

Chiral Brønsted Acid Catalyzed Enantioselective Intermolecular Allylic Aminations

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Supporting Information

General consideration:

All the solvents were purified under standard method. ^1H , ^{13}C , ^{19}F , and ^{31}P NMR spectra were recorded in CDCl_3 solution on Bruker AX-400 MHz instruments and spectral data were reported in ppm relative to tetramethylsilane (TMS) as internal standard. IR spectra were recorded on a Thermo Nicolet 6700 spectrometer. High-resolution mass spectral analysis (HRMS) data were measured on the Thermo Exactive by means of the ESI technique. Ee's were determined by high performance liquid chromatography (HPLC) analysis. All the amine nucleophiles were used directly after purchase.

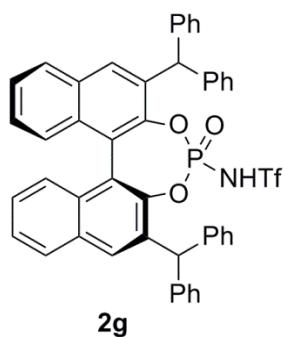
General procedure for the synthesis of Allylic alcohols: allylic alcohols were prepared according to the reported method. To a stirred solution of α,β -unsaturated ketone chalcone (10 mmol) in methanol (25 mL) was added $\text{CeCl}_3 \cdot 7\text{H}_2\text{O}$ (3.73 g, 10 mmol), the solution was cooled to 0 °C and NaBH_4 (0.418 g, 11 mmol) was added. The mixture was stirred at room temperature for 1 h before water was added and exacted with CH_2Cl_2 . The combined organic phases were dried with Na_2SO_4 , and the solvents evaporated under vacuum. The pure compound was obtained after flash chromatography on silica gel.

P. N. Chatterjee and S. Roy, *Tetrahedron*, 2012, **68**, 3776.

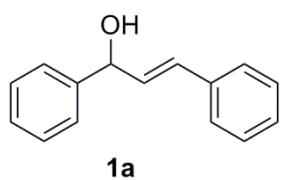
Typical procedure for the synthesis of chiral sulfonamides from allylic alcohol (Table 2, entry 1): To a dried Schlenk flask was added (*E*)-1,3-diphenylpropenol (**1a**) (0.084 g, 0.40 mmol), toluene sulfonamide (**3**) (0.1026 g, 0.60 mmol), and CHCl_3 (1.0 mL) under Ar. The reaction mixture was stirred at -60 °C for 10 min before addition of chiral phosphoramido **2g** (0.0325 g, 0.04 mmol). The reaction mixture was stirred at -60 °C for 10 h and quenched with saturated NaHCO_3 . The organic phase was directly purified by flash chromatography on silica gel (CH_2Cl_2) to afford the desired product **4a** as a white solid.

Typical procedure for the determination of the structures of regioisomers (4i** and **4i'**):** To a solution of **4i** and **4i'** (0.0782 g, 0.19 mmol) in a mixed solvent $\text{CCl}_4/\text{CH}_3\text{CN}/\text{H}_2\text{O}$ (1:1:1.6, v/v/v, 1.1 mL), NaIO_4 (0.165 g, 0.77 mmol) was added and stirred for 10 min. Then $\text{RuCl}_3 \cdot \text{H}_2\text{O}$ (0.0013 g, 0.0056 mmol) was added, and the mixture was stirred for 24 h at room temperature and quenched with 1N HCl. The mixture was extracted with CH_2Cl_2 , filtered with celite, concentrated and used directly for the GC-MS analysis.

C.-Y. Zhou, S.-F. Zhu, L.-X. Wang and Q.-L. Zhou, *J. Am. Chem. Soc.*, 2010, **132**, 10955.



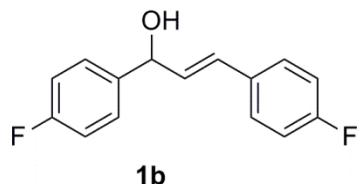
(S)-(+)-[3,3'-Bis(diphenylmethyl)-1,1'-binaphthalen-2,2'-yl]-N-triflylphosphoramide: Yellow solid; m.p. 218-220 °C; $[\alpha]_D^{20} -101.6$ (*c* 1.03, CH_2Cl_2); IR (film) 3061.1, 3027.3, 1647.5, 1598.6, 1494.3, 1450.8 cm^{-1} ; ^1H NMR (400 MHz, CDCl_3) δ 7.72 (t, *J* = 7.6 Hz, 2H), 7.50 (d, *J* = 5.0 Hz, 1H), 7.45-7.34 (m, 4H), 7.32-7.10 (m, 19H), 7.12-7.08 (m, 4H), 6.34-6.23 (m, 2H), 3.47 (brs, 1H); ^{31}P NMR (122 MHz, CDCl_3) δ -1.5; ^{19}F NMR (376 MHz, CDCl_3) δ -77.3; HRMS (FT-ICRMS) calcd for $\text{C}_{47}\text{H}_{33}\text{NO}_5\text{PSNa}$ ($\text{M}+\text{Na}$): 834.1661, found: 834.1653.



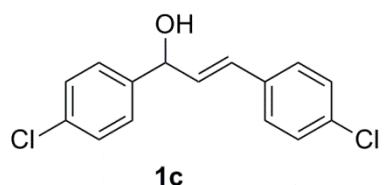
(E)-1,3-diphenylprop-2-en-1-ol (1a): 85% yield, white solid; ^1H NMR (400 MHz, CDCl_3) δ 7.42 (d, *J* = 7.2 Hz, 2H), 7.37 (t, *J* = 8.0 Hz, 4H), 7.30 (t, *J* = 7.6 Hz, 3H), 7.26-7.20 (m, 1H), 6.68 (d, *J* = 15.8 Hz, 1H), 6.37 (dd, *J* = 15.8, 6.4 Hz, 1H), 5.37 (d, *J* = 6.4 Hz, 1H), 2.12 (brs, 1H); ^{13}C NMR

(100 MHz, CDCl₃) δ 143.0, 136.7, 131.7, 130.8, 128.8, 128.7, 128.0, 127.9, 126.8, 126.6, 75.3.

J. M. Dickinson, J. A. Murphy, C. W. Patterson and N. F. Wooster, *J. Chem. Soc., Perkin Trans. 1*, 1990, 1179;

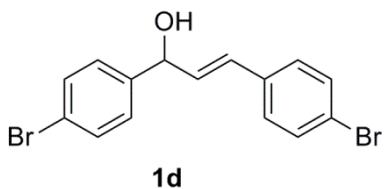


(E)-1,3-bis(4-fluorophenyl)prop-2-en-1-ol (1b): 61% yield, colorless oil; IR (film) 3371, 1602, 1508, 1227, 1157, 1085 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) δ 7.34-7.22 (m, 4H), 7.15-6.90 (m, 4H), 6.52 (d, *J* = 16.0 Hz, 1H), 6.18 (dd, *J* = 16.0, 6.0 Hz, 1H), 5.24 (d, *J* = 6.0 Hz, 1H), 2.99 (brs, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 162.6 (d, *J* = 246.0 Hz), 162.4 (d, *J* = 244.0 Hz), 138.6 (d, *J* = 3.0 Hz), 132.6 (d, *J* = 3.0 Hz), 131.2 (d, *J* = 2.0 Hz), 129.6, 128.3 (d, *J* = 9.0 Hz), 128.2 (d, *J* = 8.0 Hz), 115.7 (d, *J* = 12.0 Hz), 115.5 (d, *J* = 12.0 Hz), 74.4; HRMS (FT-ICRMS) calcd for C₁₅H₁₁F₂O (M-H): 245.0772, found: 245.0781.



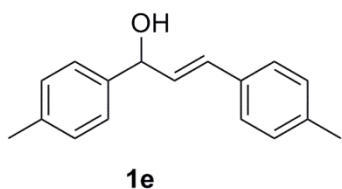
(E)-1,3-bis(4-chlorophenyl)prop-2-en-1-ol (1c): 72% yield, white solid; ¹H NMR (400 MHz, CDCl₃) δ 7.35-7.25 (m, 4H), 7.25-7.21 (m, 4H), 6.55 (d, *J* = 16.0 Hz, 1H), 6.24 (dd, *J* = 16.0, 6.4 Hz, 1H), 5.27 (d, *J* = 6.4 Hz, 1H), 2.61 (brs, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 141.1, 134.9, 133.7, 131.8, 129.8, 128.9, 128.0, 127.9, 74.4.

P. N. Chatterjee and S. Roy, *Tetrahedron*, 2012, **68**, 3776.



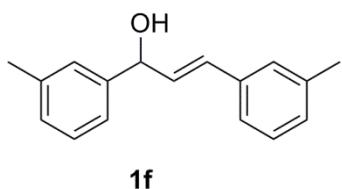
(E)-1,3-bis(4-bromophenyl)prop-2-en-1-ol (1d): 89% yield, white solid; ^1H NMR (400 MHz, CDCl_3) δ 7.49 (d, $J = 6.8$, 2H), 7.43 (d, $J = 6.8$, 2H), 7.29 (d, $J = 8.4$ Hz, 2H), 7.23 (d, $J = 6.8$ Hz, 2H), 6.59 (d, $J = 16.0$ Hz, 1H), 6.30 (dd, $J = 16.0$, 6.4 Hz, 1H), 5.36-5.30 (m, 1H), 2.15 (d, $J = 3.6$ Hz, 1H); ^{13}C NMR (100 MHz, CDCl_3) δ 141.7, 135.5, 132.0, 132.0, 130.0, 128.4, 128.3, 122.0, 74.6.

P. N. Chatterjee and S. Roy, *Tetrahedron*, 2012, **68**, 3776.



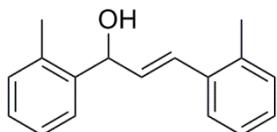
(E)-1,3-di-p-tolylprop-2-en-1-ol (1e): 81% yield, white solid; ^1H NMR (400 MHz, CDCl_3) δ 7.24 (d, $J = 8.0$ Hz, 4H), 7.20 (d, $J = 8.0$ Hz, 4H), 7.10 (d, $J = 8.0$ Hz, 2H), 7.05 (d, $J = 8.0$ Hz, 2H), 6.54 (d, $J = 14.6$ Hz, 1H), 6.24 (dd, $J = 14.6$, 6.4 Hz, 1H), 5.21 (d, $J = 6.4$ Hz, 1H), 2.63 (brs, 1H), 2.30 (s, 3H), 2.28 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 140.2, 137.5, 137.4, 134.0, 130.9, 130.3, 129.3, 129.3, 126.6, 126.5, 75.0, 21.3, 21.2.

P. N. Chatterjee and S. Roy, *Tetrahedron*, 2012, **68**, 3776.



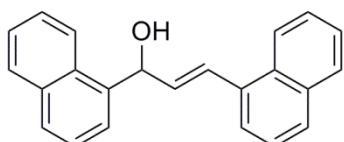
(E)-1,3-di-m-tolylprop-2-en-1-ol (1f): 79% yield, colorless oil; IR (film) 3348, 1605, 1487, 1455, 1081 cm^{-1} ; ^1H NMR (400 MHz, CDCl_3) δ 7.25-7.13 (m, 6H), 7.06 (d, $J = 7.2$ Hz, 1H), 7.02 (d, $J =$

6.0 Hz, 1H), 6.59 (d, J = 15.8 Hz, 1H), 6.28 (dd, J = 15.8, 6.4 Hz, 1H), 5.25 (d, J = 6.4 Hz, 1H), 2.44 (brs, 1H), 2.32 (s, 3H), 2.29 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 143.0, 138.4, 138.2, 136.7, 131.6, 130.6, 128.7, 128.6, 128.6, 127.5, 127.2, 124.0, 123.6, 75.2, 21.6, 21.5; HRMS (FT-ICRMS) calcd for $\text{C}_{17}\text{H}_{17}\text{O}$ (M-H): 237.1274, found: 237.1280.



1g

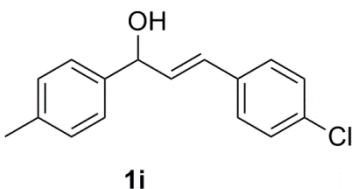
(E)-1,3-di(o-tolyl)prop-2-en-1-ol (1g): 80% yield, colorless oil; IR (film) 3336, 1602, 1486, 1460, 1011 cm^{-1} ; ^1H NMR (400 MHz, CDCl_3) δ 7.52 (d, J = 7.6 Hz, 1H), 7.44-7.37 (m, 1H), 7.40-7.10 (m, 6H), 6.86 (d, J = 16.0 Hz, 1H), 6.22 (dd, J = 16.0, 6.0 Hz, 1H), 5.57 (d, J = 6.0 Hz, 1H), 2.39 (s, 3H), 2.33 (s, 3H), 2.10 (brs, 1H); ^{13}C NMR (100 MHz, CDCl_3) δ 140.9, 135.9, 135.8, 135.4, 132.3, 130.8, 130.5, 128.7, 127.8, 126.6, 126.3, 126.0, 126.0, 72.3, 20.0, 19.4; HRMS (FT-ICRMS) calcd for $\text{C}_{17}\text{H}_{17}\text{O}$ (M-H): 237.1274, found: 237.1280.



1h

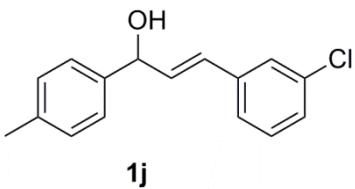
(E)-1,3-di(naphthalen-1-yl)prop-2-en-1-ol (1h): 65% yield, yellow oil; ^1H NMR (400 MHz, CDCl_3) δ 8.34 (d, J = 8.4 Hz, 1H), 8.13-8.05 (m, 1H), 7.95-7.89 (m, 1H), 7.88-7.81 (m, 2H), 7.70 (t, J = 7.2 Hz, 2H), 7.60-7.46 (m, 7H), 7.41 (t, J = 8.0 Hz, 1H), 6.63 (dd, J = 15.6, 5.6 Hz, 1H), 6.25-6.21 (m, 1H), 2.35 (d, J = 4.0 Hz, 1H); ^{13}C NMR (100 MHz, CDCl_3) δ 138.5, 134.6, 134.4, 134.3, 133.8, 131.4, 131.0, 129.1, 128.9, 128.7, 128.4, 128.3, 126.5, 126.3, 126.0, 126.0, 125.8, 125.7, 124.4, 124.2, 124.0, 124.0, 72.6.

K. Itoh, T. Ikeda, S. Tazuke and T. Shibata, *J. Phys. Chem.*, 1992, **96**, 5759–5765

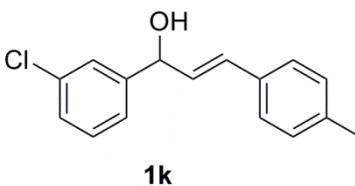


(E)-3-(4-chlorophenyl)-1-(p-tolyl)prop-2-en-1-ol (1i): 73% yield, white solid; ^1H NMR (400 MHz, CDCl_3) δ 7.33-7.23 (m, 6H), 7.18 (d, $J = 8.0$ Hz, 2H), 6.63 (d, $J = 15.8$ Hz, 1H), 6.35 (dd, $J = 15.8, 6.2$ Hz, 1H), 5.37-5.31 (m, 1H), 2.35 (s, 3H), 2.00 (d, $J = 3.0$ Hz, 1H); ^{13}C NMR (100 MHz, CDCl_3) δ 139.9, 138.0, 135.4, 133.5, 132.5, 129.6, 129.2, 128.9, 128.0, 126.5, 75.1, 21.4.

P. N. Chatterjee and S. Roy, *Tetrahedron*, 2012, **68**, 3776.

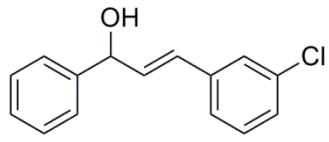


(E)-3-(3-chlorophenyl)-1-(p-tolyl)prop-2-en-1-ol (1j): 83% yield, colorless oil; IR (film) 3342, 1594, 1566, 1512, 1426 cm^{-1} ; ^1H NMR (400 MHz, CDCl_3) δ 7.27 (s, 1H), 7.25-7.20 (m, 2H), 7.16-7.08 (m, 5H), 6.49 (d, $J = 15.8$ Hz, 1H), 6.31 (dd, $J = 15.8, 6.2$ Hz, 1H), 5.21 (d, $J = 6.2$ Hz, 1H), 2.98 (brs, 1H), 2.30 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 139.7, 138.7, 137.7, 134.5, 133.3, 129.8, 129.4, 128.7, 127.7, 126.6, 126.5, 124.9, 74.7, 21.3; HRMS (FT-ICRMS) calcd for $\text{C}_{16}\text{H}_{14}\text{ClO}$ ($\text{M}-\text{H}$): 257.0728, found: 257.0736.



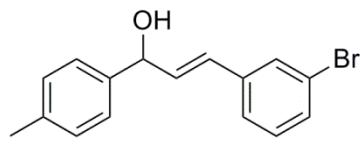
(E)-1-(3-chlorophenyl)-3-(p-tolyl)prop-2-en-1-ol (1k): 77% yield, colorless oil; IR (film) 3339, 1596, 1574, 1512, 1428, 1188 cm^{-1} ; ^1H NMR (400 MHz, CDCl_3) δ 7.33 (s, 1H), 7.20-7.13 (m, 5H),

7.08-7.01 (m, 2H), 6.49 (d, J = 15.8 Hz, 1H), 6.14 (dd, J = 15.8, 6.8 Hz, 1H), 5.14 (d, J = 6.8 Hz, 1H), 3.27 (brs, 1H), 2.27 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 145.1, 137.9, 134.4, 133.5, 131.2, 129.9, 129.4, 127.7, 126.7, 126.5, 124.6, 74.5, 21.3; HRMS (FT-ICRMS) calcd for $\text{C}_{16}\text{H}_{14}\text{ClO}$ (M-H): 257.0728, found: 257.0736.



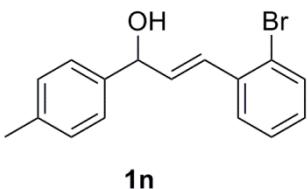
1l

(E)-3-(3-chlorophenyl)-1-phenylprop-2-en-1-ol (1l): 88%, colorless oil; IR (film) 3338, 1594, 1565, 1453, 1094, 1076 cm^{-1} ; ^1H NMR (400 MHz, CDCl_3) δ 7.40-7.26 (m, 6H), 7.20-7.15 (m, 3H), 6.57 (d, J = 16.0 Hz, 1H), 6.33 (dd, J = 16.0, 6.4 Hz, 1H), 5.33-5.28 (m, 1H), 2.47 (brs, 1H); ^{13}C NMR (100 MHz, CDCl_3) δ 142.6, 138.6, 134.7, 133.2, 129.9, 129.1, 128.9, 128.1, 127.8, 126.7, 126.5, 125.0, 75.0; HRMS (FT-ICRMS) calcd for $\text{C}_{15}\text{H}_{12}\text{ClO}$ (M-H): 243.0571, found: 243.0578.

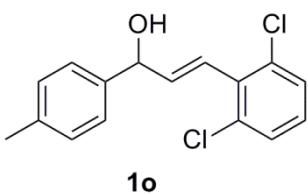


1m

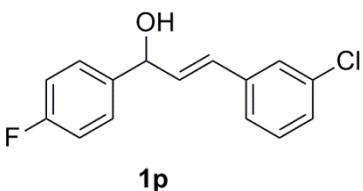
(E)-3-(3-bromophenyl)-1-(p-tolyl)prop-2-en-1-ol (1m): 75% yield, white solid; IR (film) 3340, 1590, 1562, 1473, 1072 cm^{-1} ; ^1H NMR (400 MHz, CDCl_3) δ 7.45 (s, 1H), 7.30 (d, J = 8.0 Hz, 1H), 7.26-7.18 (m, 3H), 7.16-7.05 (m, 3H), 6.50 (d, J = 16.0 Hz, 1H), 6.29 (dd, J = 16.0, 6.0 Hz, 1H), 5.24 (d, J = 6.0 Hz, 1H), 2.75 (brs, 1H), 2.31 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 139.7, 139.0, 137.7, 133.4, 130.6, 130.2, 129.5, 129.4, 128.6, 126.5, 125.4, 122.8, 74.7, 21.3; HRMS (FT-ICRMS) calcd for $\text{C}_{16}\text{H}_{14}\text{BrO}$ (M-H): 301.0222, found: 301.0231.



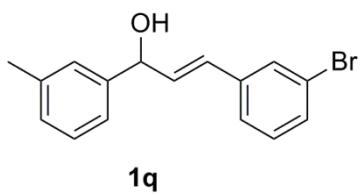
(E)-3-(2-bromophenyl)-1-(p-tolyl)prop-2-en-1-ol (1n): 67% yield, white solid; IR (film) 3355, 1512, 1466, 1436, 1022 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) δ 7.51 (d, *J* = 8.0 Hz, 1H), 7.45 (d, *J* = 8.0 Hz, 1H), 7.30 (d, *J* = 8.0 Hz, 2H), 7.23-7.13 (m, 3H), 7.10-6.98 (m, 2H), 6.28 (dd, *J* = 15.6, 6.4 Hz, 1H), 5.36-5.30 (m, 1H), 2.36 (d, *J* = 3.2 Hz, 1H), 2.33 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 139.7, 137.7, 136.7, 134.9, 133.1, 129.5, 129.1, 127.6, 127.4, 126.5, 125.6, 124.0, 75.0, 21.3; HRMS (FT-ICRMS) calcd for C₁₆H₁₄BrO (M-H): 301.0222, found: 301.0232.



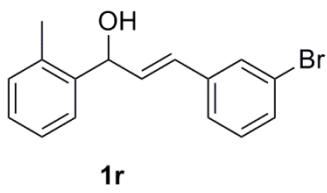
(E)-3-(2,6-dichlorophenyl)-1-(p-tolyl)prop-2-en-1-ol (1o): 80% yield, white solid; IR (film) 3294, 1555, 1512, 1428, 1180, 1090 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) δ 7.37 (d, *J* = 8.0 Hz, 2H), 7.30 (d, *J* = 8.0 Hz, 2H), 7.20 (d, *J* = 8.0 Hz, 2H), 7.09 (t, *J* = 8.0 Hz, 1H), 6.74 (d, *J* = 16.0 Hz, 1H), 6.42 (dd, *J* = 16.0, 5.6 Hz, 1H), 5.44-5.37 (m, 1H), 2.36 (s, 3H), 2.06 (d, *J* = 4.0 Hz, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 140.1, 139.4, 137.9, 134.7, 134.5, 129.6, 128.6, 128.4, 126.9, 124.0, 75.0, 21.4; HRMS (FT-ICRMS) calcd for C₁₆H₁₃Cl₂O (M-H): 291.0338, found: 291.0336.



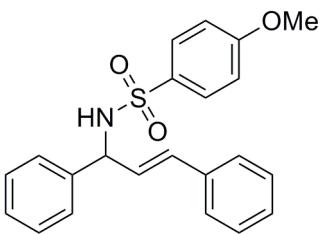
(E)-3-(3-chlorophenyl)-1-(4-fluorophenyl)prop-2-en-1-ol (1p): 66% yield, colorless oil; IR (film) 3339, 1602, 1509, 1224, 1078 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) δ 7.36-7.28 (m, 3H), 7.20-7.13 (m, 3H), 7.06-6.93 (m, 2H), 6.53 (d, *J* = 16.0 Hz, 1H), 6.28 (dd, *J* = 16.0, 6.4 Hz, 1H), 5.28-5.24 (m, 1H), 2.80 (brs, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 162.4 (d, *J* = 244.0 Hz), 138.4, 138.3 (d, *J* = 3.0 Hz), 134.7, 132.9, 130.0, 129.3, 128.2 (d, *J* = 8.0 Hz), 128.0, 126.6, 125.0, 115.6 (d, *J* = 21.0 Hz), 74.3; HRMS (FT-ICRMS) calcd for C₁₅H₁₁ClFO (M-H): 261.0477, found: 261.0485.



(E)-3-(3-bromophenyl)-1-(m-tolyl)prop-2-en-1-ol (1q): 73% yield, colorless oil; IR (film) 3339, 1562, 1474, 1072 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) δ 7.53 (s, 1H), 7.37-7.31 (m, 1H), 7.29-7.10 (m, 6H), 6.61 (d, *J* = 15.6 Hz, 1H), 6.37 (dd, *J* = 15.6, 6.0 Hz, 1H), 5.35-5.31 (m, 1H), 2.36 (s, 3H), 2.09 (d, *J* = 3.4 Hz, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 142.6, 139.0, 138.7, 133.3, 130.8, 130.3, 129.6, 129.0, 128.9, 128.8, 127.2, 125.5, 123.6, 123.0, 75.1, 21.7; HRMS (FT-ICRMS) calcd for C₁₆H₁₄BrO (M-H): 301.0222, found: 301.0231.

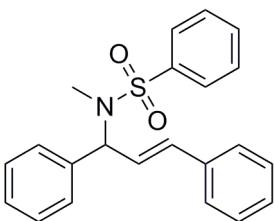


(E)-3-(3-bromophenyl)-1-(o-tolyl)prop-2-en-1-ol (1r): 79% yield, colorless oil; IR (film) 3326, 1590, 1562, 1473, 1072 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) δ 7.52-7.46 (m, 2H), 7.37-7.31 (m, 1H), 7.29-7.11 (m, 5H), 6.56 (d, *J* = 16.0 Hz, 1H), 6.37 (dd, *J* = 16.0, 6.0 Hz, 1H), 5.58-5.48 (m, 1H), 2.37 (s, 3H), 2.06 (d, *J* = 3.4 Hz, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 140.5, 139.0, 135.5, 132.5, 130.9, 130.8, 130.3, 129.6, 129.1, 128.0, 126.7, 126.1, 125.4, 123.0, 71.8, 19.4; HRMS (FT-ICRMS) calcd for C₁₆H₁₄BrO (M-H): 301.0222, found: 301.0231.

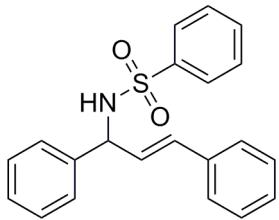


(E)-N-(1,3-diphenylallyl)-4-methoxybenzenesulfonamide(Table 1, entry 12**):** White solid; yield: 72%; $[\alpha]_D^{20}$ -22.4 (*c* 0.82, CH₂Cl₂) (72% ee); ¹H NMR (400 MHz, CDCl₃) δ 7.69 (d, *J* = 8.8 Hz, 2H), 7.29-7.14 (m, 10H), 6.78 (d, *J* = 8.8 Hz, 2H), 6.35 (d, *J* = 15.6 Hz, 1H), 6.07 (dd, *J* = 15.6, 6.6 Hz, 1H), 5.17-5.05 (m, 2H), 3.74 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 162.9, 139.8, 136.3, 132.5, 132.3, 129.6, 128.9, 128.7, 128.4, 128.1, 128.0, 127.3, 126.7, 114.2, 60.0, 55.7.

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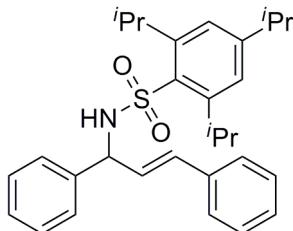


(E)-N-(1,3-diphenylallyl)-N-methylbenzenesulfonamide(Table 1, entry 13**):** Colorless oil; yield: 93%; $[\alpha]_D^{20}$ -0.5 (*c* 0.98, CH₂Cl₂) (2% ee); IR (film) 3060, 1447, 1339, 1164, 1088 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) δ 7.83 (d, *J* = 8.6 Hz, 2H), 7.50-7.46 (m, 1H), 7.45-7.38 (m, 2H), 7.35-7.18 (m, 10H), 6.40 (d, *J* = 16.0 Hz, 1H), 6.16 (dd, *J* = 16.0, 7.4 Hz, 1H), 5.87 (d, *J* = 7.4 Hz, 1H), 2.71 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 139.7, 138.6, 136.3, 134.9, 132.6, 129.1, 128.8, 128.3, 128.1, 128.0, 127.6, 126.7, 124.2, 62.6, 30.4; HRMS (ESI) calcd for C₂₂H₂₁NO₂SNa (M+Na): 386.1185, found: 386.1182.

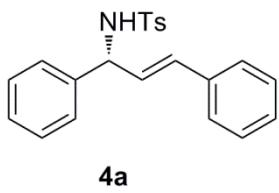


(E)-N-(1,3-diphenylallyl)benzenesulfonamide (Table 1, entry 14): White solid; yield: 78%; $[\alpha]_D^{20}$ -19.9 (*c*0.96, CH₂Cl₂) (72% ee); ¹H NMR (400 MHz, CDCl₃) δ 7.77 (d, *J* = 8.0 Hz, 2H), 7.48-7.42 (m, 1H), 7.39-7.31 (m, 2H), 7.29-7.13 (m, 10H), 6.38 (d, *J* = 15.8 Hz, 1H), 6.06 (dd, *J* = 15.8, 6.4 Hz, 1H), 5.19-5.08 (m, 2H); ¹³C NMR (100 MHz, CDCl₃) δ 140.9, 139.7, 136.2, 132.6, 132.4, 129.0, 128.9, 128.7, 128.4, 128.2, 128.1, 127.2, 126.7, 60.0.

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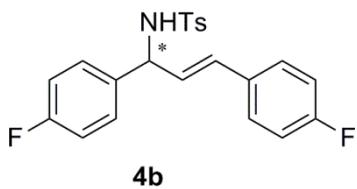


(E)-N-(1,3-diphenylallyl)-2,4,6-triisopropylbenzenesulfonamide (Table 1, entry 15): White solid; yield: 82%; m.p. 151-153 °C; $[\alpha]_D^{20}$ -22.4 (*c*0.78, CH₂Cl₂) (84% ee); IR (film) 3300, 1600, 1455, 1425, 1321, 1151 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) δ 7.28-7.21 (m, 4H), 7.21-7.15 (m, 4H), 7.15-7.10 (m, 2H), 7.09 (s, 2H), 6.39 (d, *J* = 15.8 Hz, 1H), 6.10 (dd, *J* = 15.8, 6.6 Hz, 1H), 5.19 (dd, *J* = 6.6, 6.0 Hz, 1H), 4.75 (d, *J* = 6.0 Hz, 1H), 4.16-4.04 (m, 2H), 2.90-2.80 (m, 1H), 1.25 (s, 3H), 1.24 (s, 3H), 1.23 (s, 3H), 1.21 (s, 3H), 1.15 (s, 3H), 1.14 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 152.9, 150.1, 140.2, 136.3, 134.0, 132.2, 129.0, 128.7, 128.6, 128.2, 128.1, 127.4, 126.7, 123.8, 59.9, 34.3, 30.0, 25.2, 24.8, 23.8, 23.7; HRMS (ESI) calcd for C₃₀H₃₇NO₂SNa (M+Na): 498.2437, found: 498.2432.

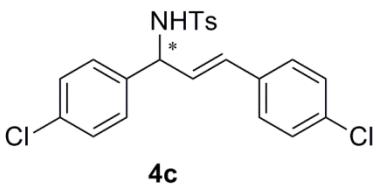


N-[1,3-diphenyl-(E)-2-propenyl]-4-methylbenzenesulfonamide (4a) (Table 2, entry 1): White solid; yield: 88%; $[\alpha]_D^{20} -23.2$ (*c* 0.98, CH_2Cl_2) (75% ee) ($[\alpha]_D^{25} -31.4$ (*c* 1.0, CHCl_3), 95% ee for *R*); ^1H NMR (400 MHz, CDCl_3) δ 7.65 (d, *J* = 8.2 Hz, 2H), 7.29-7.10 (m, 12H), 6.35 (d, *J* = 15.8 Hz, 1H), 6.07 (dd, *J* = 15.8, 6.8 Hz, 1H), 5.11 (dd, *J* = 7.0, 6.8 Hz, 1H), 4.97 (d, *J* = 7.0 Hz, 1H), 2.32 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 143.5, 139.9, 138.0, 136.3, 132.4, 129.7, 129.0, 128.7, 128.4, 128.1, 128.0, 127.6, 127.3, 126.7, 60.0, 21.6.

Bower, J. F., Jumnah, R., Williams, A. C. and Williams, J. M. J. *J. Chem. Soc., Perkin Trans. 1*, **1997**, 1411-1420.

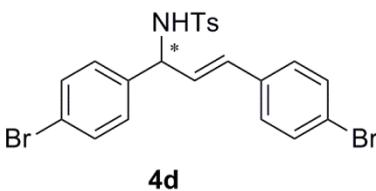


(E)-N-(1,3-bis(4-fluorophenyl)allyl)-4-methylbenzenesulfonamide (4b) (Table 2, entry 2): White solid; yield: 82%; m.p. 137-139 °C; $[\alpha]_D^{20} -20.4$ (*c* 1.04, CH_2Cl_2) (70% ee); IR (film) 3269, 1601, 1509, 1158 cm^{-1} ; ^1H NMR (400 MHz, CDCl_3) δ 7.64 (d, *J* = 8.2 Hz, 2H), 7.20-7.08 (m, 6H), 6.96-6.84 (m, 4H), 6.28 (d, *J* = 15.8 Hz, 1H), 5.98 (dd, *J* = 15.8, 6.8 Hz, 1H), 5.69 (d, *J* = 7.6 Hz, 1H), 5.07 (dd, *J* = 7.6, 6.8 Hz, 1H), 2.31 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 162.6 (d, *J* = 246.0 Hz), 162.4 (d, *J* = 245.0 Hz), 143.4, 137.8, 135.6 (d, *J* = 3 Hz), 132.3 (d, *J* = 4 Hz), 131.1, 129.6, 129.0 (d, *J* = 8 Hz), 128.3 (d, *J* = 8 Hz), 127.9, 127.4, 115.7 (d, *J* = 9 Hz), 115.5 (d, *J* = 10 Hz), 59.2, 21.5; HRMS (FT-ICRMS) calcd for $\text{C}_{22}\text{H}_{19}\text{F}_2\text{NO}_2\text{SNa}$ ($\text{M}+\text{Na}$): 422.0997, found: 422.0992.



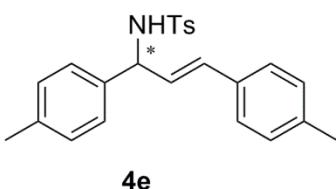
(E)-N-(1,3-bis(4-chlorophenyl)allyl)-4-methylbenzenesulfonamide (4c) (Table 2, entry 3):

White solid; yield: 61%; m.p. 170-172 °C; $[\alpha]_D^{20}$ -11.5 (*c* 1.00, CH₂Cl₂) (85% ee); IR (film) 3270, 1490, 1324, 1159 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) δ 7.62 (d, *J* = 8.2 Hz, 2H), 7.23-7.03 (m, 10H), 6.26 (d, *J* = 15.8 Hz, 1H), 6.02 (dd, *J* = 15.8, 6.6 Hz, 1H), 5.66-5.60 (m, 1H), 5.07-5.02 (m, 1H), 2.32 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 143.7, 138.1, 137.7, 134.6, 133.9, 133.8, 131.3, 129.7, 129.0, 128.8, 128.7, 128.5, 127.9, 127.4, 59.3, 21.6; HRMS (FT-ICRMS) calcd for C₂₂H₁₉Cl₂NO₂SNa (M+Na): 454.0406, found: 454.0400.



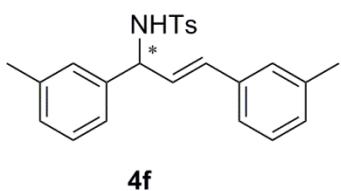
(E)-N-(1,3-bis(4-bromophenyl)allyl)-4-methylbenzenesulfonamide (4d) (Table 2, entry 4):

White solid; yield: 60%; m.p. 203-204 °C; $[\alpha]_D^{20}$ -8.3 (*c* 0.96, CH₂Cl₂) (86% ee); IR (film) 3261, 1487, 1321, 1158 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) δ 7.60 (d, *J* = 8.4 Hz, 2H), 7.40-7.30 (m, 4H), 7.14 (d, *J* = 8.0 Hz, 2H), 7.03 (d, *J* = 8.4 Hz, 4H), 6.27 (d, *J* = 15.8 Hz, 1H), 6.04 (dd, *J* = 15.8, 6.8 Hz, 1H), 5.26 (d, *J* = 7.2 Hz, 1H), 5.05 (dd, *J* = 7.2, 6.8 Hz, 1H), 2.35 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 143.8, 138.5, 137.7, 135.0, 132.0, 131.9, 131.6, 129.7, 129.0, 128.6, 128.3, 127.5, 122.2, 59.3, 21.7; HRMS (FT-ICRMS) calcd for C₂₂H₁₉Br₂NO₂SNa (M+Na): 541.9400, found: 541.9390.

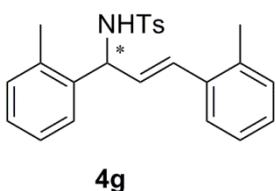


(E)-N-(1,3-di-p-tolylallyl)-4-methylbenzenesulfonamide (4e) (Table 2, entry 5): White solid;

yield: 77%; m.p. 121-123 °C; $[\alpha]_D^{20}$ -16.5 (*c* 1.00, CH₂Cl₂) (66% ee); IR (film) 3273, 1512, 1434, 1324, 1159 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) δ 7.70 (d, *J* = 8.2 Hz, 2H), 7.19 (d, *J* = 8.0 Hz, 2H), 7.19-7.05 (m, 8H), 6.35 (d, *J* = 15.8 Hz, 1H), 6.06 (dd, *J* = 15.8, 6.4 Hz, 1H), 5.11 (dd, *J* = 7.2, 6.4 Hz, 1H), 5.05 (d, *J* = 7.2 Hz, 1H), 2.38 (s, 3H), 2.36 (s, 3H), 2.35 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 143.3, 138.0, 137.9, 137.8, 137.1, 133.6, 132.0, 129.6, 129.5, 129.3, 127.5, 127.2, 126.6, 59.8, 21.6, 21.4, 21.2; HRMS (FT-ICRMS) calcd for C₂₄H₂₅NO₂SNa (M+Na): 414.1498, found: 414.1492.

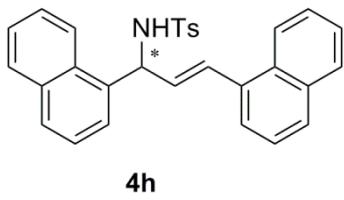


(E)-N-(1,3-di-m-tolylallyl)-4-methylbenzenesulfonamide (4f) (Table 2, entry 6): White solid; yield: 80%; m.p. 142-144 °C; $[\alpha]_D^{20}$ -19.7 (*c* 0.89, CH₂Cl₂) (62% ee); IR (film) 3274, 1605, 1326, 1159, 1092 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) δ 7.66 (d, *J* = 7.2 Hz, 2H), 7.18-7.10 (m, 4H), 7.07-6.90 (m, 6H), 6.33 (d, *J* = 16.0 Hz, 1H), 6.06 (dd, *J* = 16.0, 6.4 Hz, 1H), 5.07 (dd, *J* = 6.4, 6.4 Hz, 1H), 4.96 (brs, 1H), 2.34 (s, 3H), 2.31 (s, 3H), 2.25 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 143.4, 139.8, 138.6, 138.2, 138.1, 136.3, 132.3, 129.6, 128.8, 128.7, 128.6, 128.5, 128.3, 128.0, 127.6, 127.4, 124.3, 123.9, 60.0, 21.6, 21.5, 21.4; HRMS (FT-ICRMS) calcd for C₂₄H₂₅NO₂SNa (M+Na): 414.1498, found: 414.1492.

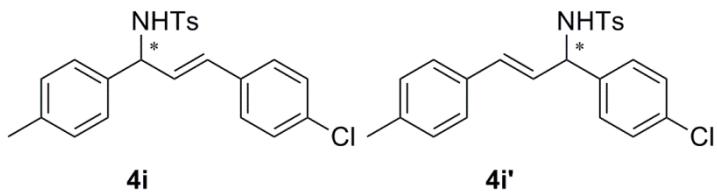


(E)-N-(1,3-di-o-tolylallyl)-4-methylbenzenesulfonamide (4g) (Table 2, entry 7): White solid; yield: 80%; m.p. 103-104 °C; $[\alpha]_D^{20}$ -7.5 (*c* 0.93, CH₂Cl₂) (22% ee); IR (film) 3275, 1603, 1489, 1459, 1325, 1159, 1092 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) δ 7.66 (d, *J* = 8.2 Hz, 2H), 7.26-7.04 (m,

10H), 6.54 (d, J = 15.6 Hz, 1H), 5.99 (dd, J = 15.6, 6.4 Hz, 1H), 5.38 (dd, J = 7.2, 6.4 Hz, 1H), 5.30 (d, J = 7.2 Hz, 1H), 2.33 (s, 6H), 2.18 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 143.3, 138.0, 137.9, 135.7, 135.6, 135.4, 130.9, 130.3, 129.9, 129.6, 129.5, 127.9, 127.8, 127.3, 127.0, 126.5, 126.1, 125.9, 56.7, 21.6, 19.8, 19.4; HRMS (FT-ICRMS) calcd for $\text{C}_{24}\text{H}_{25}\text{NO}_2\text{SNa}$ ($M+\text{Na}$): 414.1498, found: 414.1494.

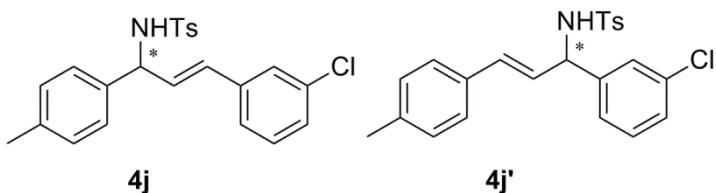


(E)-N-(1,3-di(naphthalen-1-yl)allyl)-4-methylbenzenesulfonamide (4h) (Table 2, entry 8):
 White solid; yield: 63%; m.p. 76-78 °C; $[\alpha]_D^{20} -0.8$ (c 0.80, CH_2Cl_2) (18% ee); IR (film) 3274, 1597, 1510, 1326, 1157, 1091 cm^{-1} ; ^1H NMR (400 MHz, CDCl_3) δ 8.16 (d, J = 6.8 Hz, 1H), 7.88-7.72 (m, 5H), 7.66 (d, J = 8.2 Hz, 2H), 7.54-7.30 (m, 8H), 7.24 (d, J = 15.0 Hz, 1H), 7.02 (d, J = 8.0 Hz, 2H), 6.37 (dd, J = 15.8, 6.0 Hz, 1H), 6.02 (dd, J = 7.2, 6.0 Hz, 1H), 5.44 (d, J = 7.2 Hz, 1H), 2.24 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 143.4, 137.9, 135.4, 134.2, 134.1, 133.7, 131.6, 131.2, 130.7, 129.9, 129.5, 129.1, 129.0, 128.6, 128.4, 127.4, 126.8, 126.3, 126.1, 126.0, 125.8, 125.6, 125.4, 124.1, 123.9, 123.6, 57.0, 21.5; HRMS (FT-ICRMS) calcd for $\text{C}_{30}\text{H}_{25}\text{NO}_2\text{SNa}$ ($M+\text{Na}$): 486.1498, found: 486.1493.

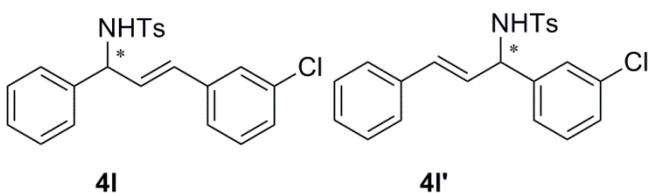


(E)-N-(3-(4-cholophenyl)-1-(p-tolyl)allyl)-4-methylbenzenesulfonamide (4i) and (E)-N-(1-(4-chlorophenyl)-3-(p-tolyl)allyl)-4-methylbenzenesulfonamide (4i') (Table 3, entry 1): White

solid; yield: 68%; m.p. 144-146 °C; $[\alpha]_D^{20}$ -17.1 (c 1.01, CH₂Cl₂) (**4i/4i'** = 4.3:1, major product **4i** 82% ee, minor product **4i'** 69% ee); IR (film) 3271, 1511, 1492, 1325, 1159, 1090 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) for major product **4i**: δ 7.64 (d, J = 7.8 Hz, 2H), 7.18 (d, J = 7.8 Hz, 2H), 7.15-7.01 (m, 8H), 6.29 (d, J = 15.8 Hz, 1H), 6.04 (dd, J = 15.8, 6.2 Hz, 1H), 5.42 (d, J = 6.8 Hz, 1H), 5.03 (dd, J = 6.8, 6.2 Hz, 1H), 2.30 (s, 3H), 2.27 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) for major product **4i**: δ 143.4, 137.9, 137.8, 136.6, 134.9, 133.5, 130.6, 130.0, 129.4, 128.7, 127.9, 127.5, 127.1, 126.6, 59.7, 21.6, 21.2; HRMS (FT-ICRMS) calcd for C₂₃H₂₂NO₂ClSNa (M+Na): 434.0952, found: 434.0944.

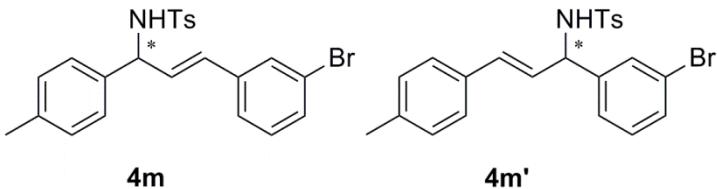


(E)-N-(3-(m-chlorophenyl)-1-(p-tolyl)allyl)-4-methylbenzenesulfonamide (4j) and (E)-N-(1-(m-chlorophenyl)-3-(p-tolyl)allyl)-4-methylbenzenesulfonamide (4j') (Table 3, entry 2): White solid; yield: 70%; m.p. 152-154 °C; $[\alpha]_D^{20}$ -18.3 (c 1.01, CH₂Cl₂) (**4j/4j'** = 8:1, major product **4j** 66% ee); IR (film) 3270, 1595, 1430, 1324, 1158, 1092 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) for major product **4j**: δ 7.66 (d, J = 7.6 Hz, 2H), 7.20-7.12 (m, 4H), 7.11-7.00 (m, 6H), 6.27 (d, J = 15.8 Hz, 1H), 6.05 (dd, J = 15.8, 6.4 Hz, 1H), 5.27-5.21 (m, 1H), 5.05 (dd, J = 6.4, 6.4 Hz, 1H), 2.33 (s, 3H), 2.28 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) for major product **4j**: δ 143.5, 138.3, 138.0, 136.5, 134.5, 130.6, 130.2, 129.8, 129.6, 127.9, 127.5, 127.1, 126.6, 125.0, 59.6, 21.6, 21.2; HRMS (FT-ICRMS) calcd for C₂₃H₂₂NO₂ClSNa (M+Na): 434.0952, found: 434.0948.

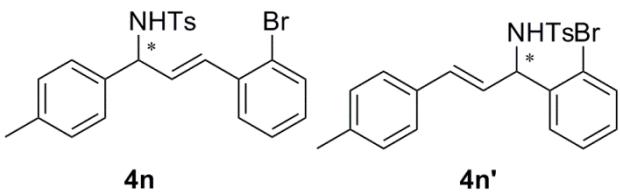


(E)-N-(3-(3-chlorophenyl)-1-phenylallyl)-4-methylbenzenesulfonamide (4l) and (E)-N-(1-(3-(3-chlorophenyl)-1-phenylallyl)-4-methylbenzenesulfonamide (4l')

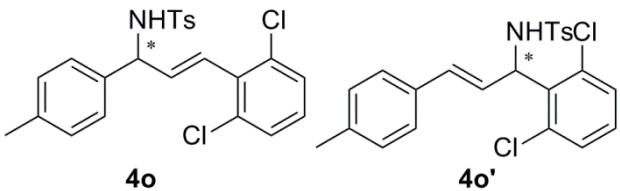
chlorophenyl)-3-phenylallyl)-4-methylbenzenesulfonamide (4l'**) (Table 3, entry 4):** White solid; yield: 56%; m.p. 123-125 °C; $[\alpha]_D^{20}$ -17.9 (*c* 0.98, CH₂Cl₂) (**4l/4l'** = 3.5:1, major product **4l** 67% ee, minor product **4l'** 58% ee); IR (film) 3269, 1596, 1325, 1159, 1092 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) for major product **4l**: δ 7.65 (d, *J* = 8.2 Hz, 2H), 7.26-7.20 (m, 3H), 7.20-7.10 (m, 6H), 7.10-7.06 (m, 1H), 7.05-7.00 (m, 1H), 6.28 (d, *J* = 16.0 Hz, 1H), 6.06 (dd, *J* = 16.0, 6.8 Hz, 1H), 5.34 (d, *J* = 7.4 Hz, 1H), 5.08 (dd, *J* = 7.4, 6.8 Hz, 1H), 2.32 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) for major product **4l**: δ 143.6, 139.5, 138.2, 137.9, 134.6, 130.8, 130.0, 129.9, 129.7, 129.0, 128.2, 127.9, 127.5, 127.2, 126.6, 125.0, 59.8, 21.6; HRMS (FT-ICRMS) calcd for C₂₂H₂₀NO₂ClSNa (M+Na): 420.0796, found: 420.0791.



(E)-N-(3-(3-bromophenyl)-1-(p-tolyl)allyl)-4-methylbenzenesulfonamide (4m**) and (E)-N-(1-(3-bromophenyl)-3-(p-tolyl)allyl)-4-methylbenzenesulfonamide (**4m'**) (Table 3, entry 5):** White solid; yield: 87%; m.p. 153-155 °C; $[\alpha]_D^{20}$ -17.9 (*c* 0.98, CH₂Cl₂) (**4m/4m'** = 6.7:1, major product **4m** 76% ee, minor product **4m'** 70% ee); IR (film) 3269, 1594, 1324, 1158, 1092 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) for major product **4m**: δ 7.66 (d, *J* = 8.2 Hz, 2H), 7.30 (d, *J* = 7.2 Hz, 1H), 7.26-7.20 (m, 1H), 7.14 (d, *J* = 8.2 Hz, 2H), 7.11-7.00 (m, 6H), 6.26 (d, *J* = 15.8 Hz, 1H), 6.04 (dd, *J* = 15.8, 6.6 Hz, 1H), 5.36-5.20 (m, 1H), 5.05 (dd, *J* = 6.8, 6.6 Hz, 1H), 2.33 (s, 3H), 2.28 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) for major product **4m**: δ 143.5, 138.5, 138.0, 136.5, 130.8, 130.5, 130.2, 130.1, 129.6, 129.5, 127.5, 127.1, 125.4, 122.7, 59.6, 21.6, 21.2; HRMS (FT-ICRMS) calcd for C₂₃H₂₂NO₂BrSNa (M+Na): 478.0447, found: 478.0441.

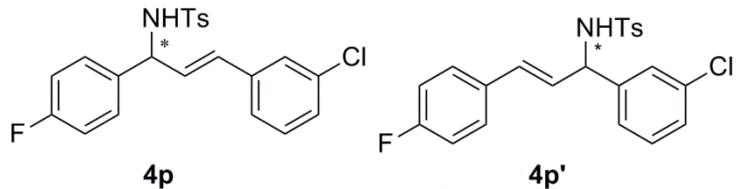


(E)-N-(3-(o-bromophenyl)-1-(p-tolyl)allyl)-4-methylbenzenesulfonamide (4n) and (E)-N-(1-(o-bromophenyl)-3-(p-tolyl)allyl)-4-methylbenzenesulfonamide (4n') (Table 3, entry 6): White solid; yield: 87%; m.p. 115-116 °C; $[\alpha]_D^{20}$ 1.3 (*c* 1.10, CH₂Cl₂) (**4n/4n'** = 3:1, major product **4n** 13% ee, minor product **4n'** 94% ee); IR (film) 3271, 1600, 1512, 1436, 1325, 1159, 1092 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) for major product **4n**: δ 7.69 (d, *J* = 7.6 Hz, 2H), 7.48 (d, *J* = 7.8 Hz, 1H), 7.42-7.01 (m, 9H), 6.69 (d, *J* = 15.8 Hz, 1H), 6.01 (dd, *J* = 15.8, 7.0 Hz, 1H), 5.46-5.41 (m, 1H), 5.06 (dd, *J* = 7.0, 6.8 Hz, 1H), 2.32-2.24 (m, 6H); ¹³C NMR (100 MHz, CDCl₃) for major product **4n**: δ 143.4, 138.0, 136.6, 136.2, 132.9, 131.5, 130.7, 129.6, 129.4, 129.3, 129.2, 127.4, 127.3, 127.1, 126.6, 123.8, 59.7, 21.5, 21.2; HRMS (FT-ICRMS) calcd for C₂₃H₂₂NO₂BrSNa (M+Na): 478.0447, found: 478.0441.

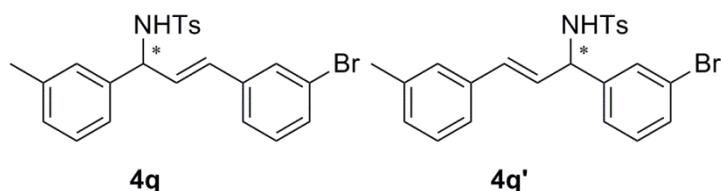


(E)-N-(3-(2,6-dichlorophenyl)-1-(p-tolyl)allyl)-4-methylbenzenesulfonamide (4o) and (E)-N-(1-(2,6-dichlorophenyl)-3-(p-tolyl)allyl)-4-methylbenzenesulfonamide (4o') (Table 3, entry 7): White solid; yield: 16%; m.p. 164-165 °C; $[\alpha]_D^{20}$ 1.8 (*c* 0.71, CH₂Cl₂) (**4o/4o'** > 20:1, major product **4o** 60% ee); IR (film) 3274, 1557, 1428, 1324, 1159, 1092 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) for major product **4o**: δ 7.76 (d, *J* = 8.2 Hz, 2H), 7.33-7.28 (m, 2H), 7.24 (d, *J* = 8.0 Hz, 2H), 7.18 (d, *J* = 8.0 Hz, 2H), 7.14-7.05 (m, 3H), 6.50 (d, *J* = 16.2 Hz, 1H), 6.28 (dd, *J* = 16.2, 6.0 Hz, 1H), 5.18 (dd, *J* = 7.2, 6.0 Hz, 1H), 5.01 (d, *J* = 7.2 Hz, 1H), 2.41 (s, 3H), 2.35 (s, 3H); ¹³C NMR (100 MHz,

CDCl_3) for major product **4o**: δ 143.5, 138.0, 137.9, 137.1, 136.5, 134.6, 133.9, 129.7, 129.6, 128.6, 127.5, 127.4, 125.8, 59.6, 21.7, 21.3; HRMS (FT-ICRMS) calcd for $\text{C}_{23}\text{H}_{21}\text{NO}_2\text{Cl}_2\text{SNa}$ ($\text{M}+\text{Na}$): 468.0562, found: 468.0556.

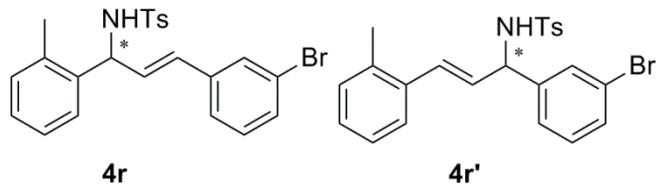


(*E*)-N-(3-(3-chlorophenyl)-1-(4-fluorophenyl)allyl)-4-methylbenzenesulfonamide (**4p**) and (*E*-N-(1-(3-chlorophenyl)-1-(3-fluorophenyl)allyl)-4-methylbenzenesulfonamide (**4p'**) (Table 3, entry 8): White solid; yield: 63%; m.p. 131-133 °C; $[\alpha]_D^{20} -17.9$ (c 0.98, CH_2Cl_2) (**4p/4p'** = 3:1, major product **4p** 59% ee, minor product **4p'** 73% ee); IR (film) 3268, 1598, 1509, 1327, 1226, 1159, 1092 cm^{-1} ; ^1H NMR (400 MHz, CDCl_3) for major product **4p**: δ 7.61 (d, J = 8.6 Hz, 2H), 7.27-7.00 (m, 8H), 6.98-6.87 (m, 2H), 6.25 (d, J = 15.8 Hz, 1H), 6.04 (dd, J = 15.8, 6.8 Hz, 1H), 5.41 (d, J = 7.4 Hz, 1H), 5.09 (dd, J = 7.4, 6.8 Hz, 1H), 2.34 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) for major product **4p**: δ 162.5 (d, J = 245 Hz), 143.8, 137.9 (d, J = 17 Hz), 135.3 (d, J = 3 Hz), 134.6, 131.1, 129.9, 129.7, 129.6 (d, J = 6 Hz), 129.0 (d, J = 8 Hz), 128.1, 127.5, 127.4, 126.6, 125.0, 115.8 (d, J = 17 Hz), 59.1, 21.6; HRMS (FT-ICRMS) calcd for $\text{C}_{22}\text{H}_{19}\text{NO}_2\text{ClFSNa}$ ($\text{M}+\text{Na}$): 438.0707, found: 438.0695.



(*E*)-N-(3-(3-bromophenyl)-1-(m-tolyl)allyl)-4-methylbenzenesulfonamide (**4q**) and (*E*-N-(1-

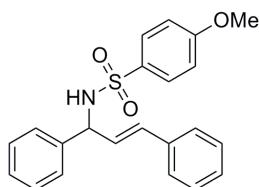
(3-bromophenyl)-3-(m-tolyl)allyl)-4-methylbenzenesulfonamide (4q'**) (Table 3, entry 9):** White solid; yield: 64%; m.p. 136-138 °C; $[\alpha]_D^{20}$ -19.2 (*c* 0.96, CH₂Cl₂) (**4q/4q'** = 6.9:1, major product **4q** 76%, minor product **4q'** 53% ee); IR (film) 3270, 1594, 1562, 1426, 1327, 1159, 1092 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) for major product **4q**: δ 7.65 (d, *J* = 8.2 Hz, 2H), 7.31 (d, *J* = 7.2 Hz, 1H), 7.26-7.22 (m, 1H), 7.16-7.06 (m, 5H), 7.01 (d, *J* = 7.4 Hz, 1H), 6.96 (d, *J* = 7.6 Hz, 1H), 6.92 (s, 1H), 6.27 (d, *J* = 15.8 Hz, 1H), 6.04 (dd, *J* = 15.8, 6.6 Hz, 1H), 5.37 (d, *J* = 7.2 Hz, 1H), 5.05 (dd, *J* = 7.2, 6.6 Hz, 1H), 2.33 (s, 3H), 2.22 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) for major product **4q**: δ 143.5, 139.3, 138.7, 138.5, 137.9, 130.8, 130.5, 130.1, 129.6, 129.5, 128.9, 127.9, 127.5, 125.4, 124.2, 122.7, 59.8, 21.6, 21.5; HRMS (FT-ICRMS) calcd for C₂₃H₂₂NO₂BrSNa (M+Na): 478.0447, found: 478.0441.



(E)-N-(3-(3-bromophenyl)-1-(o-tolyl)allyl)-4-methylbenzenesulfonamide (4r**) and (E)-N-(1-(3-bromophenyl)-3-(o-tolyl)allyl)-4-methylbenzenesulfonamide (**4r'**) (Table 3, entry 10):** White solid; yield: 52%; m.p. 137-138 °C; $[\alpha]_D^{20}$ -20.1 (*c* 0.81, CH₂Cl₂) (**4r/4r'** = 3.8:1, major product **4r** 71% ee, minor product **4r'** 55% ee); IR (film) 3273, 1594, 1324, 1158, 1092 cm⁻¹; ¹H NMR (400 MHz, CDCl₃) for major product **4r**: δ 7.63 (d, *J* = 8.0 Hz, 2H), 7.30 (d, *J* = 7.6 Hz, 1H), 7.23-7.01 (m, 9H), 6.18 (d, *J* = 15.8 Hz, 1H), 6.07 (dd, *J* = 15.8, 5.2 Hz, 1H), 5.38-5.25 (m, 2H), 2.33 (s, 3H), 2.28 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) for major product **4r**: δ 143.5, 138.5, 137.9, 137.4, 135.7, 131.0, 130.8, 130.4, 130.1, 129.8, 129.6, 129.5, 128.0, 137.4, 127.0, 126.6, 125.4, 122.7, 56.3, 21.6, 19.4; HRMS (FT-ICRMS) calcd for C₂₃H₂₂NO₂BrSNa (M+Na): 478.0447, found: 478.0441.

The chromatograms for determination of the enantiomeric excess

Table 1, entry 12



HPLC Conditions: Column: ChiralcelOD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 0.5 mL/min; Detection: UV 254 nm

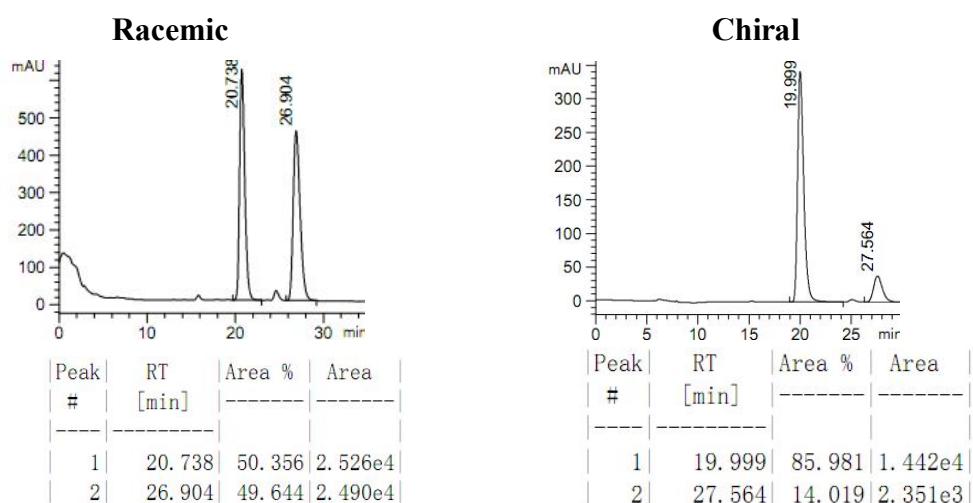
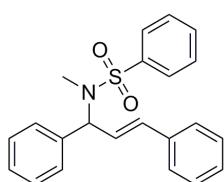


Table 1, entry 13



HPLC Conditions: Column: ChiralcelOD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 0.5 mL/min; Detection: UV 254 nm

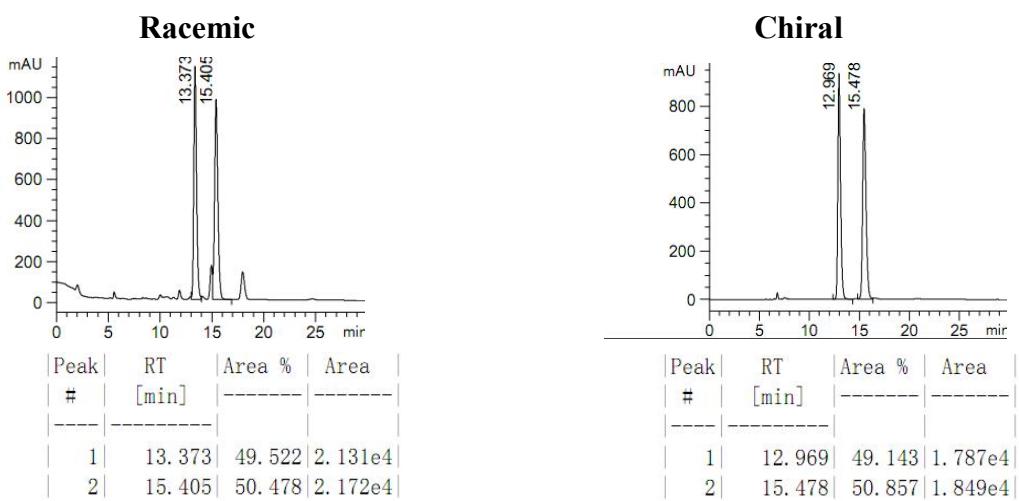
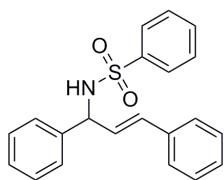
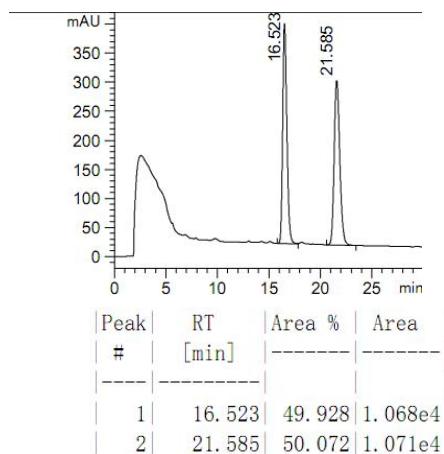
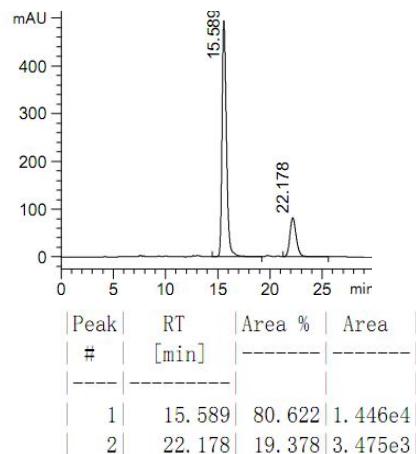
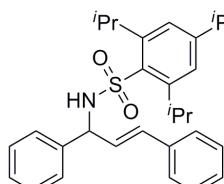


Table 1, entry 14

HPLC Conditions: Column: ChiralcelOD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 0.5 mL/min; Detection: UV 254 nm

Racemic**Chiral****Table 1, entry 15**

HPLC Conditions: Column: ChiralpakAD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (98/2); Flow rate: 1 mL/min; Detection: UV 254 nm

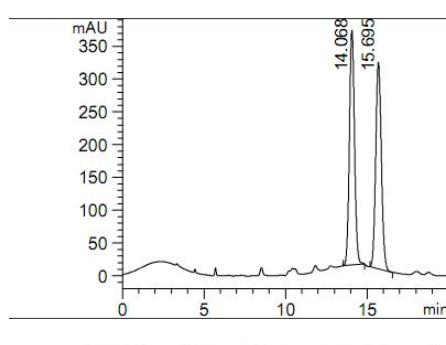
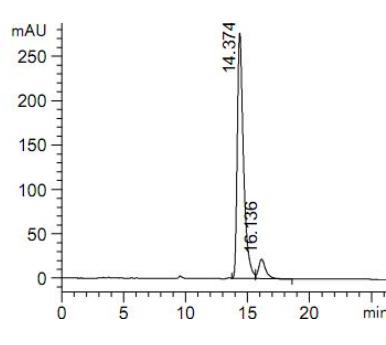
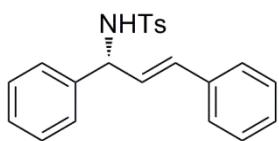
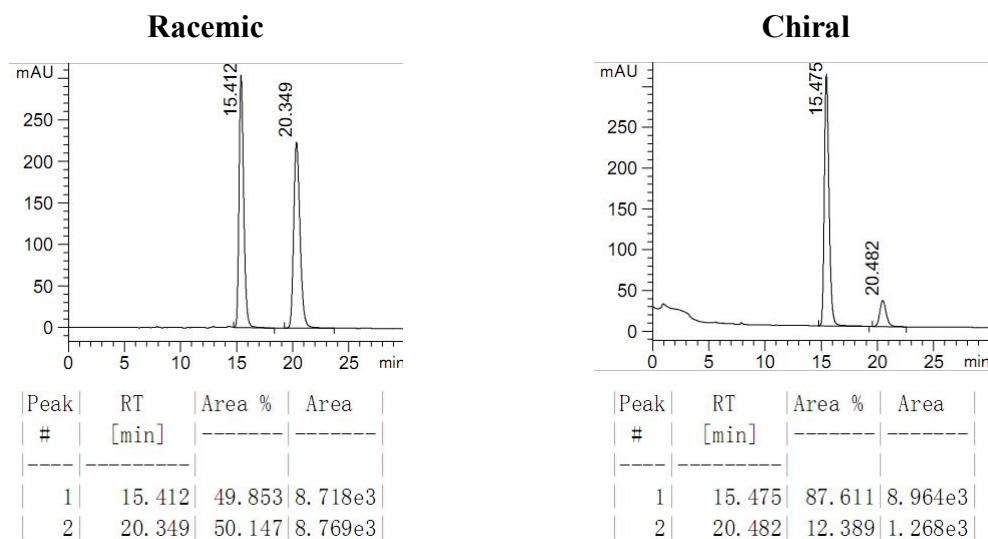
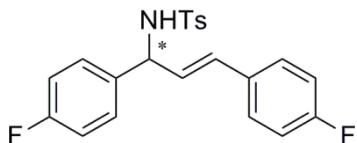
Racemic**Chiral**

Table 2, entry 1

HPLC Conditions: Column: Chiralcel OD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 0.5 mL/min; Detection: UV 254 nm

**Table 2, entry 2**

HPLC Conditions: Column: Chiralcel OD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 0.5 mL/min; Detection: UV 254 nm

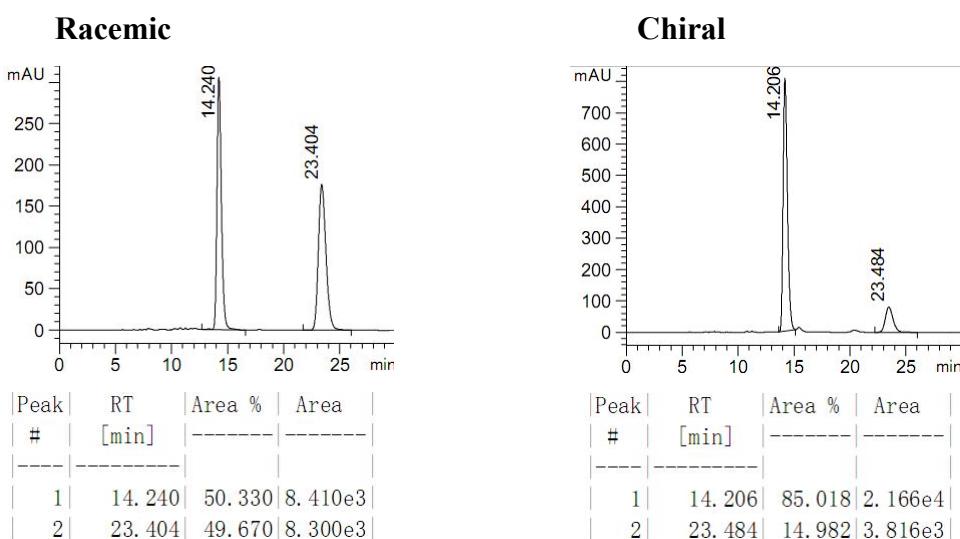
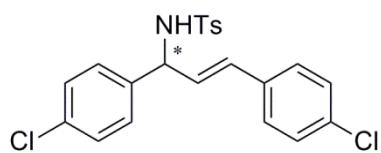
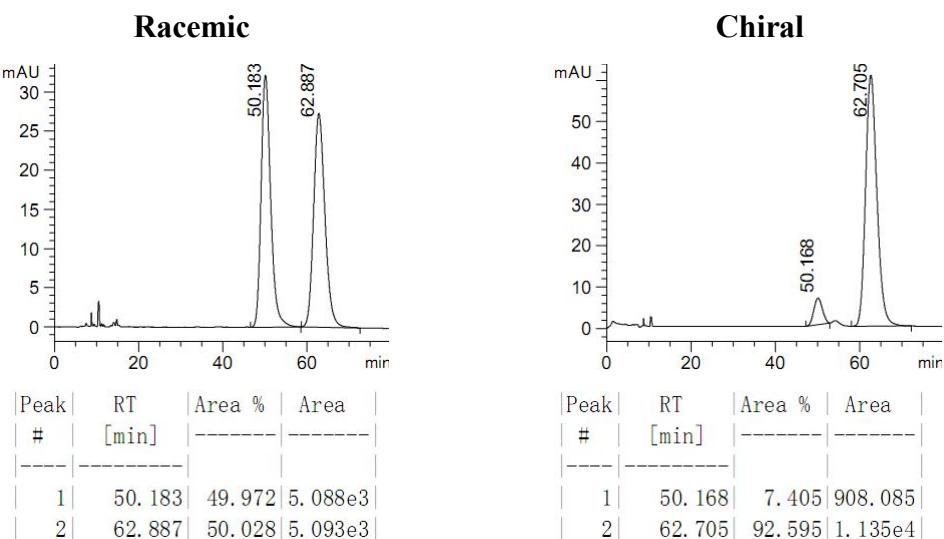
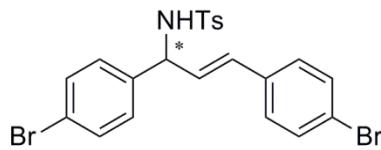


Table 2, entry 3

HPLC Conditions: Column: Chiralcel AS-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 0.5 mL/min; Detection: UV 254 nm

**Table 2, entry 4**

HPLC Conditions: Column: Chiraldak AD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 0.5 mL/min; Detection: UV 254 nm

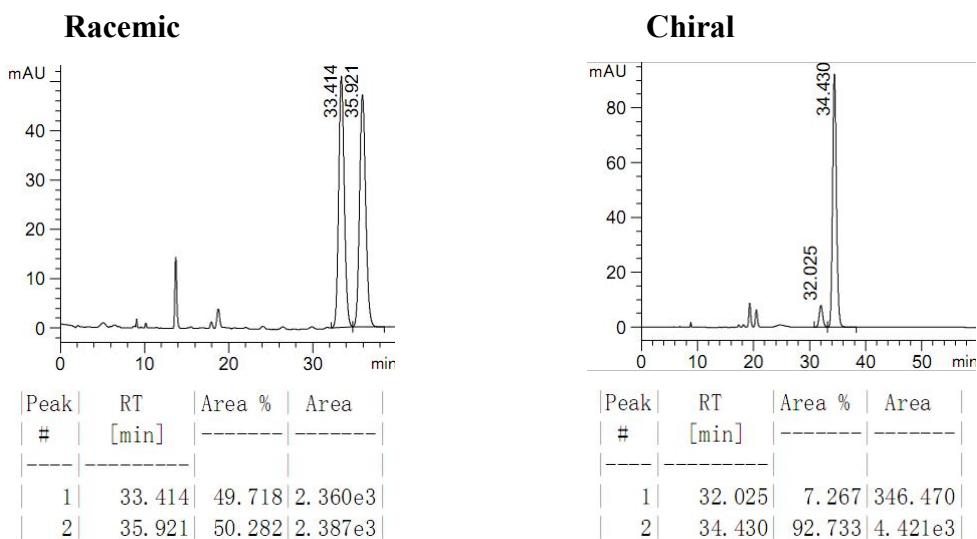
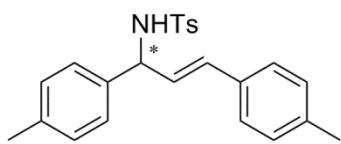
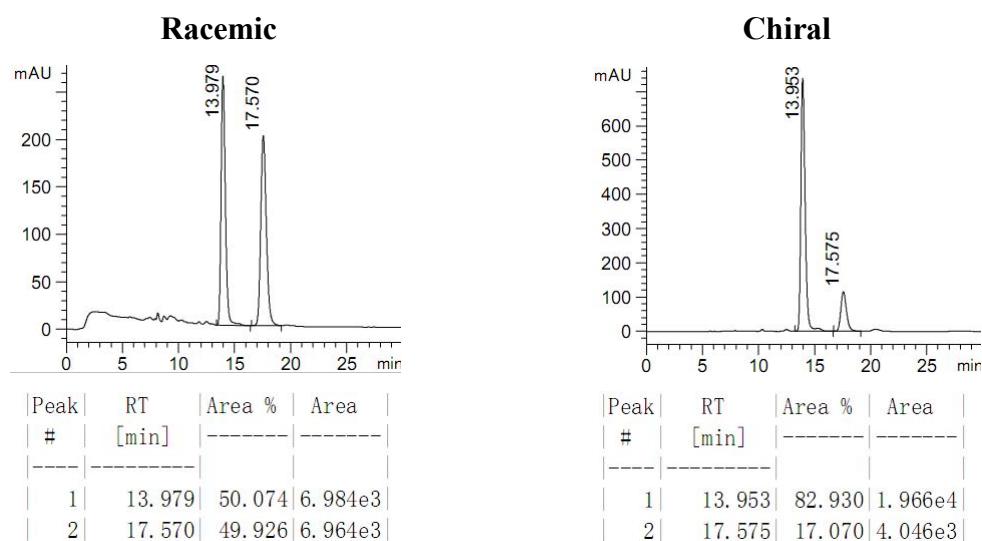
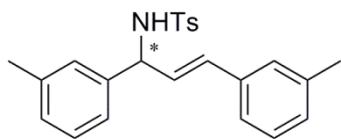


Table 2, entry 5

HPLC Conditions: Column: Chiralcel OD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 0.5 mL/min; Detection: UV 254 nm

**Table 2, entry 6**

HPLC Conditions: Column: Chiraldak AD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 1.0 mL/min; Detection: UV 254 nm

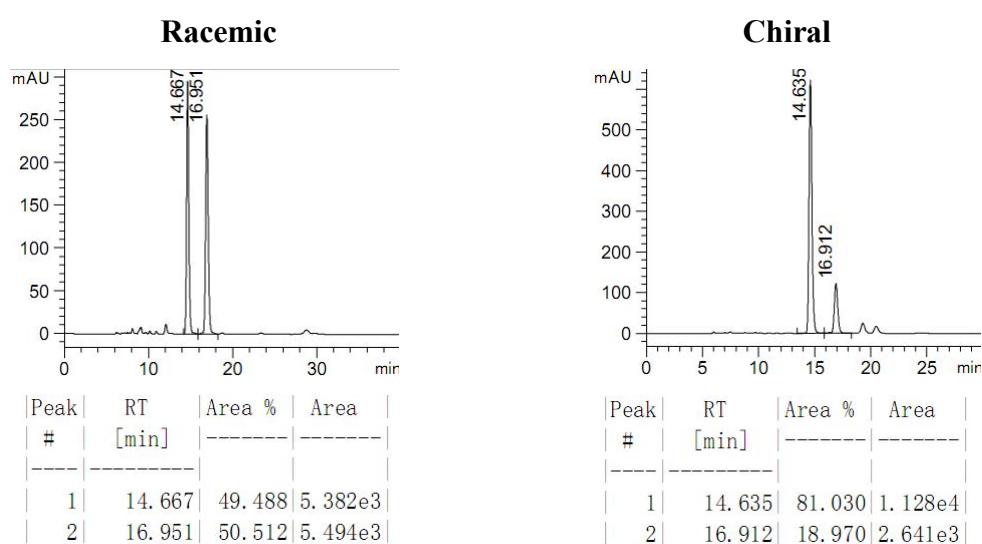
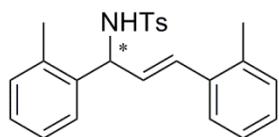
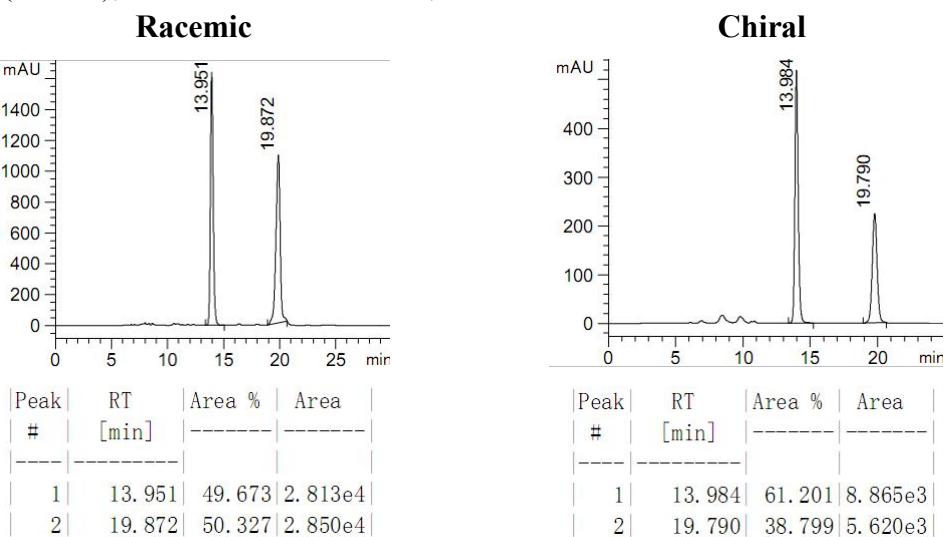
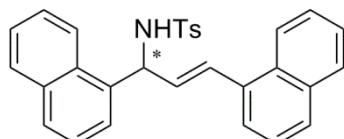


Table 2, entry 7

HPLC Conditions: Column: Chiraldak AD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 1.0 mL/min; Detection: UV 254 nm

**Table 2, entry 8**

HPLC Conditions: Column: Chiraldak AD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 0.5 mL/min; Detection: UV 310 nm

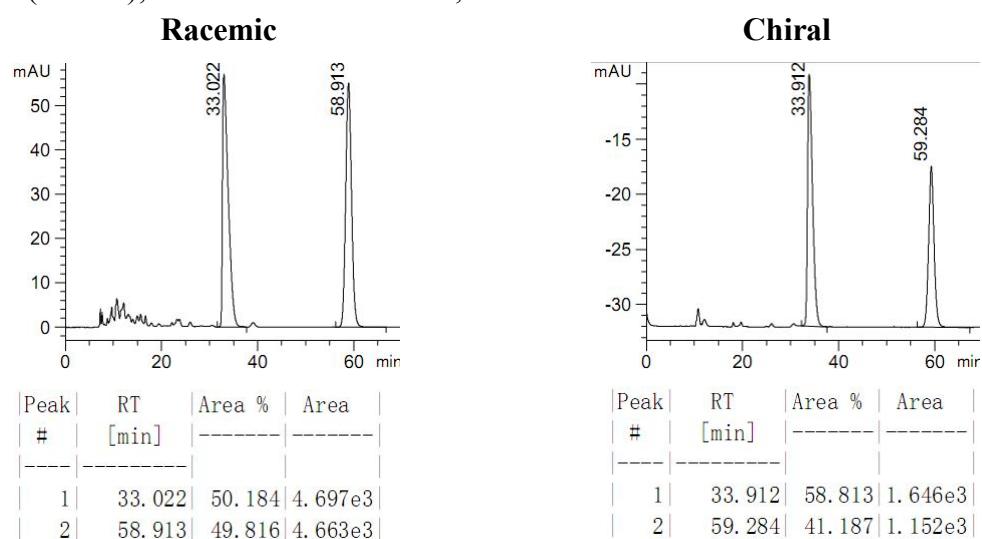
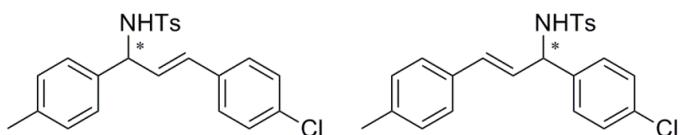
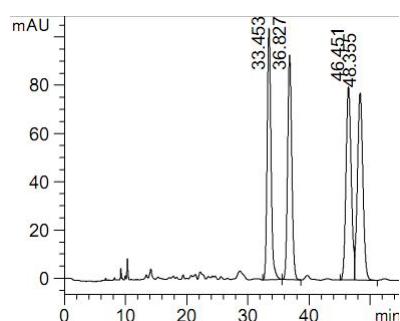
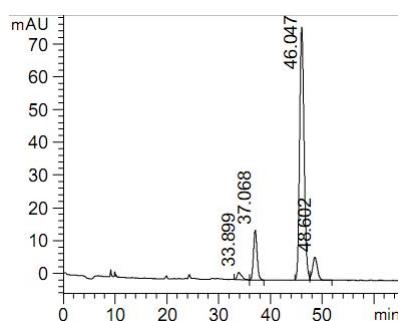


Table 3, entry 1

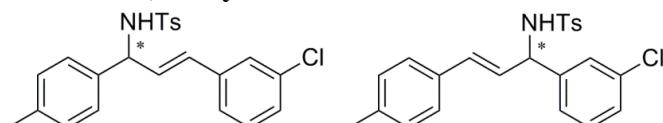
HPLC Conditions: Column: Chiraldak AD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (85/15); Flow rate: 0.5 mL/min; Detection: UV 254 nm

Racemic

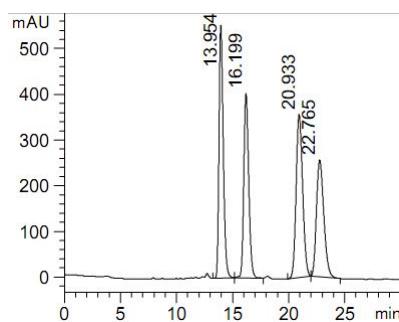
Peak #	RT [min]	Area %	Area
1	33.453	24.441	4.522e3
2	36.827	23.728	4.390e3
3	46.451	25.637	4.743e3
4	48.355	26.194	4.846e3

Chiral

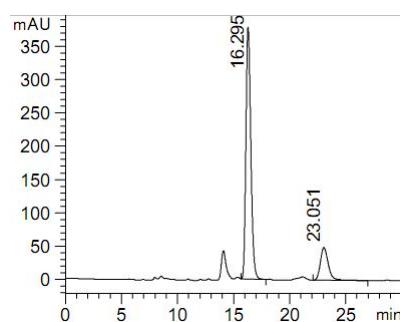
Peak #	RT [min]	Area %	Area
1	33.899	2.277	135.935
2	37.068	12.230	730.178
3	46.047	77.855	4.648e3
4	48.602	7.638	456.052

Table 3, entry 2

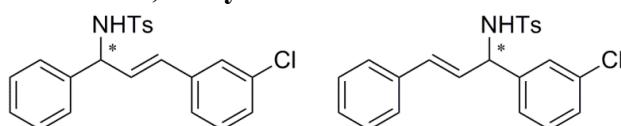
HPLC Conditions: Column: Chiralcel OD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 0.5 mL/min; Detection: UV 254 nm

Racemic

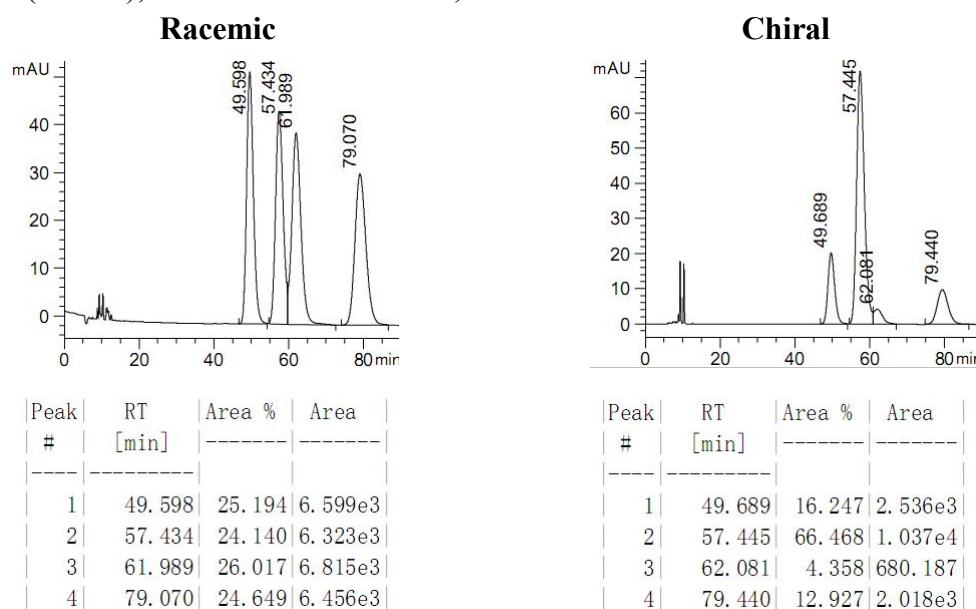
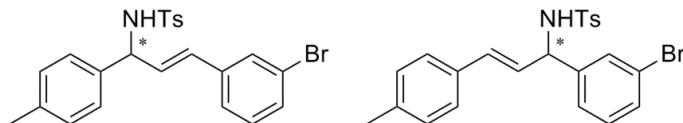
Peak #	RT [min]	Area %	Area
1	13.954	27.621	1.463e4
2	16.199	23.247	1.232e4
3	20.933	27.133	1.438e4
4	22.765	21.999	1.166e4

Chiral

Peak #	RT [min]	Area %	Area
1	16.295	82.950	1.150e4
2	23.051	17.050	2.364e3

Table 3, entry 4

HPLC Conditions: Column: Chiralcel AS-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 0.5 mL/min; Detection: UV 254 nm

**Table 3, entry 5**

HPLC Conditions: Column: Chiralcel OD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 0.5 mL/min; Detection: UV 254 nm

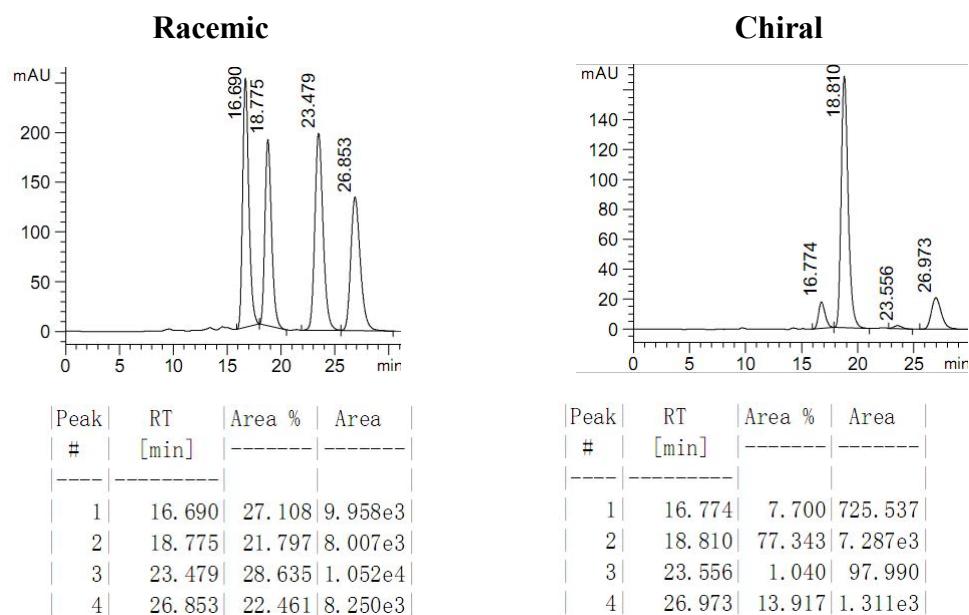
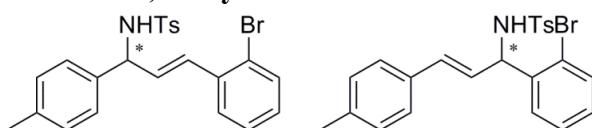
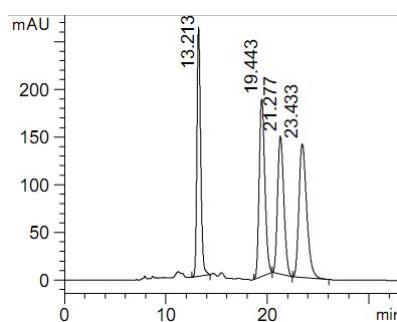
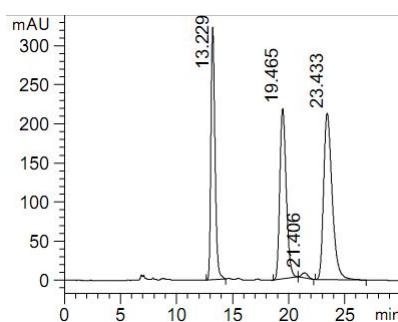


Table 3, entry 6

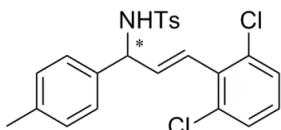
HPLC Conditions: Column: Chiralcel OD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 0.5 mL/min; Detection: UV 254 nm

Racemic

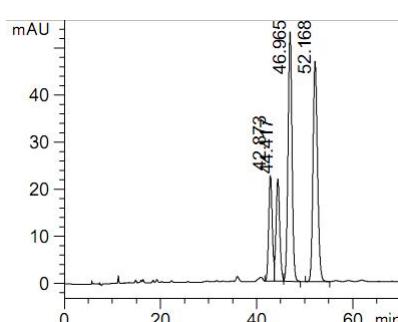
Peak #	RT [min]	Area %	Area
1	13.213	24.087	6.624e3
2	19.443	26.200	7.205e3
3	21.277	22.587	6.211e3
4	23.433	27.125	7.459e3

Chiral

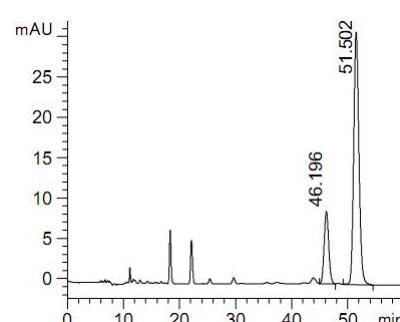
Peak #	RT [min]	Area %	Area
1	13.229	28.685	8.141e3
2	19.465	30.533	8.666e3
3	21.406	0.863	245.034
4	23.433	39.919	1.133e4

Table 3, entry 7

HPLC Conditions: Column: Chiralpak AD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (90/10); Flow rate: 0.5 mL/min; Detection: UV 254 nm

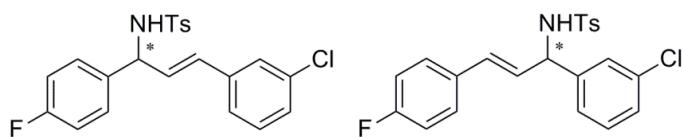
Racemic

Peak #	RT [min]	Area %	Area
1	42.873	13.523	1.141e3
2	44.417	14.001	1.182e3
3	46.965	36.200	3.055e3
4	52.168	36.276	3.062e3

Chiral

Peak #	RT [min]	Area %	Area
1	46.196	19.971	505.167
2	51.502	80.029	2.024e3

Table 3, entry 8



HPLC Conditions: Column: Chiralcel OD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (85/15); Flow rate: 0.5 mL/min; Detection: UV 254 nm

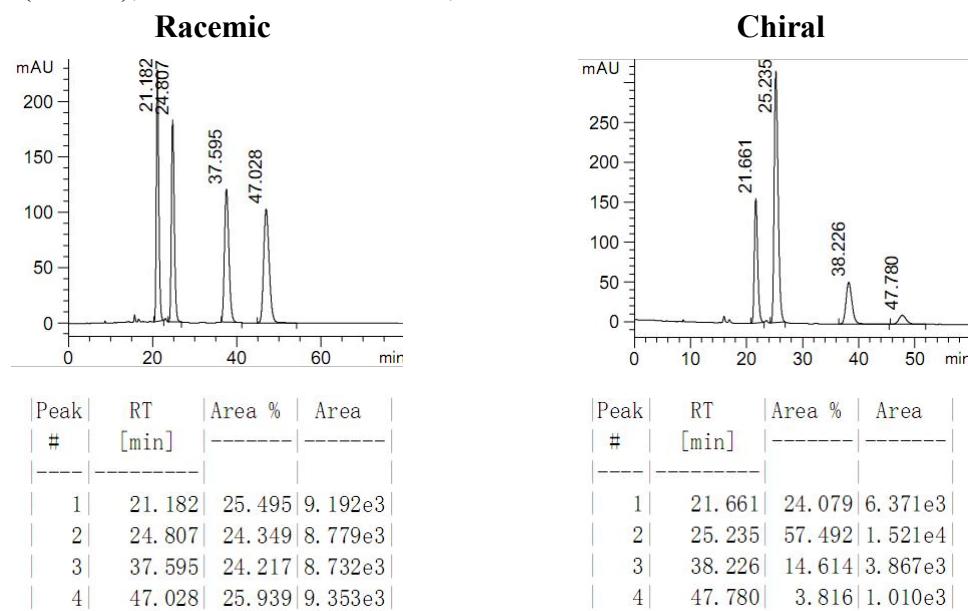
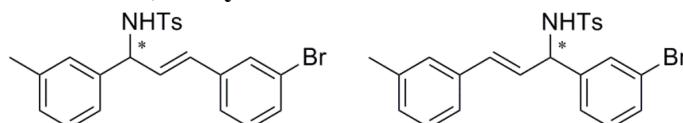


Table 3, entry 9



HPLC Conditions: Column: Chiraldak AD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (90/10); Flow rate: 0.5 mL/min; Detection: UV 254 nm

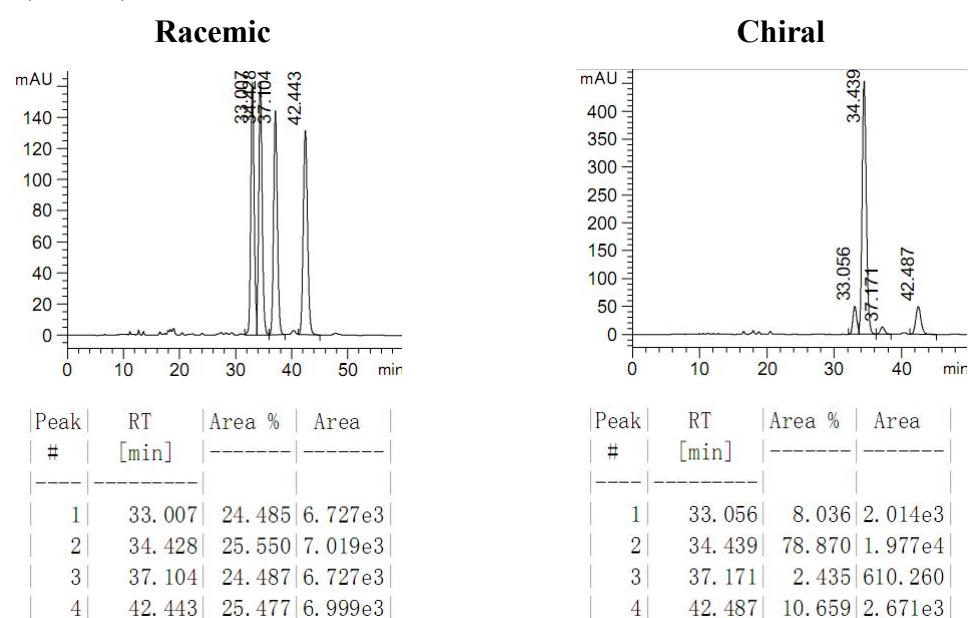
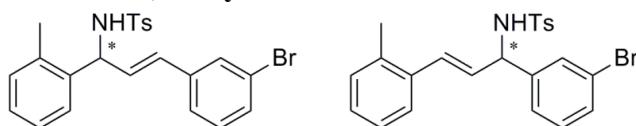
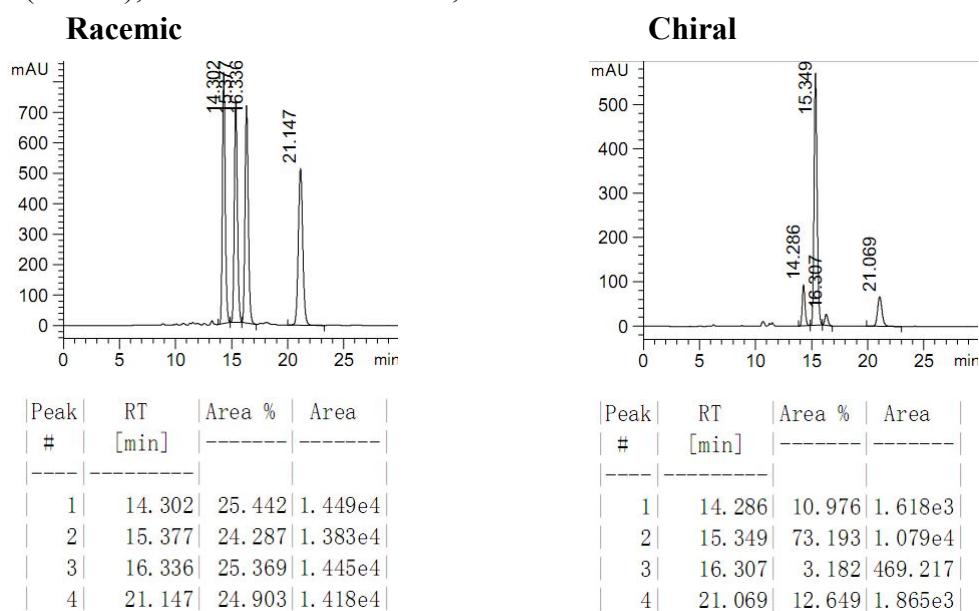
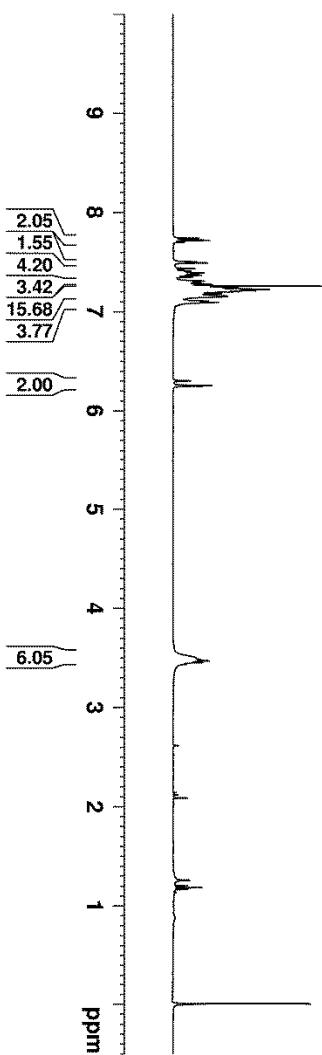
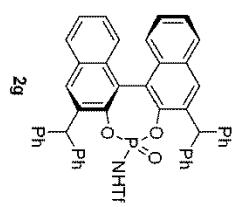


Table 3, entry 10



HPLC Conditions: Column: Chiralpak AD-H, Daicel Chemical Industries, Ltd., Eluent: Hexanes/IPA (75/25); Flow rate: 0.5 mL/min; Detection: UV 254 nm





```

NAME          zhuangminyang-4
EXPNO        29
PROCNO       1
Date_        20130911
Time         7.47
INSTRUM      spect
PROBHD      5 mm PADDL 13C
PULPROG     zg30
TD          32768
SOLVENT      CDCl3
NS           16
DS            0
SWH         12019.230 Hz
FIDRES     0.366798 Hz
AQ          1.363198 sec
RG           203
DW           41.600 usec
DE           6.500 usec
TE           300.0 K
D1          2.0000000 sec
TD0           1

===== CHANNEL f1 =====
NUC1          1H
PI           12.60 usec
SI            65536
SF          400.1300178 MHz
WDDM        EM
SSB          0
LB           0.50 Hz
GB           0
PC           1.00

```

ZMY-d1.pn

ppm

—1.5238

Current Data Parameters
NAME 310
EXPNO 1280
PROCNO 1

F2 - Acquisition Parameters

DATE_ 201308
TIME 14:48
INSTRUM spect
PROBOD 5 mm BBR BB-1H
PULPROG zg30
TD 32768
SOLVENT Acetone
NS 17
DS 0
SWH 36231.883 Hz
FIDRES 1.10509 Hz
AQ 0.45284 sec
RG 32768
DW 13.800 usec
DE 6.00 usec
TE 0.0 K
D1 2.0000000 sec
d11 0.0300000 sec
M1 0.0000000 sec
M2 0.0150000 sec
M3 0.0000000 sec

===== CHANNEL f1 =====

NUC1 31P
P1 11.40 usec
PL1 7.00 dB
SF01 121.5023628 MHz

===== CHANNEL f2 =====

NUC2 1H

P1P2 80.00 usec
PL2 2.50 dB
PL12 18.00 dB
SF02 300.1337198 MHz

F2 - Processing parameters

S1 32768
SF 121.494722 MHz
WM 4K
SSB 0
LB 10.00 Hz
GB 0
PC 0.20

10 MB plot parameters

CX 23.00 cm

CY 4.00 cm

CP 212.680 ppm

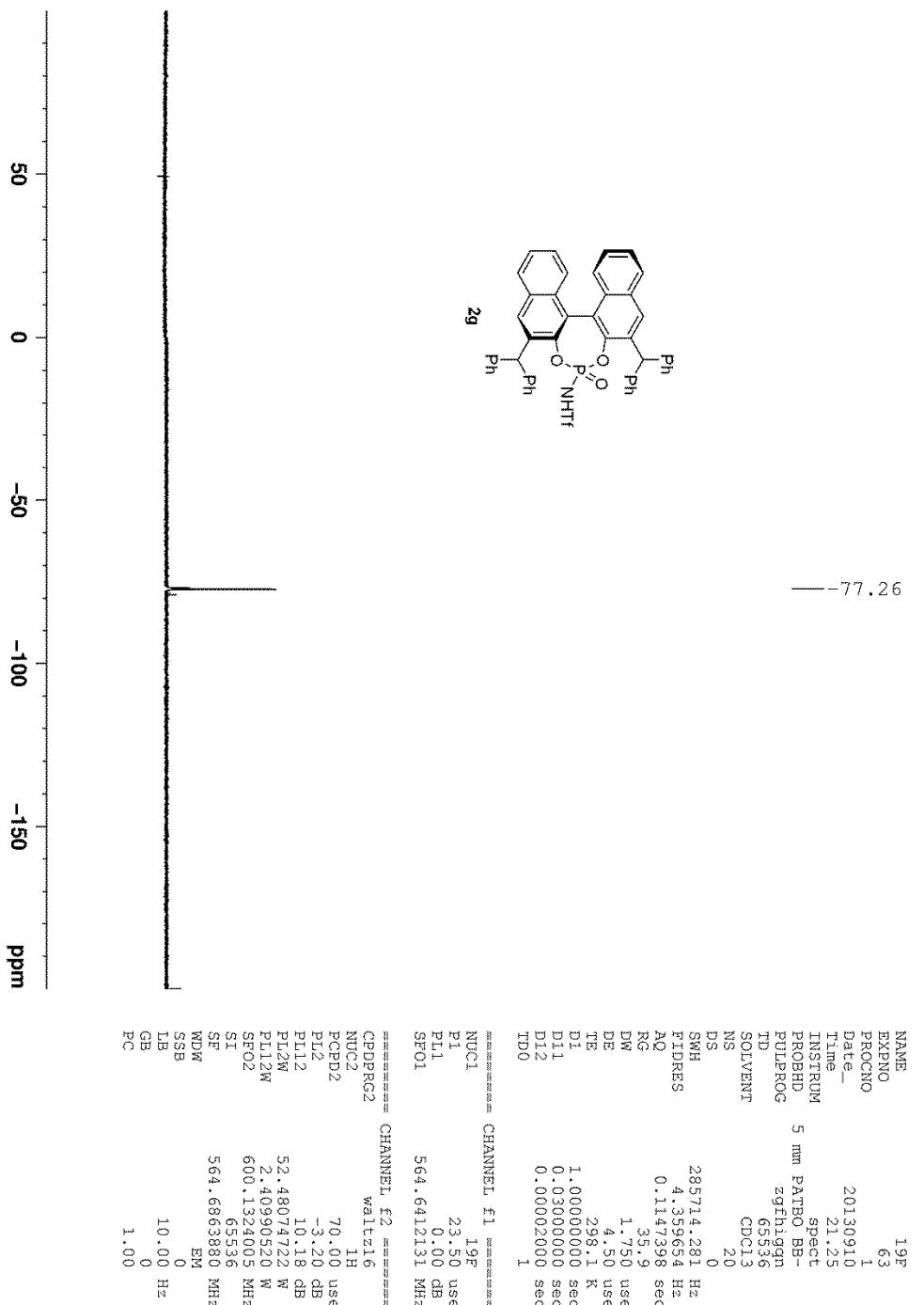
F1 28776.30 Hz

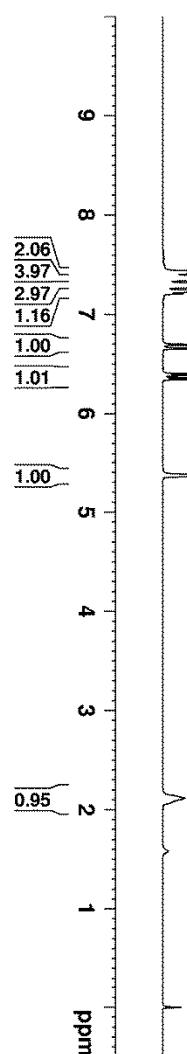
F2P -50.000 ppm

F2 6074.73 Hz

PPCM 11.39825 ppm/cm

HZCM 1384.82751 Hz/cm



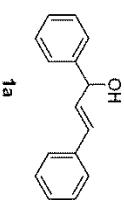


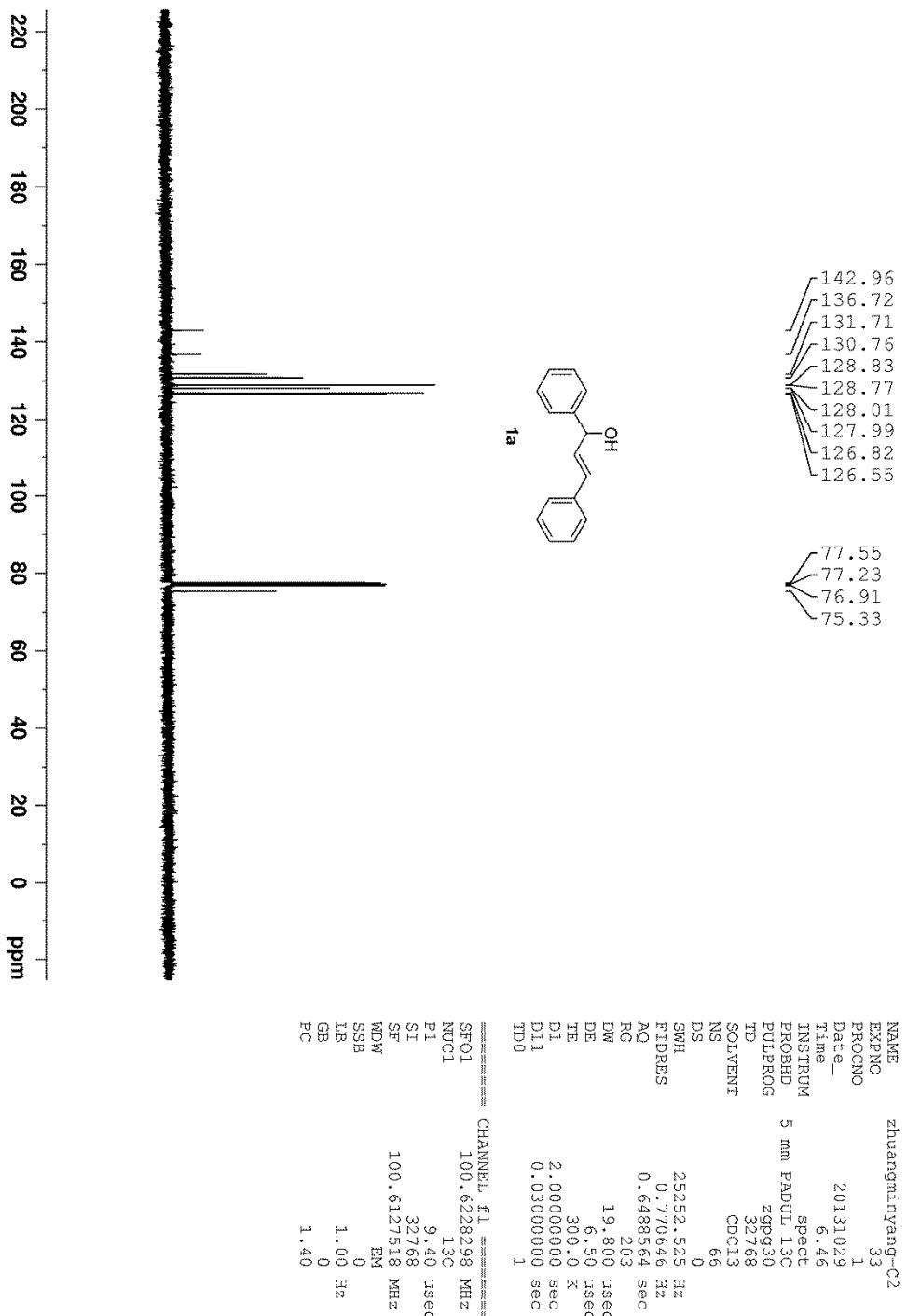
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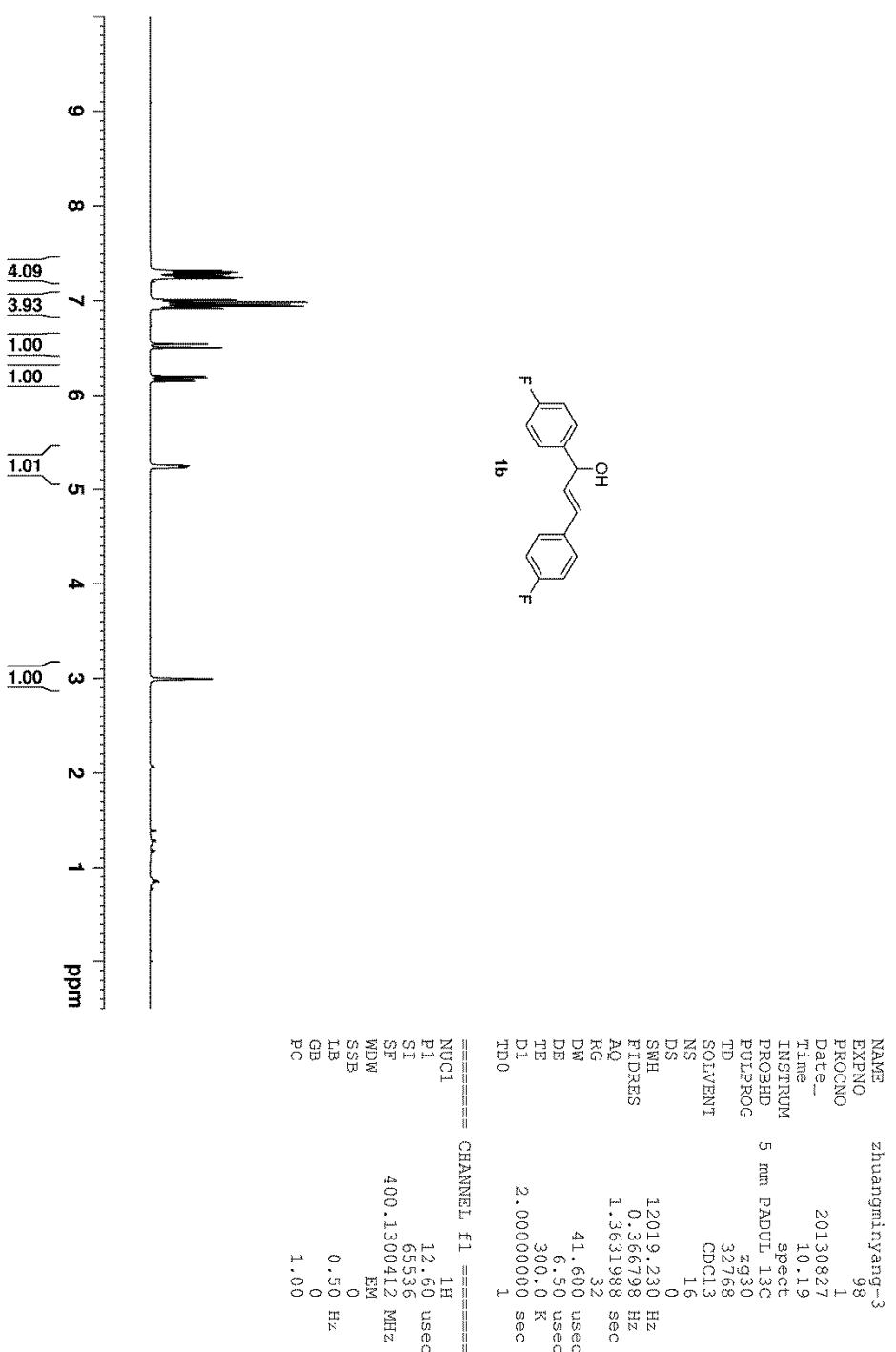
NAME          zhuangminyang-5
EXPNO        16
PROCNO       1
Date_        20131029
Time         6.45
INSTRUM     spect
PROBHD      5 mm PABUL 13C
PULPROG     zg30
TD           32768
SOLVENT      CDCl3
NS            5
DS           0
SWH         12019.230 Hz
FIDRES     0.366798 Hz
AQ          1.3631988 sec
RG           101
DW           41.600 usec
DE           6.50 usec
TE           300.0 K
TEUNIF      2.0000000 sec
DI          1
TDO          1

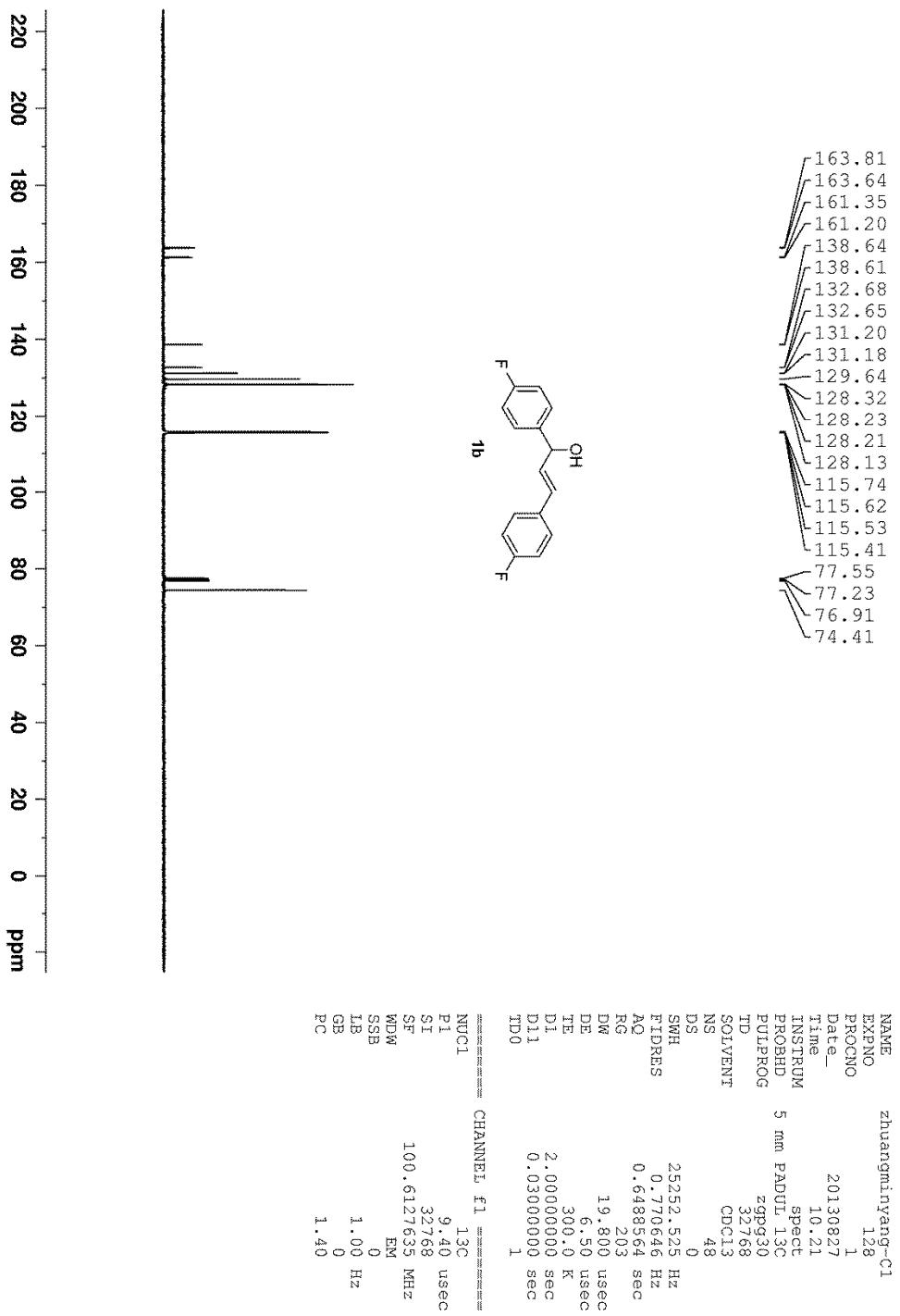
===== CHANNEL f1 =====
SFO1        400.1320007 MHz
NUC1         1H
P1          12.60 usec
SI           65536
SF          400.1300196 MHz
WDW         EM
SSB          0
LB          0.50 Hz
GB          0
PC          1.00

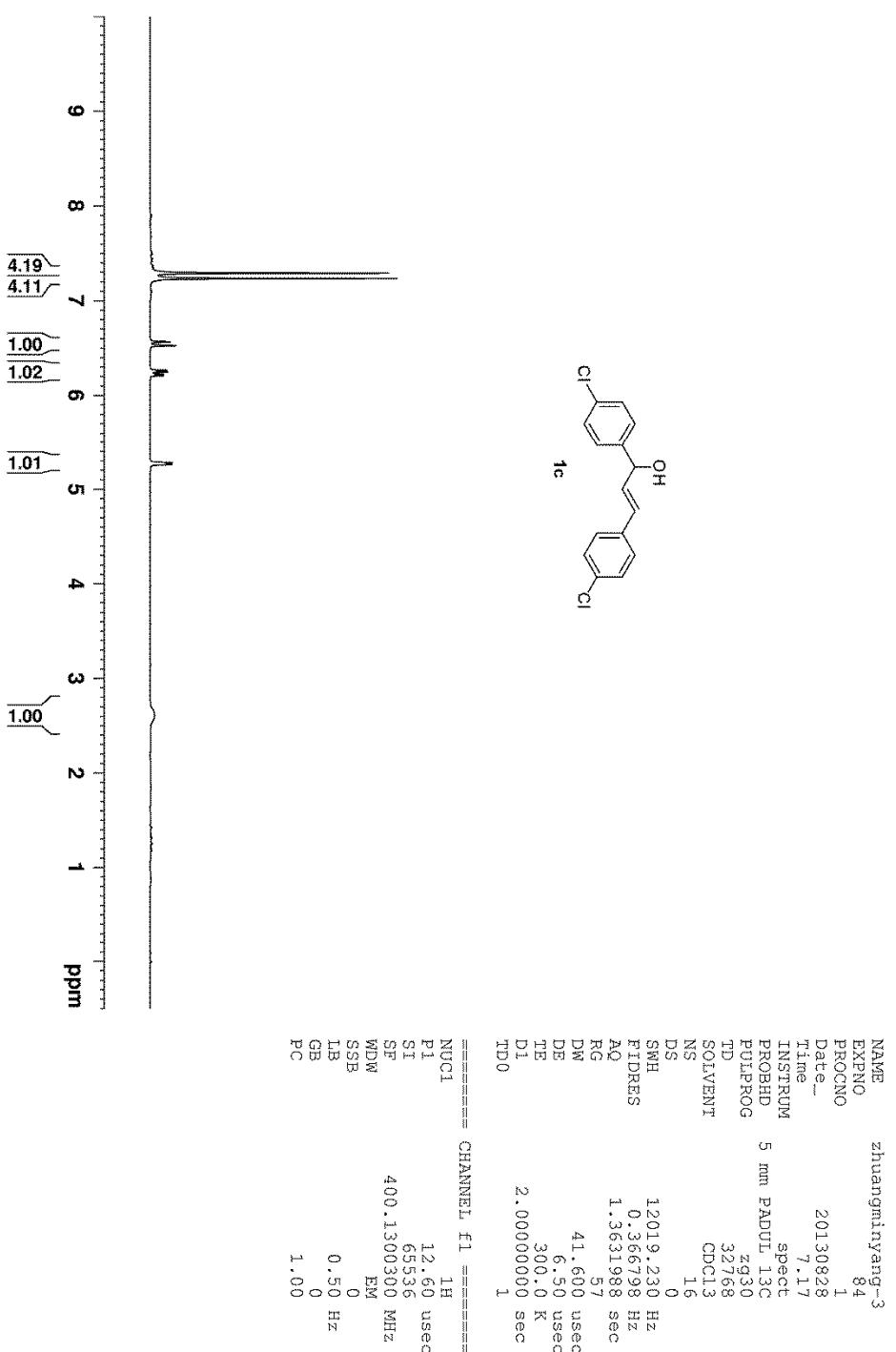
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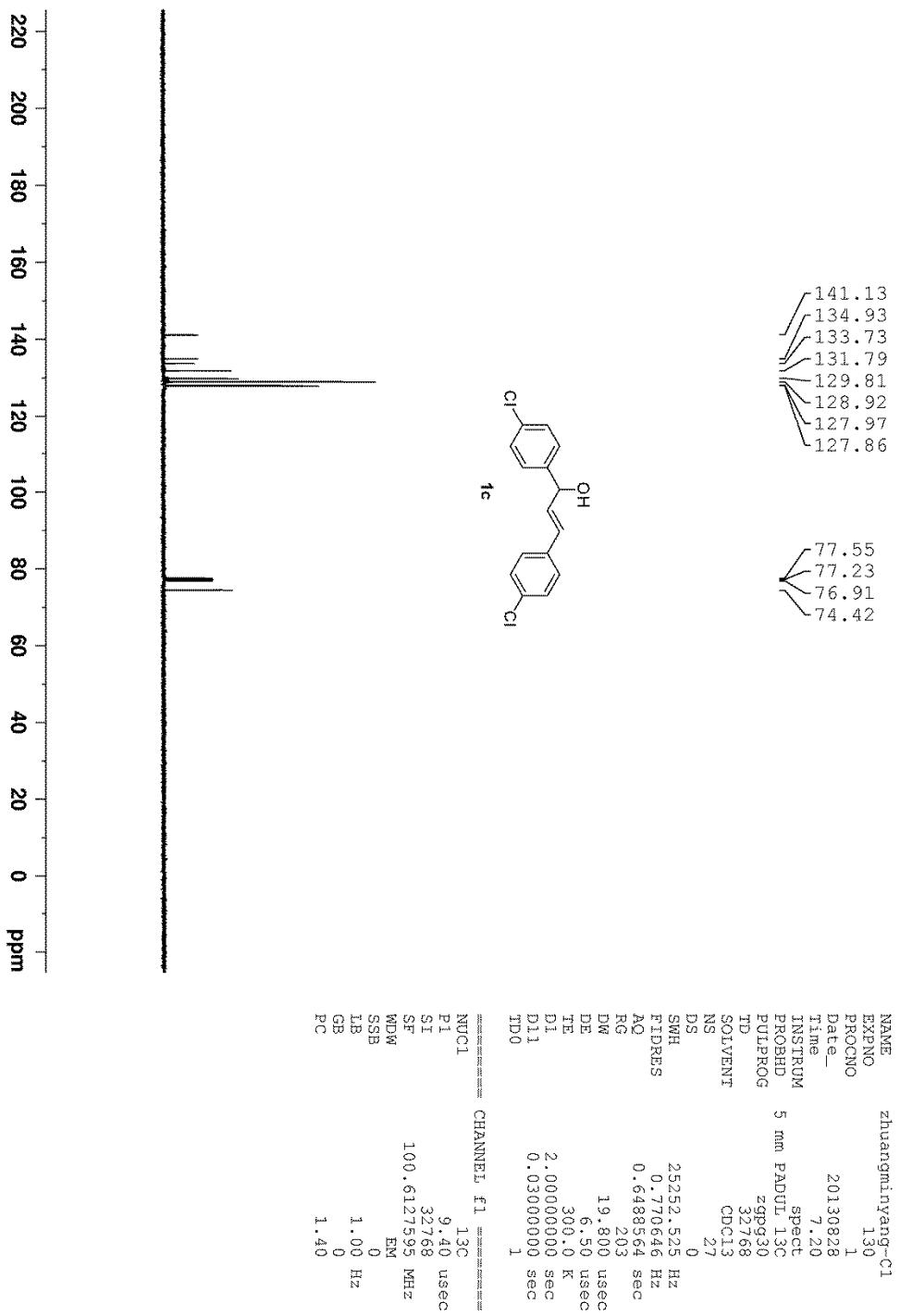


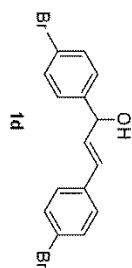
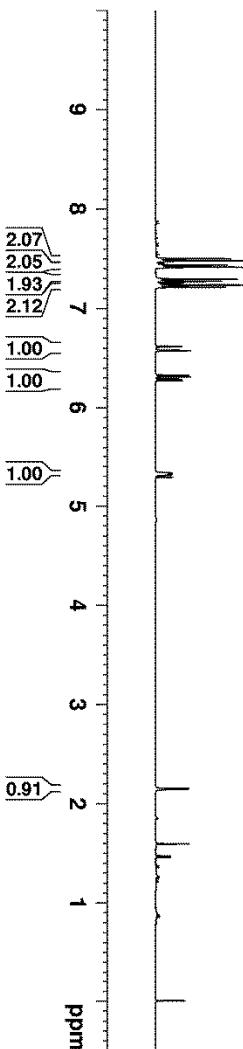










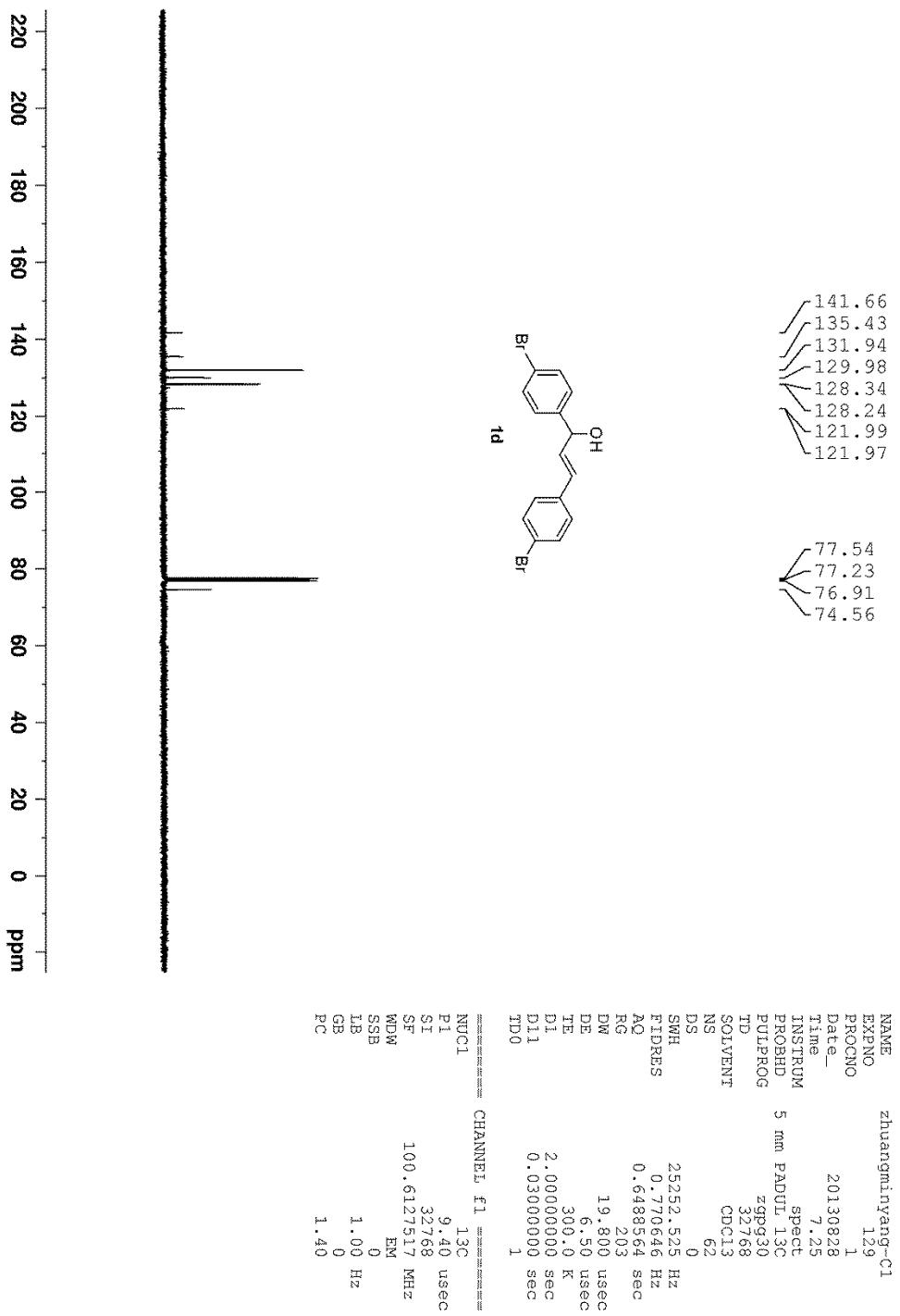


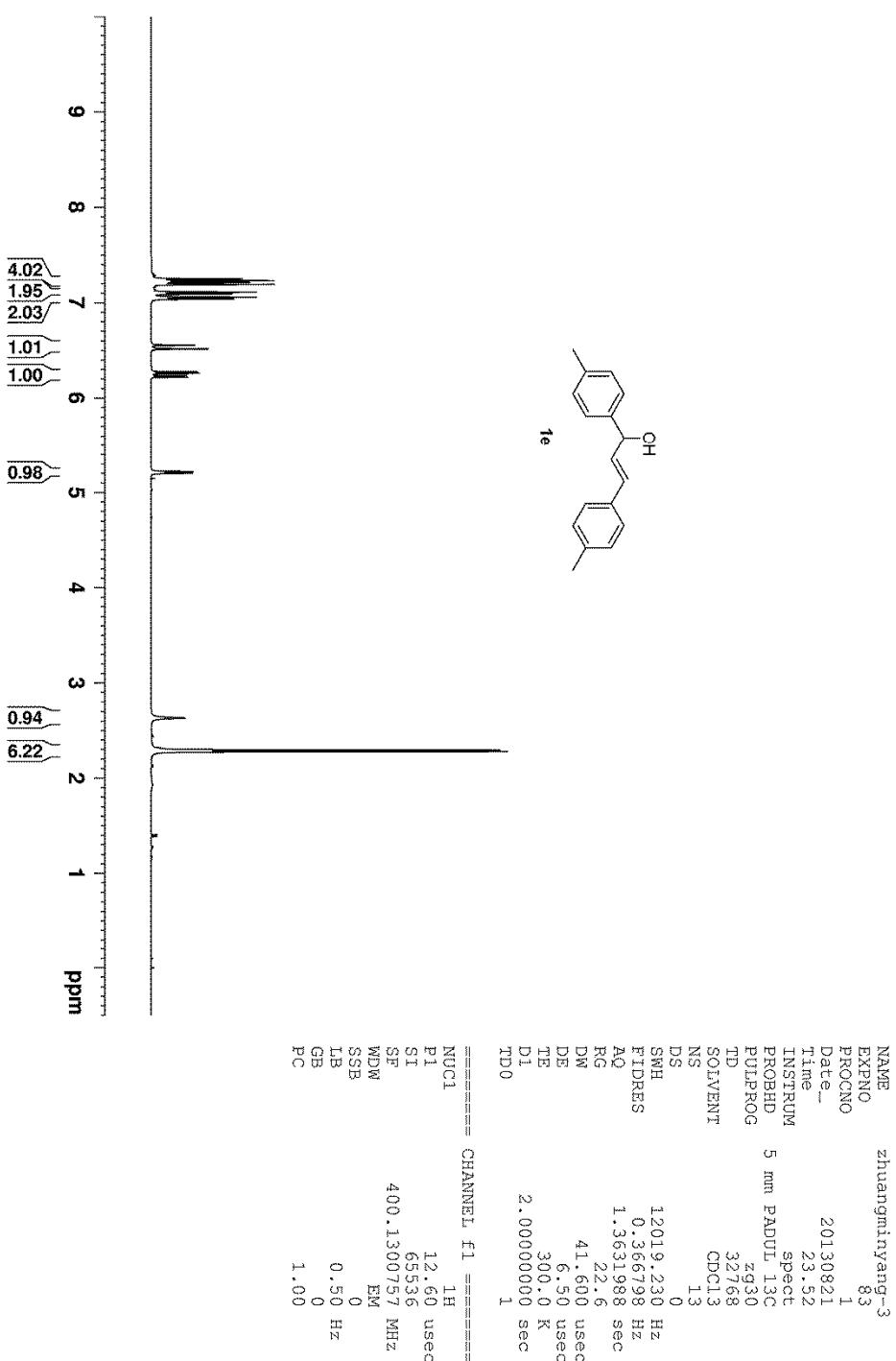
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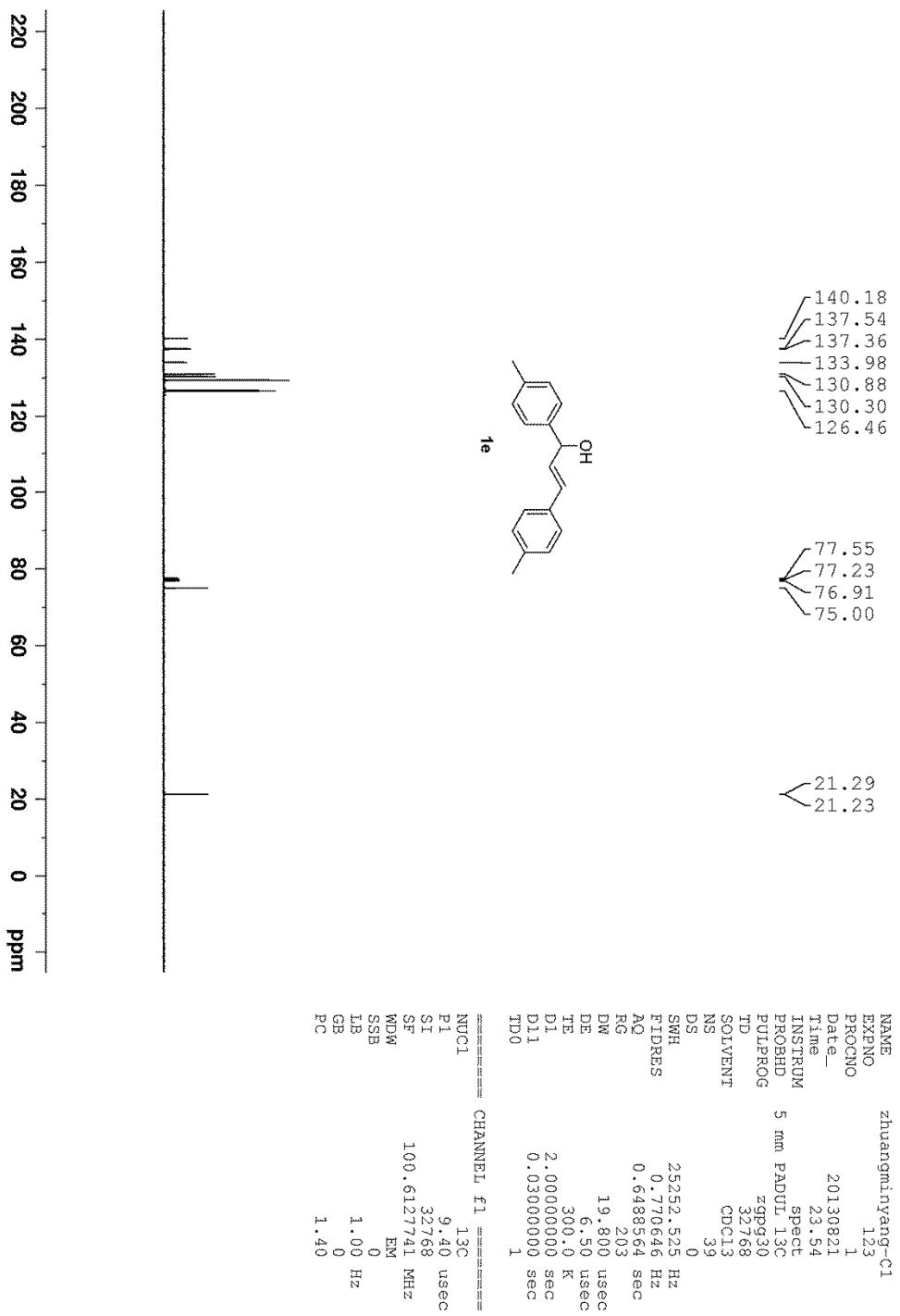
=====
NAME          zhuangminyang-3
EXPNO         100
PROCNO        1
Date_        20130828
Time       7.22
INSTRUM      spect
PROBHD      5 mm PADUL 13C
PULPROG     zg30
TD        32768
SOLVENT      CDCl3
NS           16
DS            0
SWH       0.36798 Hz
FIDRES     1.3631988 sec
AQ        203
RG          41.600 usec
DW        6.50 usec
DE        30.0 K
TE        2.0000000 sec
D1             1
TDD0

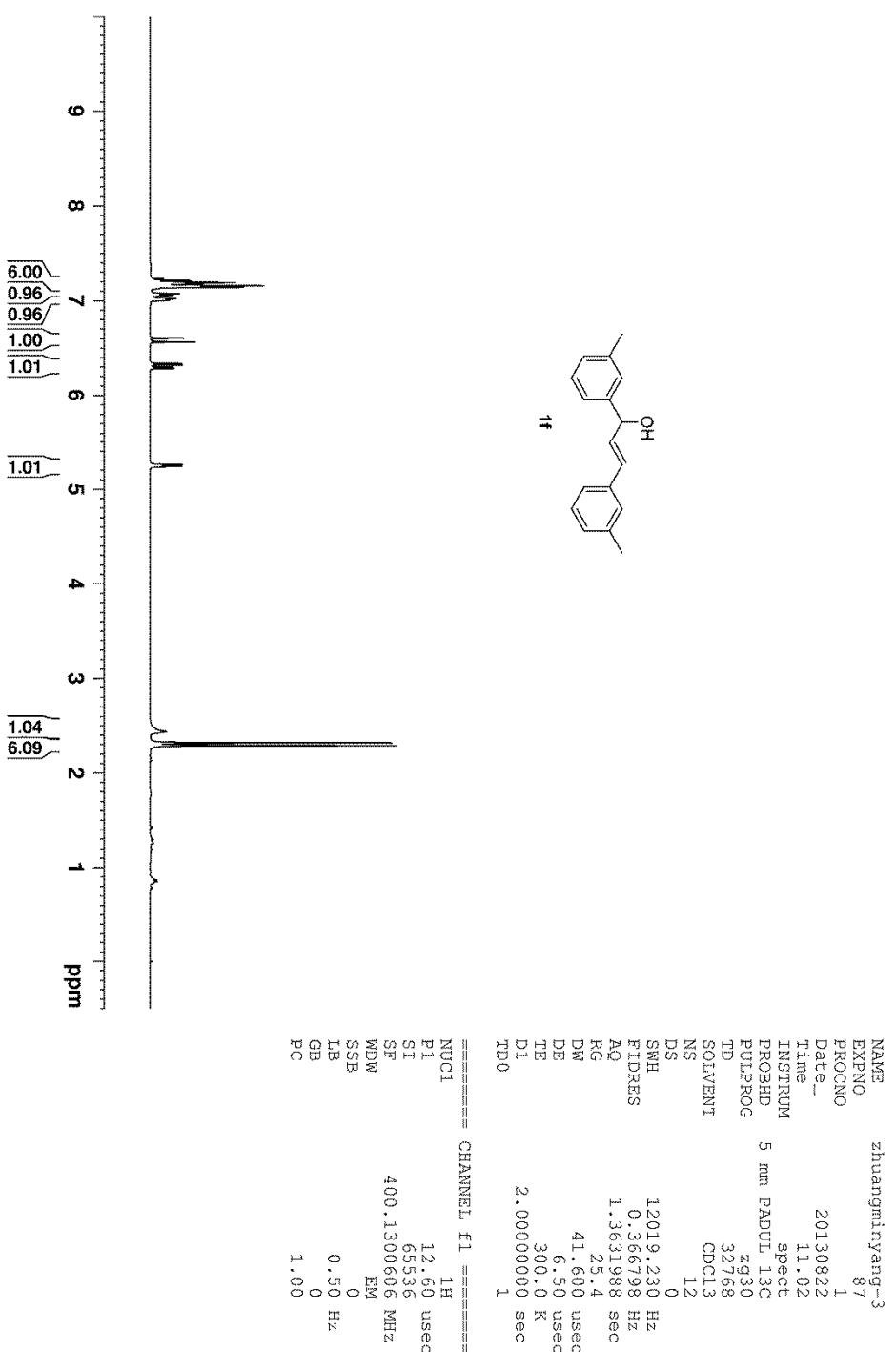
===== CHANNEL f1 =====
NUC1          1H
PI        12.60 usec
SI        65536
SF        400.1300189 MHz
WDW
SSB          0
LB        0.50 Hz
GB        0
PC        1.00

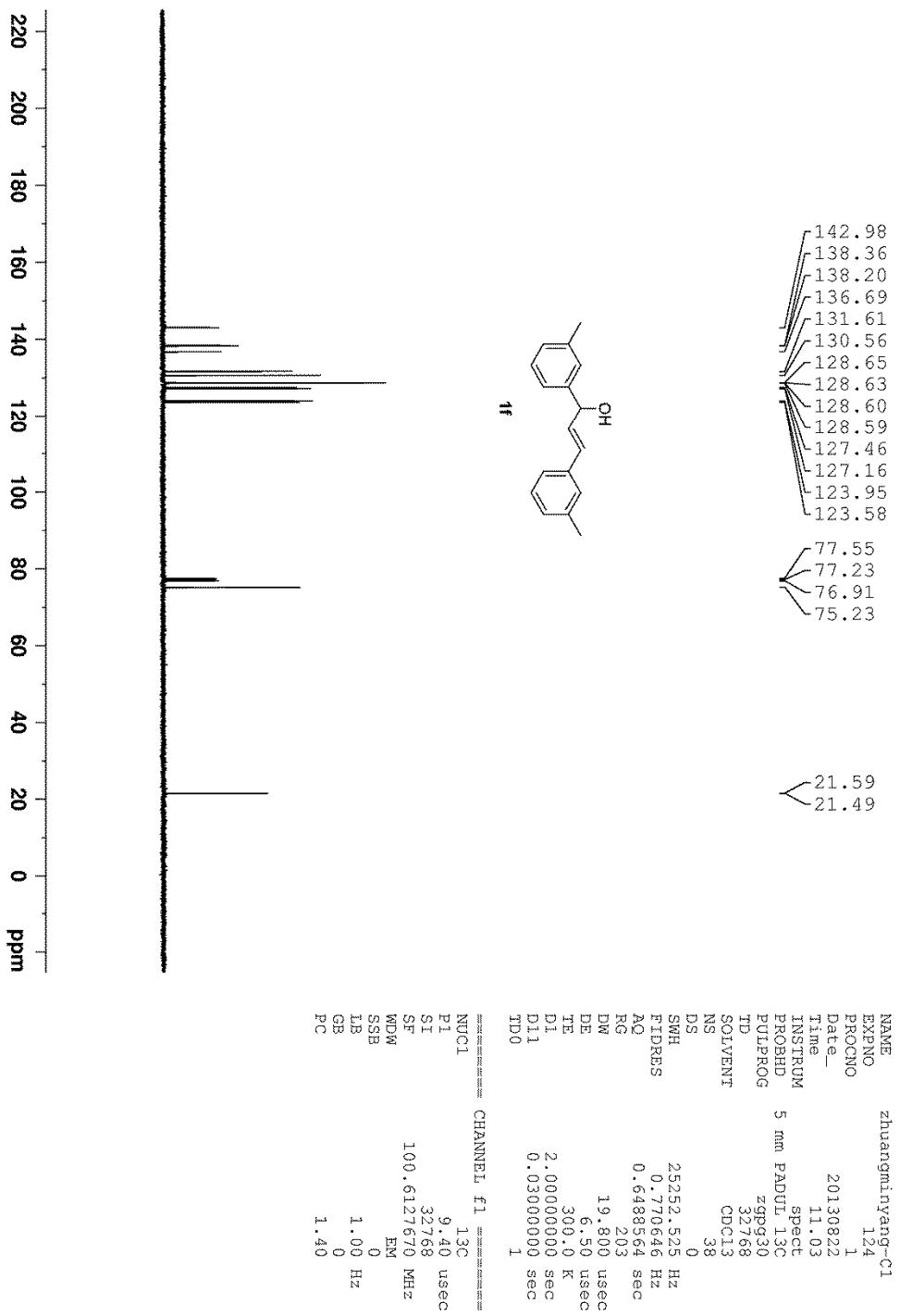
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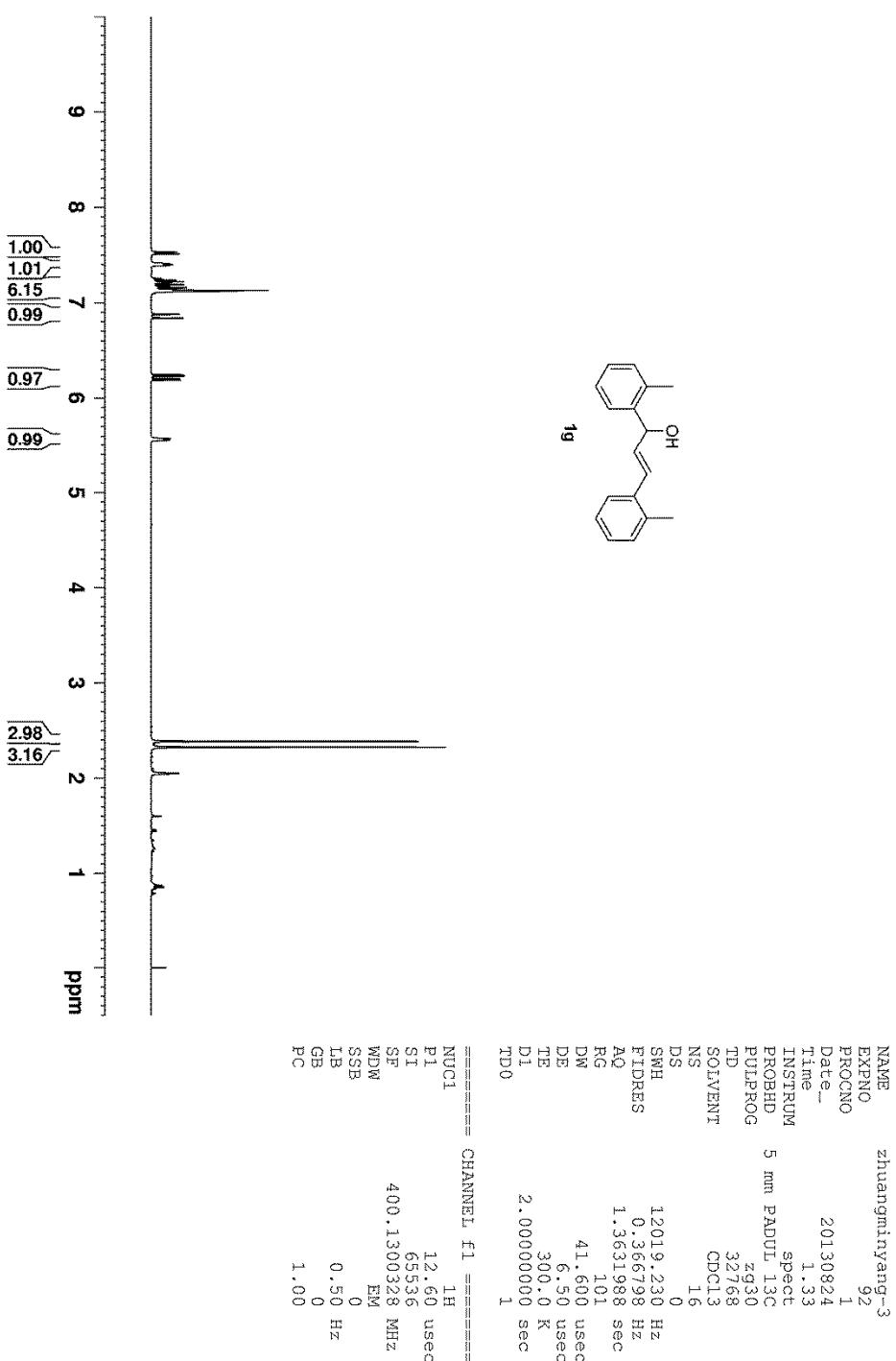


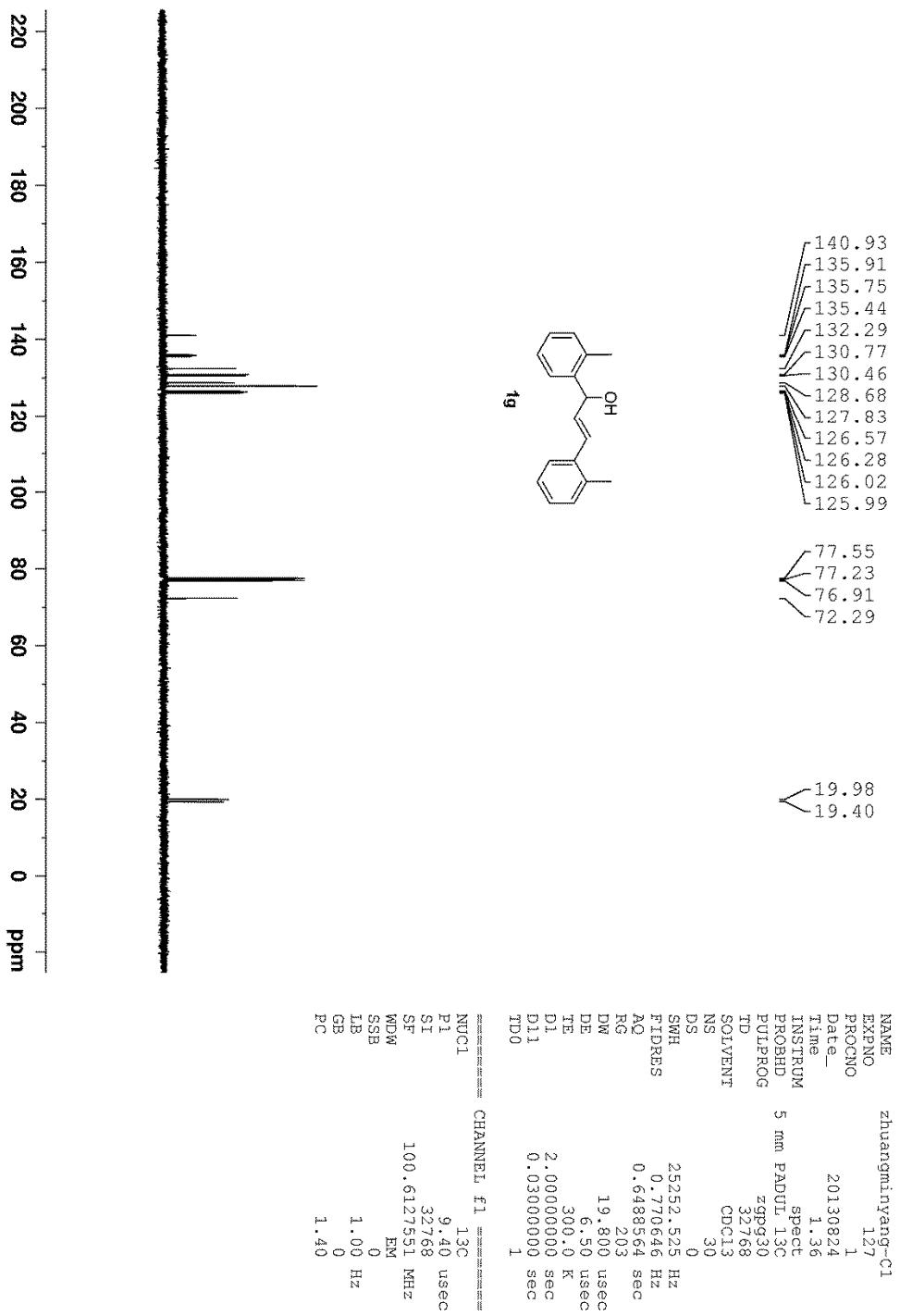


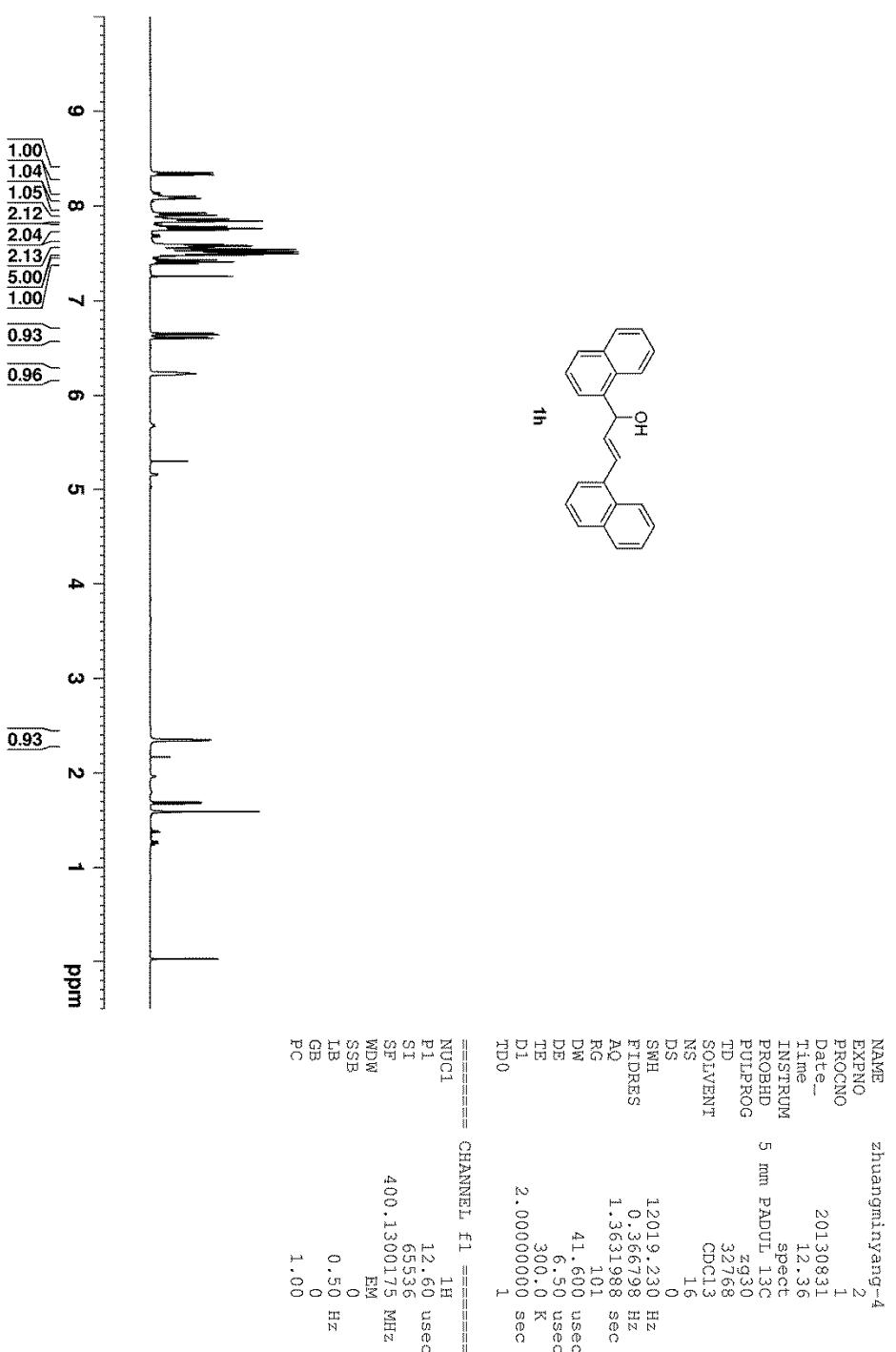


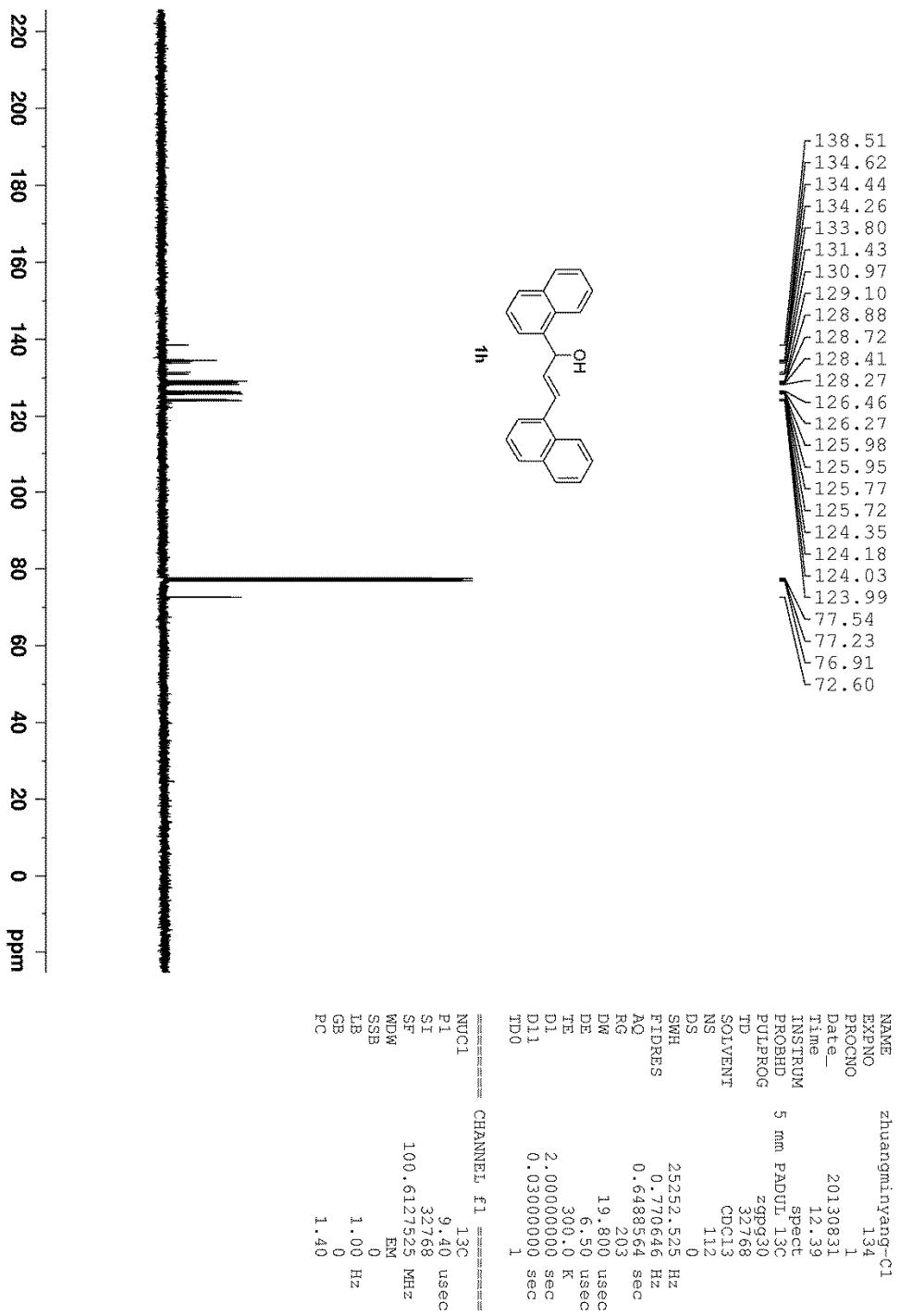


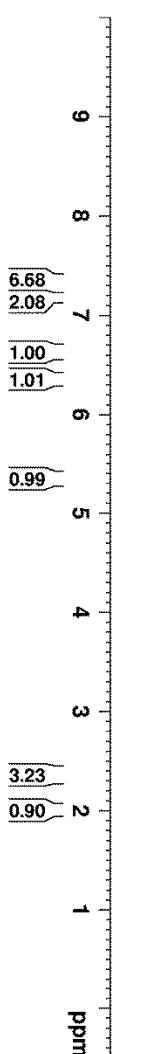
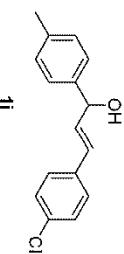










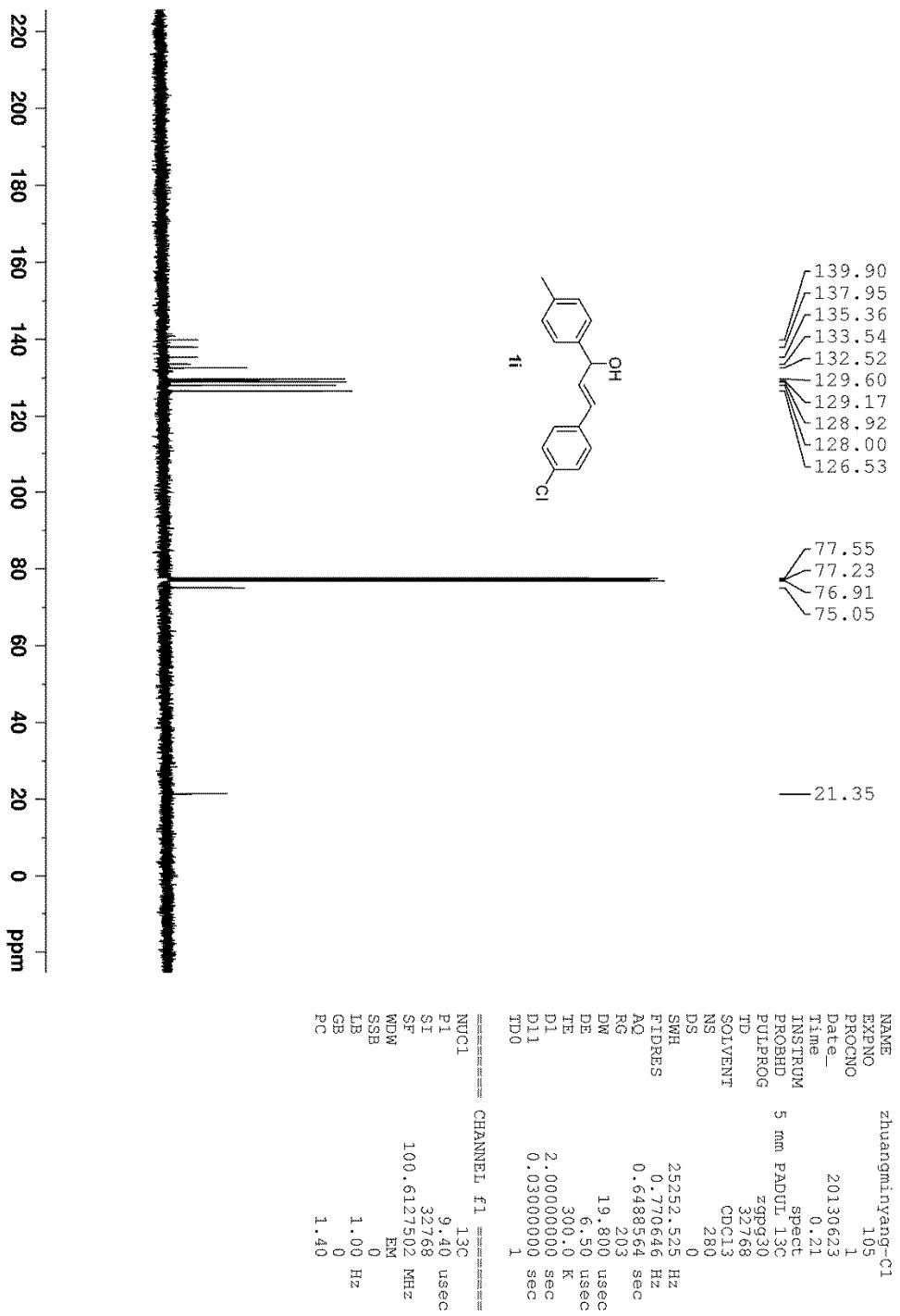


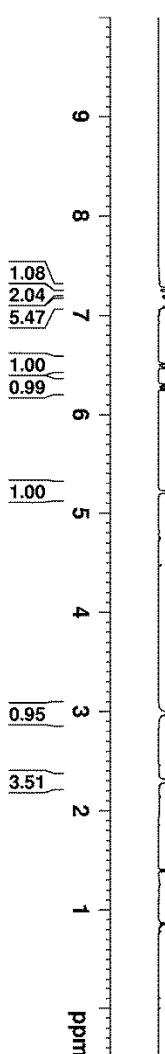
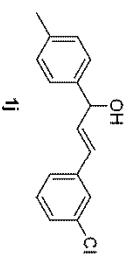
```

=====
NAME          zhuangminyang-3
EXPNO         8
PROCNO        1
Date_        20130623
Time         0.18
INSTRUM      spect
PROBHD      5 mm PADUL 13C
PULPROG     zg30
TD          32768
SOLVENT      CDCl3
NS           16
DS            0
SWH         12019.230 Hz
FIDRES     0.366798 Hz
AQ          1.3631988 sec
RG           203
DW           41.600 usec
DE           6.50 usec
TE          300.0 K
DI          2.0000000 sec
TDD0          1

=====
CHANNEL f1 =====
NUC1          1H
P1           12.60 usec
SI            65536
SF          400.1300197 MHz
MW            EM
SSB           0
LB            0.50 Hz
GB            0
PC           1.00

```



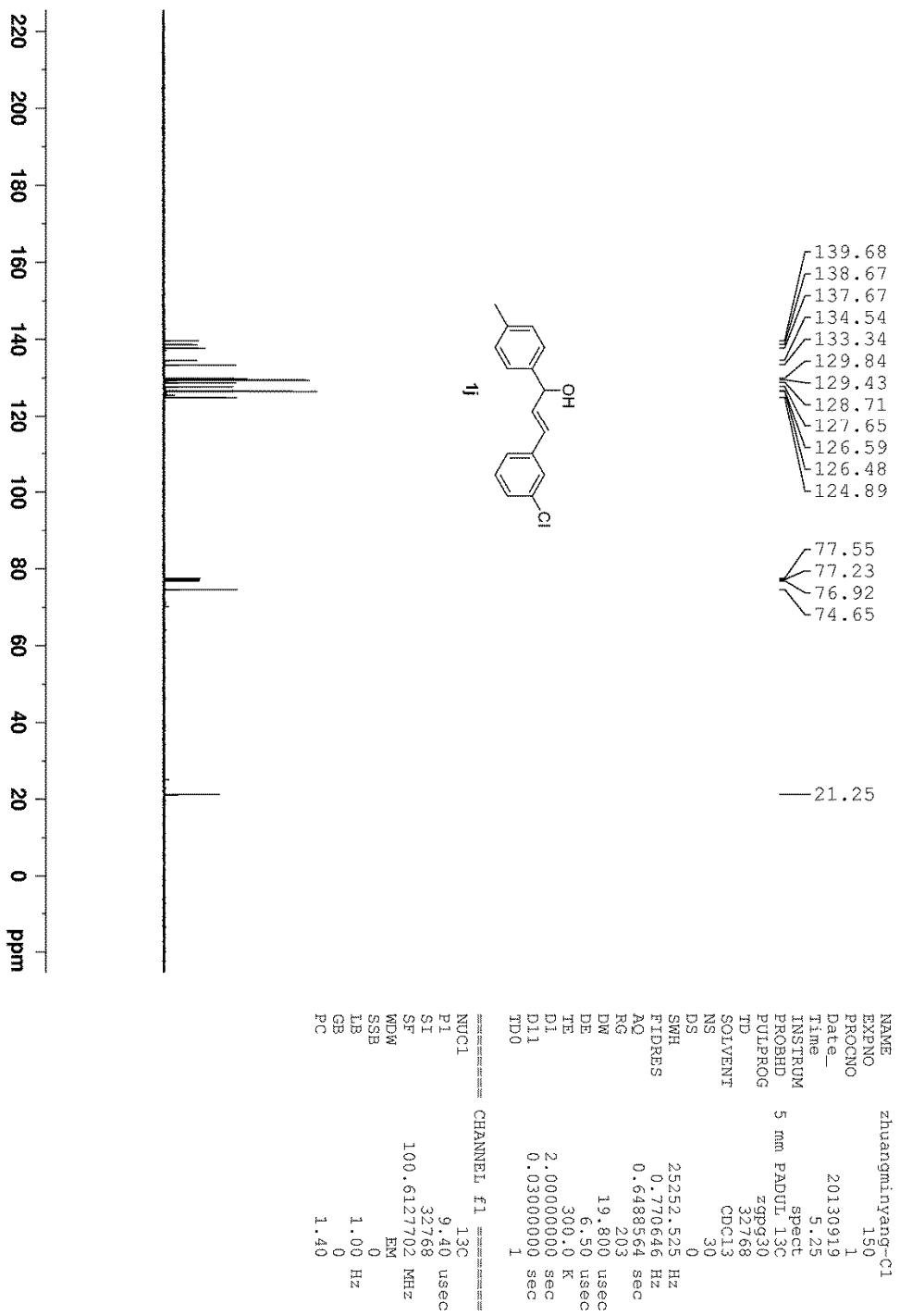


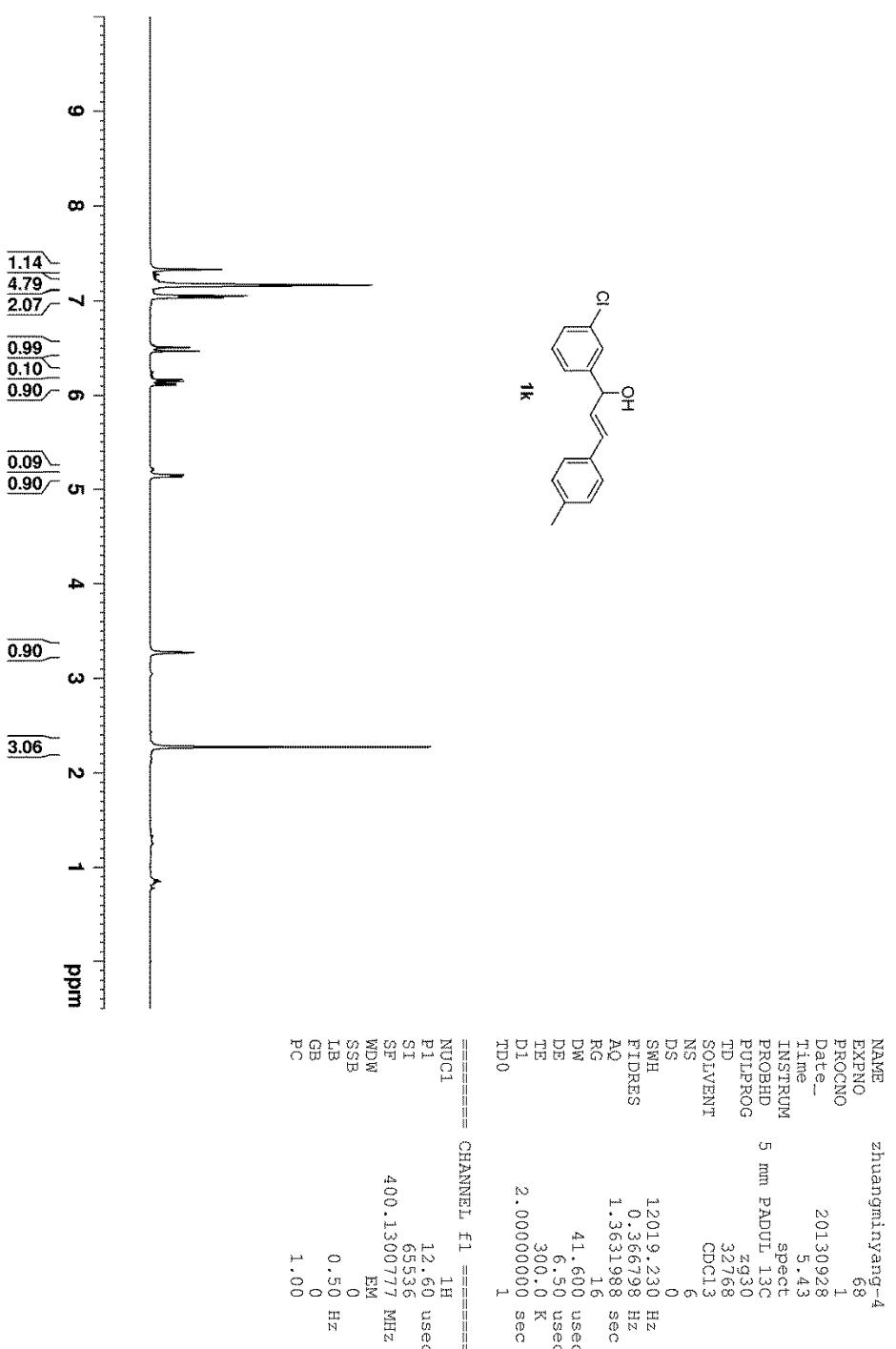
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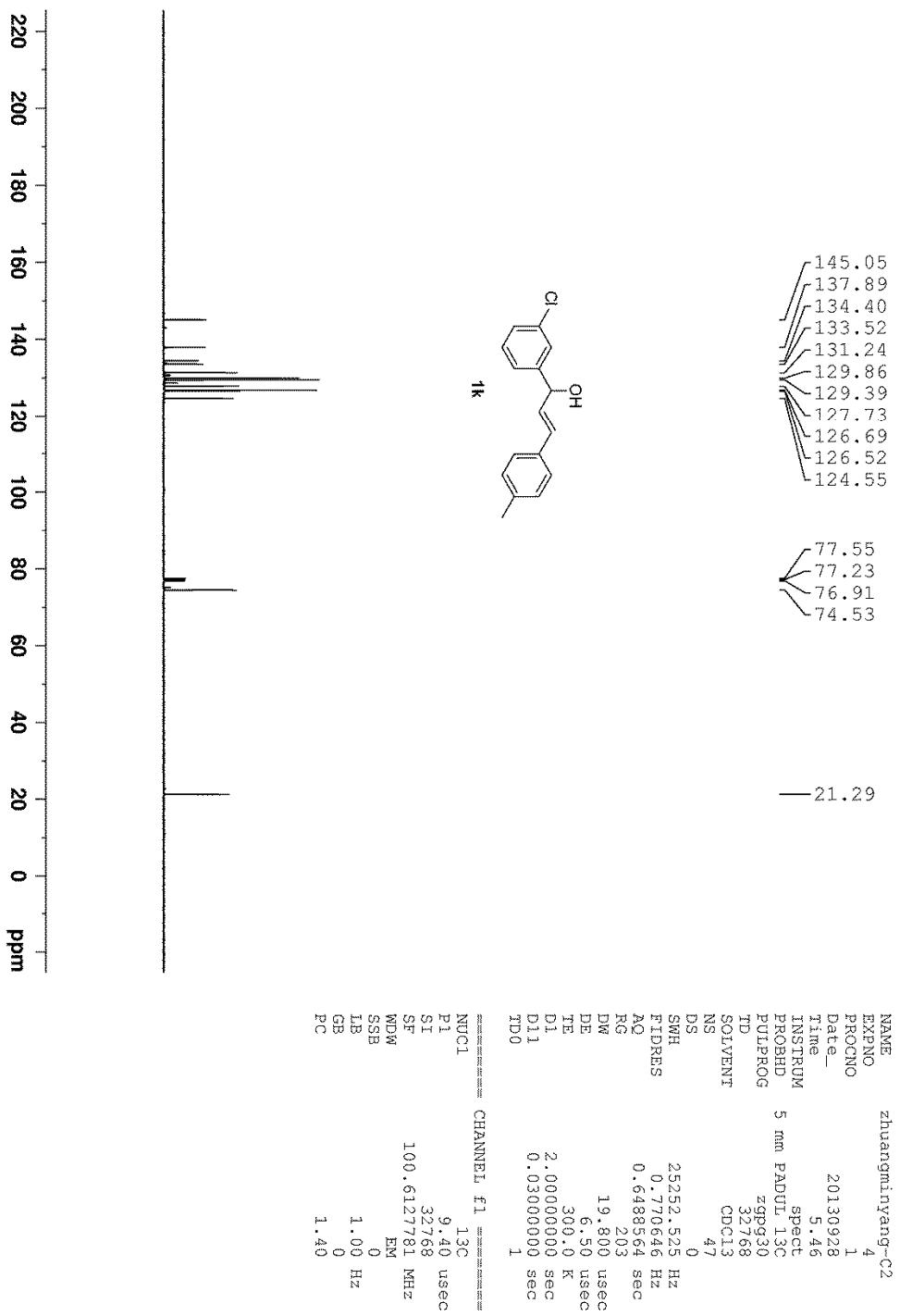
=====
NAME          zhuangminyang-4
EXPNO        50
PROCNO       1
Date_        20130919
Time         5.23
INSTRUM      spect
PROBHD      5 mm PADUL 13C
PULPROG     zg30
TD          32768
SOLVENT      CDCl3
NS           16
DS           0
SWH         12019.230 Hz
FIDRES     0.363798 Hz
AQ          1.363398 sec
RG          22.6
DW          41.600 usec
DE          6.50 usec
TE          300.0 K
DI          2.0000000 sec
TDD0         1

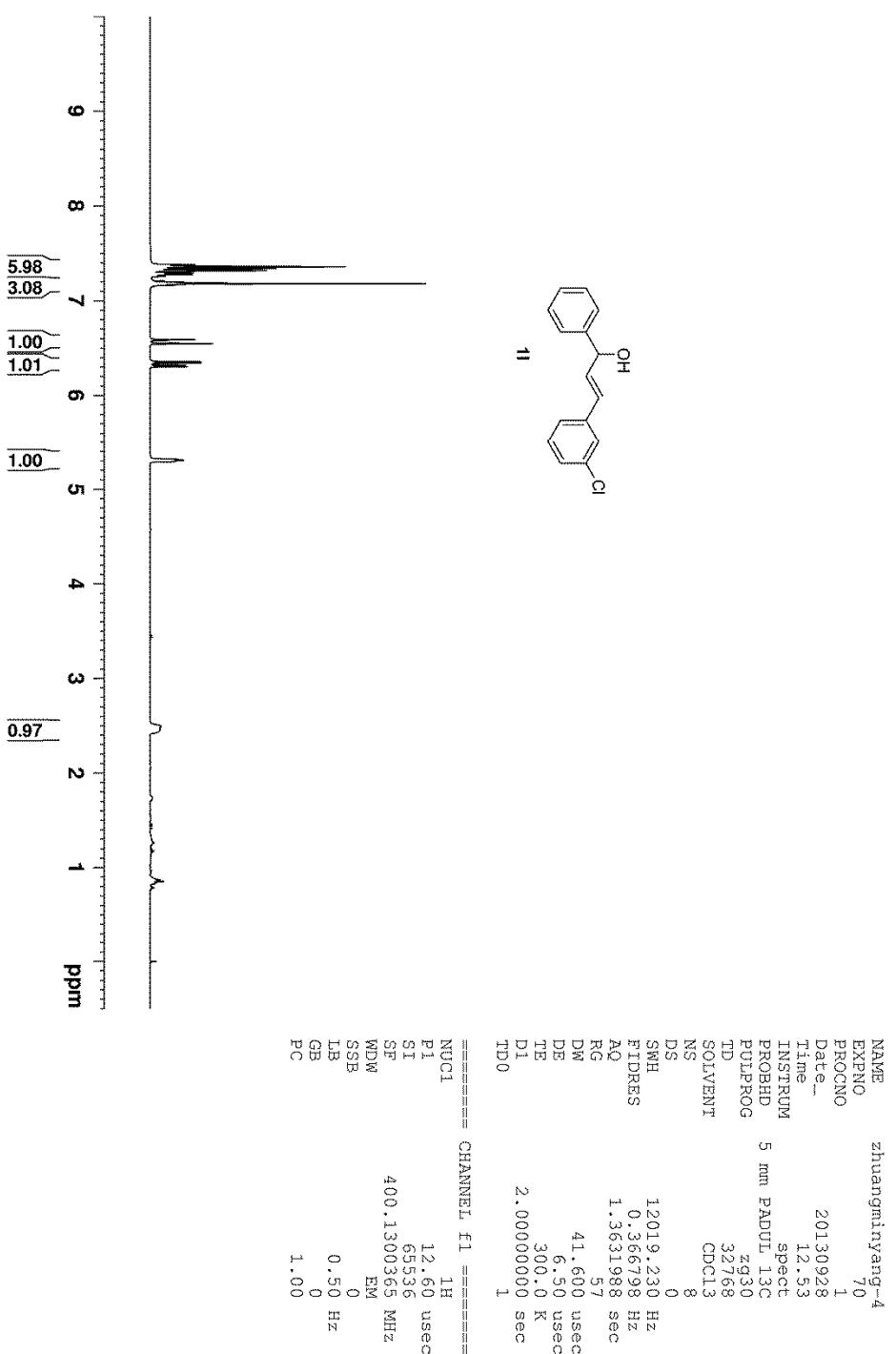
===== CHANNEL f1 =====
NUC1          1H
P1          12.60 usec
SI           65536
SF          400.1300563 MHz
WDW
SSB          0
LB          0.50 Hz
GB          0
PC          1.00

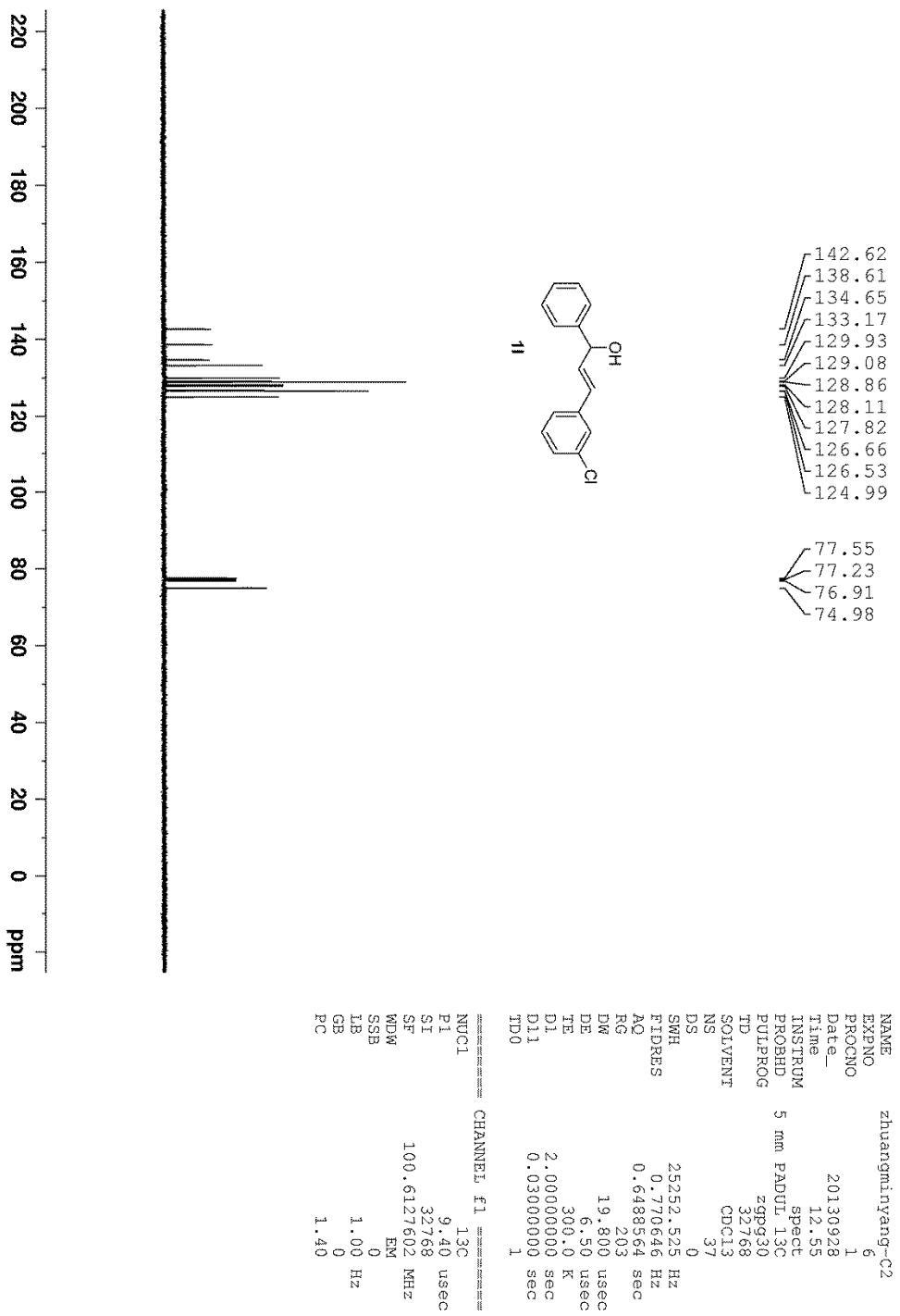
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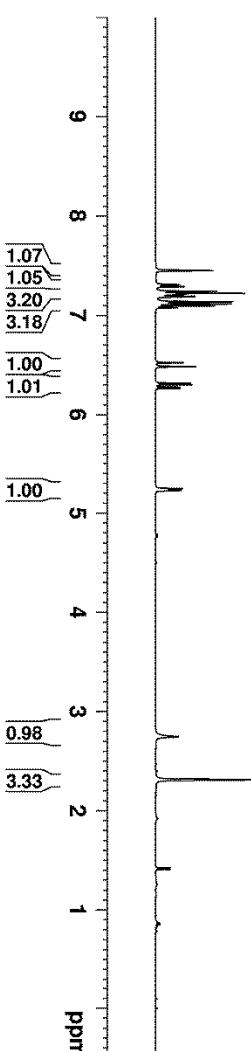
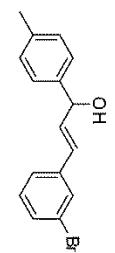










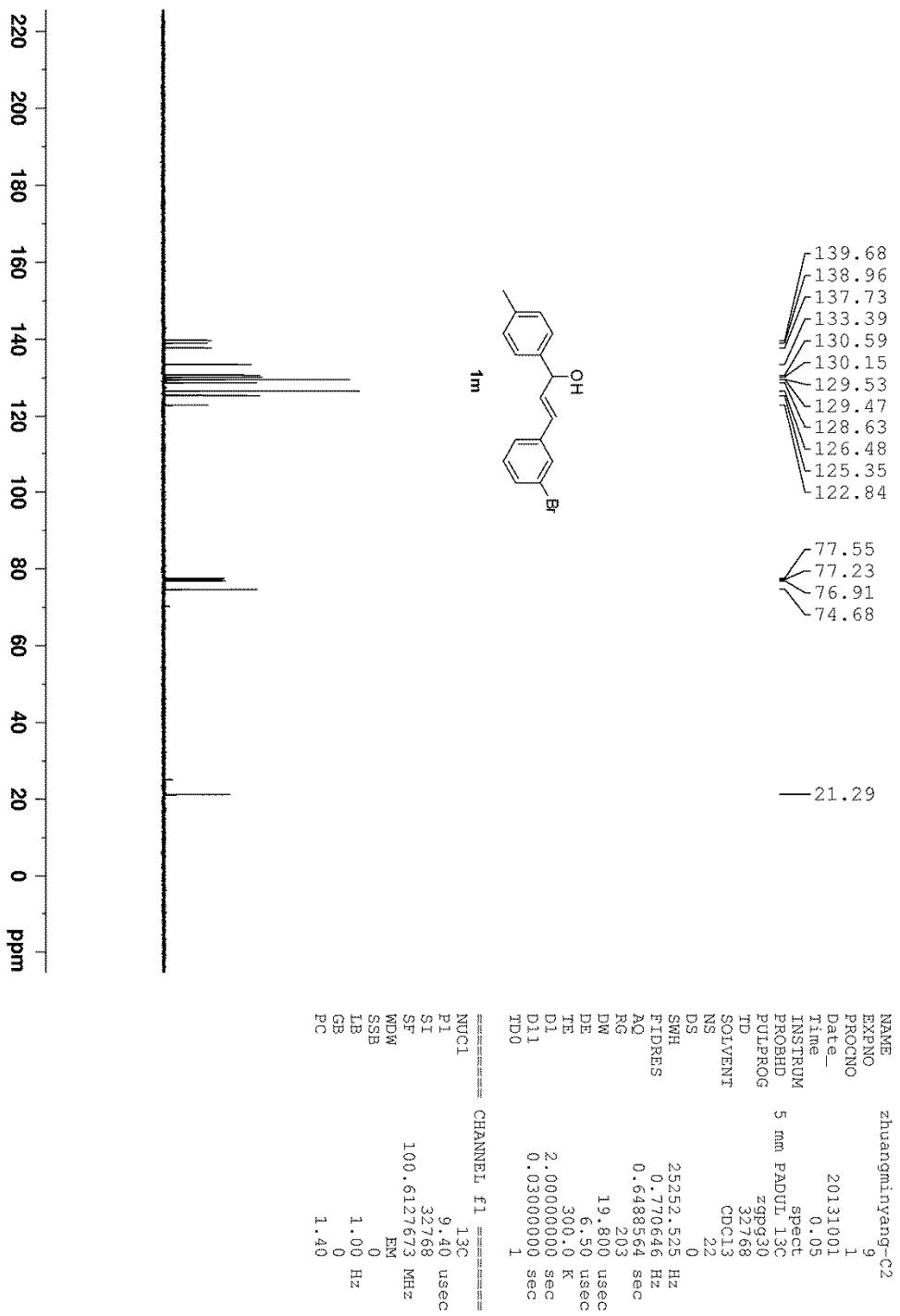


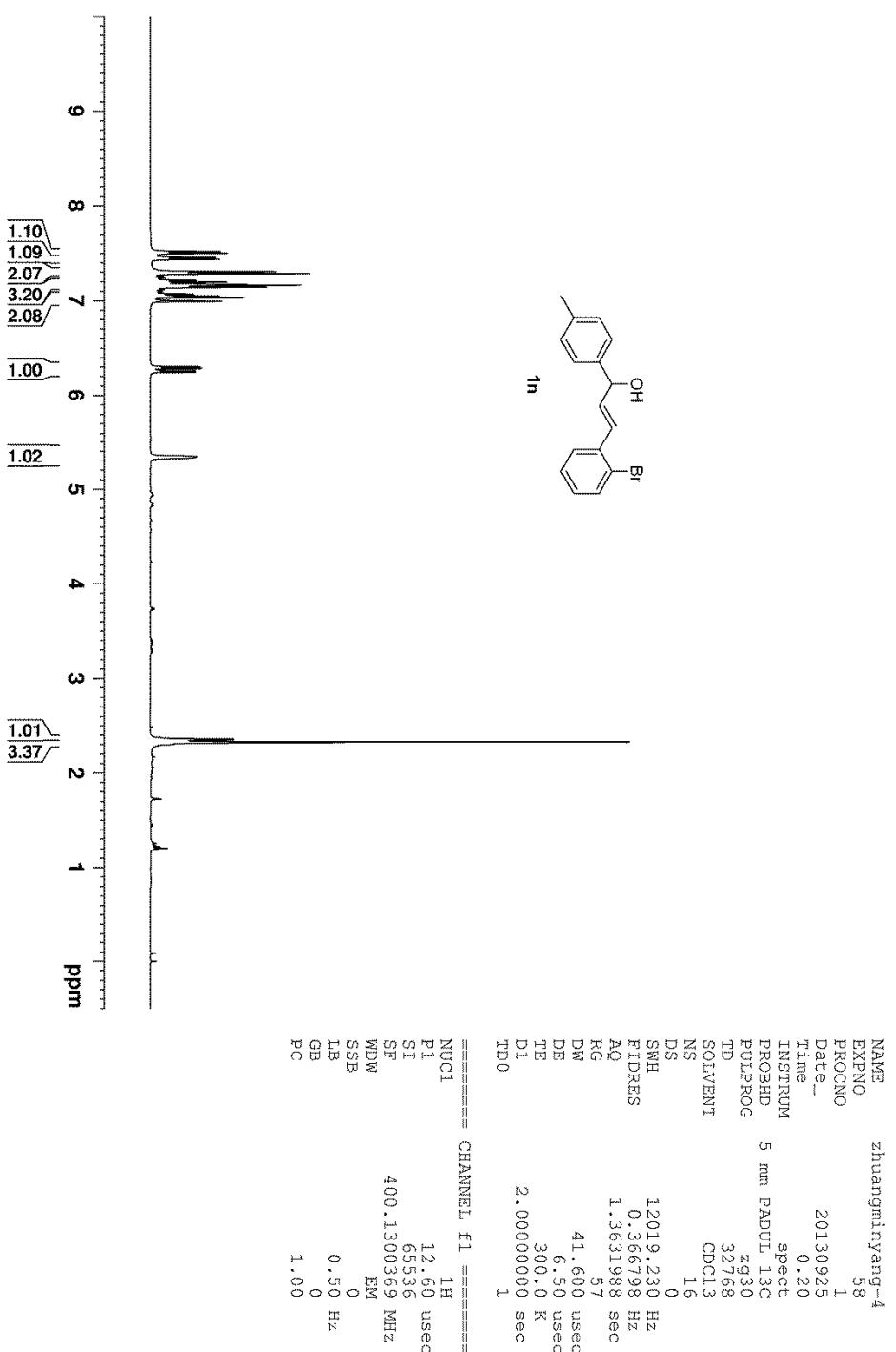
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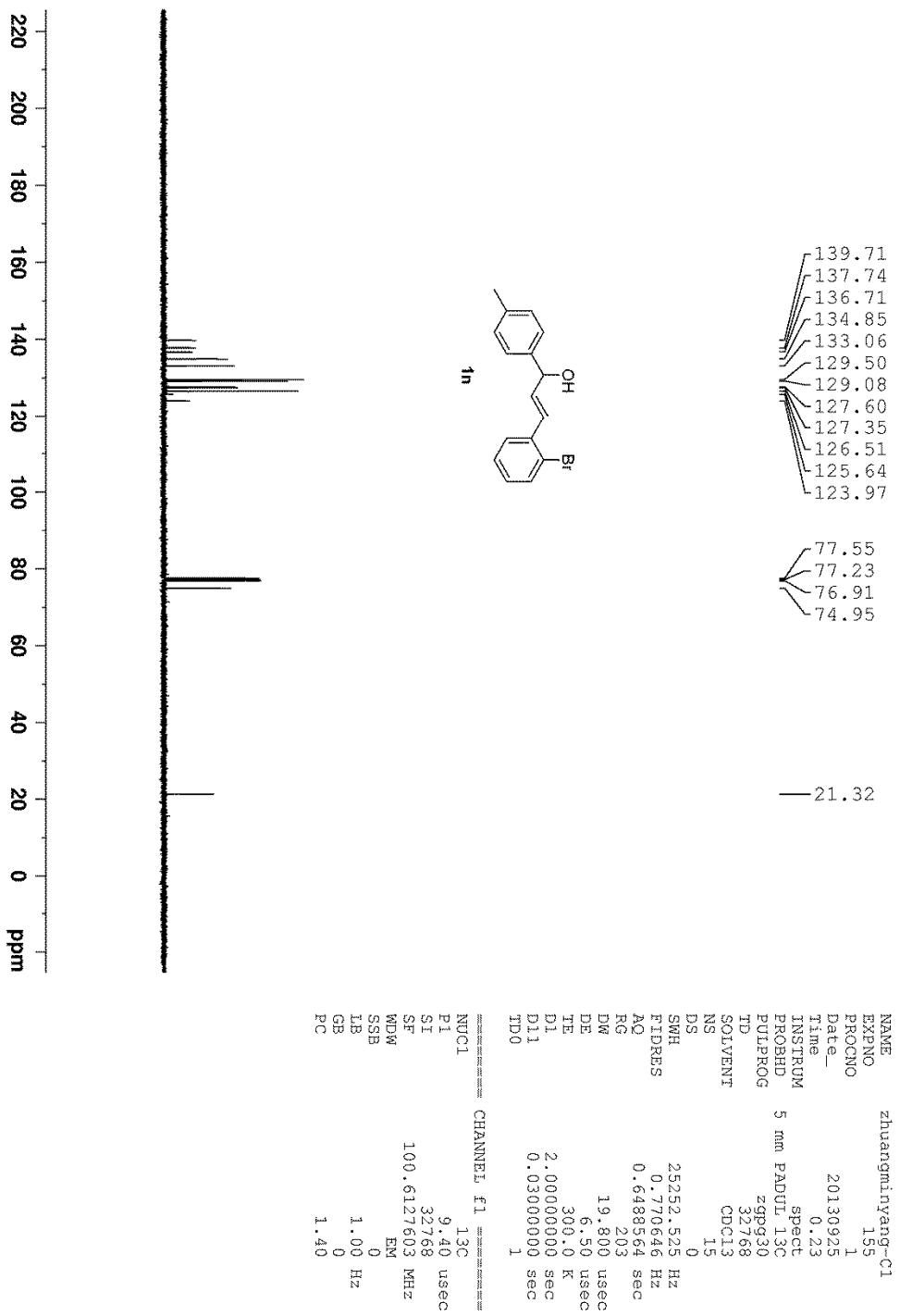
NAME          zhuangminyang-4
EXPNO        76
PROCNO       1
Date_        20131001
Time         0.03
INSTRUM      spect
PROBHD      5 mm PADUL 13C
PULPROG     zg30
TD          32768
SOLVENT      CDCl3
NS           8
DS           0
SWH         12019.230 Hz
FIDRES     0.366798 Hz
AQ          1.3633988 sec
RG          25.4
DW          41.600 usec
DE          6.50 usec
TE          310.0 K
DI          2.0000000 sec
TDD0         1

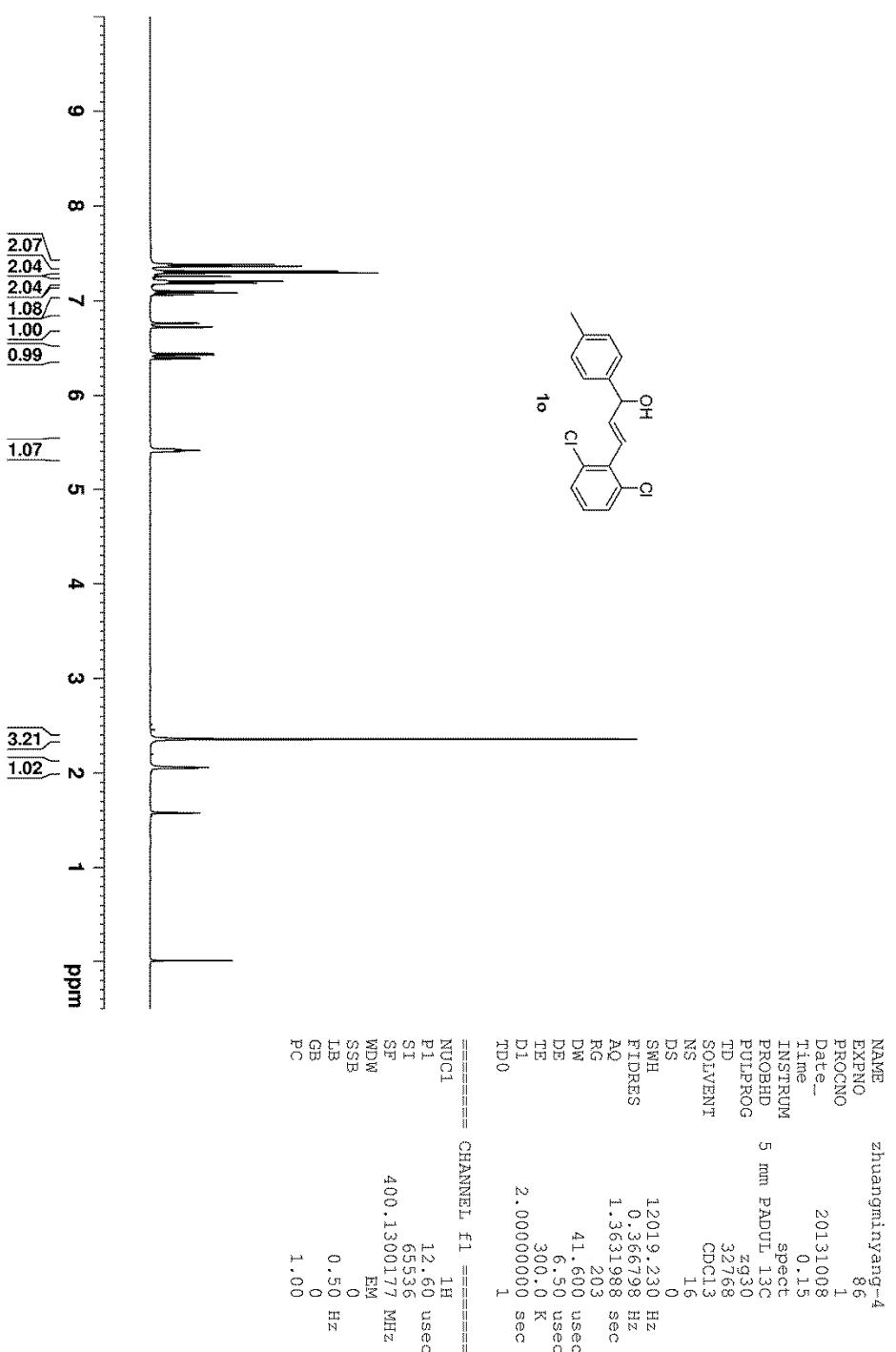
===== CHANNEL f1 =====
NUC1          1H
P1          12.60 usec
SI           65536
SF          400.1300475 MHz
WDW
SSB          0
LB          0.50 Hz
GB          0
PC          1.00

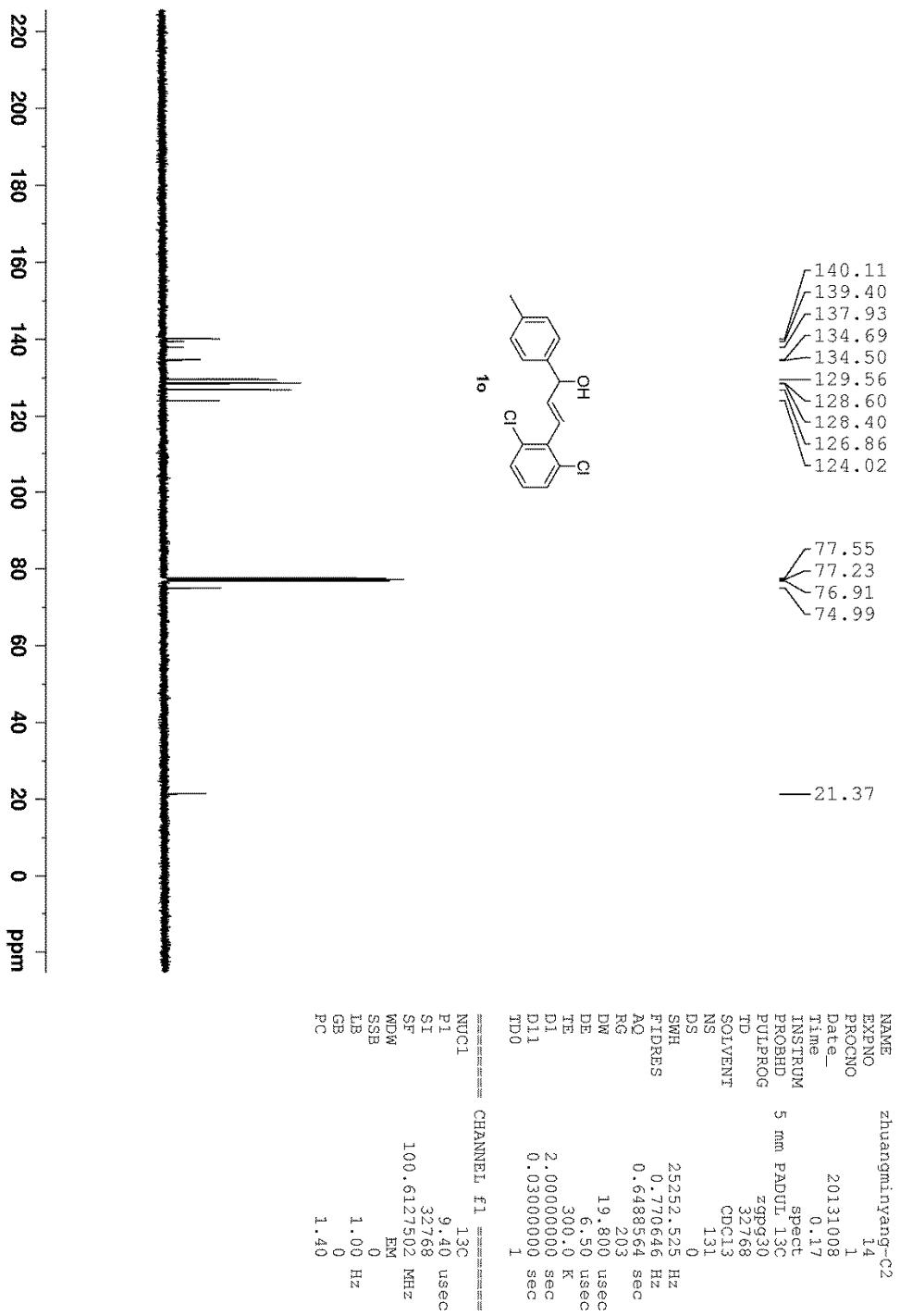
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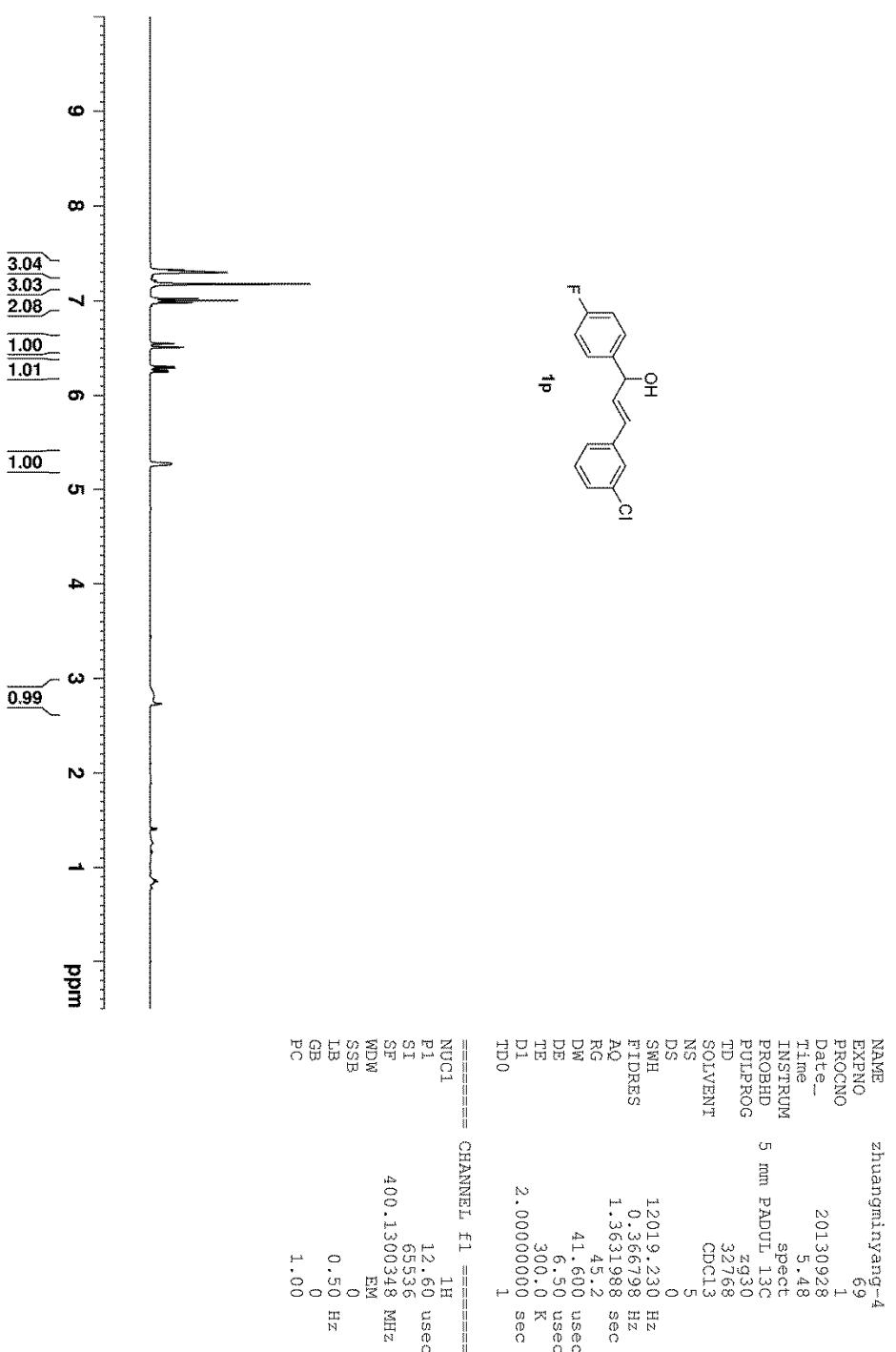


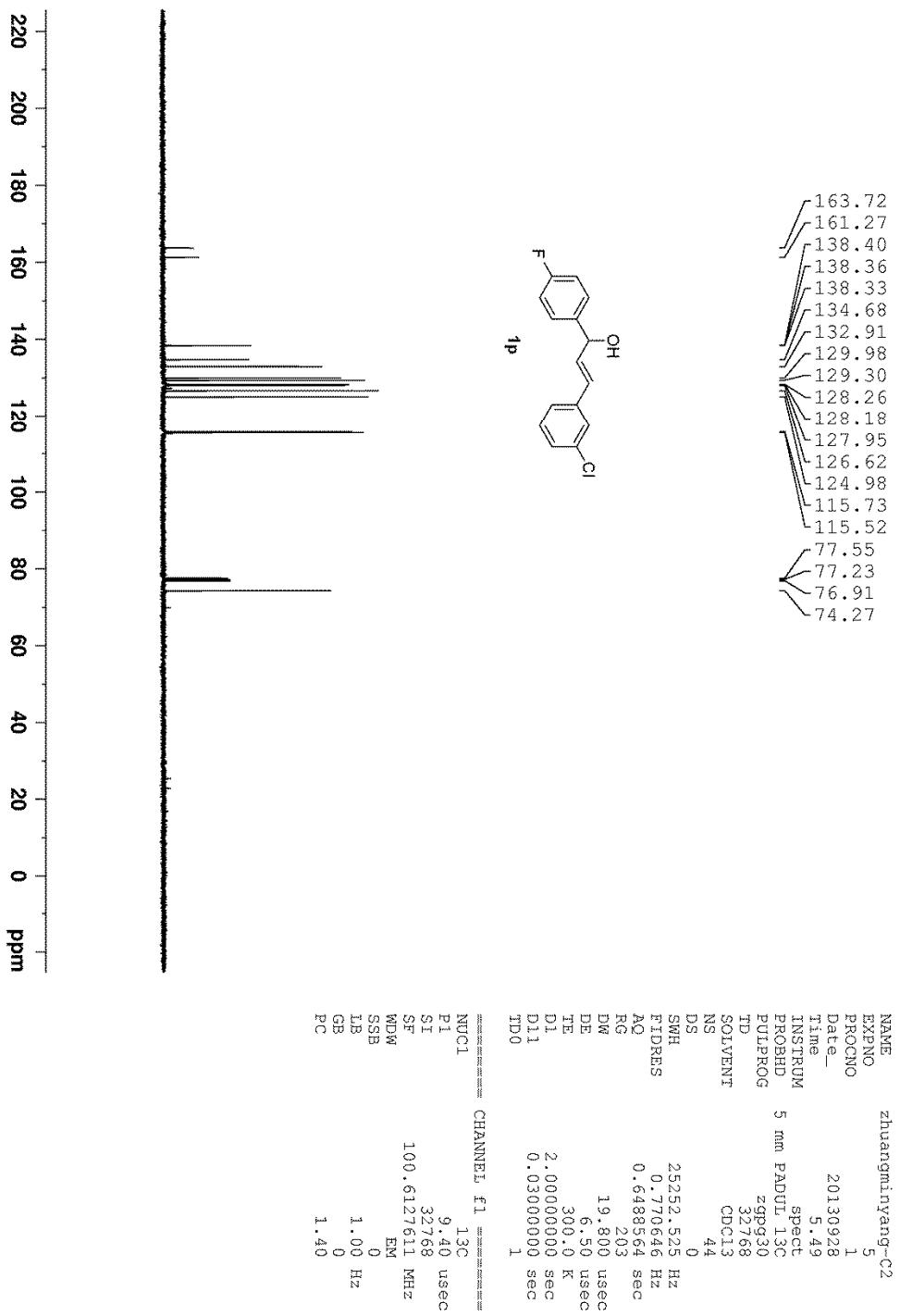


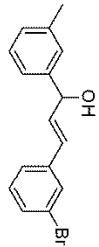
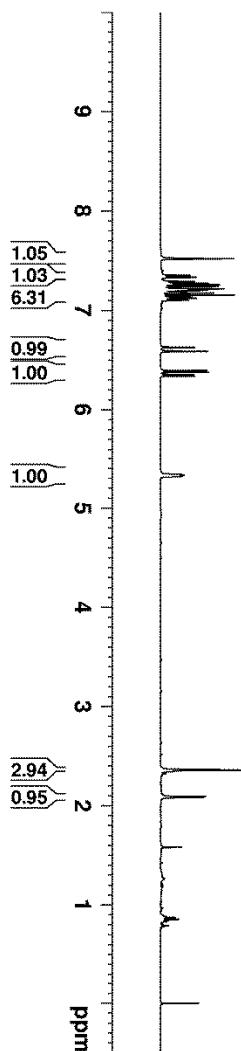










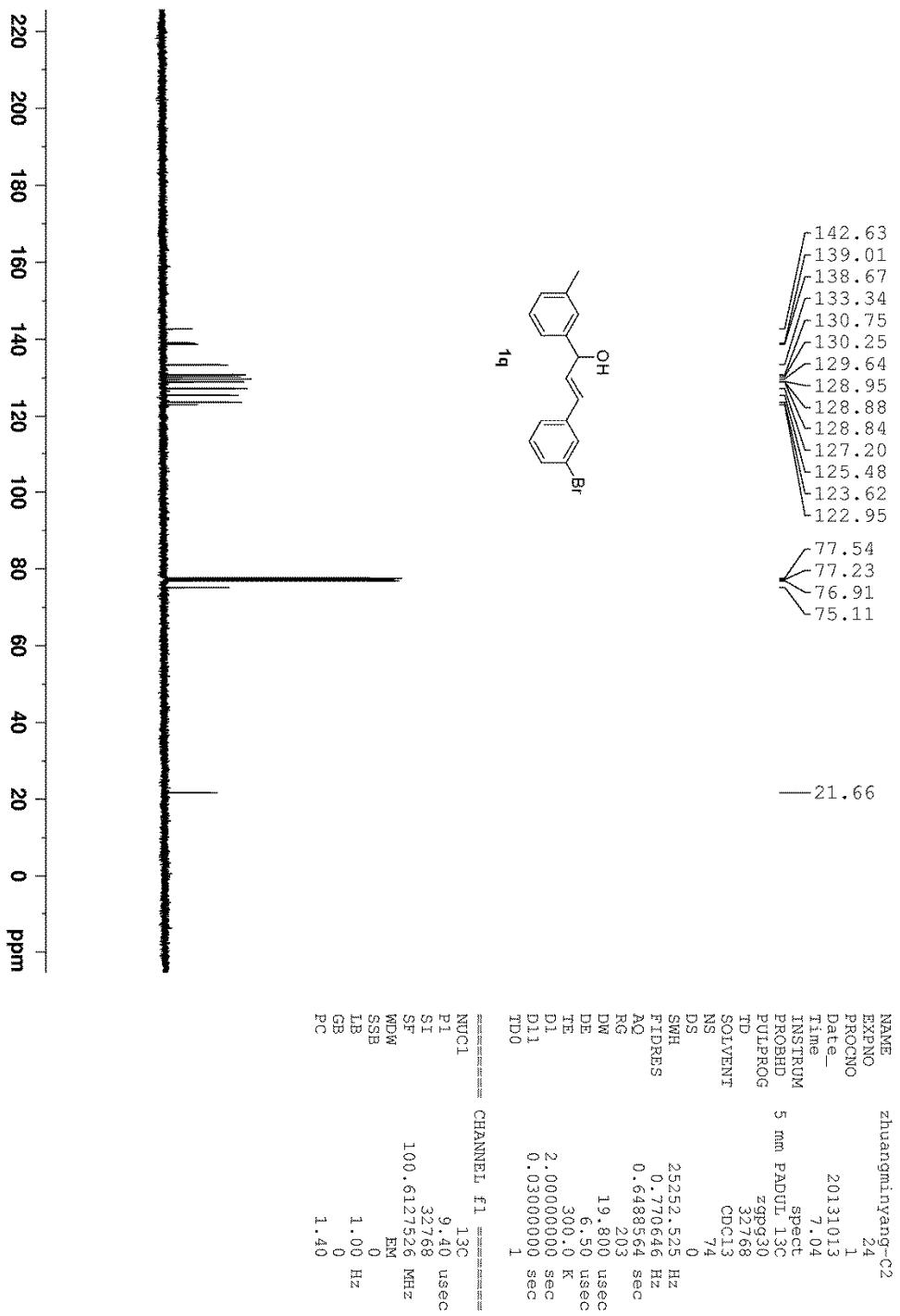


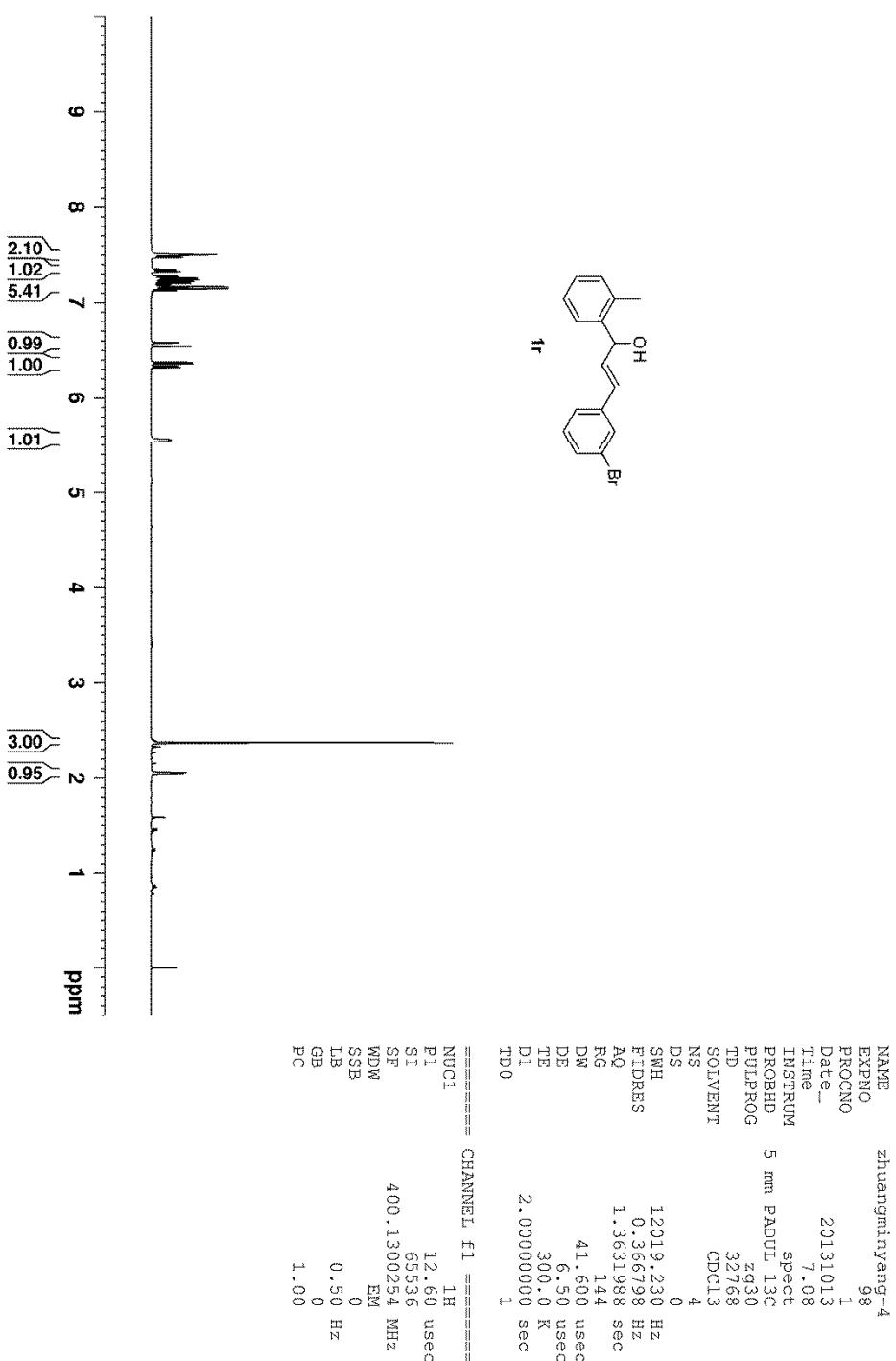
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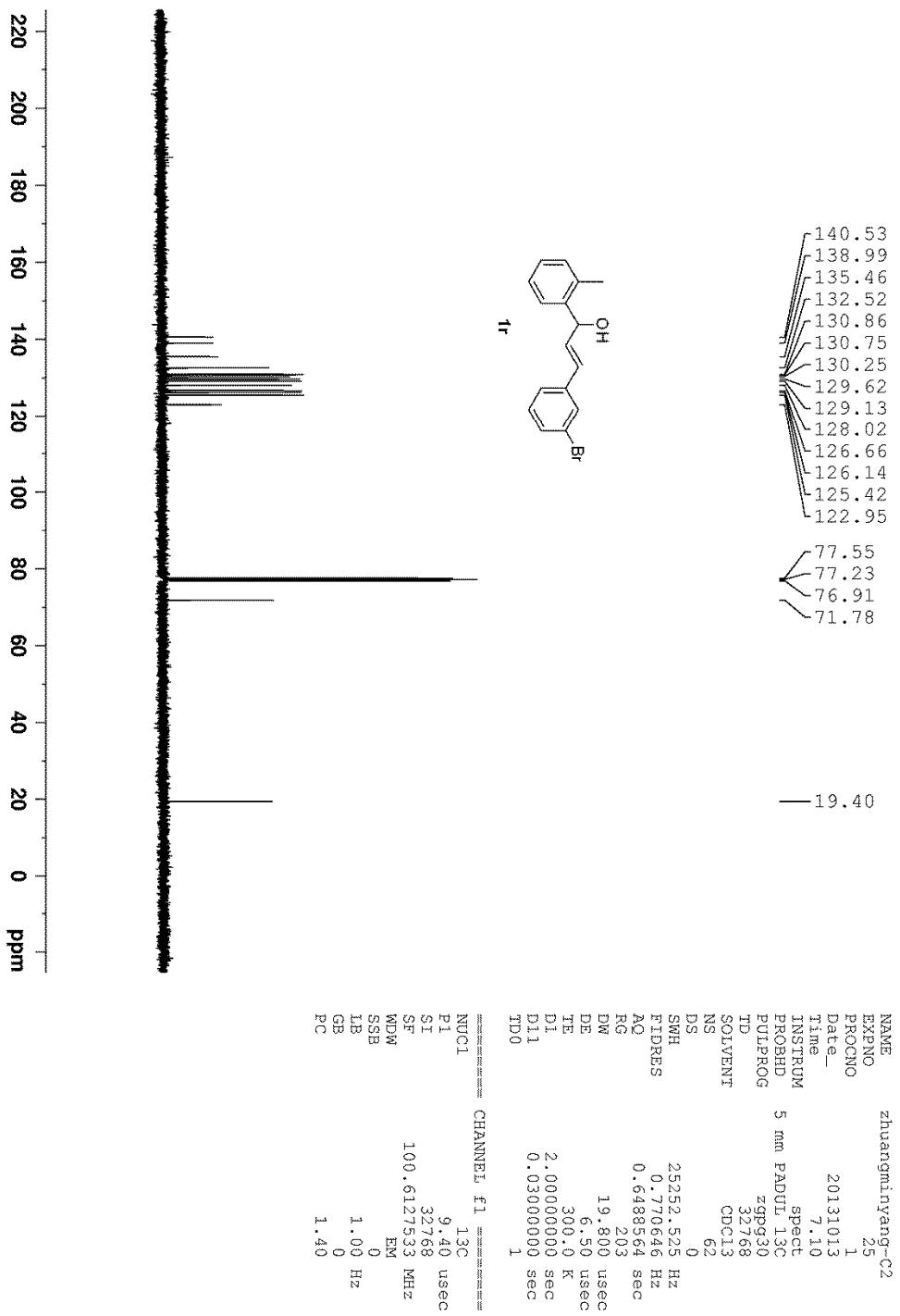
NAME          zhuangminyang-4
EXPNO         97
PROCNO        1
Date-        2013/01/13
Time-          7.02
INSTRUM     spect
PROBID       5 mm PADUL 13C
PULPROG      PULFROG
TD           32768
SOLVENT      SOLENIT
NS            6
DS            0
SWH          1201.9-230 Hz
FIDRES      0.366798 Hz
AQ            1.3631988 sec
RG           144
DW           41.600 usec
DE           6.500 usec
TE           300.0 K
D1          2.0000000 sec
TDO          1

===== CHANNEL f1 =====
NUC1          1H
P1           12.60 usec
SI            65536
SF          400.1300237 MHz
WDW          EM
SSB          0
LB           0.50 Hz
GB          0
PC          1.00

```







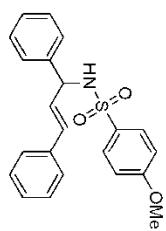
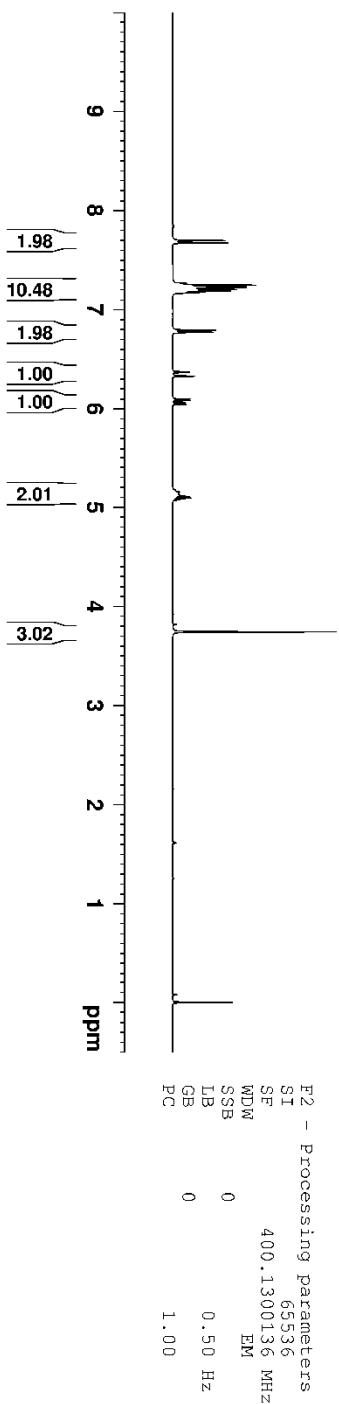


Table 1, entry 12



Current Data Parameters
NAME zhuangminyang-6
EXPNO 47
PROCNO 1

F2 - Acquisition Parameters

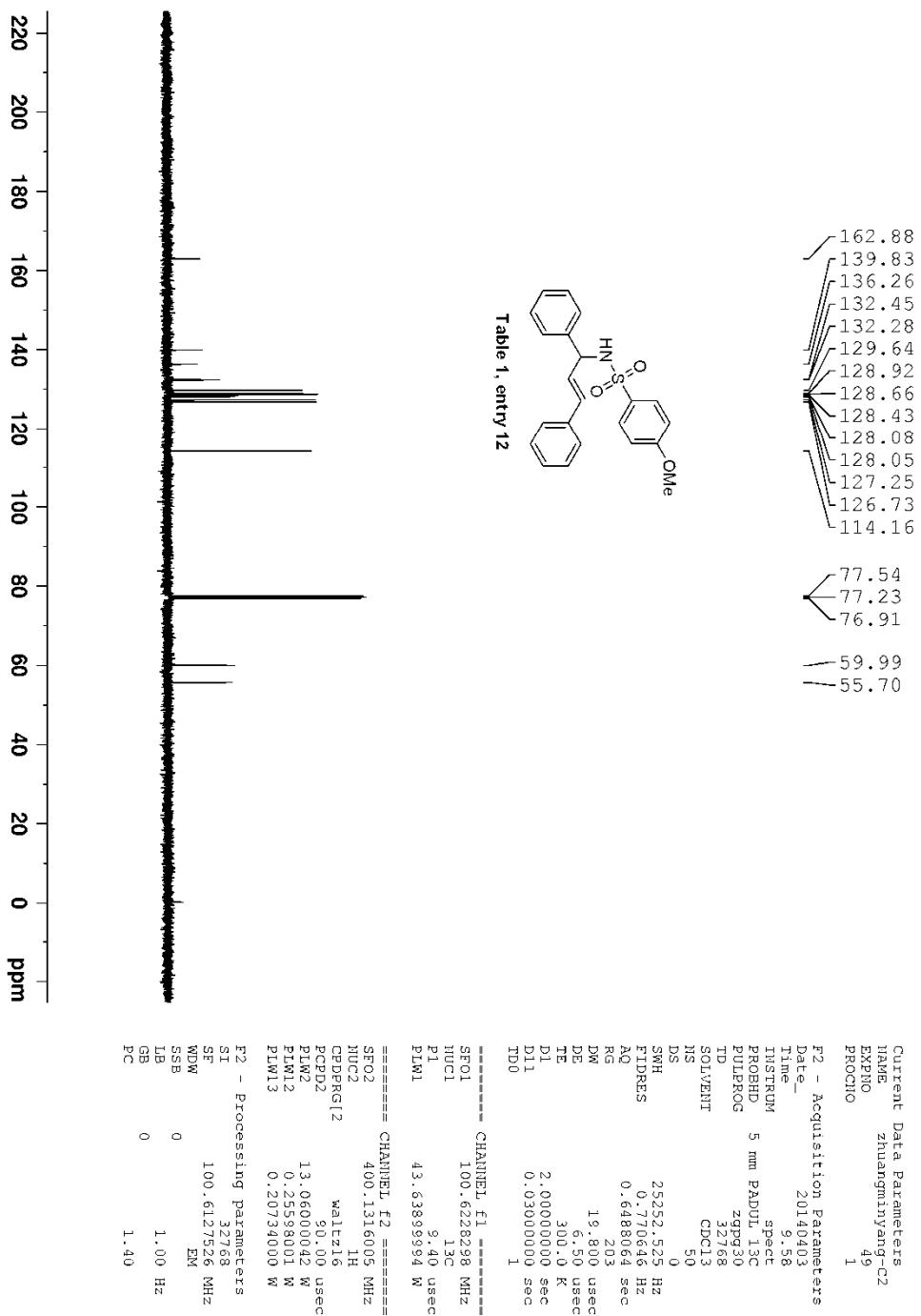
Date 20140403
Time 9.56
INSTRUM spect
PROBHD 5 mm PABUL 13C
PULPROG zg30
TD 32768
SOLVENT CDCl₃
NS 6
DS 0
SWH 12019.230 Hz
FIDRES 0.366798 Hz
AQ 1.3631488 sec
RG 114
DW 41.600 usec
DE 6.50 usec
TE 300.0 K
D1 2.0000000 sec
TDD 1

===== CHANNEL f1 =====

SP01 400.1320007 MHz
NUC1 ¹H
P1 12.60 usec
PLW1 13.06000042 W

F2 - Processing parameters

SI 65536
SF 400.1300136 MHz
WDW EM
SSB 0
LB 0.50 Hz
GB 1.00
PC



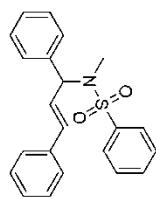
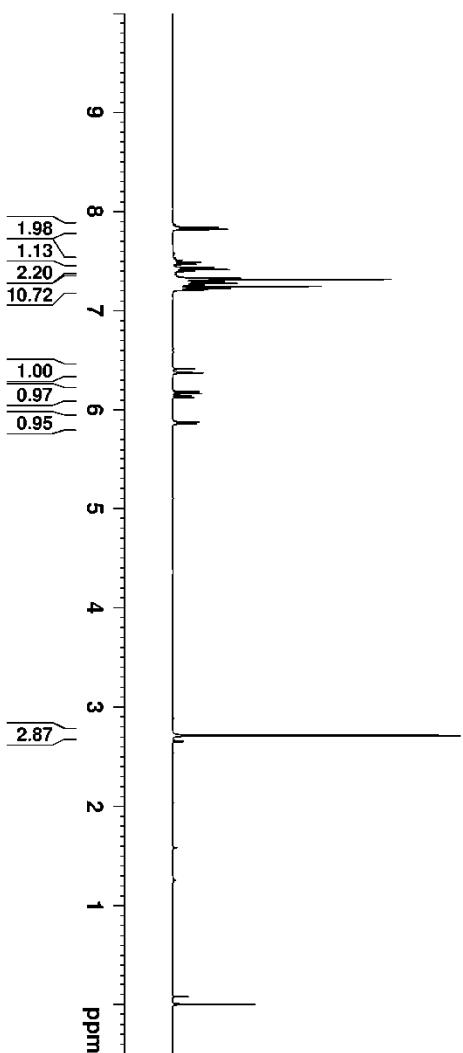


Table 1, entry 13



Current Data Parameters
NAME zhuangminyang-6
EXPNO 48
PROCNO 1

F2 - Acquisition Parameters

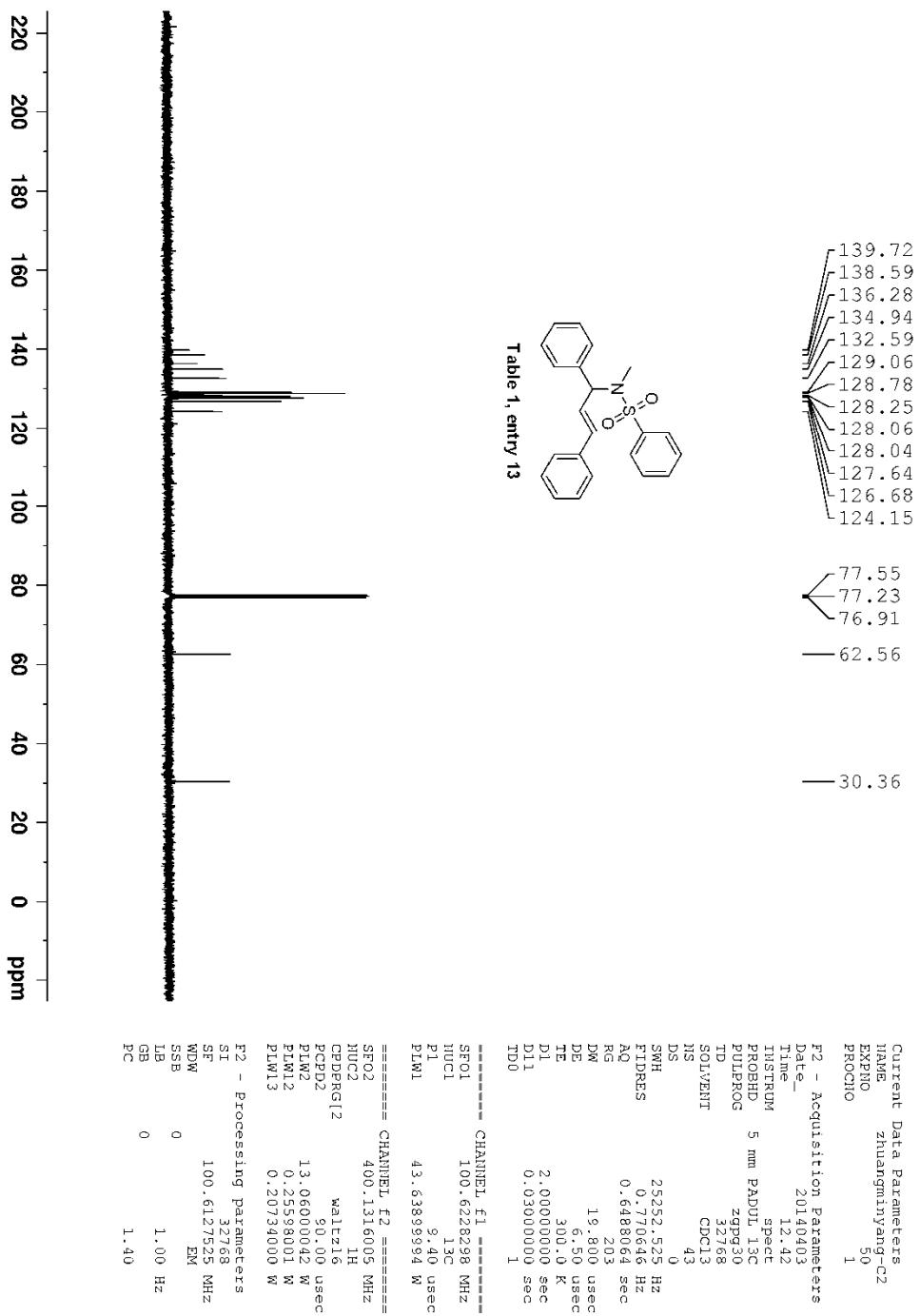
Date_ 20140403
Time 12.40
INSTRUM spect
PROBHD 5 mm PADUL 13C
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 11
DS 0
SWH 12019.230 Hz
FIDRES 0.366798 Hz
AQ 1.3631488 sec
RG 114
DW 41.600 usec
DE 6.50 usec
TE 300.0 K
D1 2.0000000 sec
TDO 1 sec

===== CHANNEL f1 =====

SFO1 400.1320007 MHz
NUC1 1H
P1 12.60 usec
PLW1 13.06000042 W

F2 - Processing parameters

SI 65536
SF 400.1300145 MHz
NDM SSB
SSB 0
LB 0.50 Hz
GB 0
PC 1.00



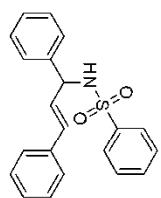
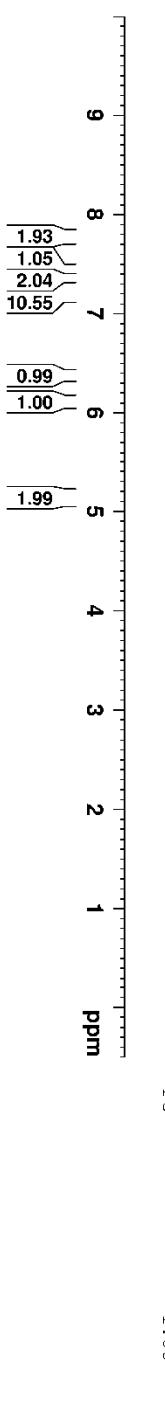
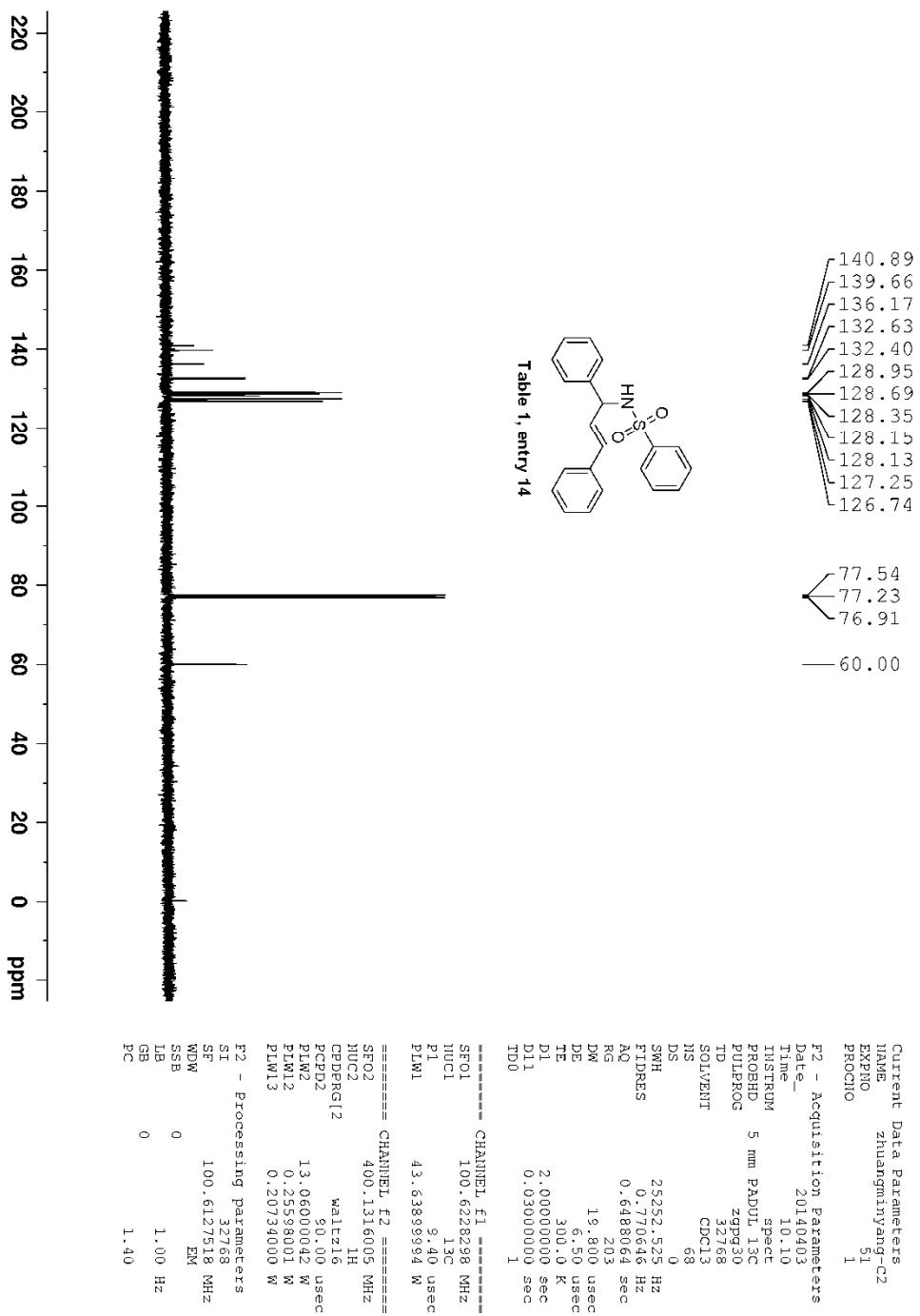


Table 1, entry 14



Current Data Parameters	
NAME	zhuangminyang-6
EXPNO	49
PROCNO	1
F2 - Acquisition Parameters	
DATE	20140403
TIME	10.09
INSTRUM	Spect
PROBHD	5 mm PABU1 13C
PULPROG	zg30
TD	32768
SOLVENT	CDCl ₃
NS	6
DS	0
SWH	12019.230 Hz
FIDRES	0.366798 Hz
AQ	1.3631488 sec
RG	128
DW	41.600 usec
DE	6.50 usec
TE	300.0 K
D1	2.0000000 sec
TDO	1 sec
===== CHANNEL f1 =====	
SFO1	400.1320007 MHz
NUC1	¹ H
P1	12.60 usec
PLW1	13.0600042 W
F2 - Processing parameters	
SI	65536
SF	400.1300136 MHz
NDM	EM
SSB	0
LB	0.50 Hz
GB	0
PC	1.00



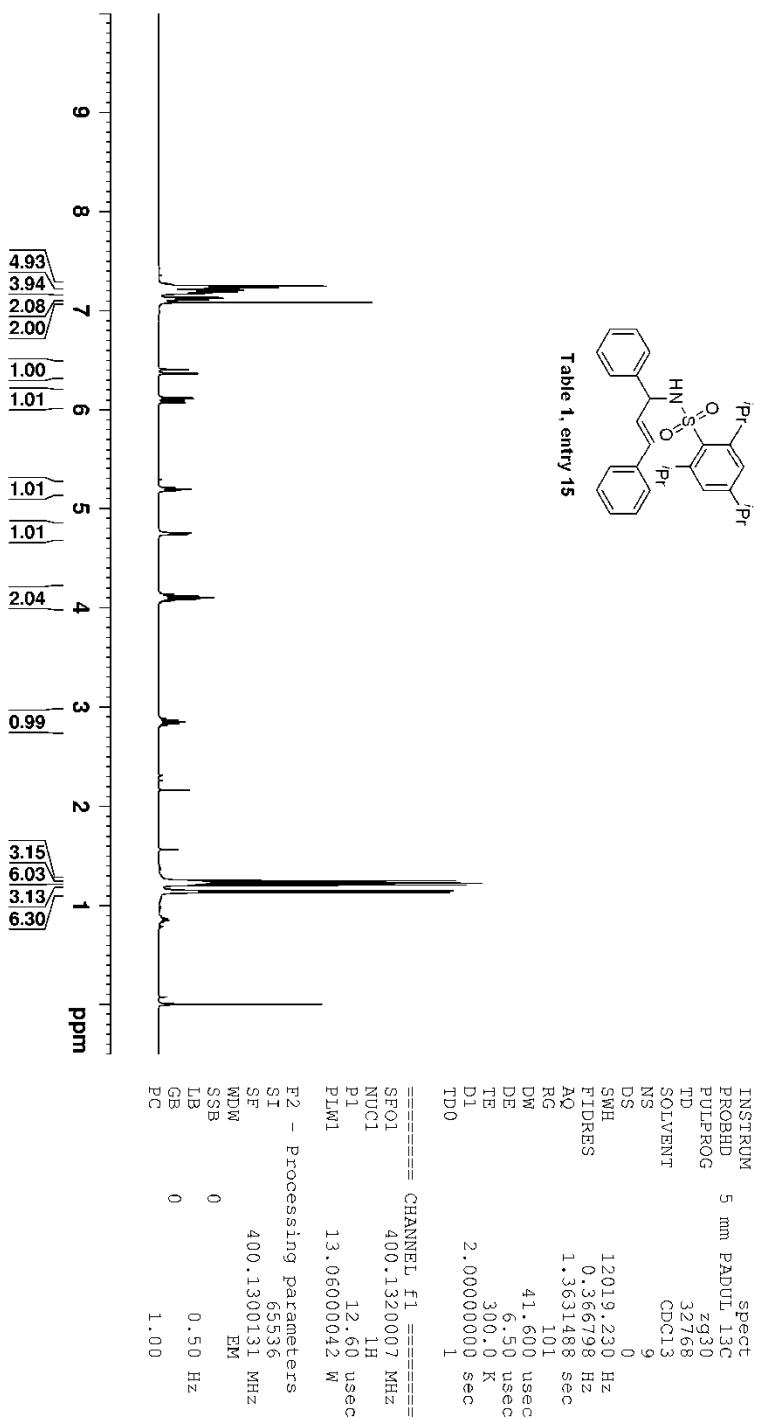
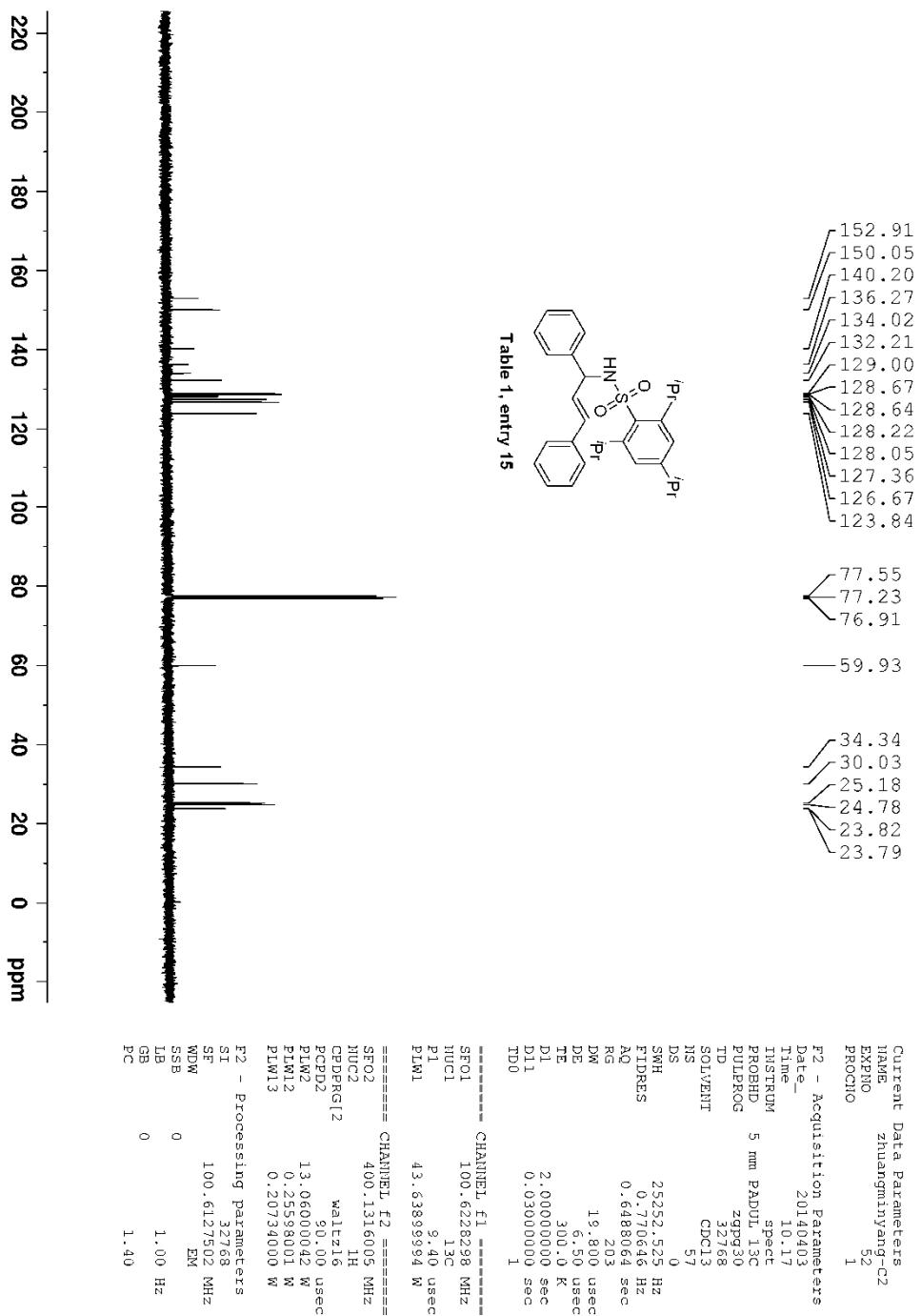
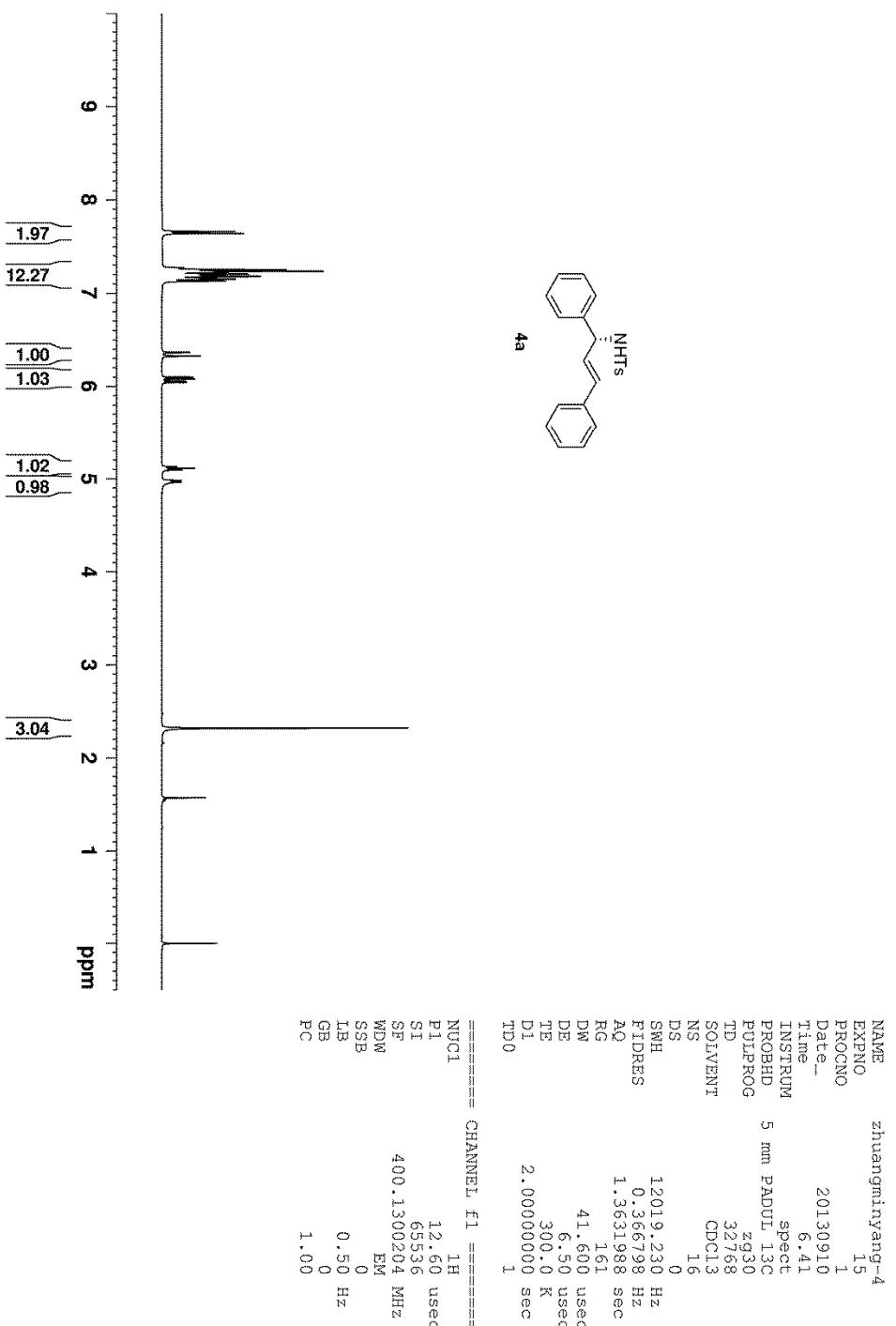
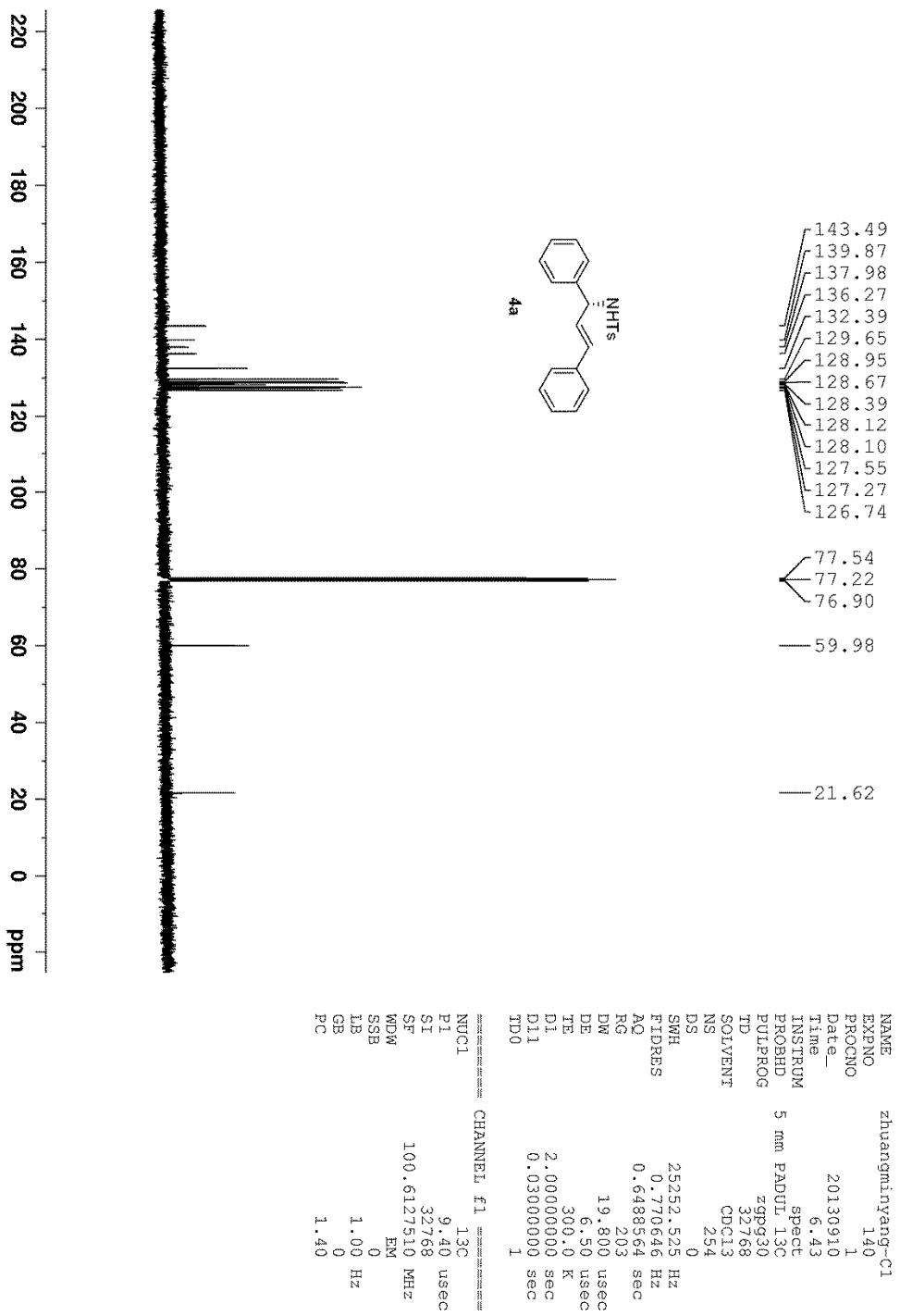
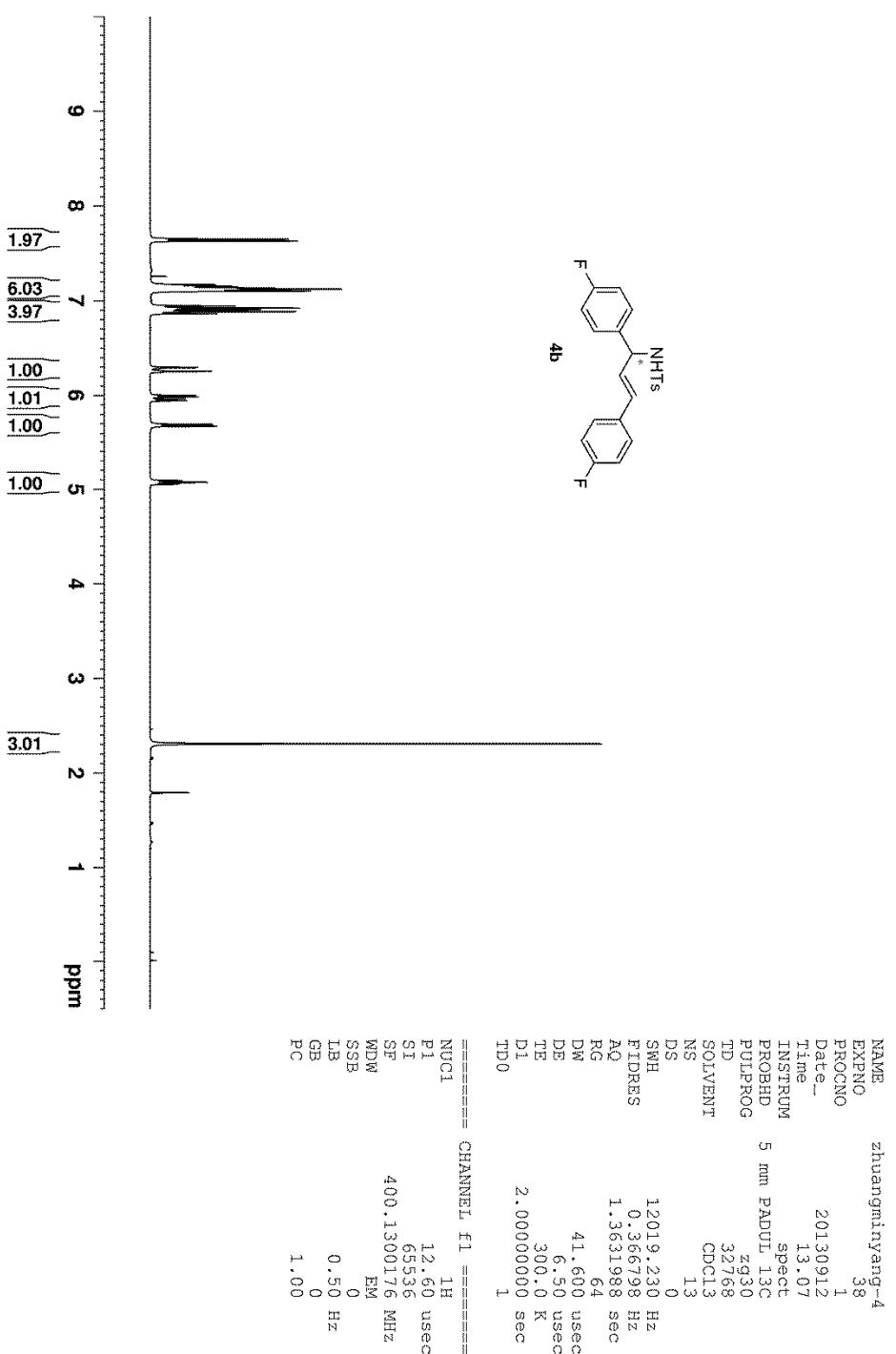


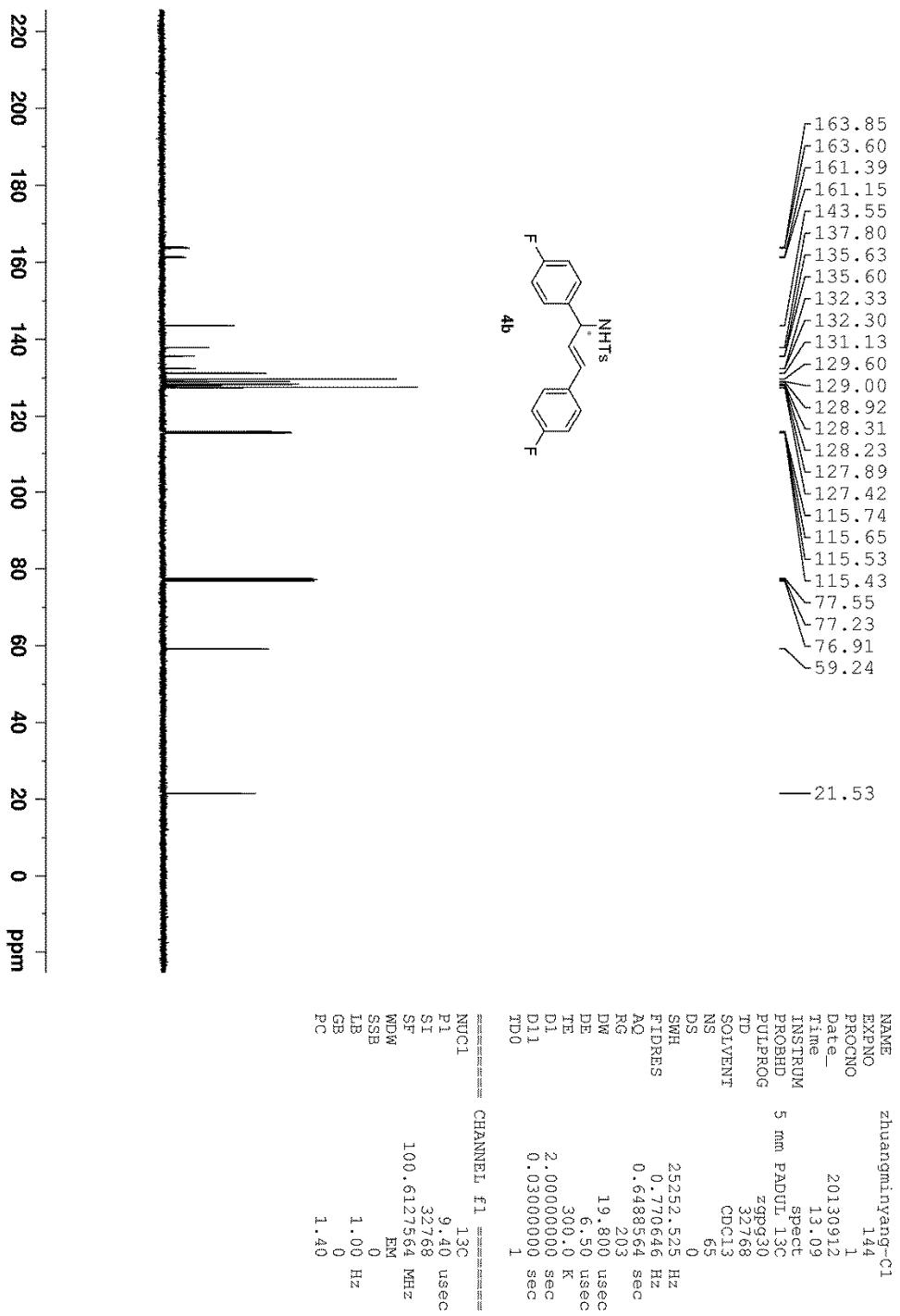
Table 1, entry 15

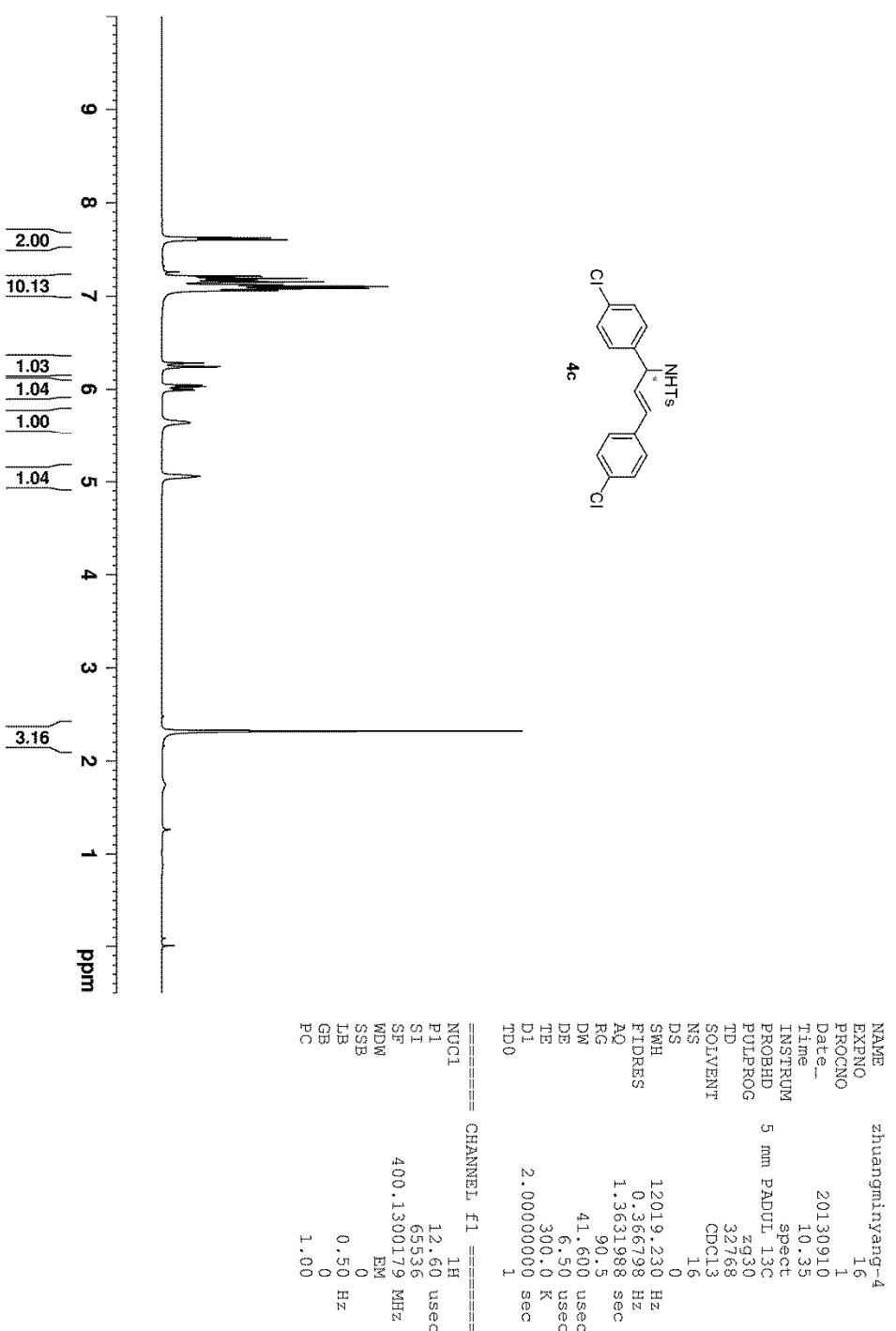


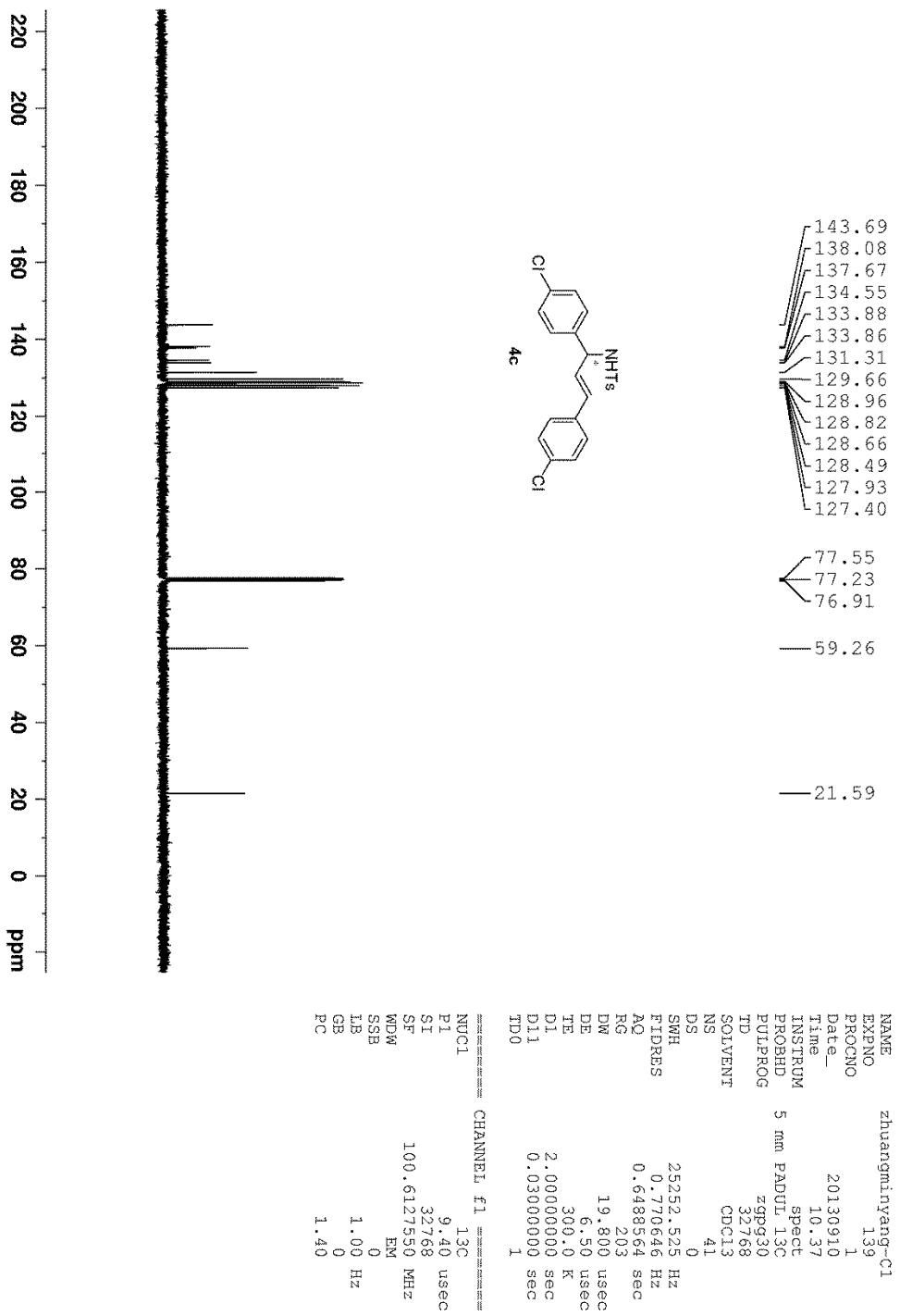


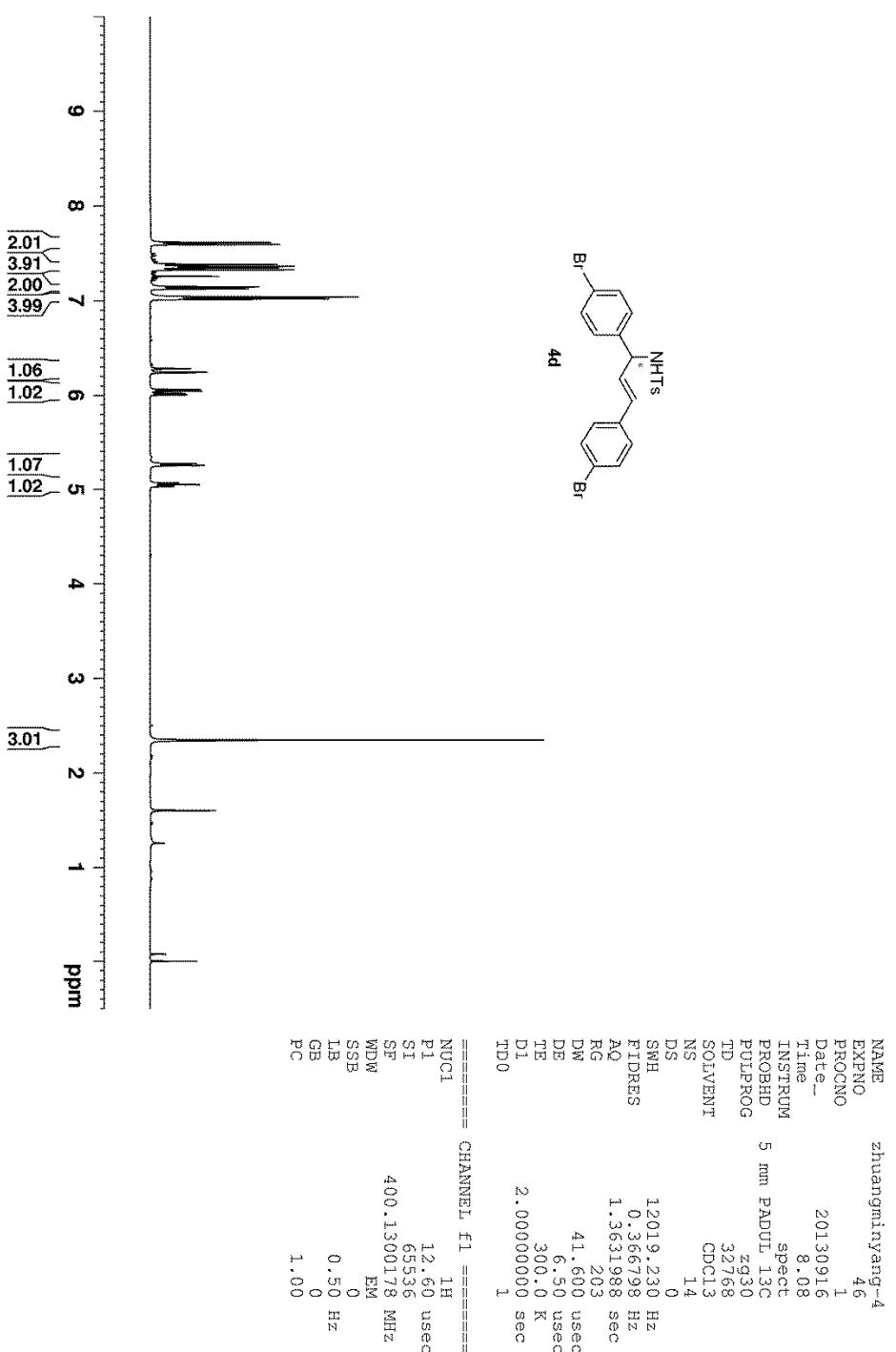


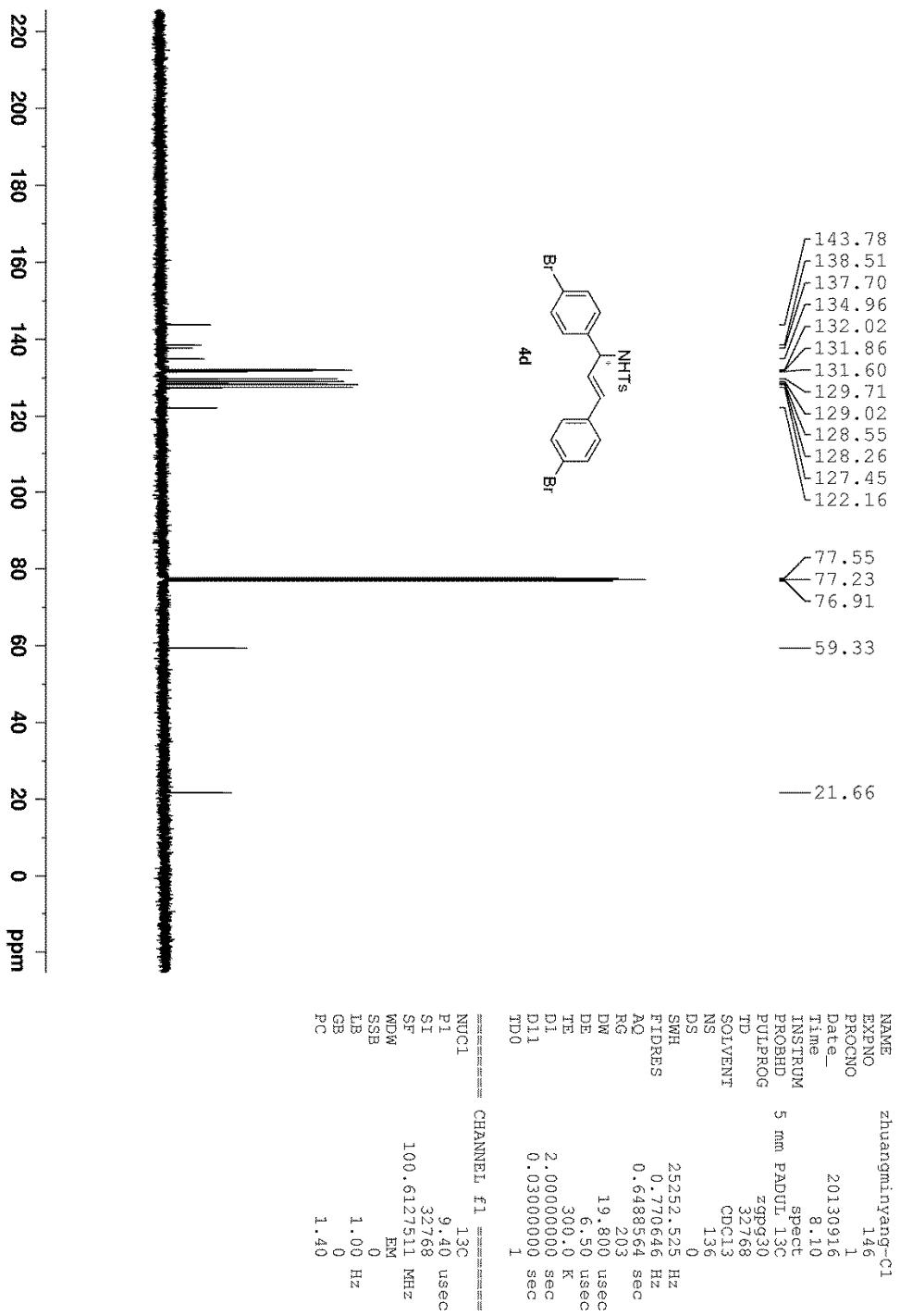


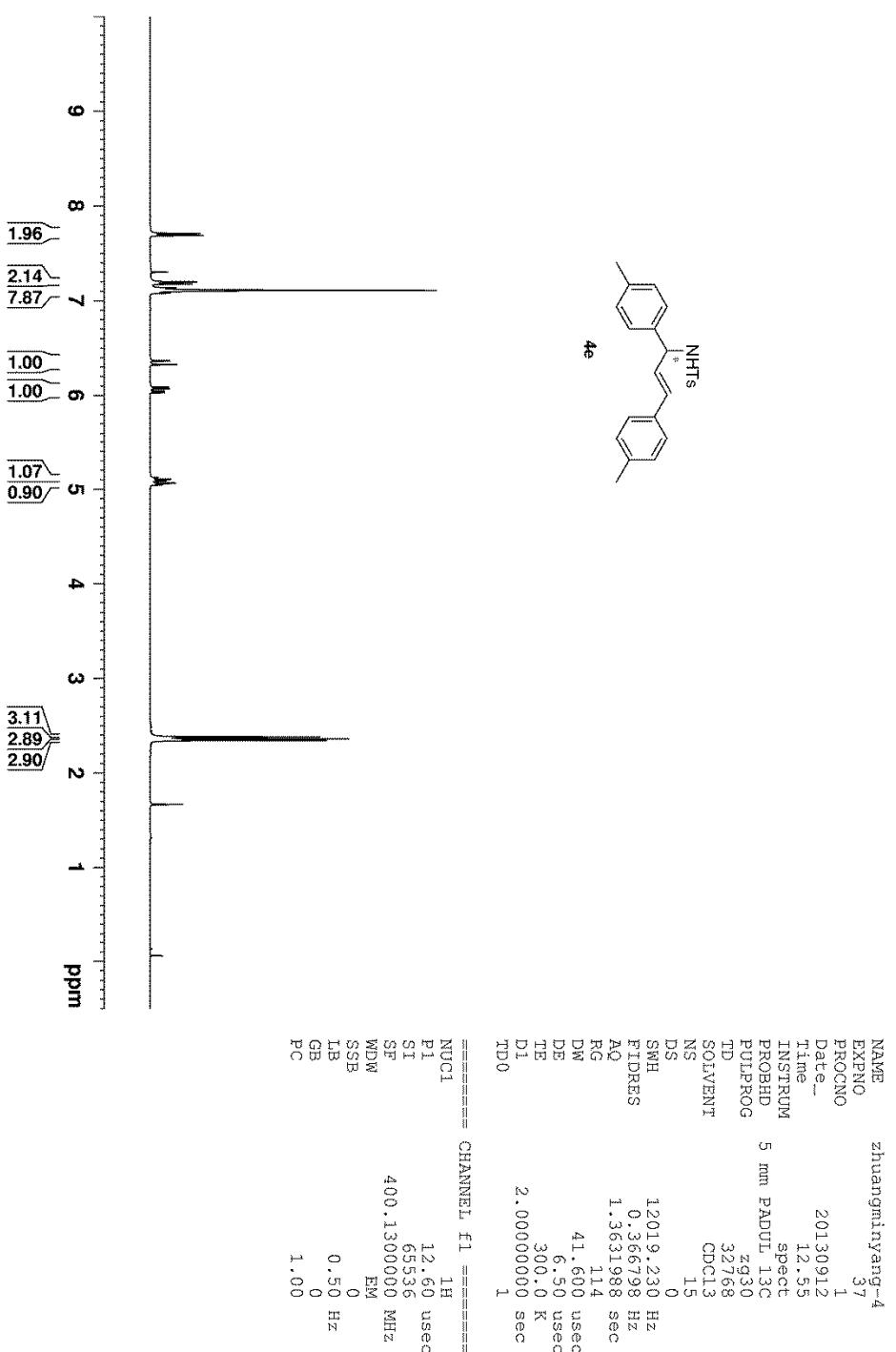


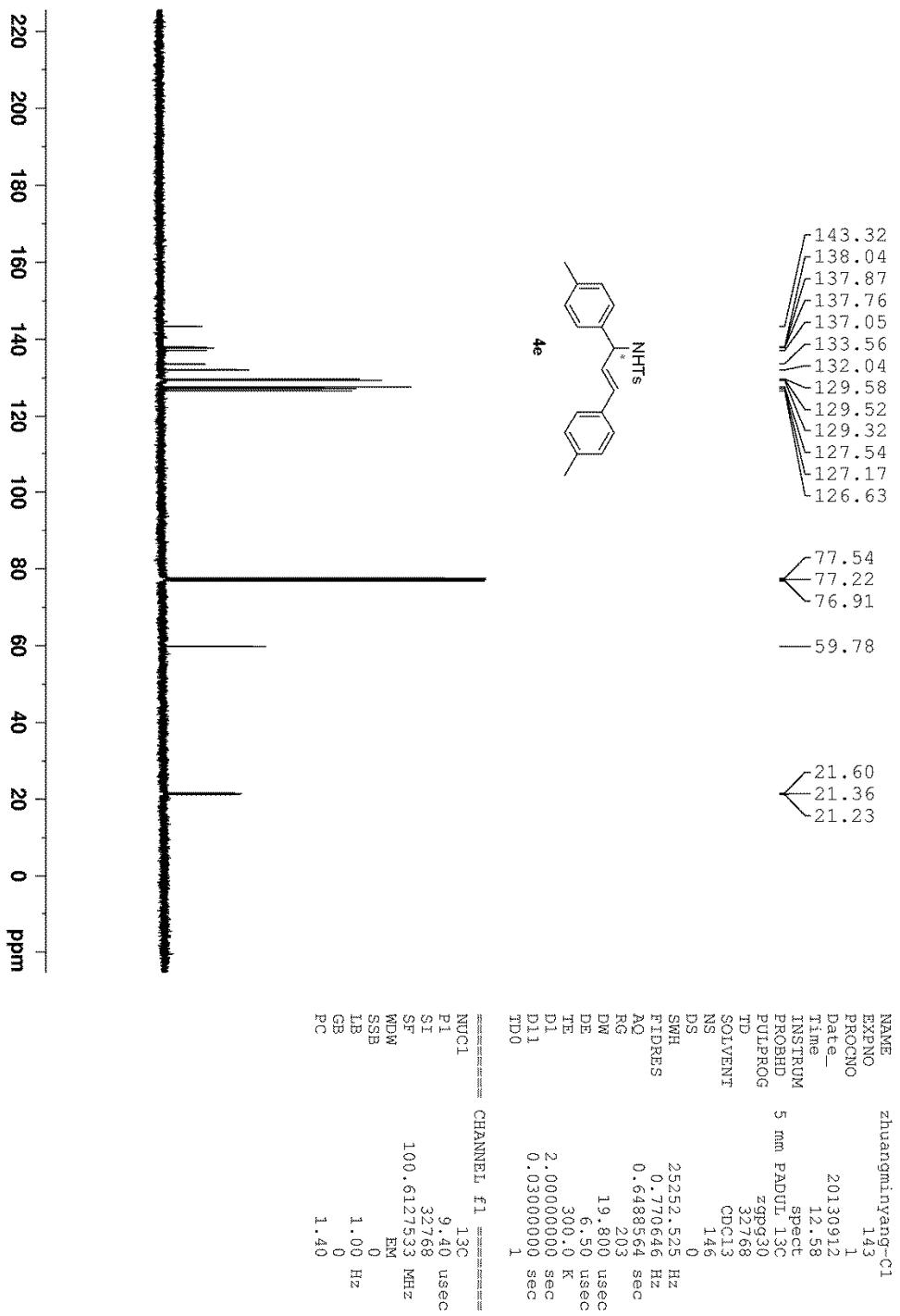


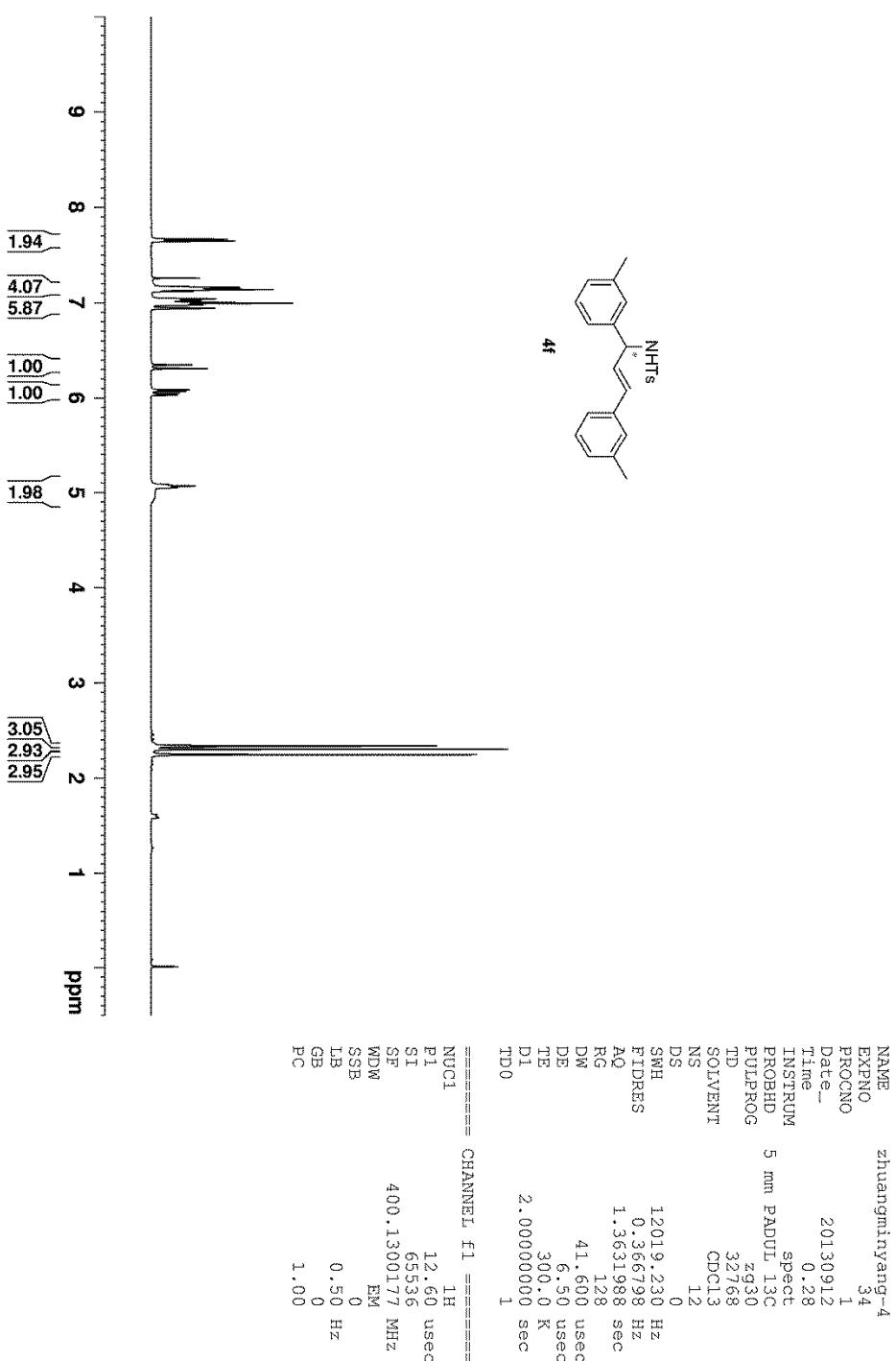




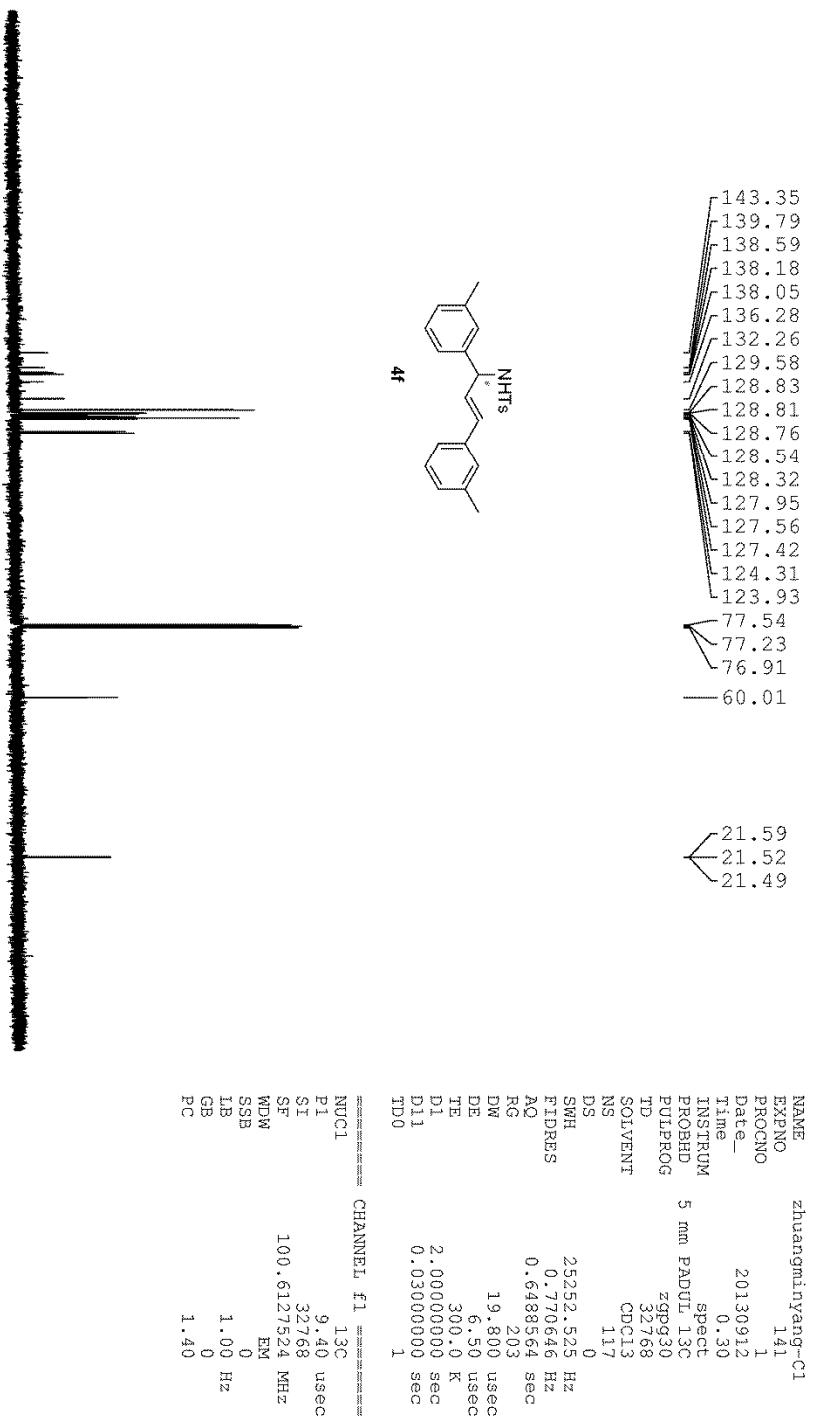


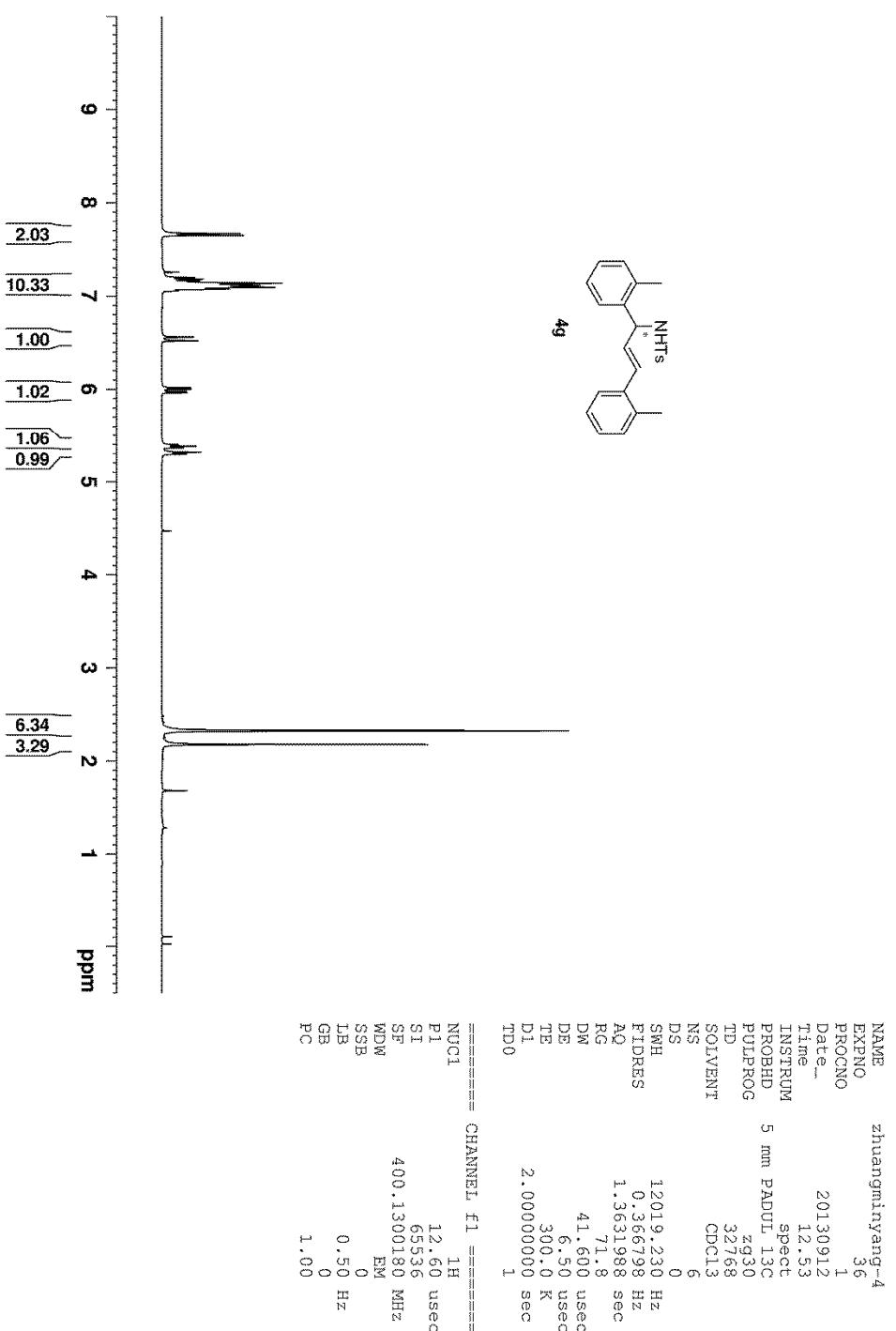


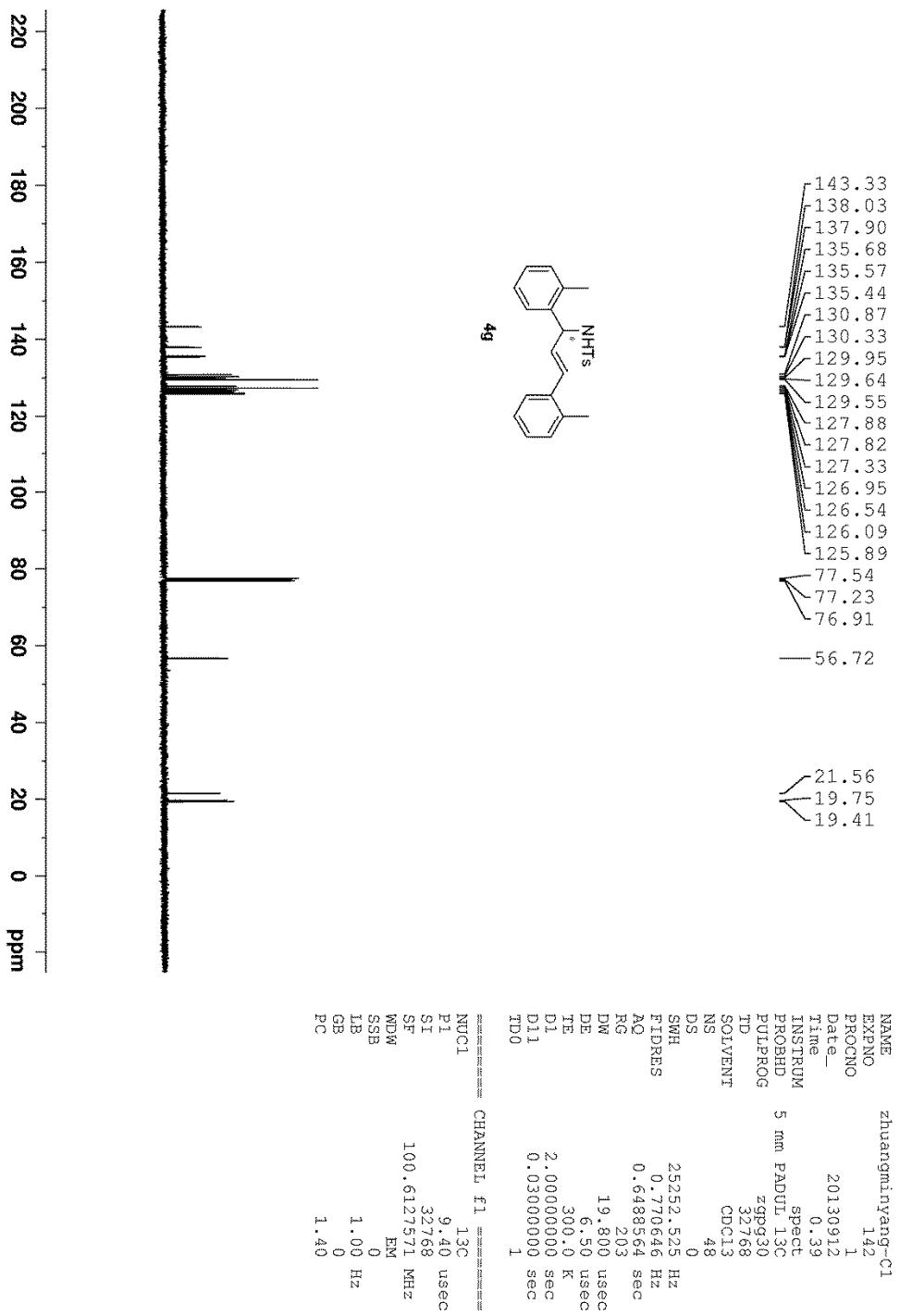


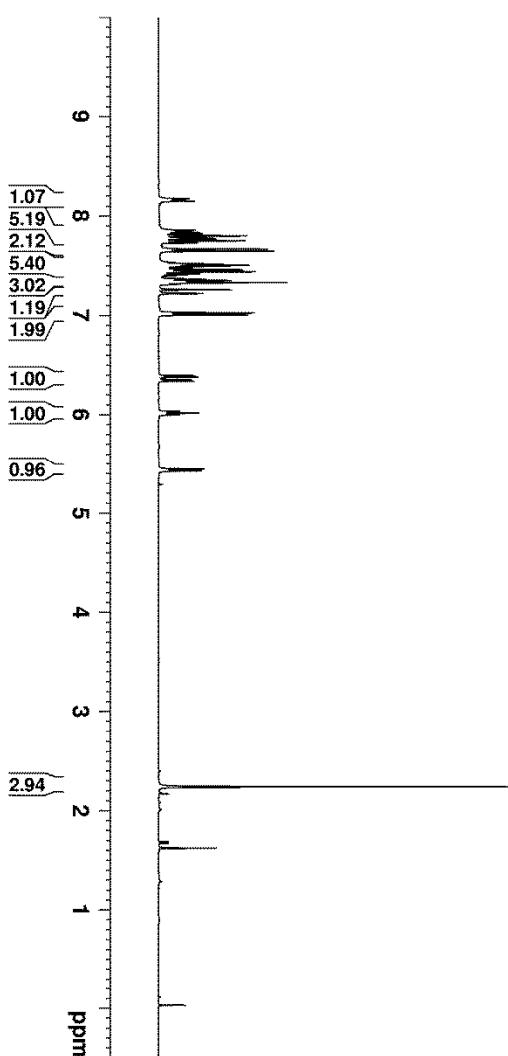
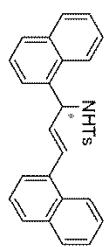


220 200 180 160 140 120 100 80 60 40 20 0 ppm







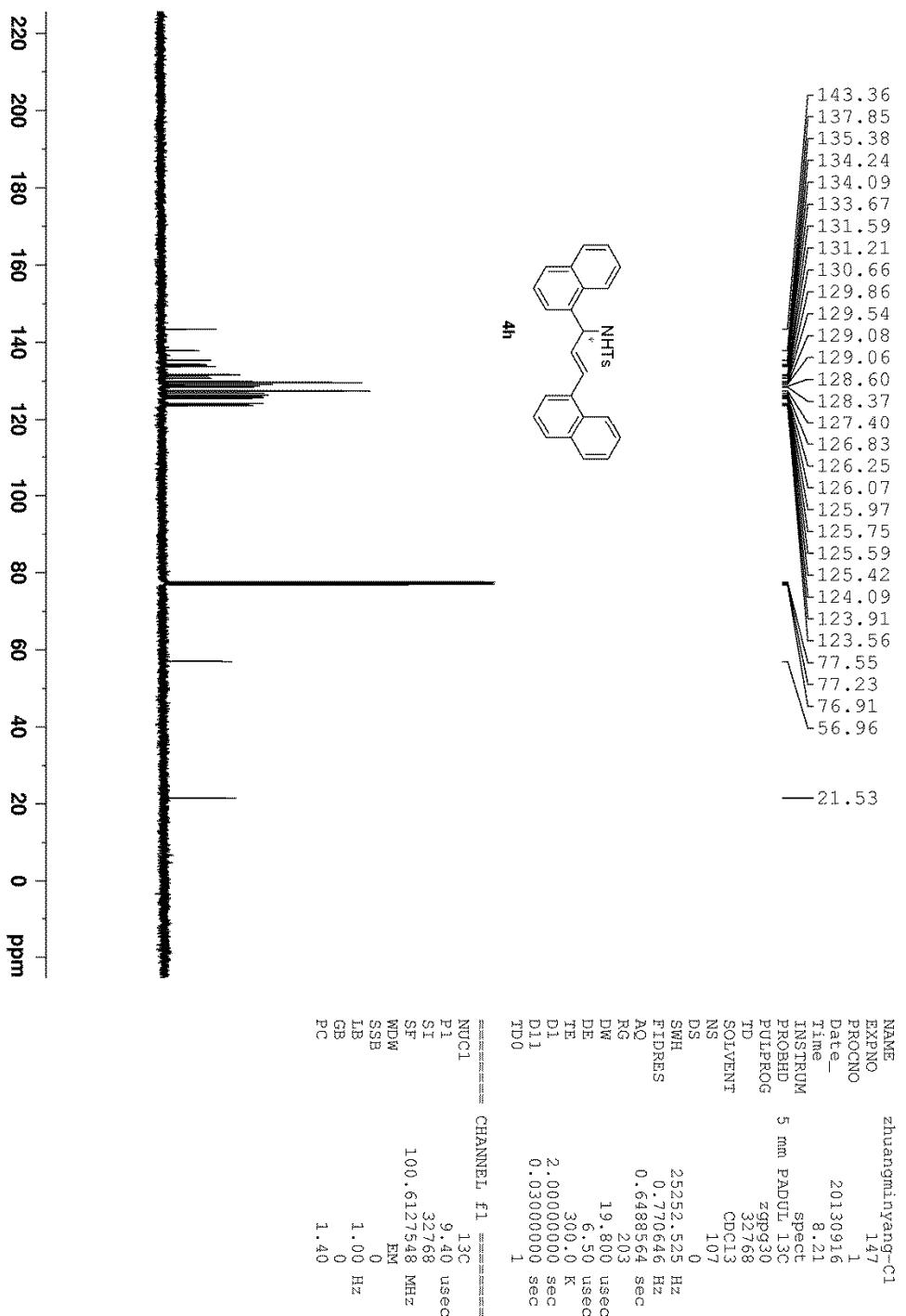


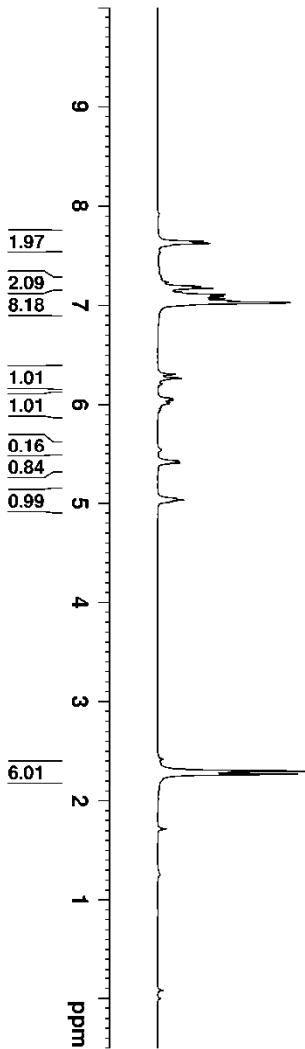
```

=====
NAME          zhuangminyang-4
EXPNO        47
PROCNO       1
Date_        20130916
Time         8.17
INSTRUM      spect
PROBHD      5 mm PADUL 13C
PULPROG     zg30
TD          32768
SOLVENT      CDCl3
NS           15
DS            0
SWH         12019.230 Hz
FIDRES     0.366798 Hz
AQ          1.3631988 sec
RG           101
DW           41.600 usec
DE           6.50 usec
TE           300.0 K
DI          2.0000000 sec
TDD0          1

===== CHANNEL f1 =====
NUC1          1H
P1           12.60 usec
SI            65536
SF          400.1300176 MHz
MW            EM
SSB           0
LB            0.50 Hz
GB            0
PC           1.00

```



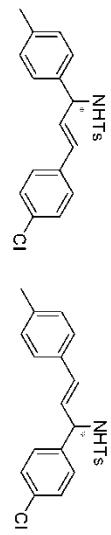


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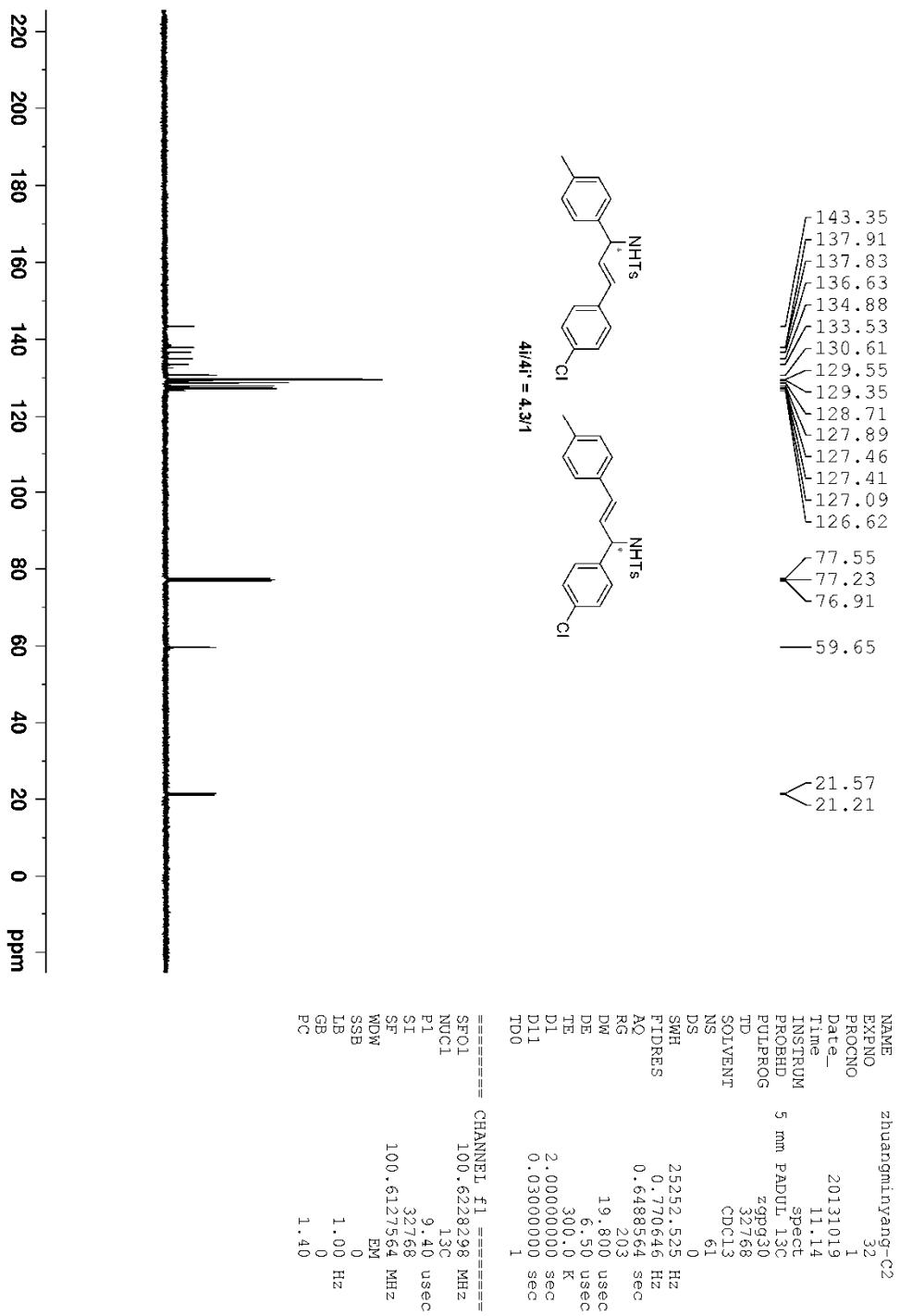
NAME          zhuangminyang-5
EXPGNO       14
PROCNO       1
Date_        20131019
Time_        11.13
INSTRUM      spect
PROBHD      5 mm PADUL 13C
PULPROG     zg30
TD          32768
SOLVENT      CDCl3
NS           12
DS           0
SWH         12019.230 Hz
F1RES       0.366798 Hz
AQ          1.363198 sec
RG           64
DW           41.600 usec
DE           6.50 usec
TE           300.0 K
D1          2.0000000 sec
TD0          1

===== CHANNEL f1 =====
SPOL        400.1320007 MHz
NUC1        1H
P1          12.60 usec
SI           65536
SF          400.1300184 MHz
WDW        EM
SSB          0
LB          0.50 Hz
GB          1.00
PC

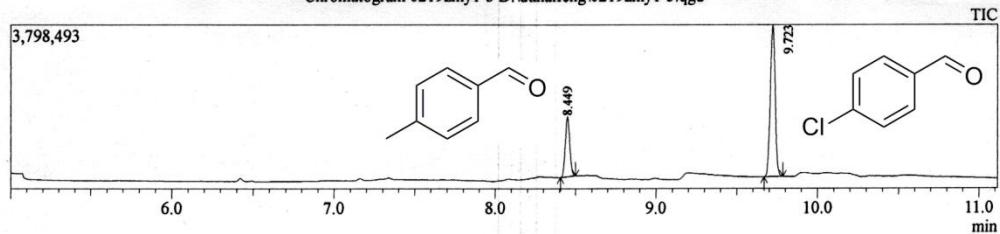
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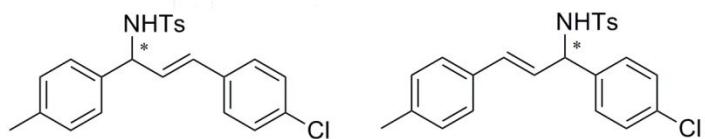
$4i/4i' = 4.3/1$



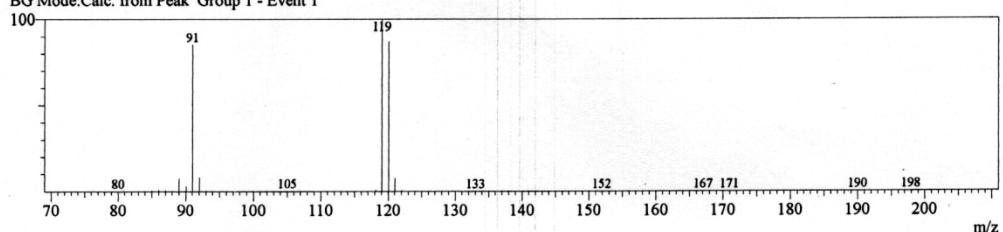
Chromatogram 0219zmy1-3 D:\duhaifeng\0219zmy1-3.qgd



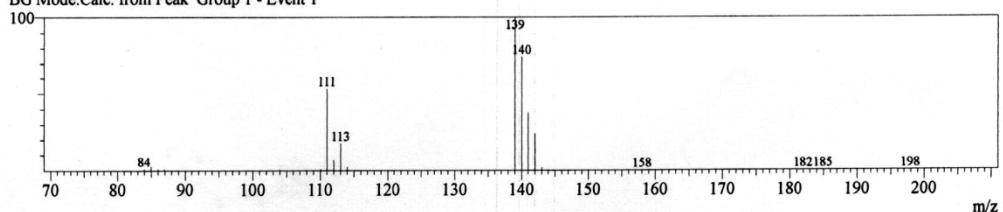
Peak#	Peak Report TIC						A/H	Mark	Name
	R.Time	I.Time	F.Time	Area	Area%	Height	Height%		
1	8.449	8.405	8.500	2661711	26.45	1399683	28.45	1.90	
2	9.723	9.670	9.785	7402449	73.55	3520924	71.55	2.10	
				10064160	100.00	4920607	100.00		

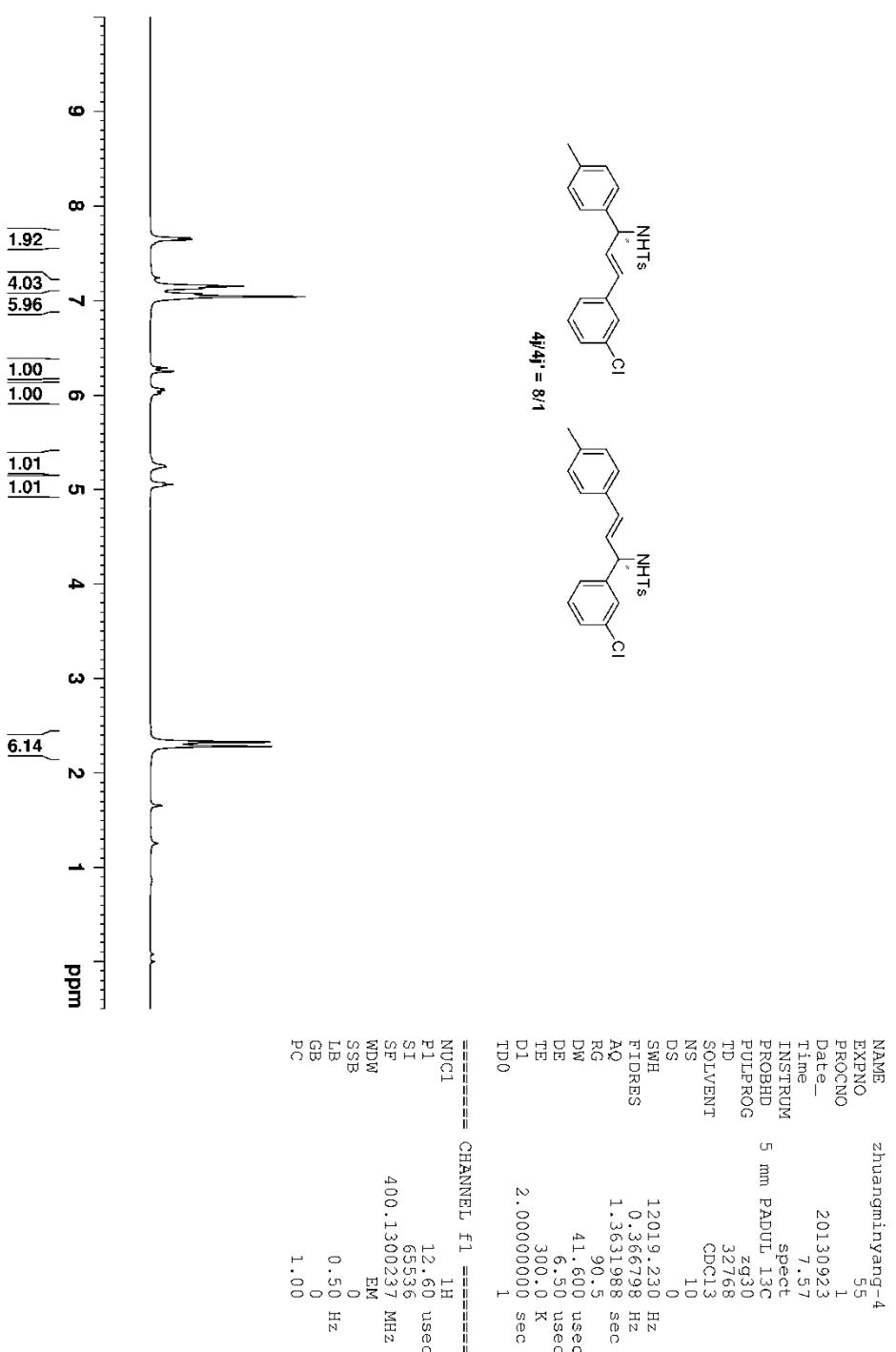


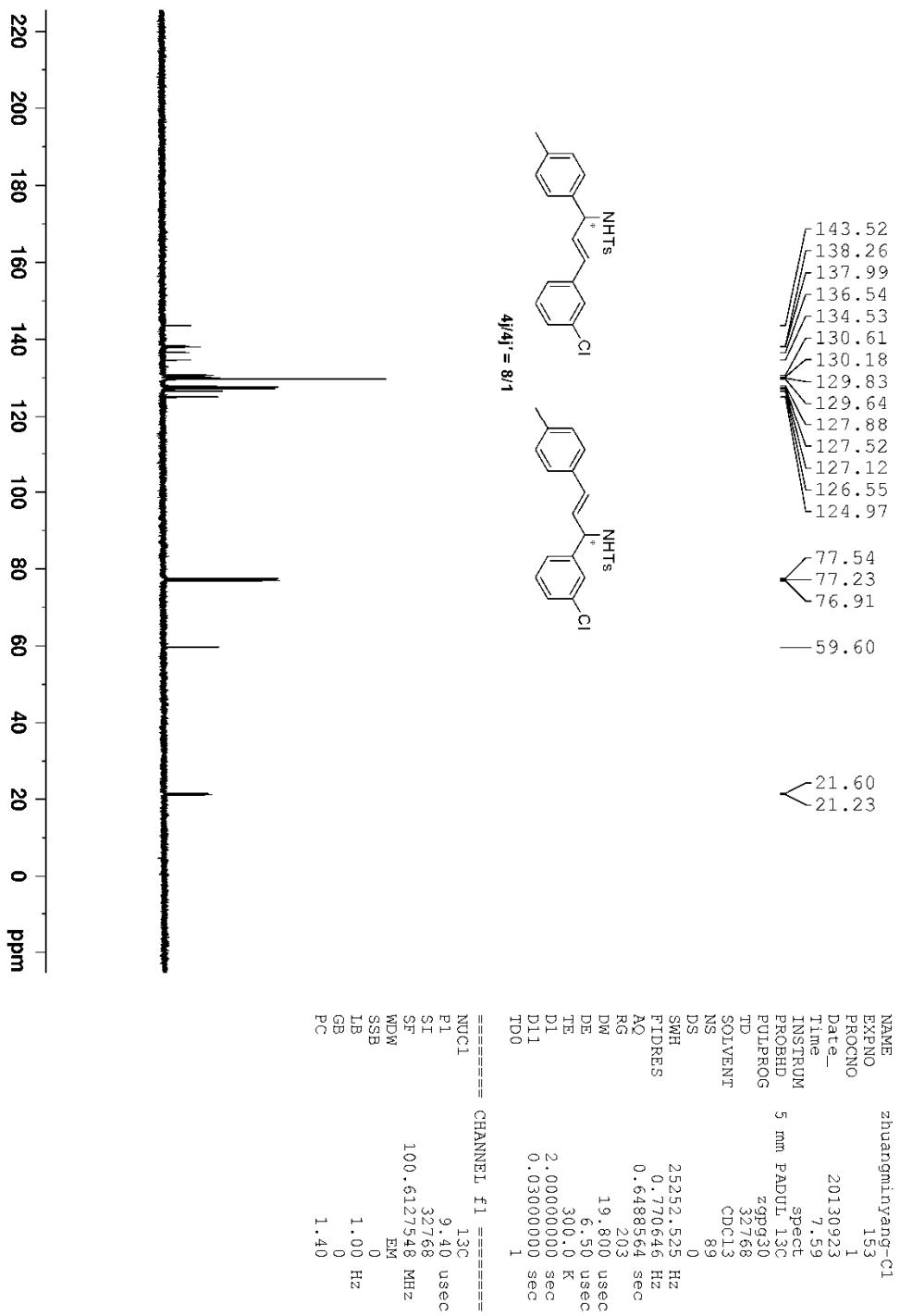
Line#1 R.Time:8.450(Scan#:691)
MassPeaks:64
RawMode:Averaged 8.445-8.455(690-692) BasePeak:119(438317)
BG Mode:Calc. from Peak Group 1 - Event 1

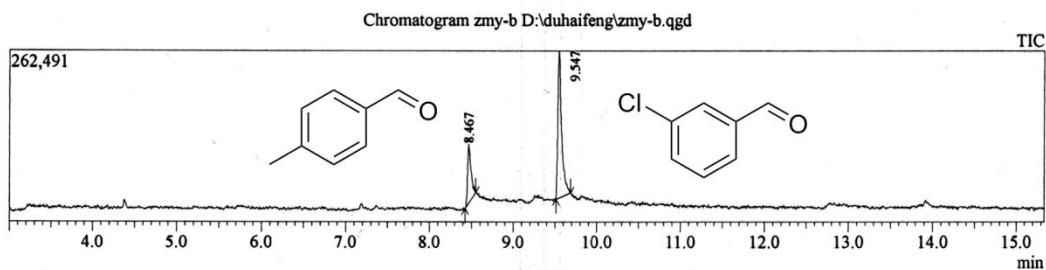


Line#2 R.Time:9.725(Scan#:946)
MassPeaks:78
RawMode:Averaged 9.720-9.730(945-947) BasePeak:139(1037101)
BG Mode:Calc. from Peak Group 1 - Event 1





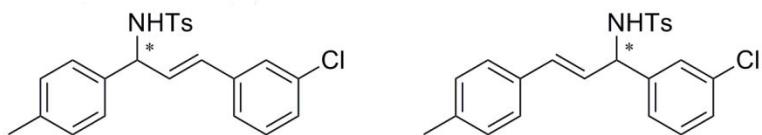




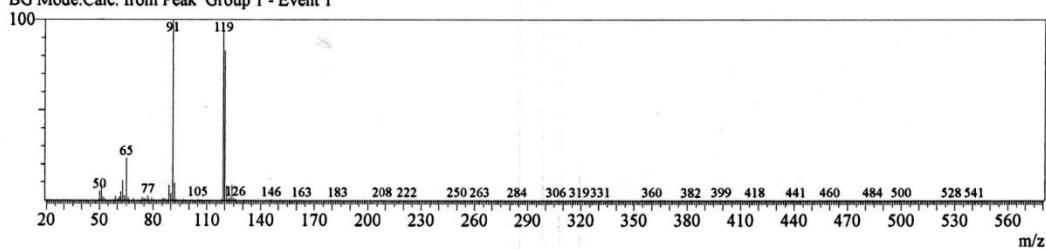
Peak#	R.Time	I.Time	F.Time	Peak Report TIC			
				Area	Area%	Height	Height%
1	8.467	8.425	8.550	262579	26.90	90925	28.76
2	9.547	9.510	9.685	713388	73.10	225185	71.24

A/H Mark Name

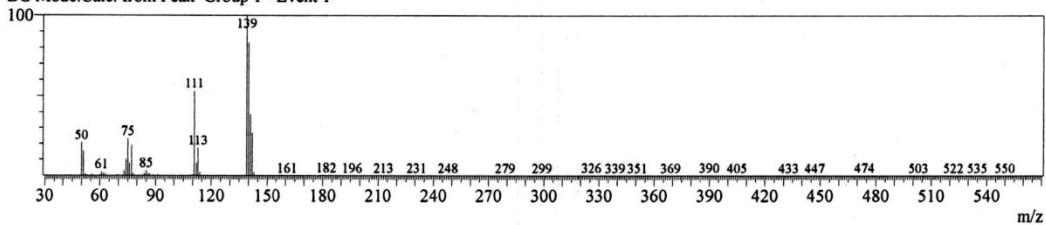
975967 100.00 316110 100.00

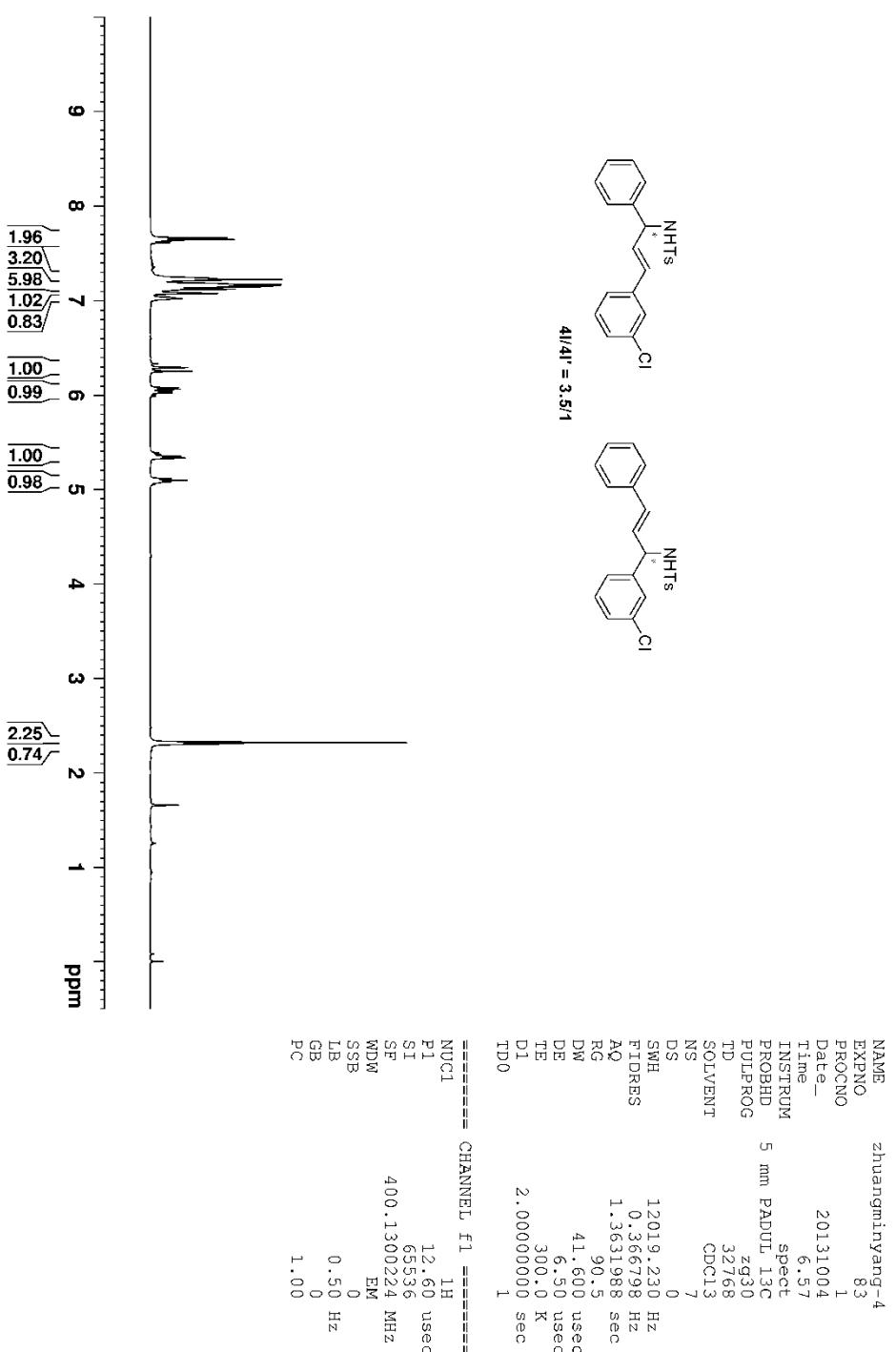


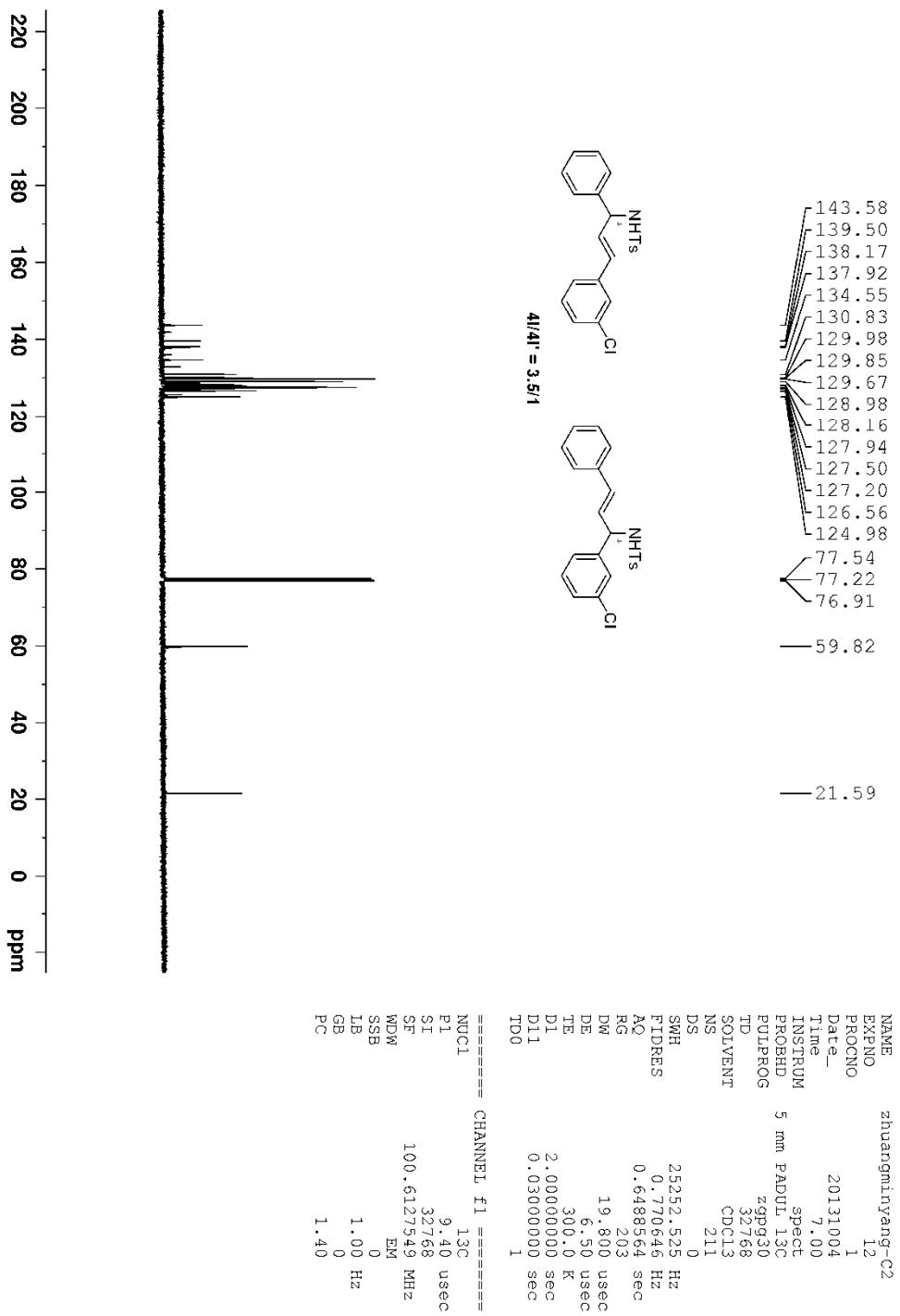
Line#:1 R.Time:8.465(Scan#:1094)
MassPeaks:295
RawMode:Averaged 8.460-8.470(1093-1095) BasePeak:91(21669)
BG Mode:Calc. from Peak Group 1 - Event 1

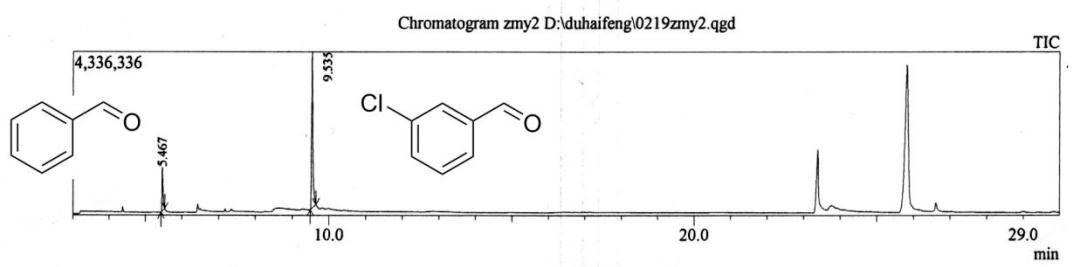


Line#:2 R.Time:9.545(Scan#:1310)
MassPeaks:280
RawMode:Averaged 9.540-9.550(1309-1311) BasePeak:139(48619)
BG Mode:Calc. from Peak Group 1 - Event 1

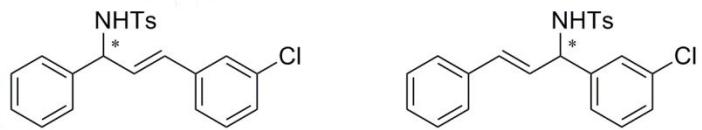






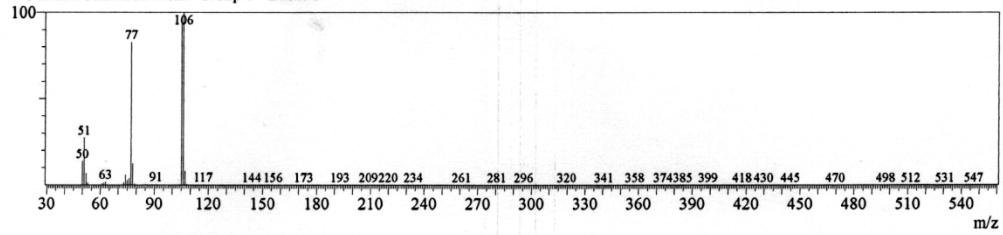


Peak#	R.Time	I.Time	F.Time	Peak Report TIC				A/H	Mark	Name
				Area	Area%	Height	Height%			
1	5.467	5.435	5.535	2330264	18.62	1161463	21.91	2.01		
2	9.535	9.490	9.645	10186921	81.38	4139118	78.09	2.46		
				12517185	100.00	5300581	100.00			

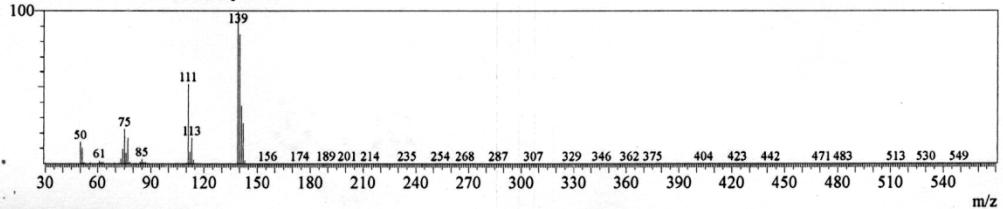


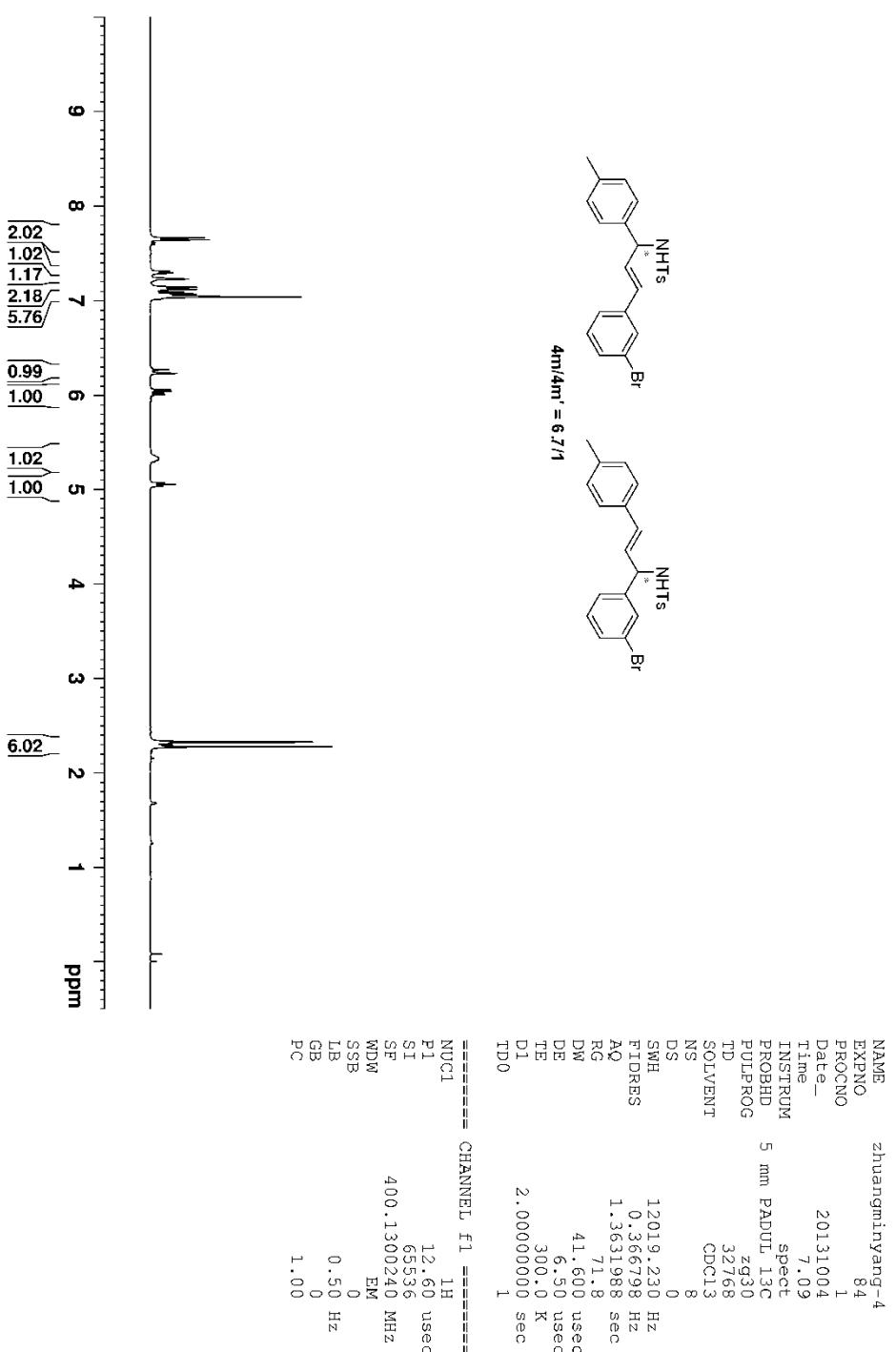
$$4I/I' = 3.5/1$$

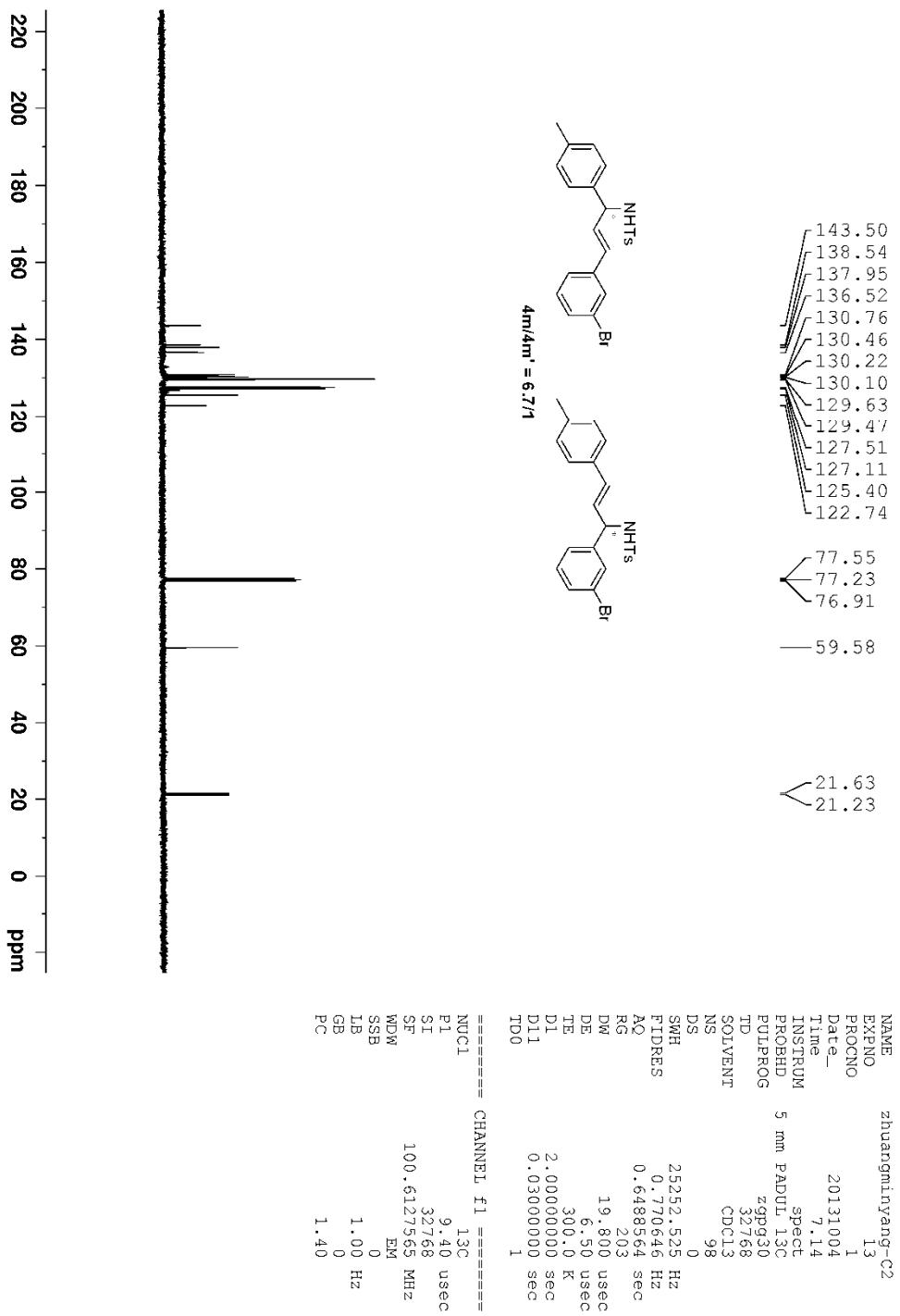
Line#:1 R.Time:5.465(Scan#:494)
MassPeaks:258
RawMode:Averaged 5.460-5.470(493-495) BasePeak:106(293239)
BG Mode:Calc. from Peak Group 1 - Event 1

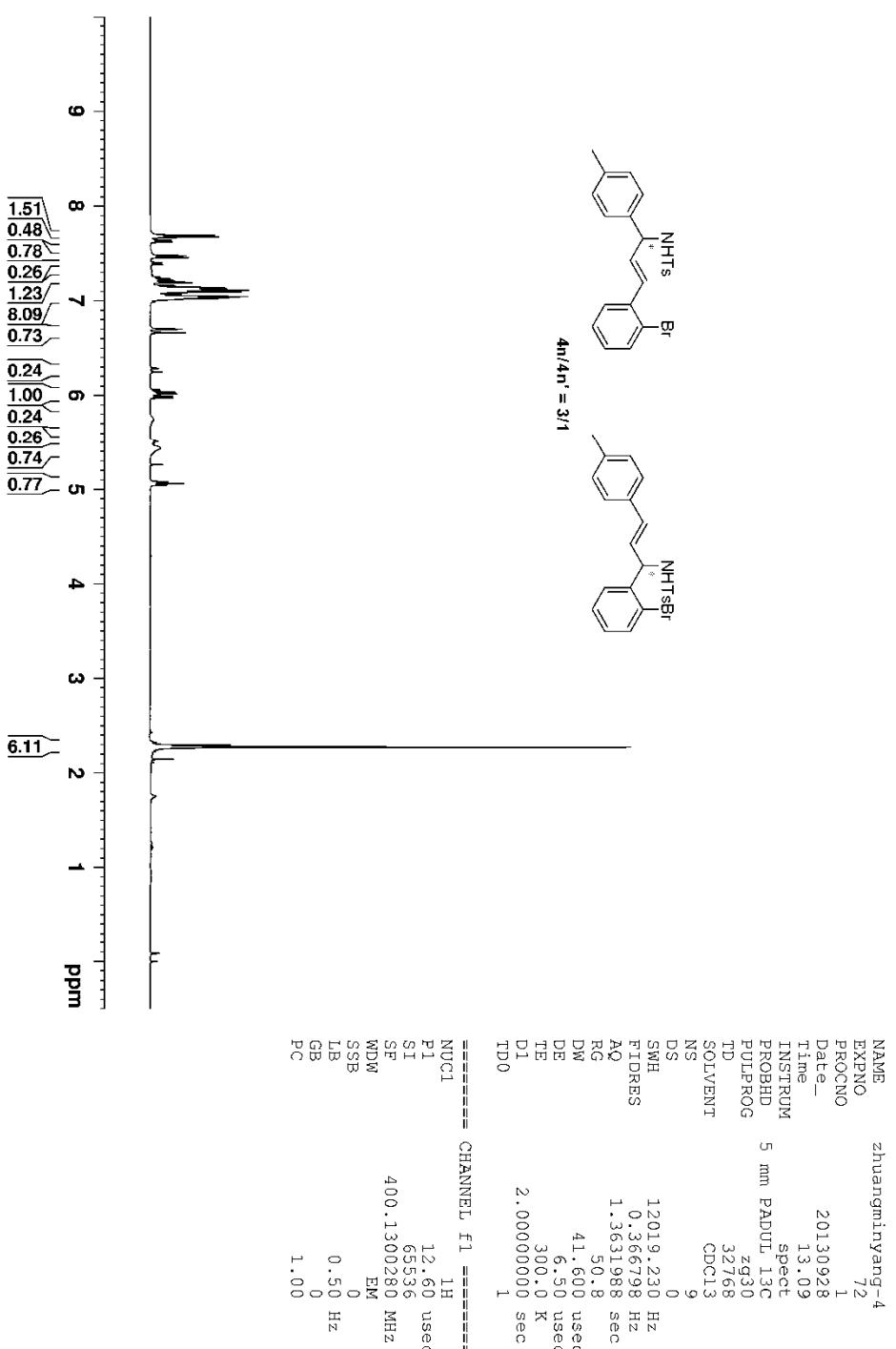


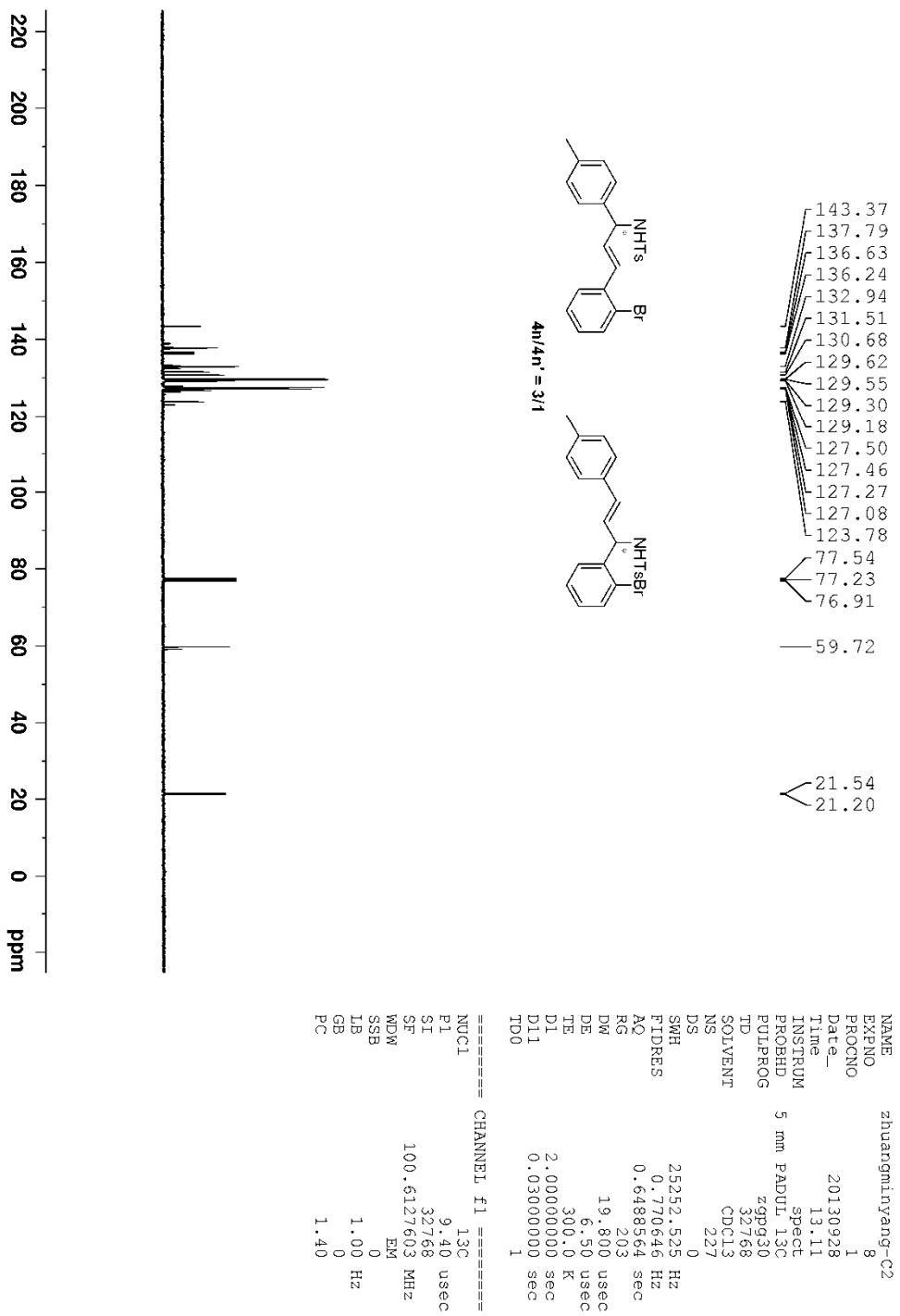
Line#:2 R.Time:9.535(Scan#:1308)
MassPeaks:305
RawMode:Averaged 9.530-9.540(1307-1309) BasePeak:139(923786)
BG Mode:Calc. from Peak Group 1 - Event 1



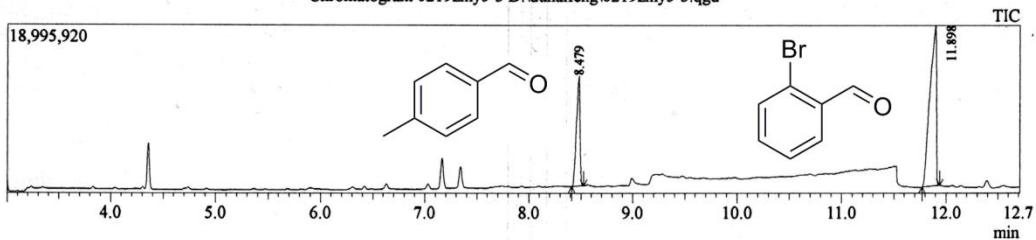








Chromatogram 0219zmy5-3 D:\duhaifeng\0219zmy5-3.qgd



Peak#	R.Time	I.Time	F.Time	Peak Report TIC			
				Area	Area%	Height	Height%
1	8.479	8.410	8.525	29242938	27.05	12468853	40.74
2	11.898	11.770	11.940	78875584	72.95	18138877	59.26
				108118522	100.00	30607730	100.00

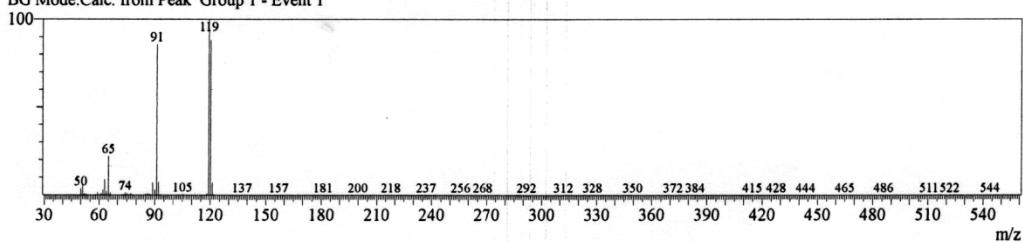


Line#:1 R.Time:8.480(Scan#:1097)

MassPeaks:254

RawMode:Averaged 8.475-8.485(1096-1098) BasePeak:119(3256888)

BG Mode:Calc. from Peak Group 1 - Event 1

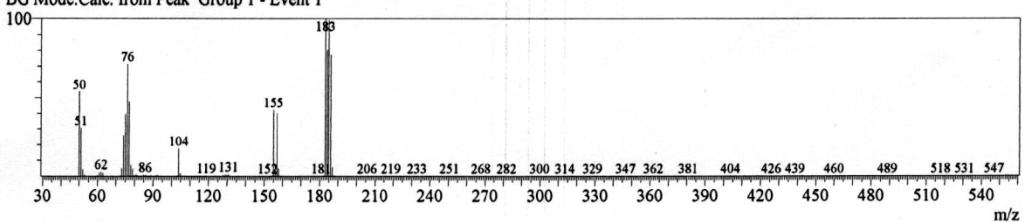


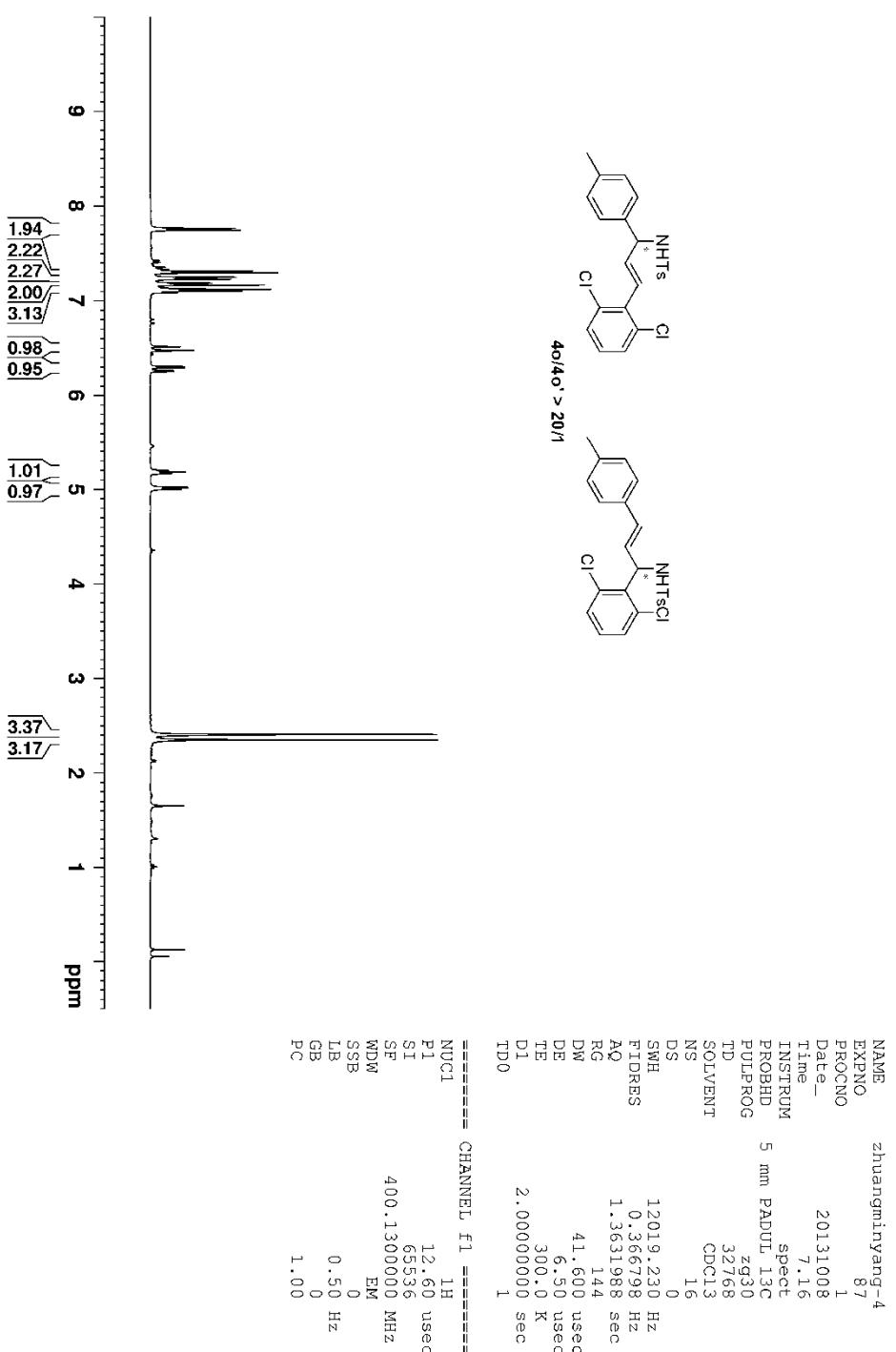
Line#:2 R.Time:11.900(Scan#:1781)

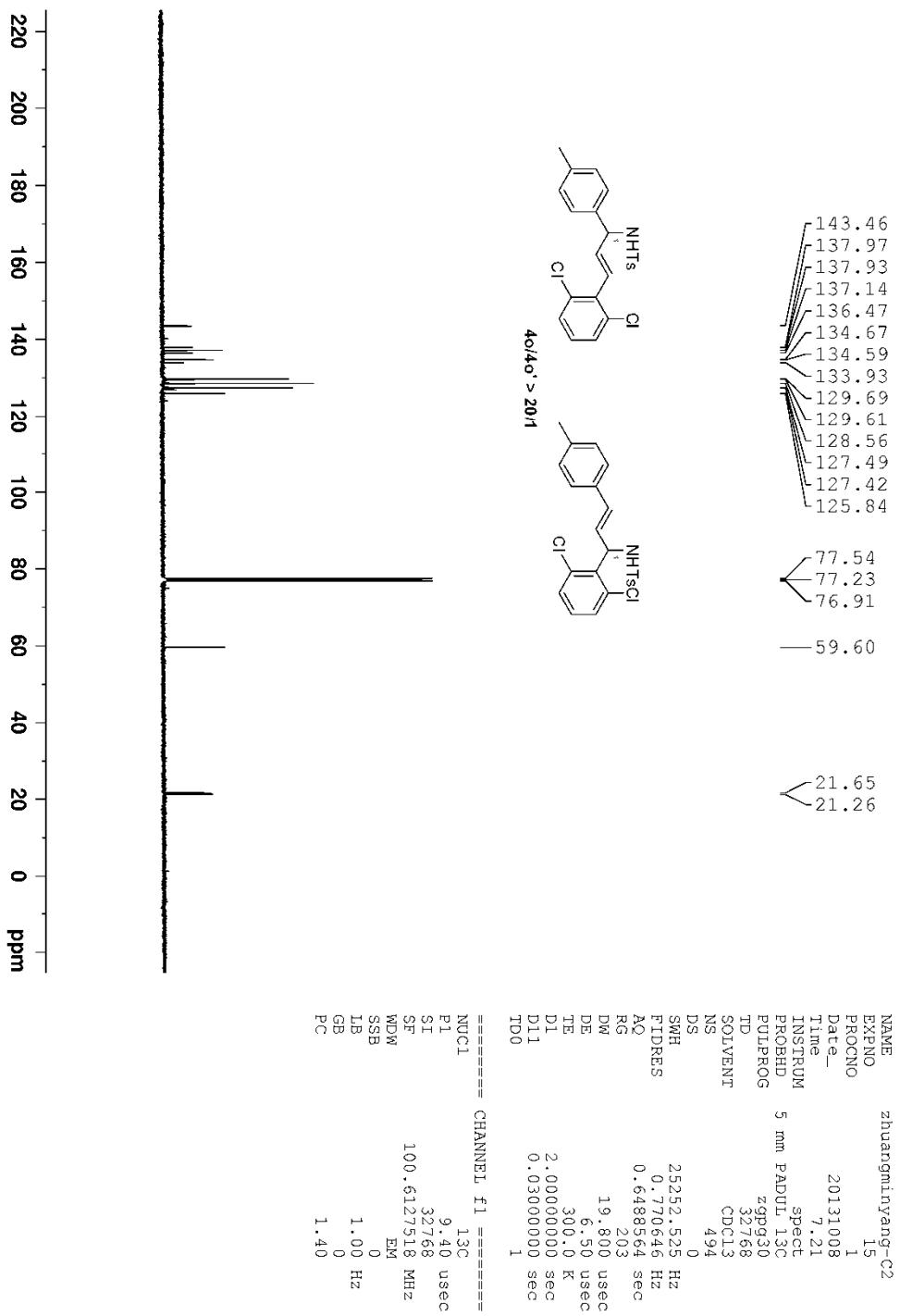
MassPeaks:279

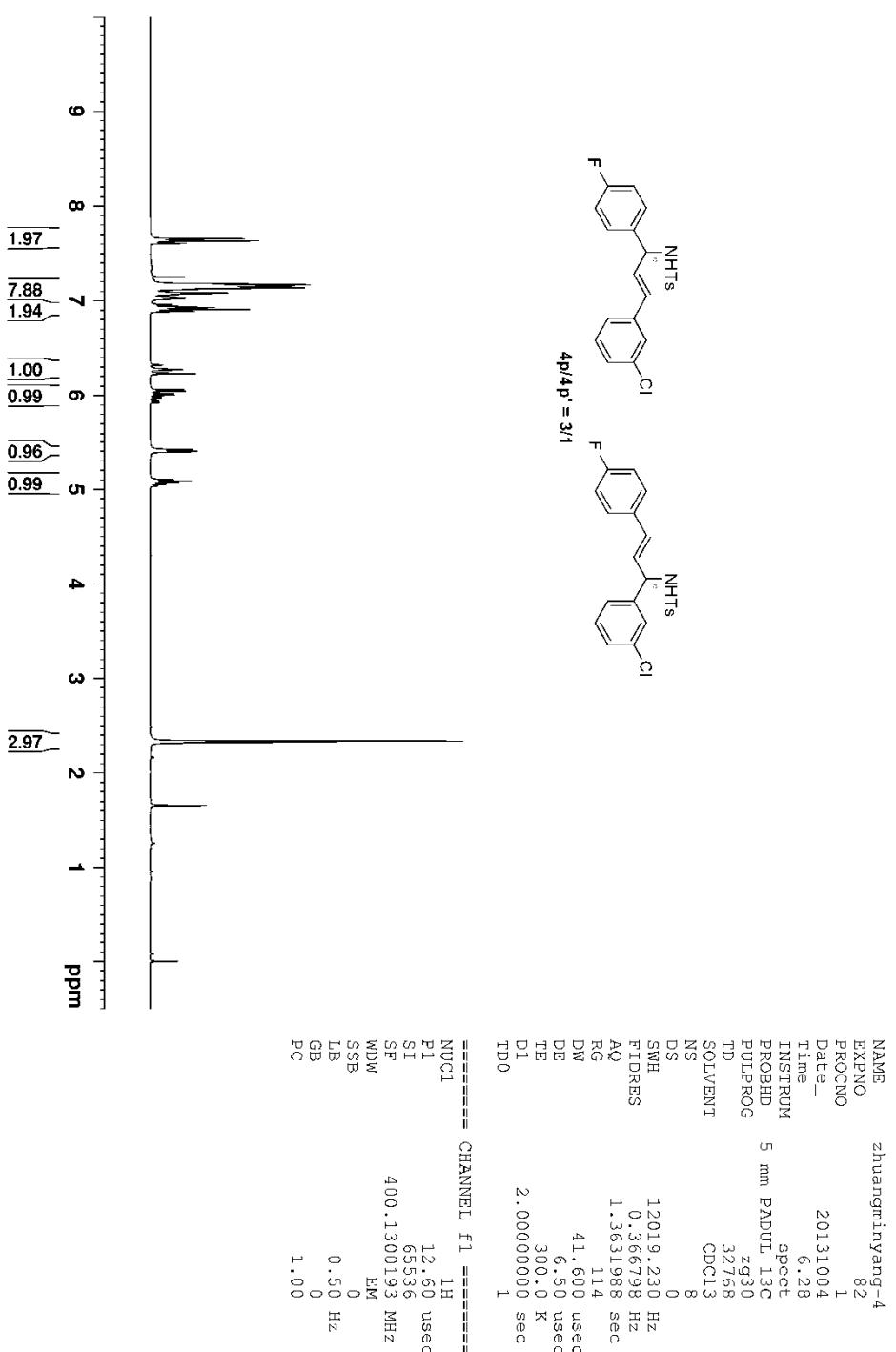
RawMode:Averaged 11.895-11.905(1780-1782) BasePeak:183(2211311)

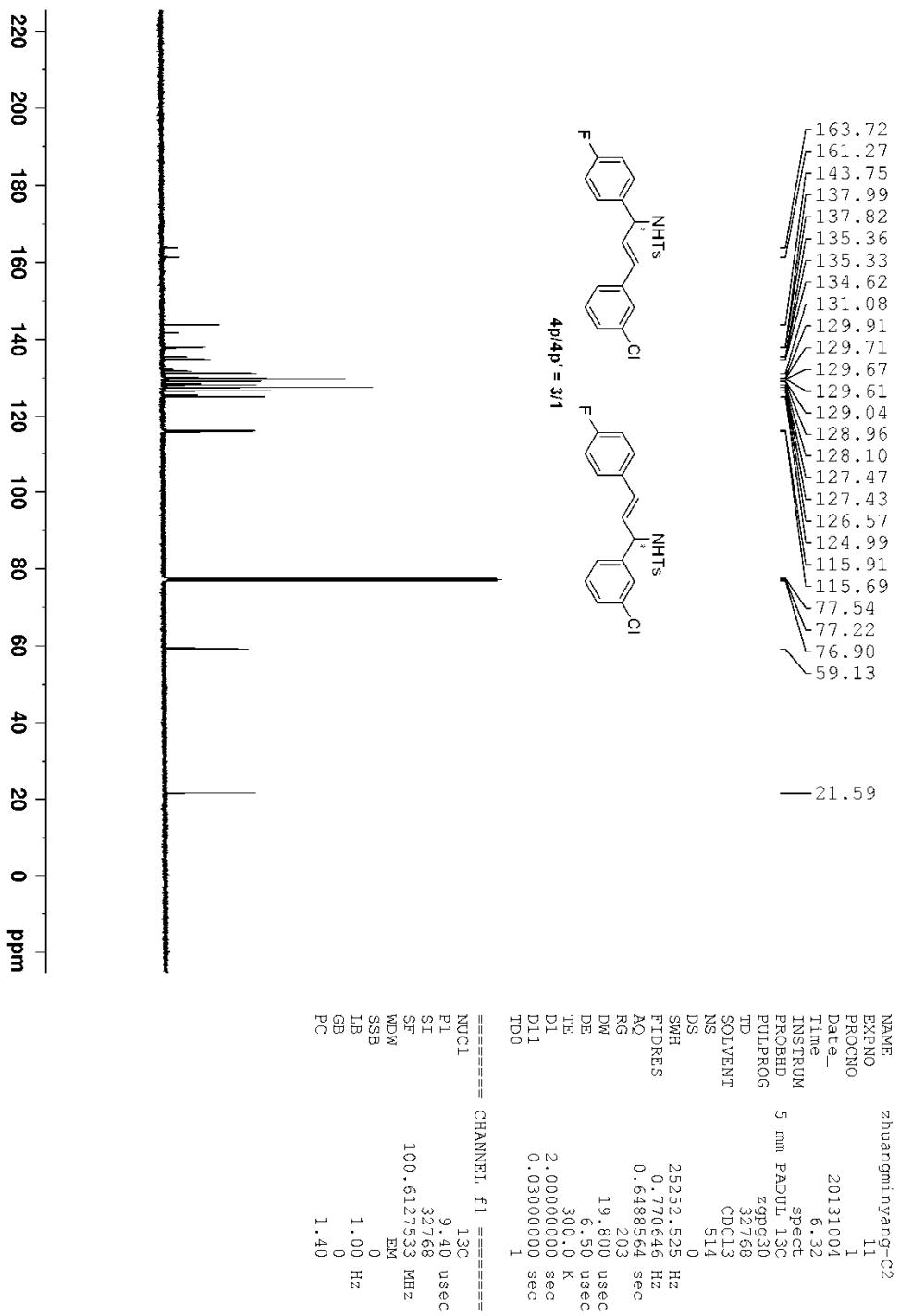
BG Mode:Calc. from Peak Group 1 - Event 1

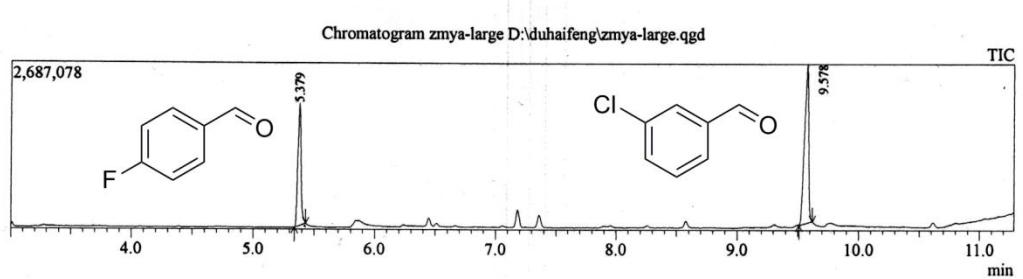




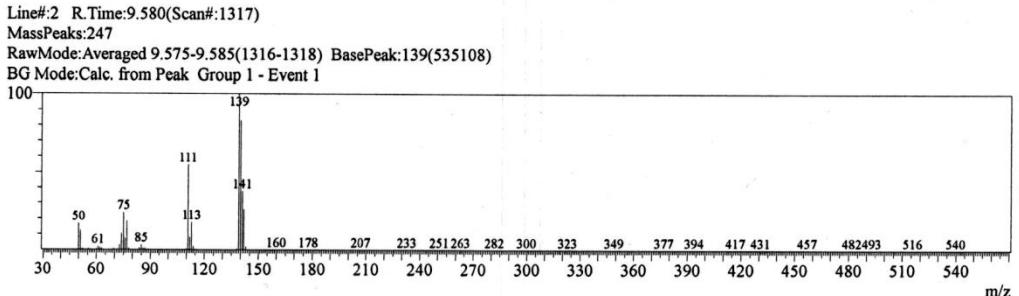
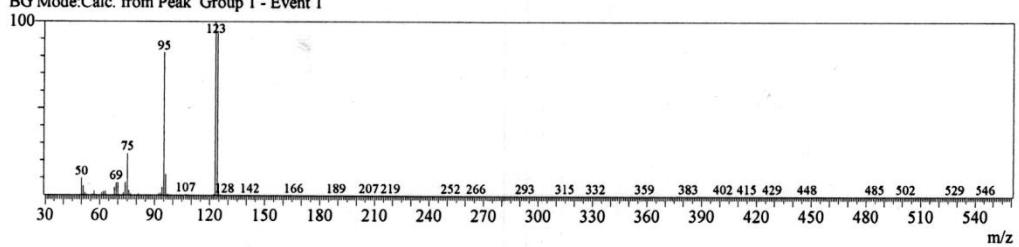
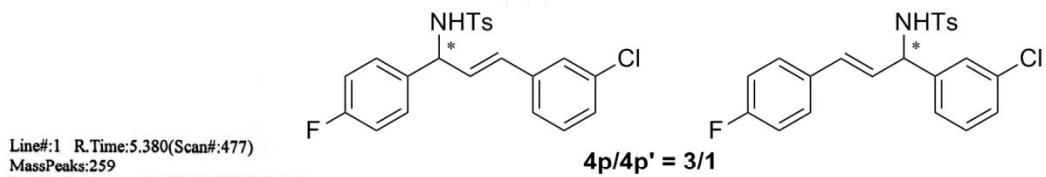


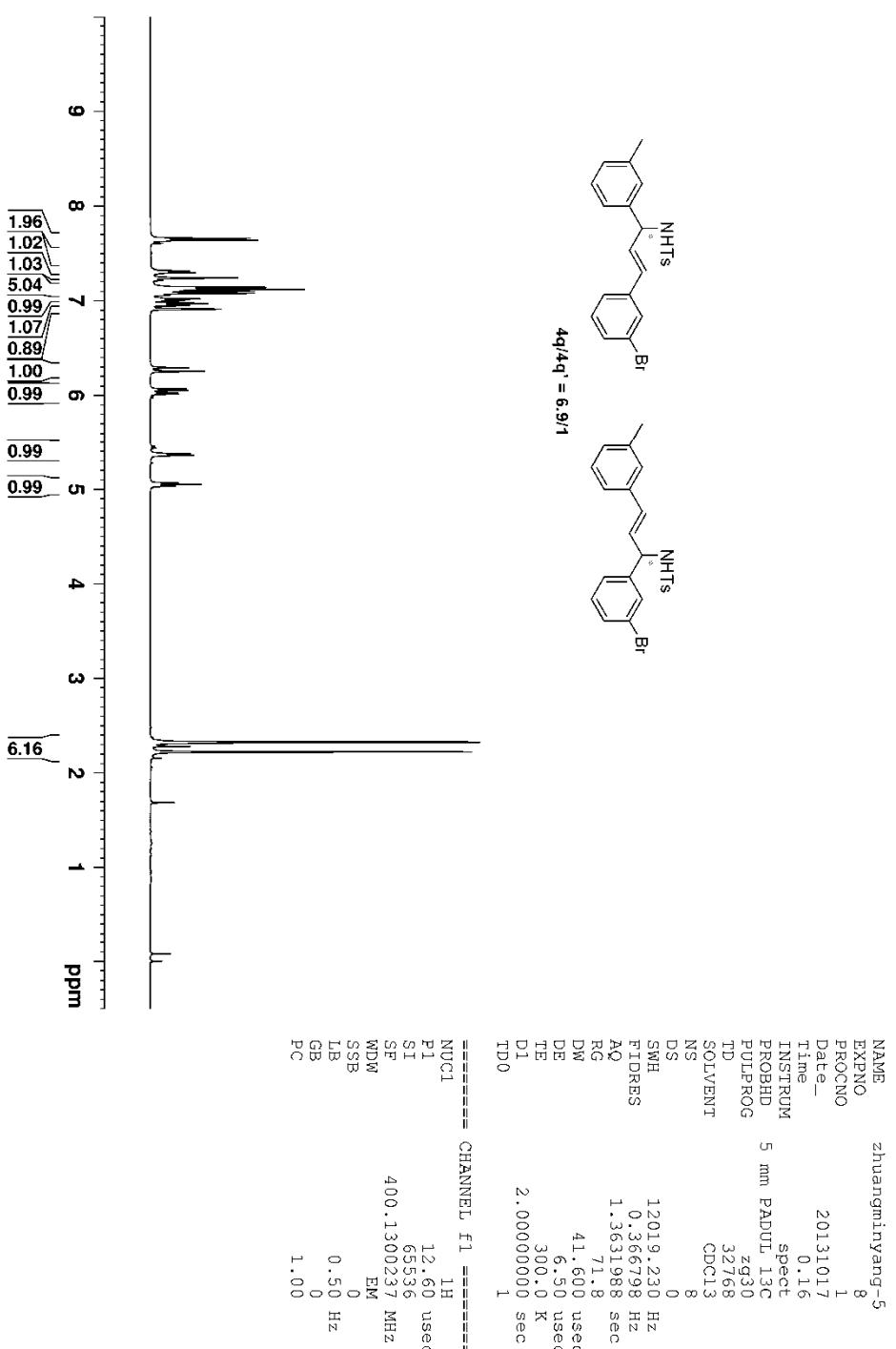


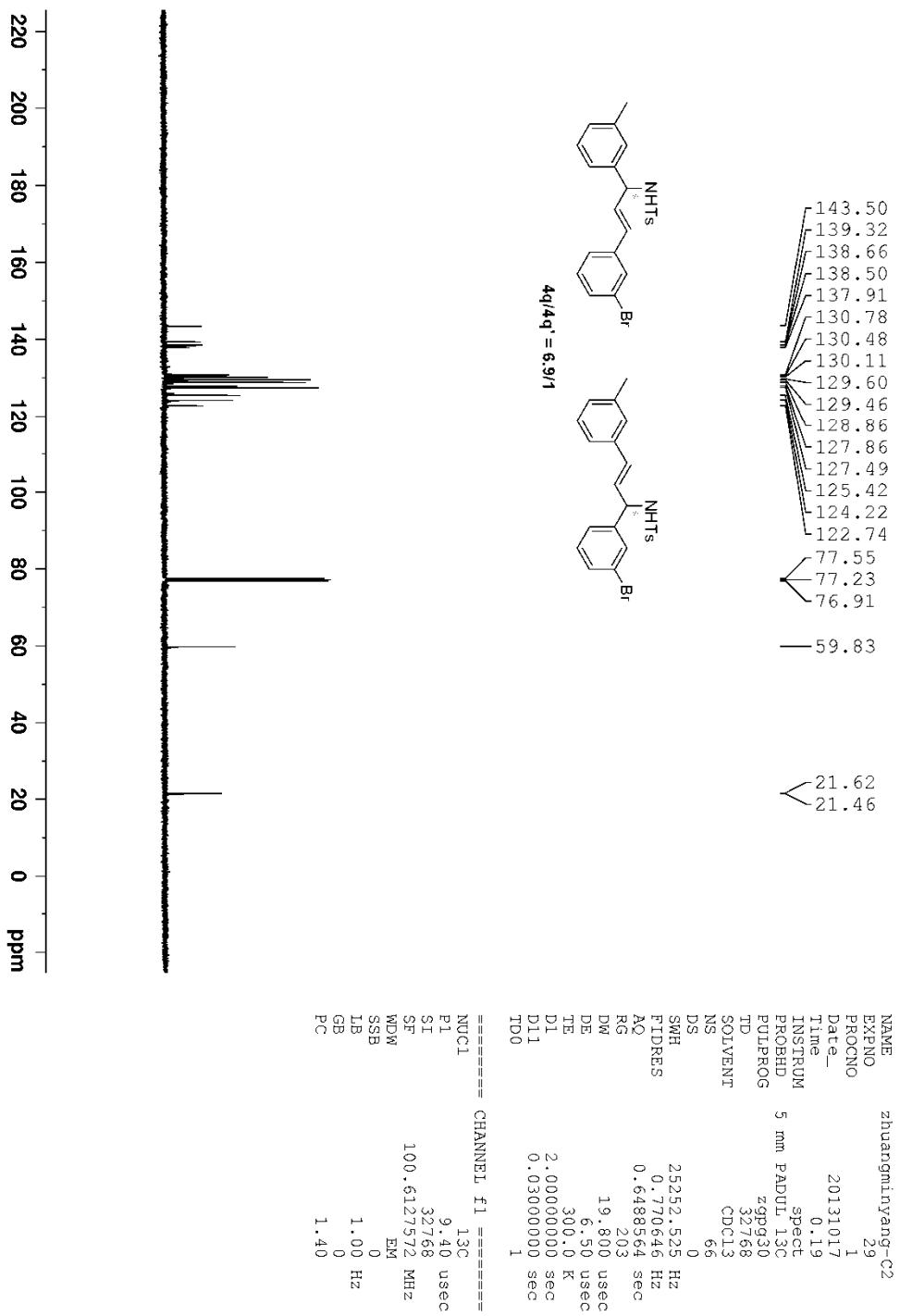




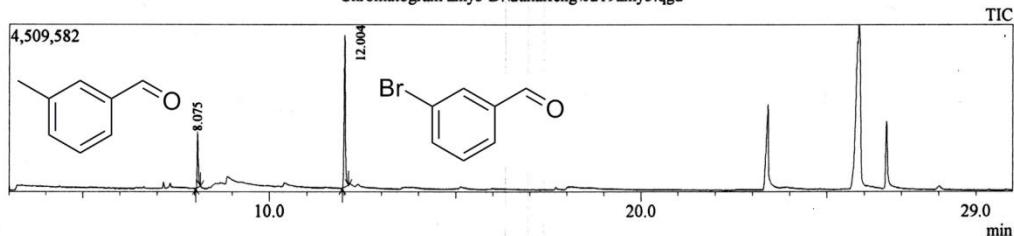
Peak#	R.Time	I.Time	F.Time	Peak Report TIC		Height%	A/H	Mark	Name
				Area	Area%				
1	5.379	5.340	5.430	3392367	35.46	1972078	43.47	1.72	
2	9.578	9.510	9.620	6173146	64.54	2564186	56.53	2.41	
				9565513	100.00	4536264			



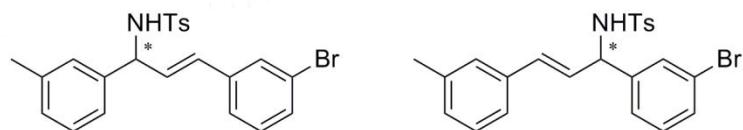




Chromatogram zmy3 D:\duhaifeng\0219zmy3.qgd



Peak#	R.Time	I.Time	F.Time	Peak Report TIC				A/H	Mark	Name
				Area	Area%	Height	Height%			
1	8.075	8.030	8.155	3370751	24.81	1504126	26.72	2.24		
2	12.004	11.950	12.110	10216488	75.19	4124529	73.28	2.48		
				13587239	100.00	5628655	100.00			

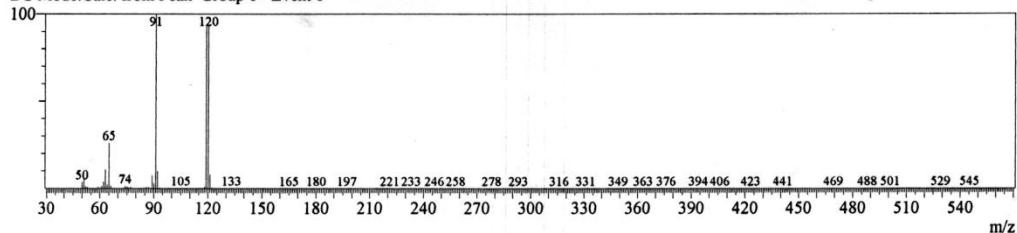


Line#:1 R.Time:8.075(Scan#:1016)

MassPeaks:300

RawMode:Averaged 8.070-8.080(1015-1017) BasePeak:91(371002)

BG Mode:Calc. from Peak Group 1 - Event 1

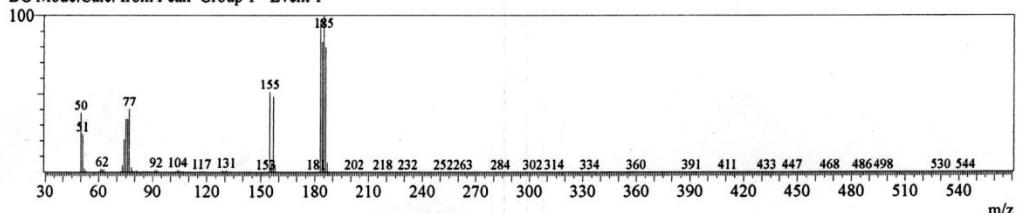


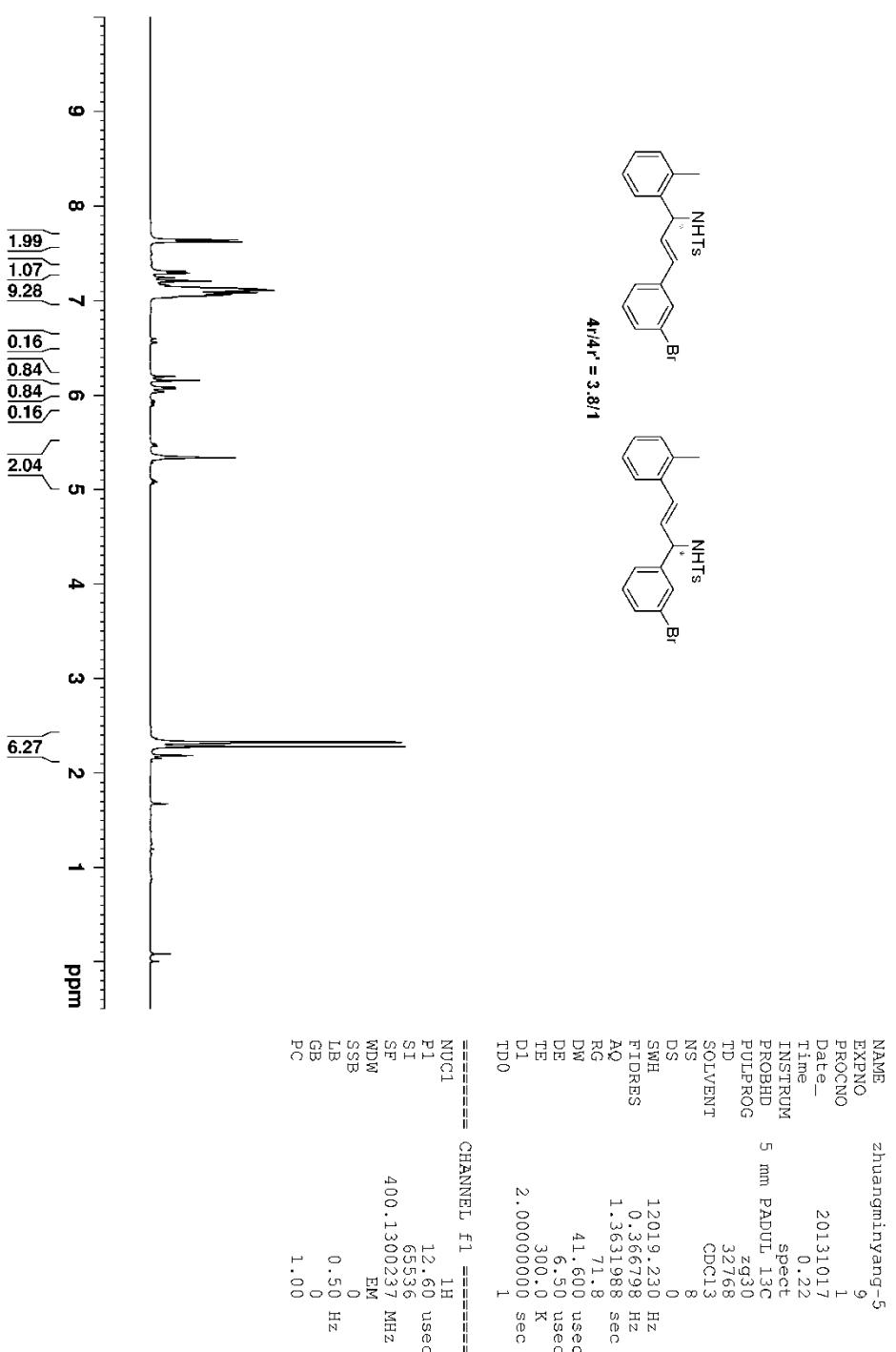
Line#:2 R.Time:12.005(Scan#:1802)

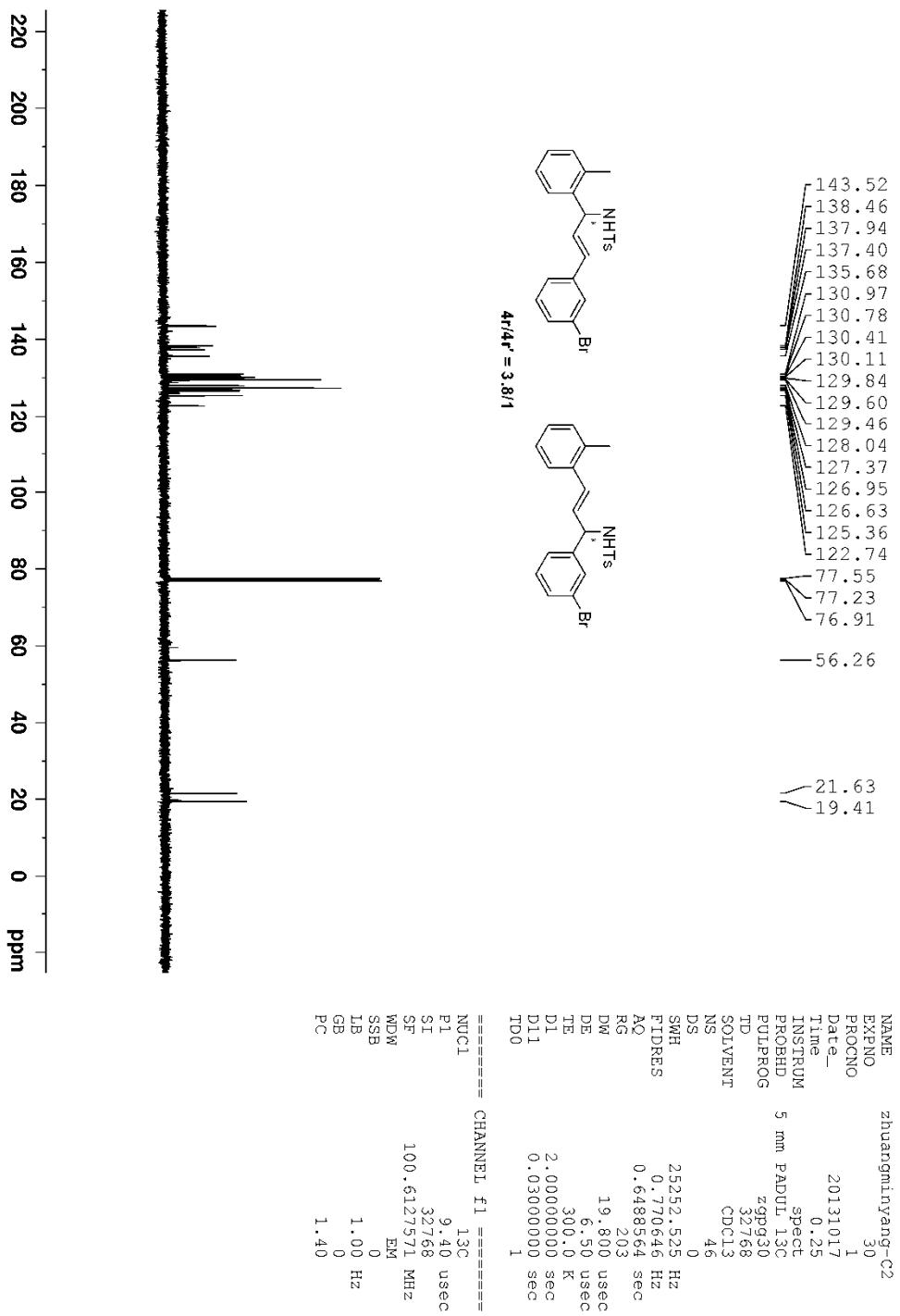
MassPeaks:320

RawMode:Averaged 12.000-12.010(1801-1803) BasePeak:185(567179)

BG Mode:Calc. from Peak Group 1 - Event 1

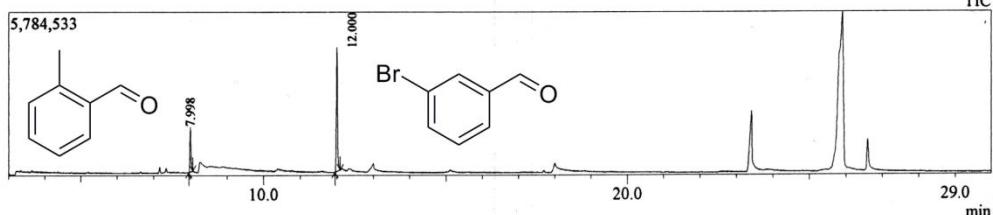




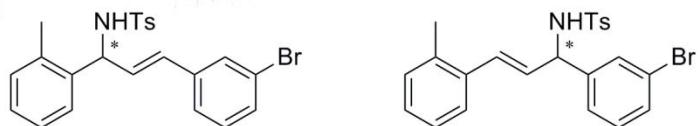


Chromatogram zmy4 D:\duhaifeng\0219zmy4.qgd

TIC



Peak#	R.Time	I.Time	F.Time	Area	Area%	Height	Height%	A/H	Mark	Name
1	7.998	7.955	8.065	3254762	24.79	1582725	26.48	2.06		
2	12.000	11.945	12.085	9876032	75.21	4393446	73.52	2.25		
				13130794	100.00	5976171	100.00			



$$4r/4r' = 3.8/1$$

Line#1 R.Time:8.000(Scan#:1001)

MassPeaks:261

RawMode:Averaged 7.995-8.005(1000-1002) BasePeak:91(395665)

BG Mode:Calc. from Peak Group 1 - Event 1

