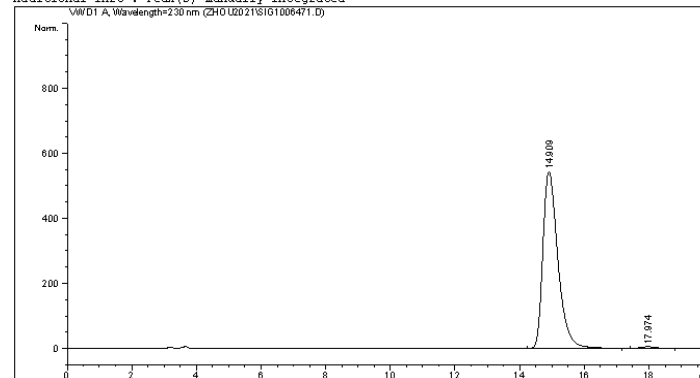
Acq. Operator :
Acq. Instrument : 仪器 1 Location : Vial 1
Injection Date : 1/20/21 16:23:56 Inj Volume : 5.000 µl

Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 1/20/21 16:22:56
(modified after loading)

Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 1/20/21 17:14:13
(modified after loading)

Sample Info : ID, n-hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

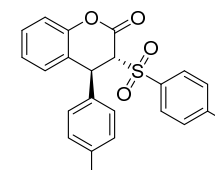


Area Percent Report

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

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	14.909	BB	0.4912	1.74835e4	543.64801	99.1906
2	17.974	BB	0.4807	142.66333	4.60158	0.8094



(+)-2j

仪器 1 1/20/21 17:14:16

Page 1 of 2

Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3001809.D
Sample Name: HZ-10-40B+-

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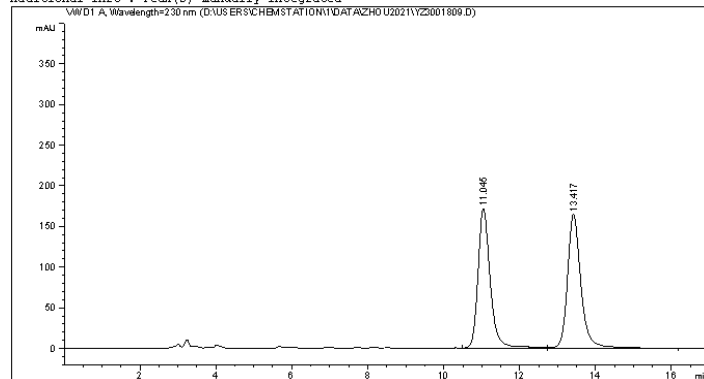
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 1/20/2021 10:29:25 AM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\def1_LC.M
Last changed : 1/20/2021 9:30:44 AM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\def1_LC.M
Last changed : 1/20/2021 11:38:37 AM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

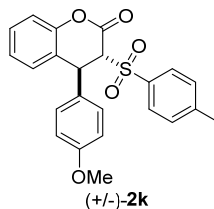


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.045	VV	0.3379	3859.35303	171.94897	48.8985
2	13.417	VB	0.3696	4033.22485	164.74094	51.1015



1260II 1/20/2021 11:38:49 AM SYSTEM

Page 1 of 2

Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3001808.D
Sample Name: HZ-10-39B

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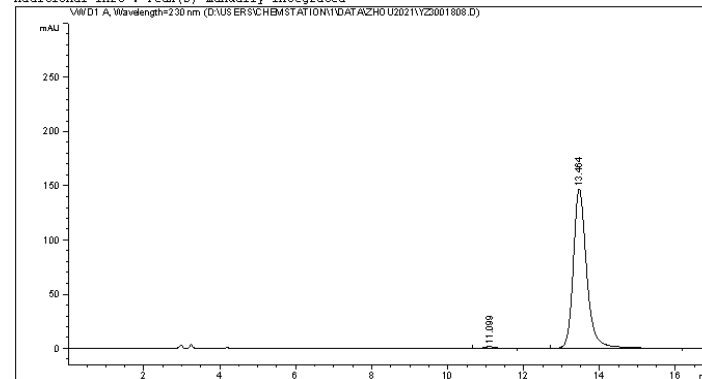
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 1/20/2021 10:06:47 AM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\def1_LC.M
Last changed : 1/20/2021 9:30:44 AM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\def1_LC.M
Last changed : 1/20/2021 11:30:22 AM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

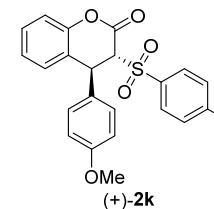


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.099	VB	0.3320	33.14118	1.52216	0.9149
2	13.464	VB	0.3688	3589.38306	147.04446	99.0851



1260II 1/20/2021 11:30:34 AM SYSTEM

Page 1 of 2

Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3001812.D
Sample Name: HZ-10-40C+-

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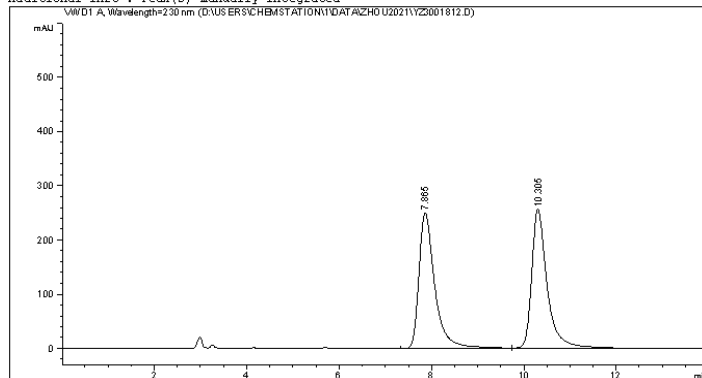
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 1/20/2021 2:46:14 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\def1_LC.M
Last changed : 1/20/2021 2:31:05 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\def1_LC.M
Last changed : 1/20/2021 4:24:19 PM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

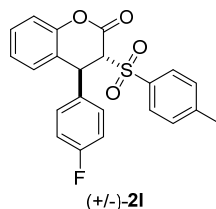


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.865	BB	0.3547	5918.61426	250.46056	49.9198
2	10.305	BB	0.3443	5937.62744	256.29160	50.0802



1260II 1/20/2021 4:24:54 PM SYSTEM

Page 1 of 2

Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3001814.D
Sample Name: HZ-10-39C

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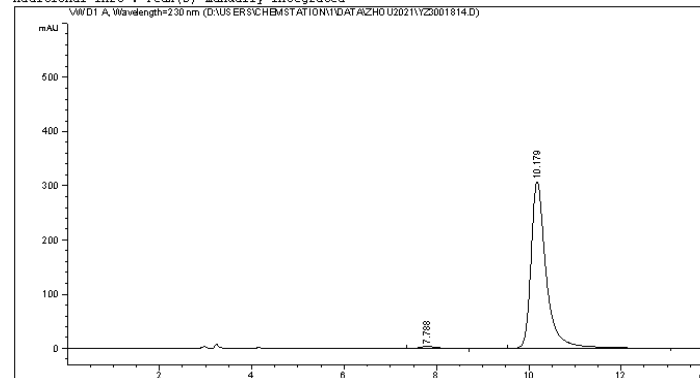
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 1/20/2021 3:35:36 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\def1_LC.M
Last changed : 1/20/2021 2:31:05 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\def1_LC.M
Last changed : 1/20/2021 4:24:19 PM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

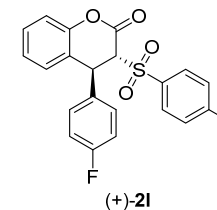


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.788	BB	0.3403	78.16211	3.45025	1.1065
2	10.179	BB	0.3369	6985.46826	307.70099	98.8935



1260II 1/20/2021 4:24:24 PM SYSTEM

Page 1 of 2

Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3001881.D
Sample Name: HZ-10-49A+-

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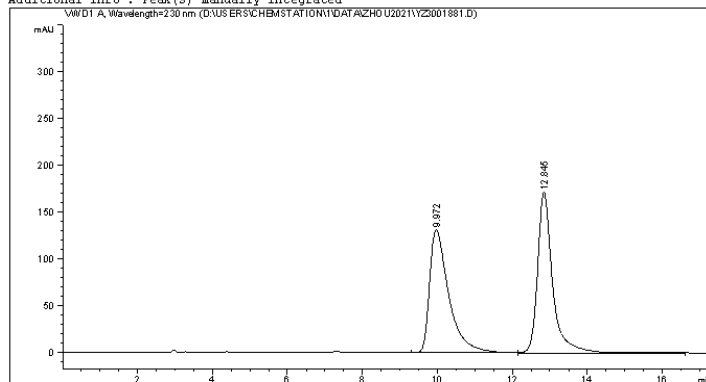
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 3/6/2021 9:17:11 AM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/6/2021 8:34:16 AM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/6/2021 9:37:31 AM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 75/25, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

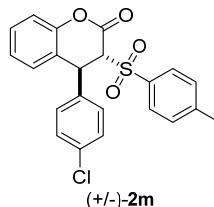


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.972	BV	0.5349	4712.80469	131.24069	49.7285
2	12.845	VB	0.4117	4764.25977	171.70033	50.2715



1260II 3/6/2021 9:37:42 AM SYSTEM

Page 1 of 2

Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3001880.D
Sample Name: HZ-10-48A

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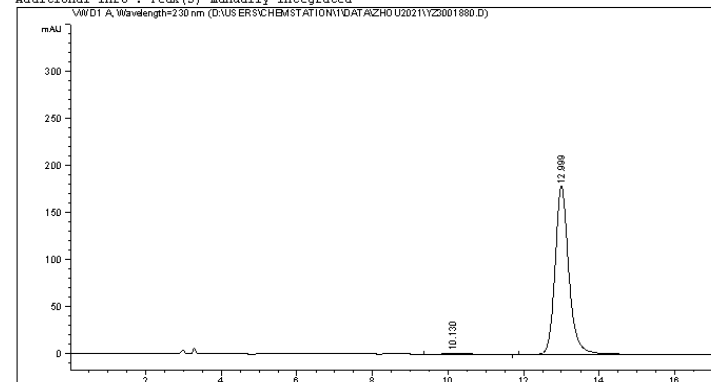
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Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 3/6/2021 8:58:21 AM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/6/2021 8:34:16 AM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/6/2021 9:22:28 AM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 75/25, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

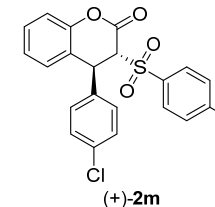


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.130	BB	0.5048	47.87781	1.43659	1.0438
2	12.999	BBA	0.3823	4539.12305	178.74617	98.9562



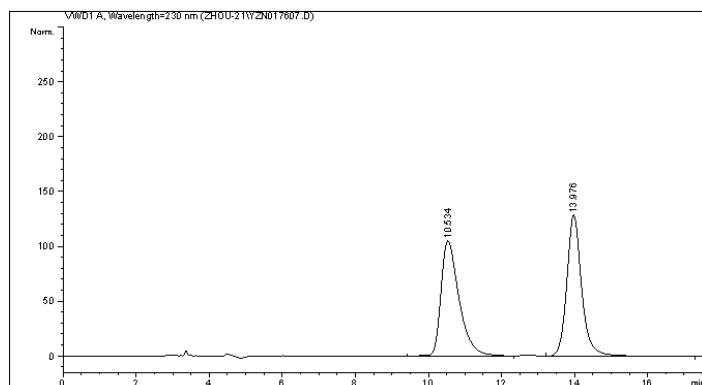
1260II 3/6/2021 9:22:41 AM SYSTEM

Page 1 of 2

Data File C:\CHEM32\1\DATA\ZHOU-21\YZN017607.D
Sample Name: HZ-10-46A+-

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : -
Injection Date : 3/3/2021 8:29:25 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/3/2021 8:22:25 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/3/2021 8:48:57 AM
(modified after loading)
Sample Info : IA, Hexane/i-PrOH = 75/25, 1.0 mL/min, 30 oC, 230 nm



Area Percent Report

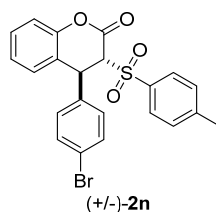
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Area %	Height [mAU]
1	10.534	BB	0.5203	3629.71680	50.0134	105.24075
2	13.976	BB	0.4222	3627.77002	49.9866	128.62926

Totals : 7257.48682 233.87000

*** End of Report ***



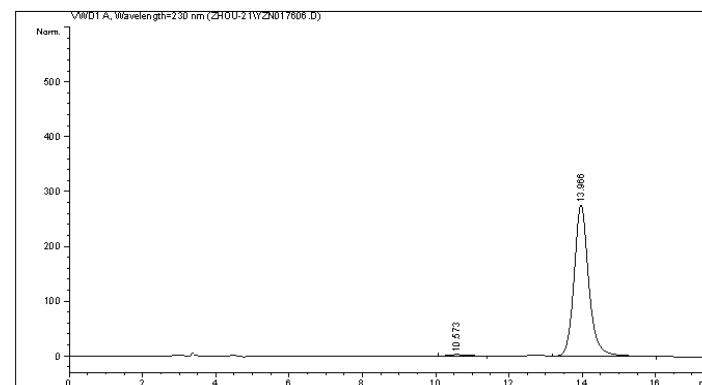
Instrument 1 3/3/2021 8:49:04 AM

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-21\YZN017606.D
Sample Name: HZ-10-45A

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Acq. Operator :
Acq. Instrument : Instrument 1 Location : -
Injection Date : 3/3/2021 8:03:11 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/3/2021 7:09:43 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/3/2021 8:50:30 AM
(modified after loading)
Sample Info : IA, Hexane/i-PrOH = 75/25, 1.0 mL/min, 30 oC, 230 nm



Area Percent Report

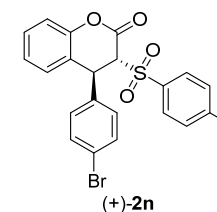
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Area %	Height [mAU]
1	10.573	BB	0.4862	79.73492	1.0263	2.49610
2	13.966	VB	0.4187	7689.28369	98.9737	275.56104

Totals : 7769.01862 278.05713

*** End of Report ***



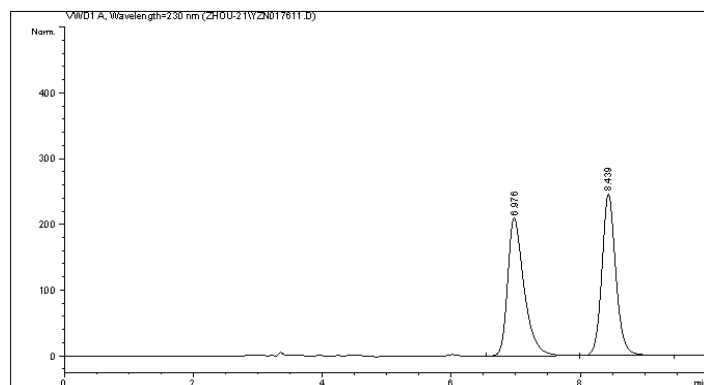
Instrument 1 3/3/2021 8:50:39 AM

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-21\YZN017611.D
Sample Name: HZ-10-46C+-

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : -
Injection Date : 3/3/2021 9:13:49 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/3/2021 8:16:42 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/3/2021 9:58:39 PM
(modified after loading)
Sample Info : IA, Hexane/i-PrOH = 75/25, 1.0 mL/min, 30 oC, 230 nm



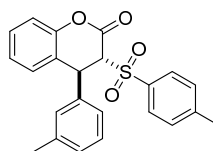
Area Percent Report

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 [mAU]	Area %
1	6.976	BB	0.2656	3693.31323	209.25581	50.0762	
2	8.439	BB	0.2274	3682.07446	245.50769	49.9238	

Totals : 7375.38770 454.76350



(+/-)-2o

*** End of Report ***

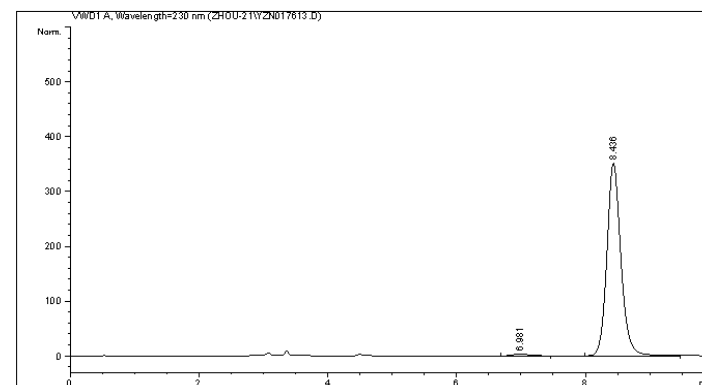
Instrument 1 3/3/2021 9:58:43 PM

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-21\YZN017613.D
Sample Name: HZ-10-45C

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Acq. Operator :
Acq. Instrument : Instrument 1 Location : -
Injection Date : 3/3/2021 9:41:12 PM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/3/2021 9:39:40 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/3/2021 9:59:28 PM
(modified after loading)
Sample Info : IA, Hexane/i-PrOH = 75/25, 1.0 mL/min, 30 oC, 230 nm



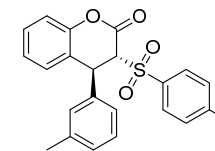
Area Percent Report

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 [mAU]	Area %
1	6.981	BB	0.2631	57.65787	3.26101	1.0773	
2	8.436	BB	0.2281	5294.17480	351.50778	98.9227	

Totals : 5351.83268 354.76879



(+)-2o

*** End of Report ***

Instrument 1 3/3/2021 9:59:31 PM

Page 1 of 1

Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3001944.D
Sample Name: HZ-10-49B+-

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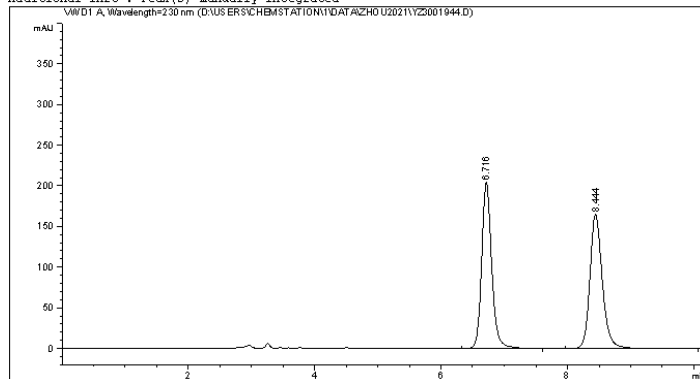
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 3/17/2021 9:07:53 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\def1_LC.M
Last changed : 3/17/2021 8:34:34 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\def1_LC.M
Last changed : 3/17/2021 9:48:47 PM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 75/25, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

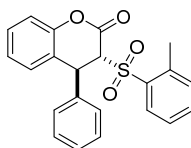


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.716	VV	0.1703	2283.18652	204.36960	49.9067
2	8.444	VV	0.2098	2291.72192	165.20837	50.0933



(+/-)-2p

1260II 3/17/2021 9:48:52 PM SYSTEM

Page 1 of 2

Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3001946.D
Sample Name: HZ-10-48B

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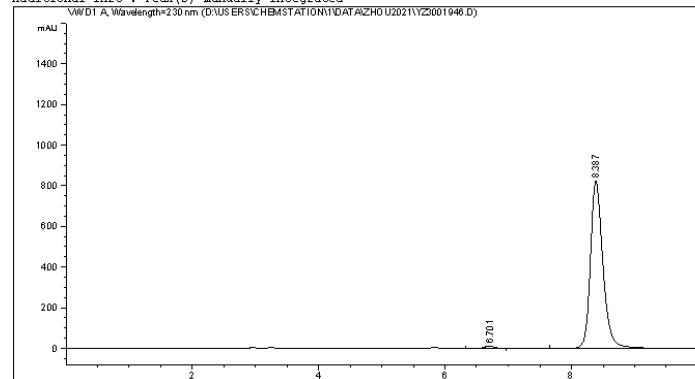
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 3/17/2021 9:35:34 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\def1_LC.M
Last changed : 3/17/2021 8:34:34 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\def1_LC.M
Last changed : 3/17/2021 9:47:39 PM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 75/25, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

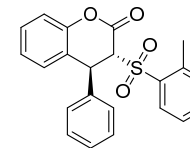


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.701	BV	0.1743	124.88231	10.53310	1.0933
2	8.387	BB	0.2083	1.12981e4	822.26489	98.9067



(+)-2p

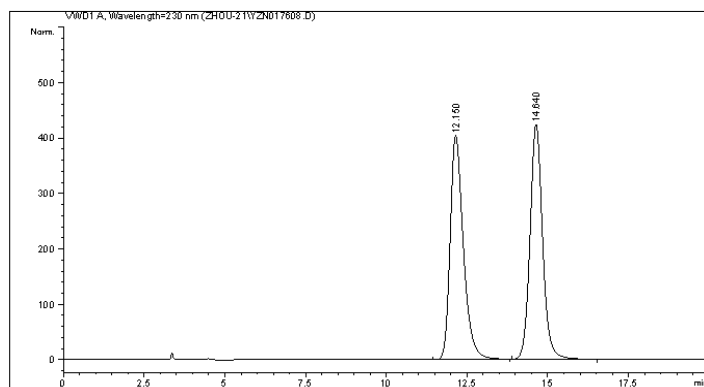
1260II 3/17/2021 9:47:46 PM SYSTEM

Page 1 of 2

Data File C:\CHEM32\1\DATA\ZHOU-21\YZN017608.D
Sample Name: HZ-10-46B+

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : -
Injection Date : 3/3/2021 8:53:07 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/3/2021 8:47:44 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/3/2021 10:09:41 AM
(modified after loading)
Sample Info : IA, Hexane/i-PrOH = 75/25, 1.0 mL/min, 30 oC, 230 nm



Area Percent Report

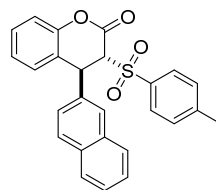
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area [mAU]	Area %
1	12.150	VB	0.4303	1.15187e4	403.88596	50.0197	
2	14.640	BB	0.4130	1.15097e4	423.76575	49.9803	

Totals : 2.30284e4 827.65170

*** End of Report ***



(+/-)-2q

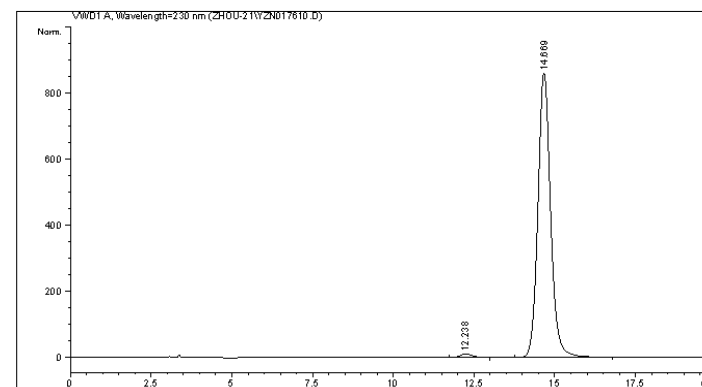
Instrument 1 3/3/2021 10:09:48 AM

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-21\YZN017610.D
Sample Name: HZ-10-45B

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : -
Injection Date : 3/3/2021 9:37:40 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/3/2021 9:36:32 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/3/2021 10:07:55 AM
(modified after loading)
Sample Info : IA, Hexane/i-PrOH = 75/25, 1.0 mL/min, 30 oC, 230 nm



Area Percent Report

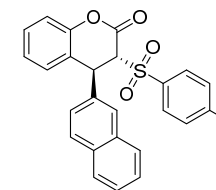
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height %s	Area [mAU]	Area %
1	12.238	BB	0.4192	327.90085	11.89184	1.3732	
2	14.669	VB	0.4151	2.35505e4	861.26312	98.6268	

Totals : 2.38784e4 873.15496

*** End of Report ***



(+)-2q

Instrument 1 3/3/2021 10:07:57 AM

Page 1 of 1

Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3001900.D
Sample Name: HZ-10-52A+-

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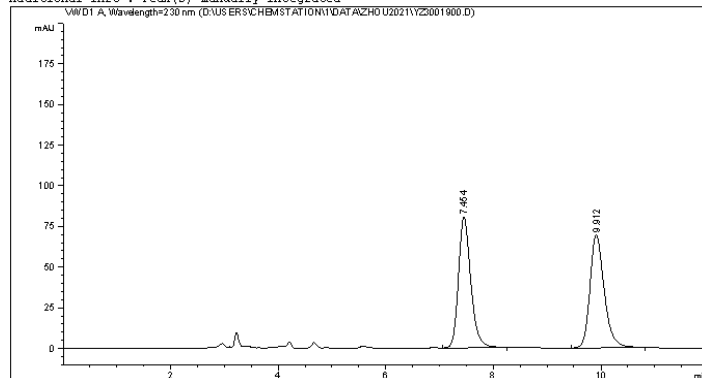
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 3/9/2021 4:10:51 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/9/2021 3:43:59 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/9/2021 4:47:03 PM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

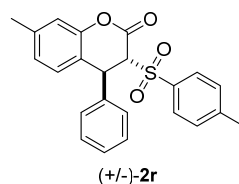


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.454	VB	0.2440	1280.01025	80.30379	49.9102
2	9.912	BB	0.2807	1284.61511	69.49427	50.0898



1260II 3/9/2021 4:47:05 PM SYSTEM

Page 1 of 2

Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3001902.D
Sample Name: HZ-10-51A

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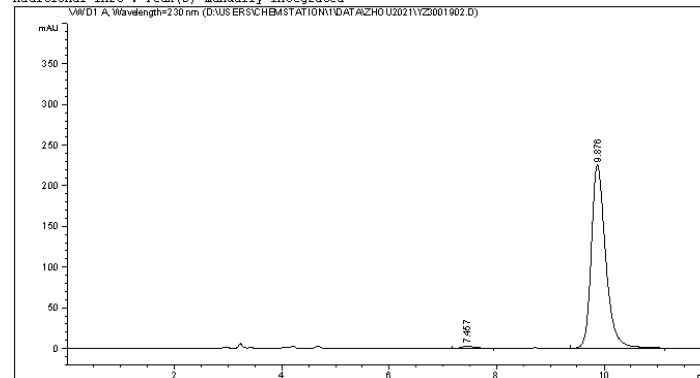
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 3/9/2021 4:42:52 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/9/2021 3:43:59 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/9/2021 5:00:16 PM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

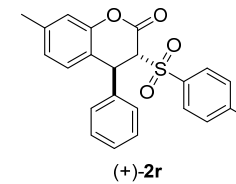


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.457	BB	0.2563	43.01477	2.51864	1.0222
2	9.876	BB	0.2806	4165.18164	225.33815	98.9778



1260II 3/9/2021 5:00:19 PM SYSTEM

Page 1 of 2

Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3001905.D
Sample Name: HZ-10-52B+-

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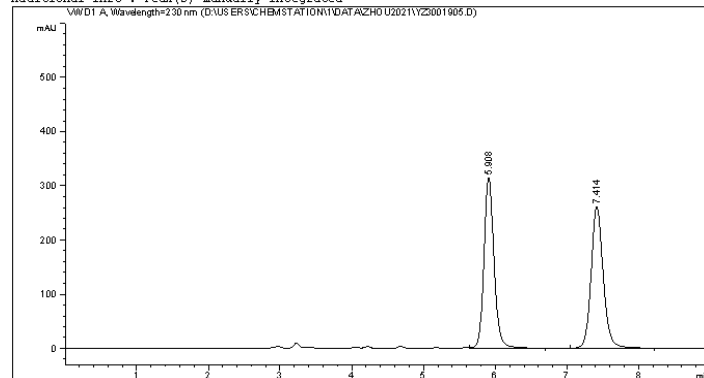
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 3/9/2021 8:25:21 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/9/2021 8:00:56 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/9/2021 9:22:36 PM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

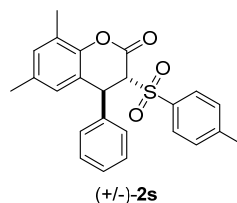


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	5.908	VB	0.1484	3068.82397	315.68811	50.0283
2	7.414	BB	0.1790	3065.34766	260.96362	49.9717



1260II 3/9/2021 9:22:40 PM SYSTEM

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Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3001907.D
Sample Name: HZ-10-51B

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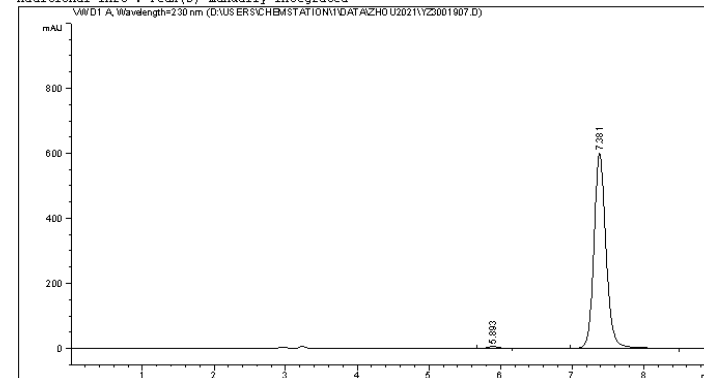
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 3/9/2021 8:59:09 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/9/2021 8:00:56 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/9/2021 9:23:02 PM by SYSTEM
(modified after loading)

Sample Info : IA, n-Hexane/i-PrOH = 70/30, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

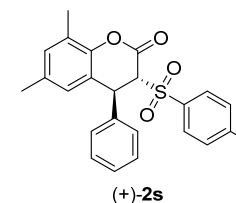


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	5.893	BB	0.1456	49.95801	5.26856	0.7086
2	7.381	BB	0.1781	7000.13770	600.20129	99.2914



1260II 3/9/2021 9:23:04 PM SYSTEM

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Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3002787.D
Sample Name: HZ-11-50D+-

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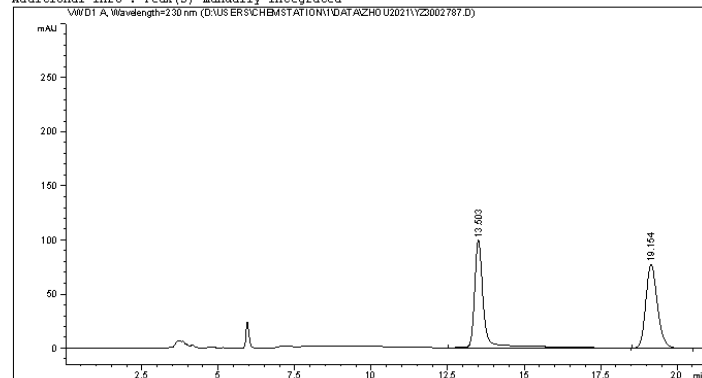
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 7/10/2021 8:06:50 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 7/10/2021 7:55:29 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 7/10/2021 9:00:48 PM by SYSTEM
(modified after loading)

Sample Info : AD-H, n-Hexane/i-PrOH = 70/30, 0.8 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

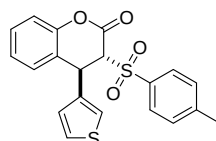


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	13.503	BB	0.3102	2101.04419	99.55585	50.8429
2	19.154	BB	0.4036	2031.38196	77.30142	49.1571



(+/-)-2t

Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3002788.D
Sample Name: HZ-11-50D

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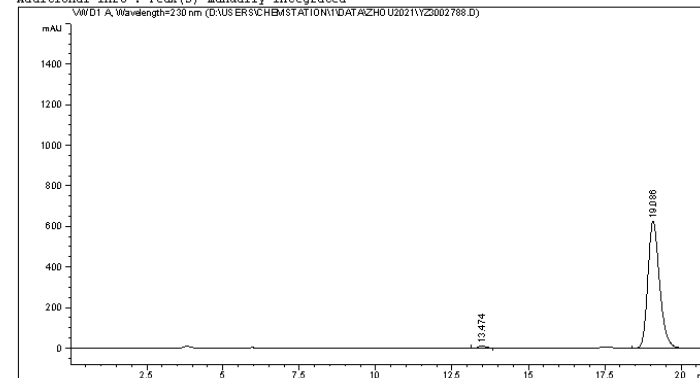
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 7/10/2021 8:39:50 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 7/10/2021 7:55:29 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 7/10/2021 9:03:26 PM by SYSTEM
(modified after loading)

Sample Info : AD-H, n-Hexane/i-PrOH = 70/30, 0.8 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

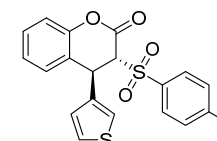


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	13.474	BB	0.2551	145.20985	8.91110	0.8698
2	19.086	BB	0.4058	1.65496e4	625.28168	99.1302



(+)-2t

Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3003713 10-32-22.D
Sample Name: HZ-11-66+

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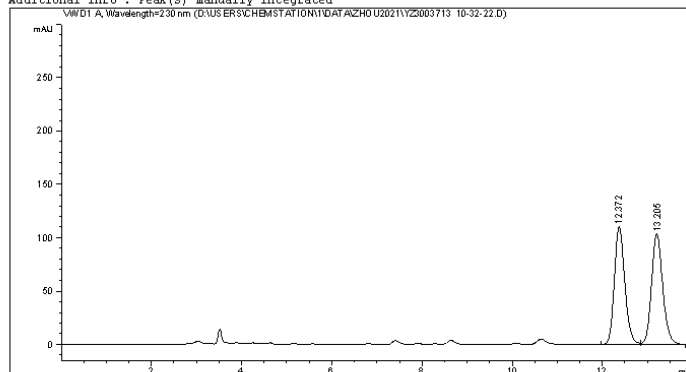
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 11/25/2021 10:32:22 AM Inj : 1
Inj Volume : No inj

Acq. Method : C:\Users\Public\Documents\ChemStation\1\Methods\def_LC.M
Last changed : 11/25/2021 10:21:18 AM by SYSTEM
(modified after loading)

Analysis Method : C:\Users\Public\Documents\ChemStation\1\Methods\def_LC.M
Last changed : 11/25/2021 10:48:35 AM by SYSTEM
(modified after loading)

Sample Info : AD-H, n-Hexane/i-PrOH = 90/10, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

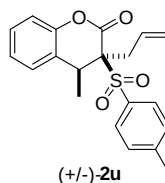


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.372	BV	0.2461	1769.60962	110.32888	49.8140
2	13.205	VB	0.2640	1782.82239	103.49602	50.1860



1260II 11/25/2021 10:48:44 AM SYSTEM

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Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3003714 10-52-55.D
Sample Name: HZ-11-65

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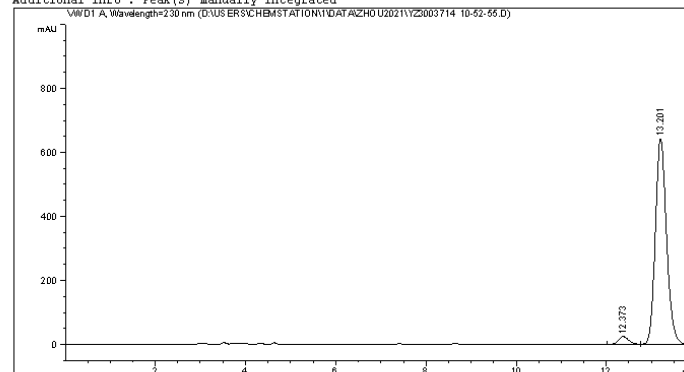
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 11/25/2021 10:52:55 AM Inj : 1
Inj Volume : No inj

Acq. Method : C:\Users\Public\Documents\ChemStation\1\Methods\def_LC.M
Last changed : 11/25/2021 10:21:18 AM by SYSTEM
(modified after loading)

Analysis Method : C:\Users\Public\Documents\ChemStation\1\Methods\def_LC.M
Last changed : 11/25/2021 11:09:31 AM by SYSTEM
(modified after loading)

Sample Info : AD-H, n-Hexane/i-PrOH = 90/10, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

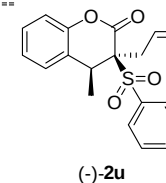


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.373	BV	0.2461	384.43906	23.96901	3.3590
2	13.201	VB	0.2647	1.10607e4	643.03699	96.6410



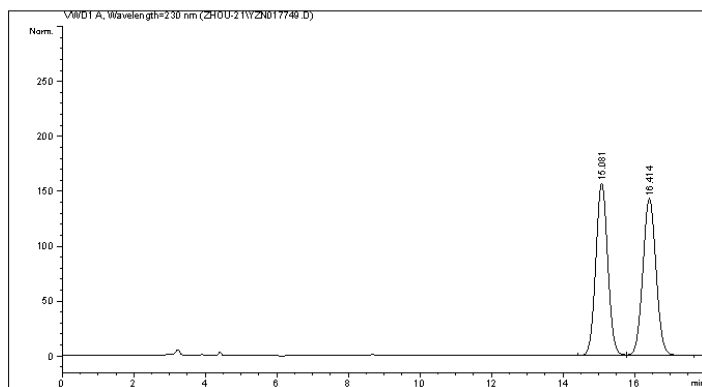
1260II 11/25/2021 11:09:39 AM SYSTEM

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Data File C:\CHEM32\1\DATA\ZHOU-21\YZN017749.D
Sample Name: HZ-10-55+-

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : -
Injection Date : 3/20/2021 8:13:58 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/20/2021 8:12:39 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/20/2021 8:37:43 AM
(modified after loading)
Sample Info : IC, Hexane/i-PrOH = 80/20, 1.0 mL/min, 30 oC, 230 nm



Area Percent Report

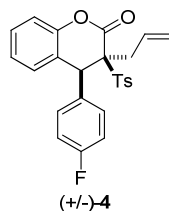
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height	Area %
1	15.081	BV	0.3649	3653.26636	156.84169	49.8823
2	16.414	VB	0.4017	3670.50830	142.85661	50.1177

Totals : 7323.77466 299.69830

*** End of Report ***



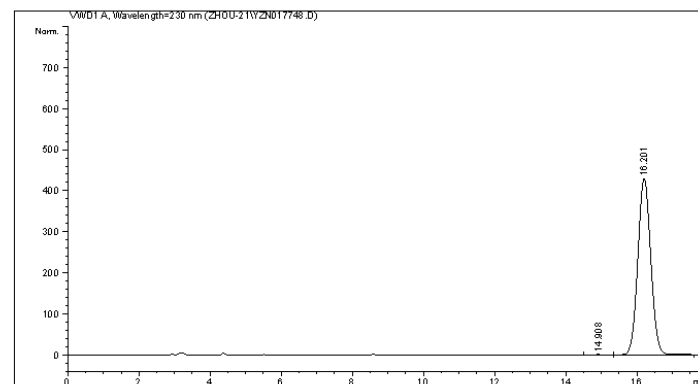
Instrument 1 3/20/2021 8:37:47 AM

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Data File C:\CHEM32\1\DATA\ZHOU-21\YZN017748.D
Sample Name: HZ-10-57

=====

Acq. Operator :
Acq. Instrument : Instrument 1 Location : -
Injection Date : 3/20/2021 7:51:59 AM
Acq. Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/20/2021 7:11:44 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\DEF_LC.M
Last changed : 3/20/2021 8:36:29 AM
(modified after loading)
Sample Info : IC, Hexane/i-PrOH = 80/20, 1.0 mL/min, 30 oC, 230 nm



Area Percent Report

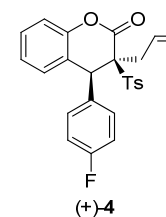
Sorted By : Signal
Multiplier: : 1.0000
Dilution: : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area mAU	Height	Area %
1	14.908	BV	0.3201	7.16760	3.30379e-1	0.0662
2	16.201	VB	0.3931	1.08167e4	429.06293	99.9338

Totals : 1.08238e4 429.39331

*** End of Report ***



Instrument 1 3/20/2021 8:36:38 AM

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Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3001970.D
Sample Name: HZ-10-59+-

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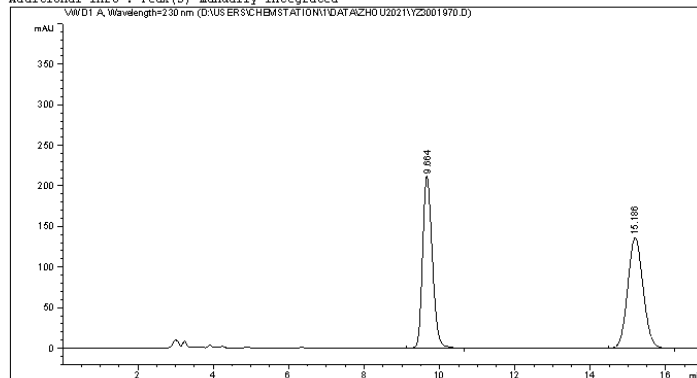
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 3/20/2021 5:04:48 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/20/2021 4:40:03 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/22/2021 10:18:49 AM by SYSTEM
(modified after loading)

Sample Info : AD-H, n-Hexane/i-PrOH = 60/40, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

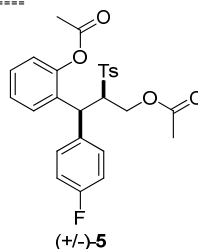


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.664	VB	0.2783	3779.10840	211.70270	49.6477
2	15.186	BB	0.4405	3832.73706	136.13089	50.3523



1260II 3/22/2021 10:18:54 AM SYSTEM

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Data File D:\USERS\CHEMSTATION\1\DATA\ZHOU2021\YZ3001969.D
Sample Name: HZ-10-61

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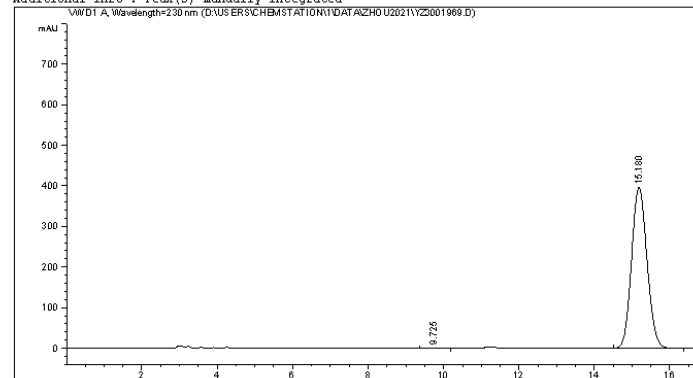
Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : 1260II Location : -
Injection Date : 3/20/2021 4:45:21 PM Inj : 1
Inj Volume : No inj

Acq. Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/20/2021 4:40:03 PM by SYSTEM
(modified after loading)

Analysis Method : C:\USERS\PUBLIC\DOCUMENTS\CHEMSTATION\1\METHODS\defl_LC.M
Last changed : 3/22/2021 10:17:49 AM by SYSTEM
(modified after loading)

Sample Info : AD-H, n-Hexane/i-PrOH = 60/40, 1.0 mL/min, 30 oC, 230 nm

Additional Info : Peak(s) manually integrated

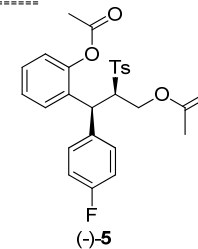


Area Percent Report

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: VWD1 A, Wavelength=230 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.725	BB	0.2672	22.31732	1.30695	0.1967
2	15.180	VB	0.4449	1.13227e4	396.83463	99.8033



1260II 3/22/2021 10:18:17 AM SYSTEM

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