

Supporting Information

Regioselective Hydrations of 1-Aryl-3-en-1-yne s with Gold and Platinum Catalysts: Selective Productions of 2-En-1-ones and 3-En-1-ones

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Content

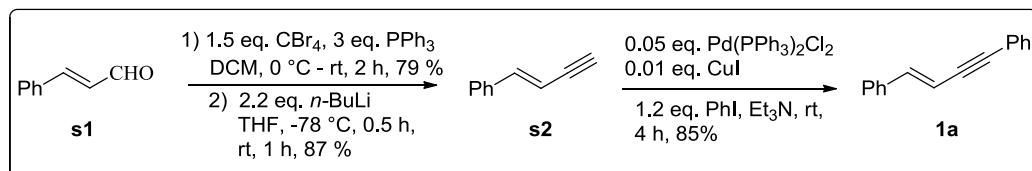
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(1) Experimental Procedures

Unless otherwise noted, all the reactions for the preparation of the substrates were performed in oven-dried glassware under nitrogen atmosphere with freshly distilled solvents. The catalytic reactions were performed in wet solvent. DCM was distilled from CaH_2 under nitrogen. THF was distilled from Na metal under nitrogen. All other commercial reagents were used without further purification, unless otherwise indicated.

^1H NMR and ^{13}C NMR spectra were recorded on a Bruker 400 MHz, Varian 500 MHz and 600 MHz Spectrometers using chloroform-d as the internal standards.

(1.1) Synthesis of (E)-but-1-en-3-yne-1,4-diyldibenzene (**1a**)



Synthesis of (*E*)-but-1-en-3-yn-1-ylbenzene (**s2**)

To a stirred solution of carbon tetrabromide (20.02 g, 60.30 mmol) in dichloromethane (200 mL) was added triphenylphosphine (23.75 g, 90.60 mmol) in DCM (100 mL) at 0 °C. The reaction mixture was stirred at 0 °C for 10 min, before a solution of cinamaldehyde (4.00 g, 30.20 mmol) in anhydrous DCM (10 mL) was added. The resulting mixture was stirred for 1 h at 0 °C before an addition of H₂O (100 mL) to partition the organic layer. The resulting mixture was extracted with DCM (3 x 50 mL); the combined organic layer was washed with brine, dried over Na₂SO₄, and concentrated under reduced pressure, to this residue was added 100 mL of pentane and the resulting suspension is filtered to remove triphenylphosphine oxide. The filtrate was concentrated in vacuo and chromatographed through a silica gel column (hexane/ether, 10:1) to afford (*E*)-(4,4-dibromobuta-1,3-dien-1-yl)benzene (8.10 g, 93.6 %) as a yellow solid.

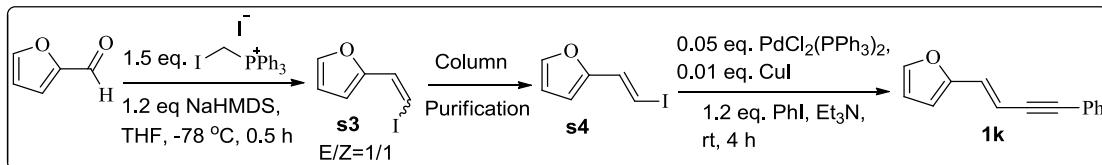
To a stirred solution of (*E*)-(4,4-dibromobuta-1,3-dien-1-yl)benzene (7.2 g, 25.08 mmol) in THF (70 mL) was added *n*BuLi (22.00 mL, 2.5 M in hexane, 50.00 mmol) dropwise at -78 °C for 30 min. The resulting solution was stirred for 30 min at -78 °C, before it was quenched with aqueous sat. NH₄Cl (5 mL) at -78 °C. The aqueous layer was extracted with (3 x 75 mL) of ether. The organic layer is washed with brine (50 mL), dried over Na₂SO₄, and concentrated under reduced pressure. The residue was eluted through a silica column to afford compound **s2** (2.80 g, 87 %) as a colorless liquid.

Synthesis of (**1a**)

To a triethylamine solution (10 mL) was added CuI (14.9 mg, 0.0780 mmol) and degassed reaction mixture using N₂ for about 20 minute before the addition of

Pd(PPh₃)₂Cl₂ (273.8 mg, 0.390 mmol) and Iodobenzene (1.9 gm 9.36 mmol). The mixture was stirred for 10 minute and was added compound **s2** (1.0 gm, 7.80 mmol) dropwise over 10 min. The resulting suspension was stirred for 4 h at room temperature. After completion of the reaction, the mixture was filtered through a short celite bed, and concentrated under reduced pressure. The residue was eluted through a silica column (hexane/ethyl acetate = 10:1) to afford compound **1a**, (1.20 gm, 80%) as a yellow solid. Follow the similar experimental procedure for the preparation of **1b-1j**, **1n**, **1q** and **1r**.

(1.2) Synthesis of (E)-2-(4-phenylbut-1-en-3-yn-1-yl)furan (**1k**)

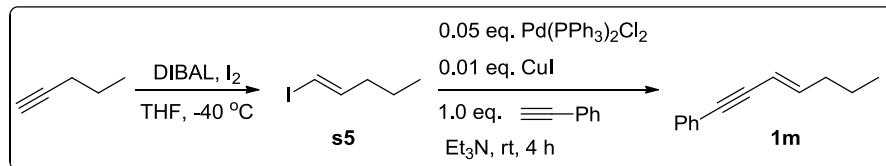


To a flame dried 250 mL round bottom flask equipped with a stir bar was added iodomethylenetriphenylphosphonium iodide (41.37 g, 78.04 mmol). The flask was then placed under vacuum for five minutes then purged with nitrogen. 50 mL of THF was added and the suspension was cooled to -20 °C and then solution of NaHMDS (62.4 mL, 0.91 M in hexane, 62.44 mmol) was added dropwise, and the solution was allowed to stir at -20 °C for 5 minutes. The flask was then cooled to -78 °C and then furan-2-carbaldehyde (5.0 g, 52.03 mmol) in 15 mL of THF was added dropwise. The reaction was allowed to stir for 10 minutes and then quenched with saturated aqueous ammonium chloride. Diethyl ether was added to the mixture and the layers were separated with a separatory funnel. The aqueous layer was washed twice with diethyl ether and the combined organic layers were dried with magnesium sulphate and concentrated to get E/Z (1:1) mixture of product **s3** (88%), which were purified using

silica column to afforded compound **s4** (42%) as yellow oil. For synthesis of **1k** from **s4** follow the similar experimental procedure as the synthesis of **1a**.

Follow the similar experimental procedure for the preparation of **1l** and **1o**.

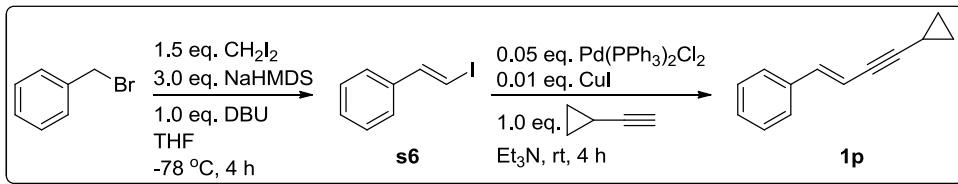
(1.3) Synthesis of (**1m**)



A 250 mL flame dried round bottom flask equipped with a stir bar was flame dried under vacuum and allowed to cool to room temperature. The flask was then purged with nitrogen and (5.0 g, 73.40 mmol) of 1-pentyne was charged into the flask. The flask was then cooled to a -40 °C bath and DIBAL (12.52 g, 88.07 mmol) was added dropwise. The reaction is then heated to 50 °C for 3 hours and then cooled to -50 °C and added iodine (22.33 g, 88.07 mmol) in 20 mL THF. The mixture is then warmed to room temperature and then cooled back down to 0 °C in an ice bath. 20% aqueous sulphuric acid was then added dropwise until there was no visible exotherm. The aqueous layer was extracted twice with pentane and the combined organic layers were washed twice with aqueous saturated sodium thiosulfate, dried with magnesium sulphate and concentrated. 12 g (83%) of crude material was isolated and used without any further purification.

For synthesis of **1m** from **s5** follow the similar experimental procedure as the synthesis of **1a**.

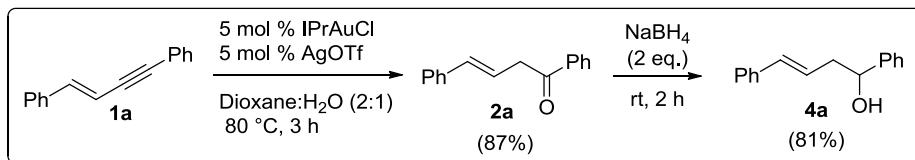
(1.4) Synthesis of (**1p**)



A solution of CH_2I_2 (1.41 mL, 17.5 mmol) in THF (10 mL) was added dropwise to a solution of NaHMDS (6.43 g, 35.1 mmol) in THF (20 mL) and ether (10 mL) at -78°C in the dark. After 20 minute, a solution of the benzyl bromide (2 g, 11.7 mmol) in THF (10 mL) was added dropwise. The reaction mixture was stirred for 90 minute then removed from the cold bath to warm to room temperature. After 30 minute, DBU (1.75 mL, 11.7 mmol) was added dropwise and the solution stirred for 1 h before ether (75 mL) was added. The mixture was filtered through a plug of celite and the solvent removed under reduced pressure. The residue was purified by flash chromatography to provide the pure vinyl iodide **s6** (80%).

For synthesis of **1p** from **s6** follow the similar experimental procedure as the synthesis of **1a**.

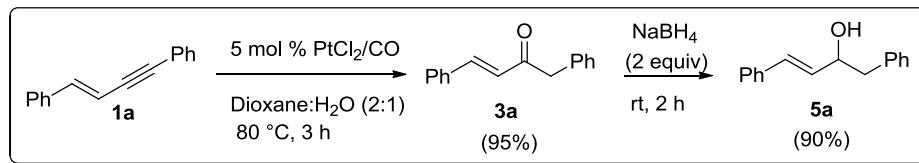
(1.5) Standard procedure for gold catalyzed one pot (E)-1,4-diphenylbut-3-en-1-ol.



A 20 mL sample vial was charged with Chloro[1,3-bis(2,6-diisopropylphenyl)imidazole-2-ylidene]gold(I) (0.0152 g, 0.0244 mmol) and Silver trifluoromethanesulfonate (0.0063 g, 0.0244 mmol) and to this mixture was added compound **1a** (0.1 g, 0.489 mmol) dissolved in 1,4-dioxane and Water (2:1

ratio). The resulting mixture was heated at 80 °C for 3 h and after the complete consumption of starting material reaction mixture was cooled to room temperature, NaBH₄ (2.0 eq.) was added slowly and reaction mixture is allowed to stir at room temperature for additional 2 h. and after complete consumption of intermediate **2a** the reaction mixture were filtered from small silica bed and dried over MgSO₄ and then purified by column chromatography to get pure compound **4a** as a off white solid (81 %).

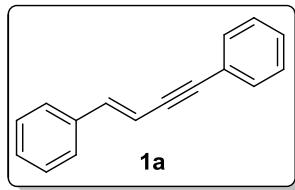
(1.6) Standard procedure for Platinum catalyzed one pot (E)-1,4-diphenylbut-3-en-2-ol.



A 20 mL catalytic sample vial was charged with platinum dichloride (0.0652 g, 0.0244 mmol) and CO balloon, before degas by using vacuum pump for about 20 minute. Compound **1a** (0.1 g, 0.489 mmol) dissolved in 1,4-dioxane (1.5 mL) and Water (5 eq) was added in to this vial. The resulting mixture was heated at 80 °C for 3 h and after the complete consumption of starting material reaction mixture was cooled to room temperature, NaBH₄ (2.0 eq.) was added slowly and reaction mixture is allowed to stir at room temperature for additional 2 h. and after complete consumption of intermediate **3a** the reaction mixture were filtered from small silica bed and dried over MgSO₄ and then purified by column chromatography to get pure compound **5a** as a yellow solid (90 %)

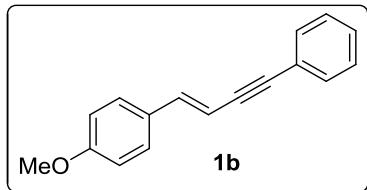
(2) Spectral data

Spectral data for (E)-but-1-en-3-yne-1,4-diylbenzene (1a)



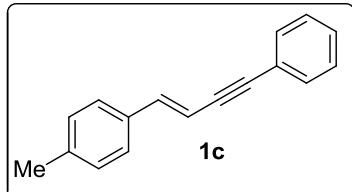
Yellow Solid; ^1H NMR (400 MHz, CDCl_3): δ 7.47 ~ 7.45 (m, 2 H), 7.44 ~ 7.40 (m, 2 H), 7.35 ~ 7.27 (m, 6 H), 7.03 (d, J = 16.2 Hz, 1 H), 6.37 (d, J = 16.2 Hz, 1 H); ^{13}C NMR (100 MHz, CDCl_3): 141.2, 136.3, 131.5, 128.7, 128.6, 128.3, 128.1, 126.3, 123.4, 108.1, 91.7, 88.9; HRMS; calcd. for $\text{C}_{16}\text{H}_{12}$: 204.0939; found: 204.0941.

Spectral data for (E)-1-methoxy-4-(4-phenylbut-1-en-3-yn-1-yl)benzene (1b)



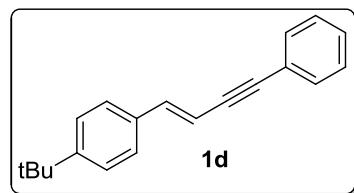
Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.46 ~ 7.45 (m, 2 H), 7.36 ~ 7.34 (m, 2 H), 7.31 ~ 7.29 (m, 3 H), 6.99 (d, J = 16.2 Hz, 1 H), 6.87 (d, J = 9 Hz, 2 H), 6.23 (d, J = 16.2 Hz, 1 H), 3.80 (s, 3 H); ^{13}C NMR (150 MHz, CDCl_3): δ 160.1, 140.8, 131.4, 129.2, 128.3, 128.0, 127.6, 123.6, 114.2, 105.7, 91.0, 89.2, 55.3; HRMS calcd. for $\text{C}_{17}\text{H}_{14}\text{O}$: 234.1045; found: 234.1048.

Spectral data for (E)-1-methyl-4-(4-phenylbut-1-en-3-yn-1-yl)benzene (1c)



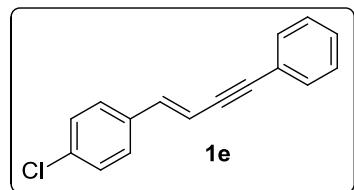
Off White Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.50 ~ 7.49 (m, 2 H), 7.35 ~ 7.31 (m, 5 H), 7.16 (d, J = 8.0 Hz, 2 H), 7.04 (d, J = 16.2 Hz, 1 H), 6.35 (d, J = 16.2 Hz, 1 H), 2.37 (s, 3 H); ^{13}C NMR (150 MHz, CDCl_3): δ 141.2, 138.7, 133.6, 131.4, 129.4, 128.3, 128.0, 126.2, 123.5, 107.0, 91.4, 89.1, 21.3; HRMS calcd. for $\text{C}_{17}\text{H}_{14}$: 218.1096; found: 218.1094.

Spectral data for (E)-1-(tert-butyl)-4-(4-phenylbut-1-en-3-yn-1-yl)benzene (1d)



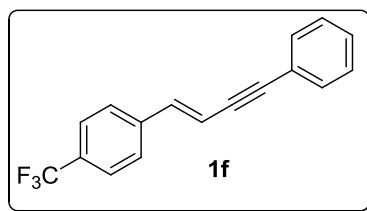
Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.48 ~ 7.46 (m, 2 H), 7.37 (s, 5 H), 7.34 ~ 7.30 (m, 2 H), 7.03 (d, J = 16.2 Hz, 1 H), 6.35 (d, J = 16.2 Hz, 1 H), 1.32 (s, 9 H); ^{13}C NMR (150 MHz, CDCl_3): δ 151.9, 141.1, 133.6, 131.5, 128.3, 128.1, 126.1, 125.7, 123.5, 107.2, 91.4, 89.1, 34.7, 31.2; HRMS: calcd. for $\text{C}_{20}\text{H}_{20}$: 260.1565; Found: 260.1563.

Spectral data for (E)-1-chloro-4-(4-phenylbut-1-en-3-yn-1-yl)benzene (1e)



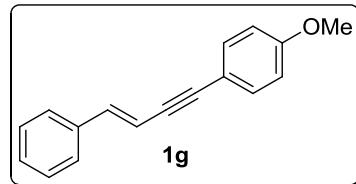
Off White Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.47 ~ 7.46 (m, 2 H), 7.34 ~ 7.29 (m, 7 H), 6.97 (d, J = 16.2 Hz, 1 H), 6.35 (d, J = 16.2 Hz, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ 139.8, 134.8, 134.3, 131.5, 128.9, 128.4, 128.3, 127.4, 123.2, 108.8, 92.3, 88.6; HRMS: calcd. for $\text{C}_{16}\text{H}_{11}\text{Cl}$: 238.0549; found: 238.0547.

Spectral data for (E)-1-(4-phenylbut-1-en-3-yn-1-yl)-4-(trifluoromethyl)benzene (1f)



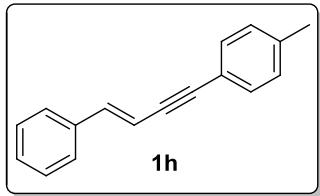
Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.58 (d, $J = 8.2$ Hz, 2 H), 7.50 ~ 7.47 (m, 4 H), 7.34 ~ 7.32 (m, 3 H), 7.03 (d, $J = 16.3$ Hz, 1 H), 6.46 (d, $J = 16.3$ Hz, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ 139.7, 139.5, 131.6, 130.2 ($J_{CF} = 33$ Hz), 128.5, 128.4, 126.4, 125.7, 125.7, 124.0 ($J_{CF} = 271.5$ Hz), 123.0, 110.9, 93.2, 88.3; HRMS: calcd. for $\text{C}_{17}\text{H}_{11}\text{F}_3$: 272.0813; found: 272.0812.

Spectral data for (E)-1-methoxy-4-(4-phenylbut-3-en-1-yn-1-yl)benzene (1g)



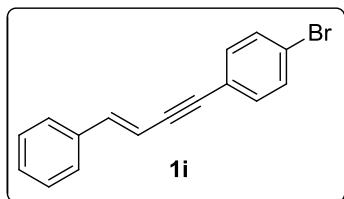
Pale Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.41 ~ 7.39 (m, 4 H), 7.34 ~ 7.31 (m, 2 H), 7.28 ~ 7.26 (m, 1 H), 6.99 (d, $J = 16.3$ Hz, 1 H), 6.85 (d, $J = 8.9$ Hz, 2 H), 6.37 (d, $J = 16.3$ Hz, 1 H), 3.81 (s, 3 H); ^{13}C NMR (150 MHz, CDCl_3): δ 159.6, 140.4, 136.5, 133.0, 128.7, 128.4, 126.2, 115.5, 114.0, 108.4, 91.8, 87.6, 55.3; HRMS: calcd. for $\text{C}_{17}\text{H}_{14}\text{O}$: 234.1045; Found: 234.1048.

Spectral data for (E)-1-methyl-4-(4-phenylbut-3-en-1-yn-1-yl)benzene (1h)



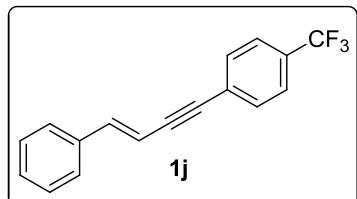
Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.41 (d, $J = 7.2$ Hz, 2 H), 7.36 (d, $J = 8.4$ Hz, 2 H), 7.33 (t, $J = 7.2$ Hz, 2 H), 7.28 ~ 7.27 (m, 1 H), 7.13 (d, $J = 7.8$ Hz, 2 H), 7.02 (d, $J = 16.2$ Hz, 1 H), 6.38 (d, $J = 16.2$ Hz, 1 H), 2.35(s, 3 H); ^{13}C NMR (150 MHz, CDCl_3): δ 140.9, 138.3, 136.4, 131.4, 129.1, 128.7, 128.5, 126.3, 120.3, 108.3, 92.0, 88.3, 21.5; HRMS: calcd. for $\text{C}_{17}\text{H}_{14}$: 218.1096; Found: 218.1092.

Spectral data for (E)-1-bromo-4-(4-phenylbut-3-en-1-yn-1-yl)benzene (1i)



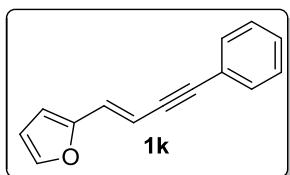
White Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.46 ~ 7.44 (m, 2 H), 7.42 ~ 7.40 (m, 2 H), 7.35 ~ 7.27 (m, 5 H), 7.04 (d, $J = 16.3$ Hz, 1 H), 6.34 (d, $J = 16.3$ Hz, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ 141.7, 136.2, 132.9, 131.6, 128.8, 126.3, 122.4, 122.4, 107.8, 90.6, 90.0; HRMS: calcd. for $\text{C}_{16}\text{H}_{11}\text{Br}$: 282.0044; Found: 282.0027.

Spectral data for (E)-1-(4-phenylbut-3-en-1-yn-1-yl)-4-(trifluoromethyl)benzene (1j)



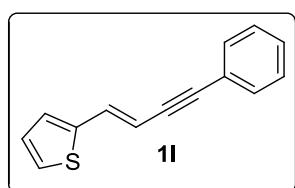
Off White Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.56 (q, $J = 8.6$ Hz, 4 H), 7.43 (d, $J = 8.8$ Hz, 2 H), 7.35 (t, $J = 7.4$ Hz, 2 H), 7.32 ~ 7.29 (m, 1 H), 7.08 (d, $J = 16.3$ Hz, 1 H), 6.37 (d, $J = 16.3$ Hz, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ 142.5, 136.0, 131.7, 129.8 ($J_{CF} = 33$ Hz), 129.0, 128.8, 127.2, 126.4, 125.3, 125.2, 123.9 ($J_{CF} = 271.5$ Hz), 107.4, 91.3, 90.2; HRMS: calcd. for $\text{C}_{17}\text{H}_{11}\text{F}_3$: 272.0813; Found: 272.0818.

Spectral data for (E)-2-(4-phenylbut-1-en-3-yn-1-yl)furan (1k)



Brown Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.46 ~ 7.45 (m, 2 H), 7.39 (d, $J = 1.7$ Hz, 1 H), 7.33 ~ 7.30 (m, 3 H), 6.79 (d, $J = 16.0$ Hz, 1 H), 6.41 (dd, $J = 1.8, 1.8$ Hz, 1 H), 6.35 (d, $J = 3.3$ Hz, 1 H), 6.29 (d, $J = 16.0$ Hz, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ 152.3, 143.0, 131.4, 128.3, 128.3, 128.1, 123.4, 111.8, 110.0, 106.2, 92.6, 88.9; HRMS: calcd. for $\text{C}_{14}\text{H}_{10}\text{O}$: 194.0732; Found: 194.0731.

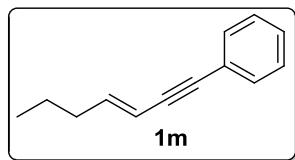
Spectral data for (E)-2-(4-phenylbut-1-en-3-yn-1-yl)thiophene (1l)



Light Brown Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.48 ~ 7.46 (m, 2 H), 7.34 ~ 7.31 (m, 3 H), 7.21 (d, $J = 5.0$ Hz, 1 H), 7.14 (d, $J = 15.9$ Hz, 1 H), 7.05 (d, $J = 3.5$ Hz, 1 H), 7.00 (dd, $J = 3.7, 3.6$ Hz, 1 H), 6.21 (d, $J = 15.9$ Hz, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ

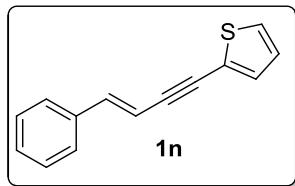
141.4, 134.0, 131.4, 128.3, 128.1, 127.7, 127.0, 125.5, 123.3, 107.3, 92.1, 88.7; HRMS:
calcd. for C₁₄H₁₀S: 210.0503; Found: 210.0507.

Spectral data for (E)-hept-3-en-1-yn-1-ylbenzene (1m)



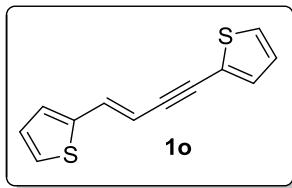
Yellow Oil; ¹H NMR (600 MHz, CDCl₃): δ 7.42 ~ 7.40 (m, 2 H), 7.30 ~ 7.26 (m, 3 H), 6.26 ~ 6.21 (m, 1 H), 5.69 (d, *J* = 16.2 Hz, 1 H), 2.15 ~ 2.11 (m, 2 H), 1.45 (q, *J* = 7.5 Hz, 2 H), 0.93 (t, *J* = 7.5 Hz, 3 H); ¹³C NMR (150 MHz, CDCl₃): δ 145.0, 131.4, 128.2, 127.8, 123.6, 109.6, 88.3, 87.8, 35.2, 22.0, 13.6; HRMS: calcd. for C₁₃H₁₄: 170.1096;
Found: 170.1091.

Spectral data for (E)-2-(4-phenylbut-3-en-1-yn-1-yl)thiophene (1n)



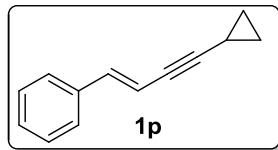
Yellow Solid; ¹H NMR (600 MHz, CDCl₃): δ 7.42 ~ 7.40 (m, 2 H), 7.35 ~ 7.32 (m, 2 H), 7.30 ~ 7.28 (m, 1 H), 7.27 ~ 7.25 (m, 1 H), 7.22 ~ 7.21 (m, 1 H), 7.02 (d, *J* = 16.2 Hz, 1 H), 6.98 (dd, *J* = 3.6, 3.6 Hz, 1 H), 6.37 (d, *J* = 16.2 Hz, 1 H); ¹³C NMR (150 MHz, CDCl₃): δ 141.2, 136.2, 131.7, 128.7, 128.7, 127.3, 127.1, 126.3, 123.5, 107.7, 92.8, 84.9; HRMS: calcd. for C₁₄H₁₀S: 210.0503; Found: 210.0502.

Spectral data for (E)-2,2'-(but-1-en-3-yne-1,4-diyl)dithiophene (1o)



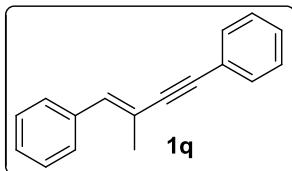
Light Brown Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.26 ~ 7.24 (m, 1 H), 7.21 ~ 7.19 (m, 2 H), 7.10 (d, $J = 15.6$ Hz, 1 H), 7.04 ~ 7.04 (d, $J = 4.2$ Hz, 1 H), 6.99 ~ 6.97 (m, 2 H), 6.18 (d, $J = 15.6$ Hz, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ 141.4, 133.9, 131.7, 127.8, 127.3, 127.2, 127.1, 125.6, 123.5, 106.9, 92.5, 85.3; HRMS: calcd. for $\text{C}_{12}\text{H}_8\text{S}_2$: 216.0067; Found: 216.0069.

Spectral data for (E)-(4-cyclopropylbut-1-en-3-yn-1-yl)benzene (1p)



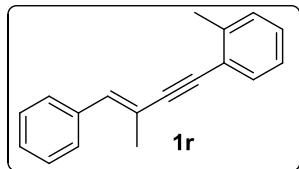
Yellow Oil; ^1H NMR (600 MHz, CDCl_3): δ 7.35 ~ 7.34 (m, 2 H), 7.31 ~ 7.29 (m, 2 H), 7.26 ~ 7.23 (m, 1 H), 6.85 (d, $J = 16.2$ Hz, 1 H), 6.12 (d, $J = 16.2$ Hz, 1 H), 1.43 ~ 1.40 (m, 1 H), 0.86 ~ 0.83 (m, 2 H), 0.80 ~ 0.75 (m, 2 H); ^{13}C NMR (150 MHz, CDCl_3): δ 140.0, 136.6, 128.6, 128.2, 126.0, 108.7, 96.7, 75.1, 8.7, 0.4; HRMS: calcd. for $\text{C}_{13}\text{H}_{12}$: 168.0939; Found: 168.0941.

Spectral data for (E)-(2-methylbut-1-en-3-yne-1,4 diyl)dibenzene (1q)



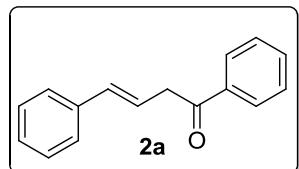
Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.49 ~ 7.48 (m, 2 H), 7.38 ~ 7.30 (m, 7 H), 7.27 ~ 7.24 (m, 1 H), 6.95 (s, 1 H), 2.17 (d, $J = 1.5$ Hz, 3 H); ^{13}C NMR (150 MHz, CDCl_3): δ 136.9, 136.1, 131.5, 129.0, 128.3, 128.3, 128.0, 127.2, 123.5, 119.9, 93.3, 88.4, 19.3; HRMS: calcd. for $\text{C}_{17}\text{H}_{14}$: 218.1096; Found: 218.1095.

Spectral data for (E)-1-methyl-2-(3-methyl-4-phenylbut-3-en-1-yn-1-yl)benzene (1r)



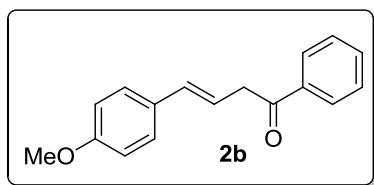
Yellow Oil; ^1H NMR (600 MHz, CDCl_3): δ 7.45 (d, $J = 7.7$ Hz, 1 H), 7.37 ~ 7.33 (m, 4 H), 7.27 ~ 7.24 (m, 1 H), 7.21 ~ 7.20 (m, 2 H), 7.16 ~ 7.14 (m, 1 H), 6.95 (s, 1 H), 2.49 (s, 3 H), 2.18 (m, 3 H); ^{13}C NMR (150 MHz, CDCl_3): δ 140.0, 136.9, 135.7, 131.7, 129.4, 129.0, 128.3, 128.1, 127.1, 125.5, 123.2, 120.1, 97.3, 87.5, 20.7, 19.4; HRMS: calcd. for $\text{C}_{18}\text{H}_{16}$: 232.1252; Found: 232.1254.

Spectral data for (E)-1,4-diphenylbut-3-en-1-one (2a)



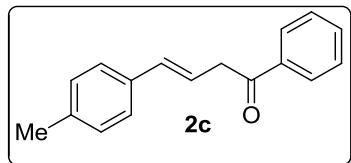
Yellow Solid; ^1H NMR (600 MHz, CDCl_3) : δ 8.0 ~ 7.98 (m, 2 H), 7.57 ~ 7.55 (m, 1 H), 7.48 ~ 7.46 (m, 2 H), 7.38 ~ 7.36 (m, 2 H), 7.30 ~ 7.27 (m, 2 H), 7.22 ~ 7.20 (m, 1 H), 6.54 (d, $J = 15.6$ Hz, 1 H), 6.49 ~ 6.44 (m, 1 H), 3.90 (dd, $J = 1.2, 1.2$ Hz, 2 H); ^{13}C NMR (150 MHz, CDCl_3): δ 198.0, 137.0, 136.6, 133.6, 133.2, 128.7, 128.5, 128.3, 127.5, 126.3, 122.6, 42.7; HRMS: calcd. for $\text{C}_{16}\text{H}_{14}\text{O}$: 222.1045; Found: 222.1040.

Spectral data for (E)-4-(4-methoxyphenyl)-1-phenylbut-3-en-1-one (2b)



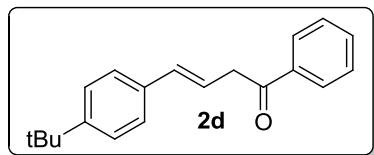
Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.99 (dd, $J = 1.2, 1.2$ Hz, 2 H), 7.57 ~ 7.54 (m, 1 H), 7.47 ~ 7.45 (m, 2 H), 7.30 (dd, $J = 2.4, 2.4$ Hz, 2 H), 6.82 (dd, $J = 1.8, 1.8$ Hz, 2 H), 6.48 (d, $J = 15.6$ Hz, 1 H), 6.33 ~ 6.28 (m, 1 H), 3.87 (dd, $J = 1.8, 1.2$ Hz, 2 H), 3.78 (s, 3 H); ^{13}C NMR (150 MHz, CDCl_3): δ 198.2, 159.1, 136.6, 133.2, 133.0, 129.9, 128.6, 128.3, 127.4, 120.3, 113.9, 55.3, 42.7; HRMS: calcd. for $\text{C}_{17}\text{H}_{16}\text{O}_2$: 252.1150; Found: 252.1145.

Spectral data for (E)-1-phenyl-4-(p-tolyl)but-3-en-1-one (2c)



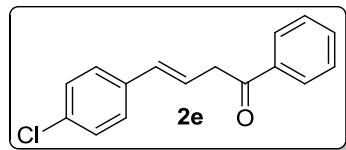
Yellow Oil; ^1H NMR (600 MHz, CDCl_3): δ 8.00 (dd, $J = 1.8, 1.2$ Hz, 2 H), 7.59 ~ 7.56 (m, 1 H), 7.49 ~ 7.47 (m, 2 H), 7.27 (t, $J = 8.1$, Hz, 2 H), 7.11 (d, $J = 7.8$ Hz, 2 H), 6.52 (d, $J = 16.2$ Hz, 1 H), 6.44 ~ 6.39 (m, 1 H), 3.90 (d, $J = 6.6$ Hz, 2 H), 2.32 (s, 3 H); ^{13}C NMR (150 MHz, CDCl_3): δ 198.1, 137.3, 136.7, 134.2, 133.4, 133.2, 129.2, 128.7, 128.3, 126.2, 121.5, 42.8, 21.2; HRMS: calcd. for $\text{C}_{17}\text{H}_{16}\text{O}$: 236.1201; Found: 236.1192.

Spectral data for (E)-4-(4-(tert-butyl)phenyl)-1-phenylbut-3-en-1-one (2d)



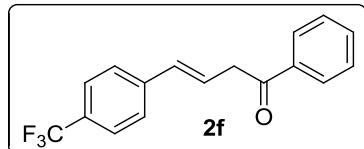
Yellow Oil; ^1H NMR (600 MHz, CDCl_3): δ 8.00 ~ 7.98 (m, 2 H), 7.57 ~ 7.54 (m, 1 H), 7.48 ~ 7.45 (m, 2 H), 7.32 (s, 4 H), 6.52 (d, J = 16.2 Hz, 1 H), 6.44 ~ 6.39 (m, 1 H), 3.89 (dd, J = 1.8, 1.2 Hz, 2 H), 1.30 (s, 9 H); ^{13}C NMR (150 MHz, CDCl_3): δ 198.1, 150.6, 136.6, 134.2, 133.3, 133.1, 128.6, 128.3, 126.0, 125.4, 121.7, 42.8, 34.5, 31.3; HRMS: calcd. for $\text{C}_{20}\text{H}_{22}\text{O}$: 278.1671; Found: 278.1669.

Spectral data for (E)-4-(4-chlorophenyl)-1-phenylbut-3-en-1-one (2e)



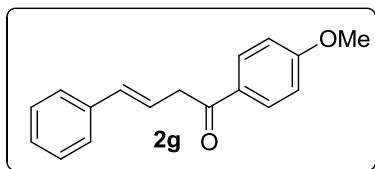
Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.99 ~ 7.97 (m, 2 H), 7.58 ~ 7.56 (m, 1 H), 7.47 (t, J = 7.8 Hz, 2 H), 7.30 ~ 7.24 (m, 5 H), 6.49 ~ 6.42 (m, 1 H), 3.89 (d, J = 5.4 Hz, 2 H); ^{13}C NMR (150 MHz, CDCl_3): δ 197.7, 136.5, 135.5, 133.3, 133.1, 132.3, 128.7, 128.6, 128.3, 127.5, 123.3, 42.5; HRMS: calcd. for $\text{C}_{16}\text{H}_{13}\text{ClO}$: 256.0655; Found: 256.0647.

Spectral data for (E)-1-phenyl-4-(4-(trifluoromethyl)phenyl)but-3-en-1-one (2f)



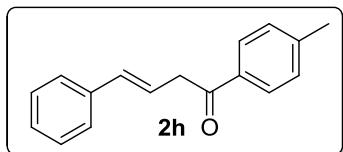
Off White Solid; ^1H NMR (600 MHz, CDCl_3): δ 8.00 ~ 7.98 (m, 2 H), 7.59 ~ 7.56 (m, 1 H), 7.54 (d, J = 8.2 Hz, 2 H), 7.49 ~ 7.45 (m, 4 H), 6.61 ~ 6.54 (m, 2 H), 3.94 (d, J = 5.8 Hz, 2 H); ^{13}C NMR (150 MHz, CDCl_3): δ 197.5, 140.4, 136.5, 133.4, 132.2, 129.3 (J_{CF} = 32.5 Hz), 128.7, 128.3, 126.4, 125.5, 124.2, (J_{CF} = 270.9 Hz), 42.5; HRMS: calcd. for $\text{C}_{17}\text{H}_{13}\text{F}_3\text{O}$: 290.0918; Found: 290.0913.

Spectral data for (E)-1-(4-methoxyphenyl)-4-phenylbut-3-en-1-one (2g)



Pale Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.97 (m, J = 2.1 Hz, 2 H), 7.37 ~ 7.35 (m, 2 H), 7.29 ~ 7.27 (m, 2 H), 7.21 ~ 7.19 (m, 1 H), 6.95 ~ 6.93 (m, 2 H), 6.52 (d, J = 16.2 Hz, 1 H), 6.48 ~ 6.43 (m, 1 H), 3.86 (s, 3 H), 3.84 (dd, J = 1.2, 1.2 Hz, 2 H); ^{13}C NMR (150 MHz, CDCl_3): δ 196.5, 163.6, 137.0, 133.3, 130.6, 129.7, 128.5, 127.4, 126.2, 123.0, 113.8, 55.5, 42.5; HRMS: calcd. for $\text{C}_{17}\text{H}_{16}\text{O}_2$: 252.1150; Found: 252.1145.

Spectral data for (E)-4-phenyl-1-(p-tolyl)but-3-en-1-one (2h)

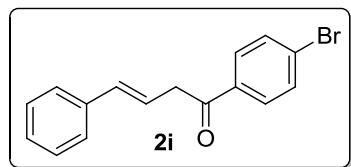


Pale Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 8.00 (d, J = 8.4 Hz, 2 H), 7.47 (d, J = 8.4 Hz, 2 H), 7.40 ~ 7.35 (m, 4 H), 7.32 ~ 7.29 (m, 1 H), 6.64 (d, J = 15.6 Hz, 1 H), 6.58 ~ 6.53 (m, 1 H), 3.97 (t, J = 3.3 Hz, 2 H), 2.51 (s, 3 H); ^{13}C NMR (150 MHz, CDCl_3): δ

197.6, 144.0, 137.1, 134.2, 133.4, 129.3, 128.5, 128.4, 127.4, 126.3, 122.8, 42.6, 21.6;

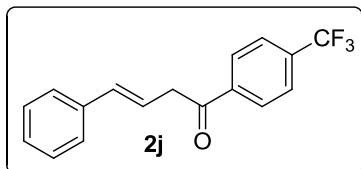
HRMS: calcd. for C₁₇H₁₆O: 236.1201; Found: 236.1204.

Spectral data for (E)-1-(4-bromophenyl)-4-phenylbut-3-en-1-one (2i)



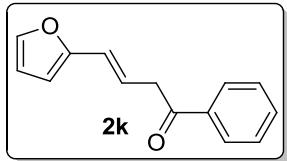
Off White Solid; ¹H NMR (600 MHz, CDCl₃): δ 7.85 (d, *J* = 8.4 Hz, 2 H), 7.60 (d, *J* = 9.0 Hz, 2 H), 7.36 (d, *J* = 7.2 Hz, 2 H), 7.29 (t, *J* = 7.7 Hz, 2 H), 7.21 (t, *J* = 7.3 Hz, 1 H), 6.54 (d, *J* = 16.2 Hz, 1 H), 6.44 ~ 6.39 (m, 1 H), 3.85 (dd, *J* = 7.2, 1.8 Hz, 2 H); ¹³C NMR (150 MHz, CDCl₃): δ 196.9, 136.9, 135.3, 133.9, 132.0, 129.8, 128.5, 128.4, 127.6, 126.3, 122.1, 42.6; HRMS calcd. for C₁₆H₁₃BrO: 300.0150; Found: 300.0155.

Spectral data for (E)-4-phenyl-1-(4-(trifluoromethyl)phenyl)but-3-en-1-one (2j)



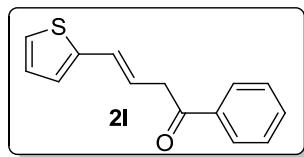
Off White Solid; ¹H NMR (600 MHz, CDCl₃): δ 8.09 (d, *J* = 8.1 Hz, 2 H), 7.74 (d, *J* = 8.2 Hz, 2 H), 7.37 (d, *J* = 8.8 Hz, 2 H), 7.30 (t, *J* = 7.7 Hz, 2 H), 7.24 ~ 7.21 (m, 1 H), 6.55 (d, *J* = 16.0 Hz, 1 H), 6.46 ~ 6.41 (m, 1 H), 3.92 (dd, *J* = 1.4, 1.4 Hz, 2 H); ¹³C NMR (150 MHz, CDCl₃): δ 196.9, 139.2, 136.7, 134.5 (*J*_{CF} = 33 Hz), 134.1, 128.7, 128.6, 127.7, 126.3, 125.8, 125.8, 123.6 (*J*_{CF} = 271.5 Hz), 121.6, 42.9; HRMS: calcd. for C₁₇H₁₃F₃O: 290.0918; Found: 290.0918.

Spectral data for (E)-4-(furan-2-yl)-1-phenylbut-3-en-1-one (2k)



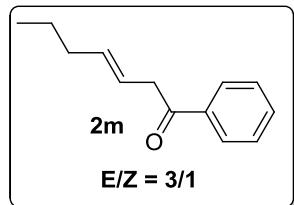
Brown Oil; ^1H NMR (600 MHz, CDCl_3): δ 7.97 (dd, $J = 1.2, 1.2$ Hz, 2 H), 7.56 ~ 7.54 (m, 1 H), 7.45 (t, $J = 7.8$ Hz, 2 H), 7.30 (d, $J = 1.4$ Hz, 1 H), 6.41 ~ 6.32 (m, 3 H), 6.19 (d, $J = 3.2$ Hz, 1 H), 3.84 (d, $J = 6.0$ Hz, 2 H); ^{13}C NMR (150 MHz, CDCl_3): δ 197.6, 152.4, 141.8, 136.5, 133.2, 128.6, 128.3, 122.0, 121.2, 111.1, 107.4, 42.4; HRMS: calcd. for $\text{C}_{14}\text{H}_{12}\text{O}_2$: 212.0837; Found: 212.0832.

Spectral data for (E)-1-phenyl-4-(thiophen-2-yl)but-3-en-1-one (2l)



Brown Oil; ^1H NMR (600 MHz, CDCl_3): δ 7.99 ~ 7.97 (m, 2 H), 7.58 ~ 7.55 (m, 1 H), 7.48 ~ 7.45 (m, 2 H), 7.12 ~ 7.11 (m, 1 H), 6.93 ~ 6.92 (m, 2 H), 6.67 ~ 6.64 (m, 1 H), 6.32 ~ 6.27 (m, 1 H), 3.86 (dd, $J = 6.6, 1.5$, Hz, 2 H); ^{13}C NMR (150 MHz, CDCl_3): δ 197.6, 142.0, 136.5, 133.2, 128.7, 128.3, 127.2, 126.6, 125.3, 124.0, 122.1, 42.4; HRMS: calcd. for $\text{C}_{14}\text{H}_{12}\text{OS}$: 228.0609; Found: 228.0611.

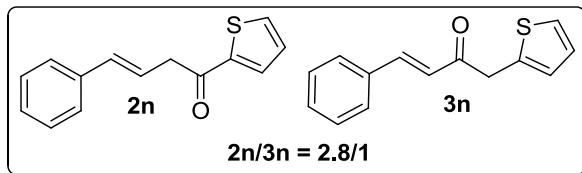
Spectral data for (E/Z)-1-phenylhept-3-en-1-one (2m)



Colorless Oil; ^1H NMR (600 MHz, CDCl_3) (*E* isomer): δ 7.96 ~ 7.94 (m, 2 H), 7.55 ~ 7.52 (m, 1 H), 7.46 ~ 7.42 (m, 2 H), 5.67 ~ 5.59 (m, 2 H), 3.68 ~ 3.66 (m, 2 H), 2.03 ~ 2.00 (m, 2 H), 1.41 ~ 1.36 (m, 2 H), 0.86 (t, $J = 7.5$ Hz, 3 H); ^1H NMR (600 MHz, CDCl_3) (*Z* isomer, major peaks): δ 3.73 ~ 3.72 (m, 2 H), 2.09 ~ 2.05 (m, 2 H), 0.91 (t, $J = 7.5$ Hz, 3 H); ^{13}C NMR (150 MHz, CDCl_3) (*E* isomer): δ 198.7, 136.7, 134.9, 133.0, 128.5, 128.3, 122.4, 42.6, 34.7, 22.3, 13.6; ^{13}C NMR (150 MHz, CDCl_3) (*Z* isomer, major peaks): δ 198.2, 133.3, 128.3, 126.9, 121.4, 37.5, 34.4, 22.5, 13.8; HRMS: calcd. for $\text{C}_{13}\text{H}_{16}\text{O}$: 188.1201; Found: 188.1198.

Spectral data for

(E)-4-phenyl-1-(thiophen-2-yl)but-3-en-1-one/(E)-4-phenyl-1-(thiophen-2-yl)but-3-en-2-one (2n/3n)

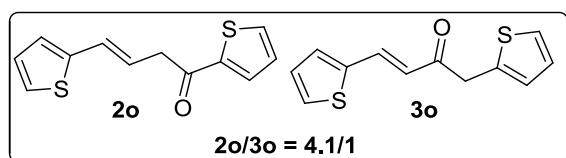


Pale Yellow Oil; ^1H NMR (600 MHz, CDCl_3) (major isomer): δ 7.78 (dd, $J = 3.8, 1.1$ Hz, 1 H), 7.64 (dd, $J = 5.0, 1.2$ Hz, 1 H), 7.38 ~ 7.36 (m, 2 H), 7.31 ~ 7.28 (m, 2 H), 7.23 ~ 7.20 (m, 1 H), 7.14 (dd, $J = 5.0, 3.8$ Hz, 1 H), 6.56 (d, $J = 15.9$ Hz, 1 H), 6.46 ~ 6.41 (m, 1 H), 3.82 (dd, $J = 7.0, 1.4$ Hz, 2 H); ^1H NMR (600 MHz, CDCl_3) (minor isomer, major peaks): δ 7.67 (s, 1 H), 7.53 ~ 7.52 (m, 2 H), 6.98 (dd, $J = 3.5, 3.4$ Hz, 1 H), 6.94 ~ 6.94 (m, 1 H), 6.81 (d, $J = 16.1$ Hz, 1 H), 4.12 (s, 2 H); ^{13}C NMR (150 MHz, CDCl_3) (major isomer): δ 190.6, 143.7, 136.8, 133.9, 133.7, 132.3, 128.5, 128.1, 127.5, 126.2, 122.2, 43.5; ^{13}C NMR (150 MHz, CDCl_3) (minor isomer, major peaks): δ 195.6, 143.8, 135.5,

134.3, 130.6, 128.9, 128.4, 127.0, 126.8, 125.1, 124.5, 42.1; HRMS: calcd. for C₁₄H₁₂OS: 228.0609; Found: 228.0608.

Spectral data for

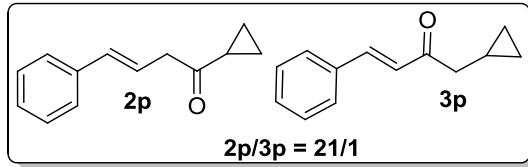
(E)-1,4-di(thiophen-2-yl)but-3-en-1-one/(E)-1,4-di(thiophen-2-yl)but-3-en-1-one (2o/3o)



Light Brown Oil; ¹H NMR (600 MHz, CDCl₃) (*major isomer*): δ 7.77 ~ 7.74 (m, 1 H), 7.64 ~ 7.63 (m, 1 H), 7.14 ~ 7.11 (m, 2 H), 6.93 ~ 6.92 (m, 2 H), 6.67 (d, *J* = 15.9 Hz, 1 H), 6.27 ~ 6.22 (m, 1 H), 3.78 ~ 3.76 (m, 2 H); ¹H NMR (600 MHz, CDCl₃) (*minor isomer, major peaks*): δ 7.38 (d, *J* = 5.1 Hz, 1 H), 7.29 ~ 7.28 (m, 1 H), 7.24 ~ 7.21 (m, 1 H), 7.05 ~ 7.04 (m, 2 H), 6.98 ~ 6.96 (m, 2 H), 6.60 (d, *J* = 15.7 Hz, 1 H), 4.07 (s, 2 H); ¹³C NMR (150 MHz, CDCl₃) (*major isomer*): δ 190.3, 143.7, 141.8, 134.0, 132.3, 128.2, 127.2, 126.9, 125.4, 124.2, 121.8, 43.3; ¹³C NMR (150 MHz, CDCl₃) (*minor isomer, major peaks*): δ 136.2, 132.0, 129.1, 128.3, 127.0, 126.8, 125.1, 123.2, 42.2; HRMS: calcd. for C₁₂H₁₀OS₂: 234.0173; Found: 234.0172.

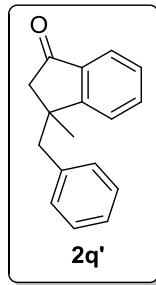
Spectral data for

(E)-1-cyclopropyl-4-phenylbut-3-en-1-one/(E)-1-cyclopropyl-4-phenylbut-3-en-2-one (2p/3p)



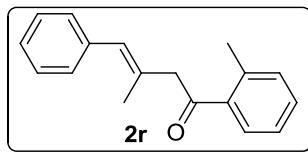
Pale Yellow Oil; ^1H NMR (600 MHz, CDCl_3) (*major isomer*): δ 7.38 ~ 7.36 (m, 2 H), 7.31 ~ 7.28 (m, 2 H), 7.23 ~ 7.20 (m, 1 H), 6.49 (d, J = 15.9 Hz, 1 H), 6.37 ~ 6.32 (m, 1 H), 3.46 ~ 3.44 (m, 2 H), 2.03 ~ 1.99 (m, 1 H), 1.07 ~ 1.04 (m, 2 H), 0.91 ~ 0.87 (m, 2 H); ^1H NMR (600 MHz, CDCl_3) (*minor isomer, major peaks*): δ 7.55 ~ 7.52 (m, 2 H), 6.79 (d, J = 16.2 Hz, 1 H), 2.55 (d, J = 6.9 Hz, 1 H), 0.60 ~ 0.57 (m, 2 H), 0.19 ~ 0.16 (m, 2 H); ^{13}C NMR (150 MHz, CDCl_3) (*major isomer*): δ 208.6, 137.0, 133.6, 128.5, 127.4, 126.2, 122.2, 47.5, 20.1, 11.1; HRMS: calcd. for $\text{C}_{13}\text{H}_{14}\text{O}$: 186.1045; Found: 186.1044.

Spectral data for 3-benzyl-3-methyl-2,3-dihydro-1H-inden-1-one (2q')



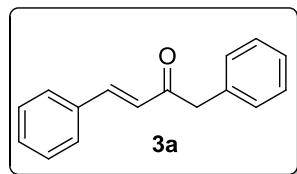
Yellow Oil; ^1H NMR (400 MHz, CDCl_3): δ 7.64 ~ 7.57 (m, 2 H), 7.48 ~ 7.42 (m, 1 H), 7.39 ~ 7.32 (m, 1 H), 7.16 ~ 7.13 (m, 3 H), 6.87 ~ 6.84 (m, 2 H), 3.00 (d, J = 13.3 Hz, 1 H), 2.90 (d, J = 13.3 Hz, 1 H), 2.81 (d, J = 18.7 Hz, 1 H), 2.37 (d, J = 18.8 Hz, 1 H), 1.48 (s, 3 H); ^{13}C NMR (100 MHz, CDCl_3): δ 205.2, 161.8, 137.3, 136.2, 134.5, 130.2, 127.9, 127.6, 126.5, 124.3, 123.2, 49.7, 48.2, 43.0, 28.1; HRMS: calcd. for $\text{C}_{17}\text{H}_{16}\text{O}$: 236.1201; Found: 236.1204.

Spectral data for (E)-3-methyl-1,4-diphenylbut-3-en-1-one (2r)



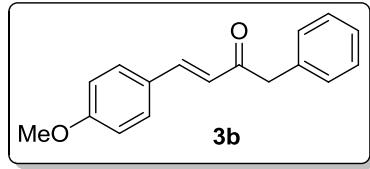
Yellow Oil; ^1H NMR (600 MHz, CDCl_3): δ 7.56 (d, $J = 7.8$ Hz, 1 H), 7.33 (t, $J = 7.5$ Hz, 1 H), 7.27 ~ 7.26 (m, 4 H), 7.24 ~ 7.22 (m, 2 H), 7.20 ~ 7.19 (m, 1 H), 6.52 (s, 1 H), 3.99 (s, 2 H), 2.49 (s, 3 H), 1.84 (d, $J = 1.2$ Hz, 3 H); ^{13}C NMR (150 MHz, CDCl_3): δ 195.6, 156.4, 140.4, 139.0, 137.3, 131.5, 130.6, 129.1, 128.4, 128.4, 126.3, 125.8, 125.6, 39.3, 24.9, 20.6; HRMS: calcd. for $\text{C}_{18}\text{H}_{18}\text{O}$: 250.1358; Found: 250.1359.

Spectral data for (E)-1,4-diphenylbut-3-en-2-one (3a)



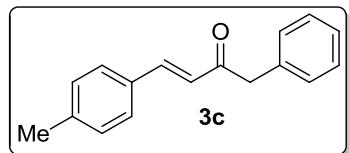
Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.63 (d, $J = 16.1$ Hz, 1 H), 7.51 ~ 7.50 (m, 2 H), 7.37 ~ 7.33 (m, 5 H), 7.28 ~ 7.26 (m, 3 H), 6.78 (d, $J = 16.1$ Hz, 1 H), 3.94 (s, 2 H); ^{13}C NMR (150 MHz, CDCl_3): δ 197.2, 143.3, 134.4, 130.5, 129.4, 128.9, 128.7, 128.3, 126.9, 125.1, 48.3; HRMS: calcd. for $\text{C}_{16}\text{H}_{14}\text{O}$: 222.1045; Found: 222.1040.

Spectral data for (E)-4-(4-methoxyphenyl)-1-phenylbut-3-en-2-one (3b)



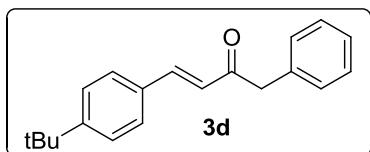
Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.59 (d, $J = 16.0$ Hz, 1 H), 7.45 (d, $J = 8.7$ Hz, 2 H), 7.34 ~ 7.31 (m, 2 H), 7.27 ~ 7.25 (m, 3 H), 6.88 (d, $J = 8.8$ Hz, 2 H), 6.66 (d, $J = 16.0$ Hz, 1 H), 3.91 (s, 2 H), 3.81 (s, 3 H); ^{13}C NMR (150 MHz, CDCl_3): δ 197.2, 161.6, 143.1, 134.6, 130.0, 129.4, 128.6, 127.0, 126.8, 122.9, 114.3, 55.3, 48.2; HRMS: calcd. for $\text{C}_{17}\text{H}_{16}\text{O}_2$: 252.1150; Found: 252.1153.

Spectral data for (E)-1-phenyl-4-(p-tolyl)but-3-en-2-one (3c)



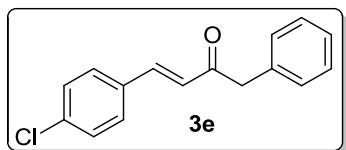
Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.60 (d, $J = 16.1$ Hz, 1 H), 7.40 (d, $J = 8.2$ Hz, 2 H), 7.34 ~ 7.32 (m, 2 H), 7.26 (d, $J = 7.2$ Hz, 3 H), 7.17 (d, $J = 8.0$ Hz, 2 H), 6.73 (d, $J = 16.1$ Hz, 1 H), 3.92 (s, 2 H), 2.35 (s, 3 H); ^{13}C NMR (150 MHz, CDCl_3): δ 197.4, 143.5, 141.1, 134.5, 131.6, 129.6, 129.4, 128.7, 128.4, 126.9, 124.3, 48.3, 21.5; HRMS: calcd. for $\text{C}_{17}\text{H}_{16}\text{O}$: 236.1201; Found: 236.1193.

Spectral data for (E)-4-(4-(tert-butyl)phenyl)-1-phenylbut-3-en-2-one (3d)



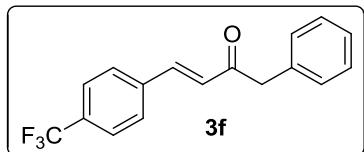
Yellow Oil; ^1H NMR (600 MHz, CDCl_3): δ 7.60 (d, $J = 16.2$ Hz, 1 H), 7.45 (m, $J = 8.4$ Hz, 2 H), 7.39 ~ 7.37 (m, 2 H), 7.35 ~ 7.31 (m, 2 H), 7.26 ~ 7.24 (m, 3 H), 6.73 (d, $J = 16.2$ Hz, 1 H), 3.92 (s, 2 H), 1.30 (s, 9 H); ^{13}C NMR (150 MHz, CDCl_3): δ 197.4, 154.2, 143.4, 134.6, 131.6, 129.5, 128.7, 128.2, 126.9, 125.9, 124.5, 48.3, 34.9, 31.1; HRMS: calcd. for $\text{C}_{20}\text{H}_{22}\text{O}$: 278.1671; Found: 278.1681.

Spectral data for (E)-4-(4-chlorophenyl)-1-phenylbut-3-en-2-one (3e)



Light Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.55 (d, $J = 16.2$ Hz, 1 H), 7.42 ~ 7.41 (m, 2 H), 7.33 ~ 7.31 (m, 4 H), 7.26 ~ 7.24 (m, 3 H), 6.73 (d, $J = 16.2$ Hz, 1 H), 3.92 (s, 2 H); ^{13}C NMR (150 MHz, CDCl_3): δ 197.0, 141.8, 136.4, 134.2, 132.9, 129.5, 129.4, 129.2, 128.8, 127.0, 125.4, 48.5; HRMS: calcd. for $\text{C}_{16}\text{H}_{13}\text{ClO}$: 256.0655; Found: 256.0651.

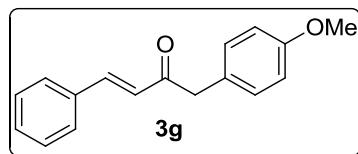
Spectral data for (E)-1-phenyl-4-(4-(trifluoromethyl)phenyl)but-3-en-2-one (3f)



Off White Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.62 ~ 7.58 (m, 5 H), 7.34 (t, $J = 7.8$ Hz, 2 H), 7.28 ~ 7.25 (m, 3 H), 6.82 (d, $J = 16.2$ Hz, 1 H), 3.94 (s, 2 H); ^{13}C NMR (150 MHz, CDCl_3): δ 196.8, 141.3, 137.8, 134.0, 131.9 ($J_{CF} = 33$), 129.5, 128.9, 128.4, 127.2, 127.1,

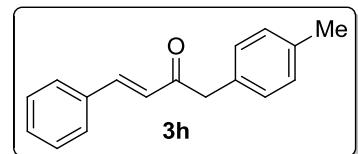
125.8, 125.8, 123.7 ($J_{CF} = 271.5$), 48.7; HRMS: calcd. for $C_{17}H_{13}F_3O$: 290.0918; Found: 290.0913.

Spectral data for (E)-1-(4-methoxyphenyl)-4-phenylbut-3-en-2-one (3g)



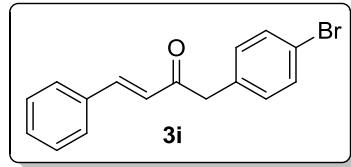
Yellow Solid; 1H NMR (600 MHz, $CDCl_3$): δ 7.61 (d, $J = 16.2$ Hz, 1 H), 7.50 (dd, $J = 6.0, 2.4$ Hz, 2 H), 7.37 ~ 7.35 (m, 3 H), 7.17 (d, $J = 8.5$ Hz, 2 H), 6.88 ~ 6.86 (m, 2 H), 6.76 (d, $J = 16.2$ Hz, 1 H), 3.86 (s, 2 H), 3.78 (s, 3 H); ^{13}C NMR (150 MHz, $CDCl_3$): δ 197.6, 158.6, 143.2, 134.4, 130.5, 128.9, 128.3, 126.4, 125.1, 114.2, 55.2, 47.5; HRMS: calcd. for $C_{17}H_{16}O_2$: 252.1150; Found: 252.1152.

Spectral data for (E)-4-phenyl-1-(p-tolyl)but-3-en-2-one (3h)



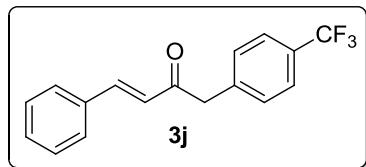
Pale Yellow Solid; 1H NMR (600 MHz, $CDCl_3$): δ 7.60 (d, $J = 16.1$ Hz, 1 H), 7.50 ~ 7.48 (m, 2 H), 7.36 ~ 7.35 (dd, $J = 5.5, 2.3$ Hz, 3 H), 7.14 (s, 4 H), 6.75 (d, $J = 16.2$ Hz, 1 H), 3.88 (s, 2 H), 2.32 (s, 3 H); ^{13}C NMR (150 MHz, $CDCl_3$): δ 197.5, 143.3, 136.6, 134.5, 131.3, 130.5, 129.5, 129.3, 128.9, 128.4, 125.2, 48.10, 21.0; HRMS: calcd. for $C_{17}H_{16}O$: 236.1201; Found: 236.1208.

Spectral data for (E)-1-(4-bromophenyl)-4-phenylbut-3-en-2-one (3i)



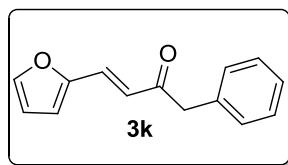
Off White Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.61 (d, $J = 16.1$ Hz, 1 H), 7.51 ~ 7.50 (m, 2 H), 7.45 (d, $J = 8.3$ Hz, 2 H), 7.38 (m, 3 H), 7.12 (d, $J = 8.2$ Hz, 2 H), 6.75 (d, $J = 16.1$ Hz, 1 H), 3.88 (s, 2 H); ^{13}C NMR (150 MHz, CDCl_3): δ 196.5, 143.7, 134.3, 133.4, 131.8, 131.2, 130.7, 128.9, 128.4, 125.0, 121.0, 47.5; HRMS: calcd. for $\text{C}_{16}\text{H}_{13}\text{BrO}$: 300.0150; Found: 300.0147.

Spectral data for (E)-4-phenyl-1-(4-(trifluoromethyl)phenyl)but-3-en-2-one (3j)



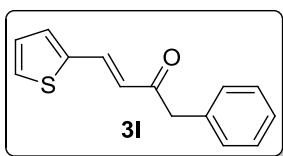
Off White Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.63 (d, $J = 16.1$ Hz, 1 H), 7.59 (d, $J = 8.3$ Hz, 2 H), 7.53 ~ 7.52 (m, 2 H), 7.40 ~ 7.36 (m, 5 H), 6.77 (d, $J = 16.1$ Hz, 1 H), 4.01 (s, 2 H); ^{13}C NMR (150 MHz, CDCl_3): δ 196.2, 143.9, 138.4, 134.2, 130.8, 129.9, 129.3 ($J_{CF} = 31.5$ Hz), 129.0, 128.4, 125.6, 125.6, 125.1, 124.1 ($J_{CF} = 271.5$ Hz), 47.7; HRMS: calcd. for $\text{C}_{17}\text{H}_{13}\text{F}_3\text{O}$: 290.0918; Found: 290.0913.

Spectral data for (E)-4-(furan-2-yl)-1-phenylbut-3-en-2-one (3k)



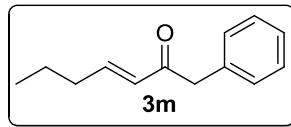
Brown Oil; ^1H NMR (600 MHz, CDCl_3): δ 7.45 ~ 7.45 (m, 1 H), 7.36 (d, $J = 15.7$ Hz, 1 H), 7.33 ~ 7.30 (m, 2 H), 7.26 ~ 7.23 (m, 3 H), 6.66 (d, $J = 15.8$ Hz, 1 H), 6.63 (d, $J = 3.4$ Hz, 1 H), 6.45 ~ 6.44 (m, 1 H), 3.86 (s, 2 H); ^{13}C NMR (150 MHz, CDCl_3): δ 196.9, 151.1, 145.0, 134.4, 129.5, 129.4, 128.7, 127.0, 122.4, 116.0, 112.5, 48.9; HRMS: calcd. for $\text{C}_{14}\text{H}_{12}\text{O}_2$: 212.0837; Found: 212.0826.

Spectral data for (E)-1-phenyl-4-(thiophen-2-yl)but-3-en-2-one (3l)



Brown Oil; ^1H NMR (600 MHz, CDCl_3): δ 7.74 (d, $J = 15.7$ Hz, 1 H), 7.36 ~ 7.31 (m, 3 H), 7.26 ~ 7.25 (m, 4 H), 7.03 (dd, $J = 3.7, 3.6$ Hz, 1 H), 6.58 (d, $J = 15.7$ Hz, 1 H), 3.88 (s, 2 H); ^{13}C NMR (150 MHz, CDCl_3): δ 196.7, 139.8, 135.7, 134.4, 131.8, 129.4, 128.9, 128.7, 128.2, 126.9, 123.9, 48.4; HRMS: calcd. for $\text{C}_{14}\text{H}_{12}\text{OS}$: 228.0609; Found: 228.0613.

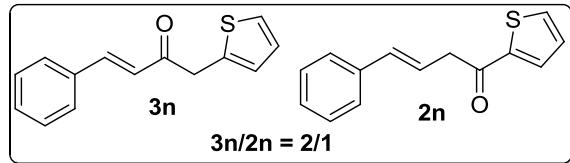
Spectral data for (E)-1-phenylhept-3-en-2-one (3m)



Colorless Oil; ^1H NMR (600 MHz, CDCl_3): δ 7.31 ~ 7.29 (m, 2 H), 7.24 ~ 7.22 (m, 1 H), 7.20 ~ 7.18 (m, 2 H), 6.92 ~ 6.87 (m, 1 H), 6.12 (d, $J = 16.2$ Hz, 1 H), 3.81 (s, 2 H), 2.17 ~ 2.13 (m, 2 H), 1.45 (q, $J = 7.5$ Hz, 2 H), 0.90 ~ 0.88 (m, 3 H); ^{13}C NMR (150 MHz,

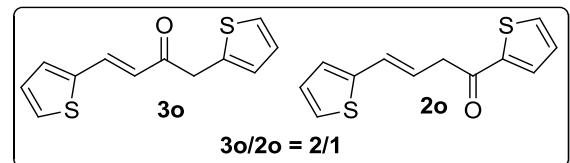
CDCl_3): δ 197.6, 148.4, 134.6, 129.5, 129.4, 128.6, 126.8, 47.6, 34.5, 21.3, 13.6; HRMS: calcd. for $\text{C}_{13}\text{H}_{16}\text{O}$: 188.1201; Found: 188.1198.

Spectral data for
(E)-4-phenyl-1-(thiophen-2-yl)but-3-en-2-one/(E)-4-phenyl-1-(thiophen-2-yl)but-3-en-1-one (3n/2n)



Yellow Oil; ^1H NMR (600 MHz, CDCl_3) (*major isomer*): δ 7.66 ~ 7.63 (m, 2 H), 7.53 ~ 7.52 (m, 2 H), 7.39 ~ 7.36 (m, 4 H), 7.30 ~ 7.28 (m, 2 H), 7.23 ~ 7.20 (m, 2 H), 6.81 (d, J = 15.6 Hz, 1 H), 4.12 (s, 2 H); ^1H NMR (600 MHz, CDCl_3) (*minor isomer, major peaks*): δ 7.78 ~ 7.77 (m, 1 H), 7.14 ~ 7.13 (m, 2 H), 6.99 ~ 6.97 (m, 2 H), 6.94 ~ 6.93 (m, 1 H), 6.56 (d, J = 16.2 Hz, 1 H), 6.45 ~ 6.40 (m, 1 H), 3.82 (q, J = 1.5 Hz, 2 H); ^{13}C NMR (150 MHz, CDCl_3) (*major isomer*): δ 195.6, 143.8, 135.5, 134.3, 130.7, 128.9, 128.4, 127.1, 126.8, 125.1, 124.5, 42.1; ^{13}C NMR (150 MHz, CDCl_3) (*minor isomer, major peaks*): δ 190.6, 136.8, 133.9, 133.8, 132.3, 128.5, 128.1, 127.5, 126.3, 122.2, 43.6; HRMS: calcd. for $\text{C}_{14}\text{H}_{12}\text{OS}$: 228.0609; Found: 228.0608.

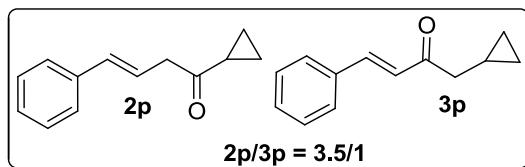
Spectral data for
(E)-1,4-di(thiophen-2-yl)but-3-en-2-one/(E)-1,4-di(thiophen-2-yl)but-3-en-1-one (3o/2o)



Brown Oil; ^1H NMR (600 MHz, CDCl_3) (*major isomer*): δ 7.75 ~ 7.74 (m, 2 H), 7.38 (d, J = 5.4 Hz, 1 H), 7.28 (d, J = 4.2 Hz, 1 H), 7.22 ~ 7.21 (m, 1 H), 7.05 ~ 7.04 (m, 1 H), 6.98 ~ 6.96 (m, 1 H), 6.60 (d, J = 15.6 Hz, 1 H), 4.07 (s, 2 H); ^1H NMR (600 MHz, CDCl_3) (*minor isomer, major peaks*): δ 7.64 ~ 7.63 (m, 1 H), 7.13 ~ 7.11 (m, 2 H), 6.93 ~ 6.92 (m, 3 H), 6.67 (d, J = 16.2 Hz, 1 H), 6.27 ~ 6.22 (m, 1 H), 3.77 (dd, J = 7.2, 1.8 Hz, 2 H); ^{13}C NMR (150 MHz, CDCl_3) (*major isomer*): δ 195.1, 139.6, 136.1, 135.5, 132.0, 129.1, 128.3, 127.0, 126.8, 125.1, 123.2, 43.3; ^{13}C NMR (150 MHz, CDCl_3) (*minor isomer, major peaks*): δ 190.3, 143.6, 141.8, 133.9, 132.3, 128.1, 127.2, 126.9, 125.4, 124.1, 121.7, 42.1; HRMS: calcd. for $\text{C}_{12}\text{H}_{10}\text{OS}_2$: 234.0173; Found: 234.0172.

Spectral data for

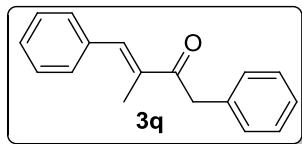
(E)-1-cyclopropyl-4-phenylbut-3-en-1-one/(E)-1-cyclopropyl-4-phenylbut-3-en-2-one (2p/3p)



Pale Yellow Oil; ^1H NMR (600 MHz, CDCl_3) (*major isomer*): δ 7.38 ~ 7.35 (m, 2 H), 7.31 ~ 7.27 (m, 2 H), 7.22 ~ 7.18 (m, 1 H), 6.48 (d, J = 15.9 Hz, 1 H), 6.36 ~ 6.31 (m, 1 H), 3.45 ~ 3.44 (m, 2 H), 2.02 ~ 1.98 (m, 1 H), 1.09 ~ 1.04 (m, 2 H), 0.90 ~ 0.86 (m, 2 H); ^1H NMR (600 MHz, CDCl_3) (*minor isomer, major peaks*): δ 7.55 ~ 7.52 (m, 2 H), 6.79 (d, J = 16.2 Hz, 1 H), 2.54 (d, J = 6.9 Hz, 1 H), 0.60 ~ 0.57 (m, 2 H), 0.19 ~ 0.16 (m, 2 H); ^{13}C NMR (150 MHz, CDCl_3) (*major isomer*): δ 208.6, 137.0, 133.6, 128.5, 127.5, 126.2, 122.2, 47.5, 20.2, 11.1; ^{13}C NMR (150 MHz, CDCl_3) (*minor isomer*): δ 142.5,

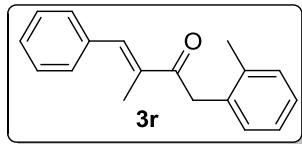
130.4, 128.9, 128.7, 128.2, 125.9, 46.1, 6.6, 4.6; HRMS: calcd. for C₁₃H₁₄O: 186.1045; Found: 186.1044.

Spectral data for (E)-3-methyl-1,4-diphenylbut-3-en-2-one (3q)



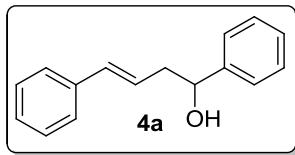
Yellow Oil; ¹H NMR (500 MHz, CDCl₃): δ 7.65 (s, 1 H), 7.41 (d, *J* = 5.0 Hz, 4 H), 7.36 ~ 7.32 (m, 3 H), 7.27 ~ 7.26 (m, 3 H), 4.14 (s, 2 H), 2.08 (d, *J* = 1.5 Hz, 3 H); ¹³C NMR (110 MHz, CDCl₃): δ 199.7, 139.8, 136.9, 135.7, 135.3, 129.7, 129.2, 128.6, 128.4, 126.7, 44.6, 13.3; HRMS: calcd. for C₁₇H₁₆O: 236.1201; Found: 236.1201.

Spectral data for (E)-3-methyl-4-phenyl-1-(o-tolyl)but-3-en-2-one (3r)



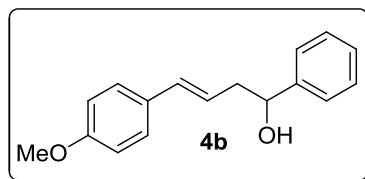
White Solid; ¹H NMR (600 MHz, CDCl₃): δ 7.66 (s, 1 H), 7.45 ~ 7.42 (m, 4 H), 7.38 ~ 7.36 (m, 1 H), 7.23 ~ 7.17 (m, 3 H), 7.15 ~ 7.13 (m, 1 H), 4.17 (s, 2 H), 2.30 (s, 3 H), 2.12 (d, *J* = 1.2 Hz, 3 H); ¹³C NMR (150 MHz, CDCl₃): δ 199.7, 139.0, 137.4, 136.7, 135.9, 134.2, 130.3, 130.1, 129.7, 128.6, 128.5, 127.0, 126.1, 42.6, 19.8, 13.4; HRMS: calcd. for C₁₈H₁₈O: 250.1358; Found: 250.1359

Spectral data for (E)-1,4-diphenylbut-3-en-1-ol (4a)



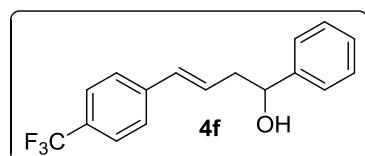
Light Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.39 ~ 7.32 (m, 6 H), 7.30 ~ 7.27 (m, 3 H), 7.20 (t, J = 7.3 Hz, 1 H), 6.49 (d, J = 15.9 Hz, 1 H), 6.22 ~ 6.17 (m, 1 H), 4.80 (t, J = 6.4 Hz, 1 H), 2.67 ~ 2.64 (m, 2 H), 2.06 (s, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ 143.9, 137.2, 133.4, 128.5, 128.5, 127.6, 127.3, 126.2, 125.9, 125.8, 73.8, 43.1; HRMS: calcd. for $\text{C}_{16}\text{H}_{14}\text{O}$: 224.1201; Found: 224.1192.

Spectral data for (E)-4-(4-methoxyphenyl)-1-phenylbut-3-en-1-ol (4b)



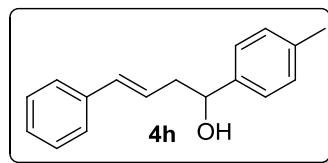
Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.39 ~ 7.34 (m, 4 H), 7.29 ~ 7.26 (m, 3 H), 6.82 (dd, J = 1.8, 2.4 Hz, 2 H), 6.44 (d, J = 15.6 Hz, 1 H), 6.07 ~ 6.02 (m, 1 H), 4.78 (q, J = 5.4 Hz, 1 H), 3.79 (s, 3 H), 2.65 ~ 2.59 (m, 2 H), 2.13 (bs, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ 159.0, 143.9, 132.9, 130.0, 128.4, 127.5, 127.3, 125.8, 123.5, 113.9, 73.7, 55.3, 43.1; HRMS: calcd. for $\text{C}_{17}\text{H}_{18}\text{O}_2$: 254.1307; Found: 254.1308.

Spectral data for (E)-1-phenyl-4-(4-(trifluoromethyl)phenyl)but-3-en-1-ol (4f)



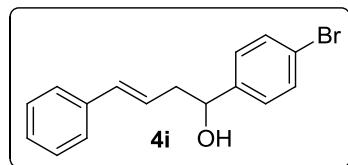
White Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.53 (d, $J = 7.8$ Hz, 2 H), 7.40 (d, $J = 8.4$ Hz, 2 H), 7.37 ~ 7.28 (m, 5 H), 6.49 (d, $J = 15.6$ Hz, 1 H), 6.33 ~ 6.28 (m, 1 H), 4.82 (t, $J = 6.3$ Hz, 1 H), 2.69 ~ 2.67 (m, 2 H), 2.17 (bs, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ 143.8, 140.7, 131.8, 129.5, 129.1, 128.5, 127.8, 126.2, 125.8, 125.5, 125.4, 124.2 ($J_{CF} = 271.5$ Hz), 73.8, 42.8; HRMS: calcd. for $\text{C}_{17}\text{H}_{15}\text{F}_3\text{O}$: 292.1075; Found: 292.9802.

Spectral data for (E)-4-phenyl-1-(p-tolyl)but-3-en-1-ol (4h)



Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.37 ~ 7.35 (m, 2 H), 7.32 ~ 7.29 (m, 4 H), 7.23 ~ 7.22 (m, 1 H), 7.20 (d, $J = 7.2$ Hz, 1 H), 6.52 (d, $J = 16.2$ Hz, 1 H), 6.25 ~ 6.20 (m, 1 H), 4.79 (t, $J = 5.4$ Hz, 1 H), 2.69 ~ 2.67 (m, 2 H), 2.38 (s, 3 H), 2.10 (bs, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ 141.0, 137.3, 137.2, 133.2, 129.1, 128.5, 127.2, 126.1, 126.1, 125.7, 73.6, 43.0, 21.1; HRMS: calcd. for $\text{C}_{17}\text{H}_{18}\text{O}$: 238.1358; Found: 238.1915.

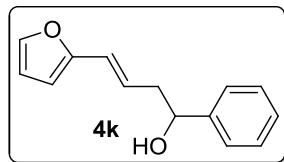
Spectral data for (E)-1-(4-bromophenyl)-4-phenylbut-3-en-1-ol (4i)



Off White Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.47 (d, $J = 8.4$ Hz, 2 H), 7.33 ~ 7.28 (m, 4 H), 7.25 ~ 7.22 (m, 3 H), 6.48 (d, $J = 15.6$ Hz, 1 H), 6.17 ~ 6.12 (m, 1 H), 4.75 (q, $J = 5.1$ Hz, 1 H), 2.64 ~ 2.58 (m, 2 H), 2.19 (bs, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ

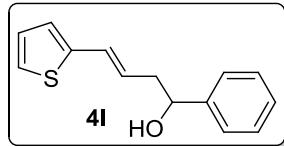
142.8, 137.0, 133.8, 131.5, 128.5, 127.5, 127.4, 126.2, 125.2, 121.3, 73.0, 43.0; HRMS: calcd. for C₁₆H₁₅BrO: 302.0306; Found: 301.9871.

Spectral data for (E)-4-(furan-2-yl)-1-phenylbut-3-en-1-ol (4k)



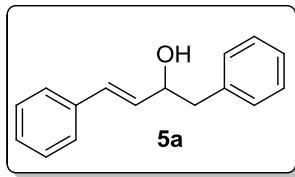
Brown Oil; ¹H NMR (600 MHz, CDCl₃): δ 7.38 ~ 7.27 (m, 6 H), 6.34 ~ 6.33 (m, 1 H), 6.30 (d, *J* = 16.2 Hz, 1 H), 6.16 ~ 6.11 (m, 2 H), 4.78 (q, *J* = 5.1 Hz, 1 H), 2.63 ~ 2.60 (m, 2 H), 2.15 (bs, 1 H); ¹³C NMR (150 MHz, CDCl₃): δ 152.7, 143.8, 141.6, 128.5, 127.6, 125.8, 124.8, 121.8, 111.1, 107.0, 73.7, 42.9; HRMS: calcd. for C₁₄H₁₄O₂: 214.0994; Found: 214.0991.

Spectral data for (E)-1-phenyl-4-(thiophen-2-yl)but-3-en-1-ol (4l)



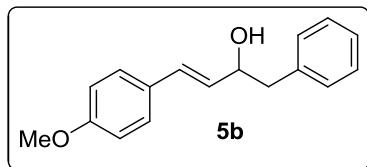
Yellow Oil; ¹H NMR (600 MHz, CDCl₃): δ 7.39 ~ 7.34 (m, 4 H), 7.30 ~ 7.27 (m, 1 H), 7.10 (d, *J* = 5.0 Hz, 1 H), 6.94 ~ 6.92 (m, 1 H), 6.89 (d, *J* = 3.3 Hz, 1 H), 6.61 (d, *J* = 15.7 Hz, 1 H), 6.06 ~ 6.01 (m, 1 H), 4.78 (q, *J* = 5.5 Hz, 1 H), 2.63 ~ 2.58 (m, 2 H), 2.15 (bs, 1 H); ¹³C NMR (150 MHz, CDCl₃): δ 143.8, 142.3, 128.5, 127.6, 127.2, 126.4, 125.8, 125.0, 123.7, 73.7, 42.9; HRMS: calcd. for C₁₄H₁₄OS: 230.0765; Found: 230.9852.

Spectral data for (E)-1,4-diphenylbut-3-en-2-ol (5a)



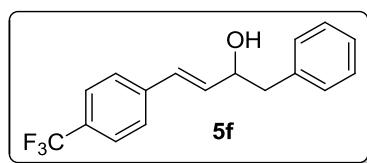
Yellow Solid; ¹H NMR (600 MHz, CDCl₃): δ 7.38 ~ 7.37 (m, 2 H), 7.34 ~ 7.31 (m, 4 H), 7.27 ~ 7.25 (m, 4 H), 6.59 (d, *J* = 15.6 Hz, 1 H), 6.29 (dd, *J* = 15.6, 6.0 Hz, 1 H), 4.54 ~ 4.51 (m, 1 H), 2.98 ~ 2.87 (m, 2 H), 1.92 (bs, 1 H); ¹³C NMR (150 MHz, CDCl₃): δ 137.6, 136.6, 131.4, 130.3, 129.5, 128.5, 128.5, 127.6, 126.6, 126.4, 73.4, 44.1; HRMS: calcd. for C₁₆H₁₆O: 224.1201; Found: 224.1202.

Spectral data for (E)-4-(4-methoxyphenyl)-1-phenylbut-3-en-2-ol (5b)



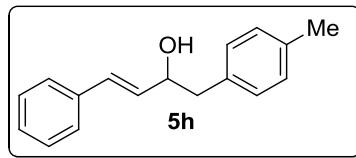
Yellow Solid; ¹H NMR (600 MHz, CDCl₃): δ 7.32 ~ 7.23 (m, 4 H), 7.25 ~ 7.23 (m, 3 H) 6.84 (dd, *J* = 2.4, 1.8 Hz, 2 H), 6.52 (d, *J* = 16.2 Hz, 1 H), 6.13 (dd, *J* = 16.2, 6.6 Hz, 1 H), 4.49 (q, *J* = 6.6 Hz, 1 H), 3.79 (s, 3 H), 2.95 (dd, *J* = 4.8, 5.4 Hz, 1 H), 2.86 (dd, *J* = 8.4, 7.8 Hz, 1 H), 1.74 (bs, 1 H); ¹³C NMR (150 MHz, CDCl₃): δ 159.3, 137.8, 130.0, 129.6, 129.4, 129.3, 128.5, 127.6, 126.5, 114.0, 73.6, 55.3, 44.2; HRMS: calcd. for C₁₇H₁₈O₂: 254.1307; Found: 254.1302.

Spectral data for (E)-1-phenyl-4-(4-(trifluoromethyl)phenyl)but-3-en-2-ol (5f)



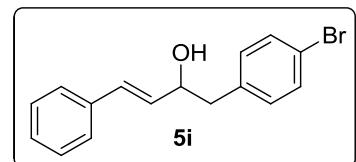
Off White Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.56 (d, $J = 7.8$ Hz, 2 H), 7.44 (d, $J = 8.4$ Hz, 2 H), 7.35 ~ 7.32 (m, 2 H), 7.27 ~ 7.25 (m, 3 H), 6.62 (d, $J = 16.2$ Hz, 1 H), 6.37 (dd, $J = 16.2, 6.0$ Hz, 1 H), 4.54 (d, $J = 6.6$ Hz, 1 H), 2.99 ~ 2.86 (m, 2 H), 1.95 (s, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ 140.2, 137.3, 134.1, 129.5, 129.4, 129.2, 128.7, 128.6, 126.7, 125.5, 125.5, 125.2, 125.1, 73.1, 44.1; HRMS: calcd. for $\text{C}_{17}\text{H}_{15}\text{F}_3\text{O}$: 292.1075; Found: 292.1081.

Spectral data for (E)-4-phenyl-1-(p-tolyl)but-3-en-2-ol (5h)



Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.37 ~ 7.36 (m, 2 H), 7.31 (t, $J = 7.5$ Hz, 2 H), 7.24 ~ 7.22 (m, 1 H), 7.16 ~ 7.12 (m, 4 H), 6.59 (d, $J = 15.6$ Hz, 1 H), 6.28 (dd, $J = 15.6, 6.0$ Hz, 1 H), 4.51 ~ 4.48 (m, 1 H), 2.93 (dd, $J = 13.8, 4.8$ Hz, 1 H), 2.83 (dd, $J = 13.8, 8.4$ Hz, 1 H), 2.33 (s, 3 H), 1.80 (bs, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ 136.7, 136.1, 134.5, 131.6, 130.3, 129.4, 129.2, 128.5, 127.6, 126.5, 73.5, 43.7, 21.0; HRMS: calcd. for $\text{C}_{17}\text{H}_{18}\text{O}$: 238.1358; Found: 238.1915.

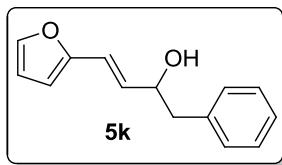
Spectral data for (E)-1-(4-bromophenyl)-4-phenylbut-3-en-2-ol (5i)



Light Yellow Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.42 (d, $J = 8.4$ Hz, 2 H), 7.35 ~ 7.34 (m, 2 H), 7.32 ~ 7.29 (m, 2 H), 7.25 ~ 7.22 (m, 1 H), 7.12 (d, $J = 8.4$ Hz, 2 H), 6.56 (d, J

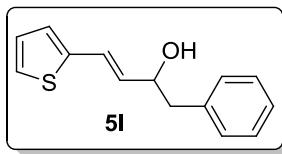
= 16.2 Hz, 1 H), 6.23 (dd, J = 16.2, 6.6 Hz, 1 H), 4.48 (q, J = 6.3 Hz, 1 H), 2.91 ~ 2.82 (m, 2 H), 1.76 (bs, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ 136.7, 136.5, 131.5, 131.3, 131.1, 130.8, 128.6, 127.8, 126.5, 120.5, 73.3, 43.4; HRMS: calcd. for $\text{C}_{16}\text{H}_{15}\text{BrO}$: 302.0306; Found: 301.9871.

Spectral data for (E)-4-(furan-2-yl)-1-phenylbut-3-en-2-ol (5k)



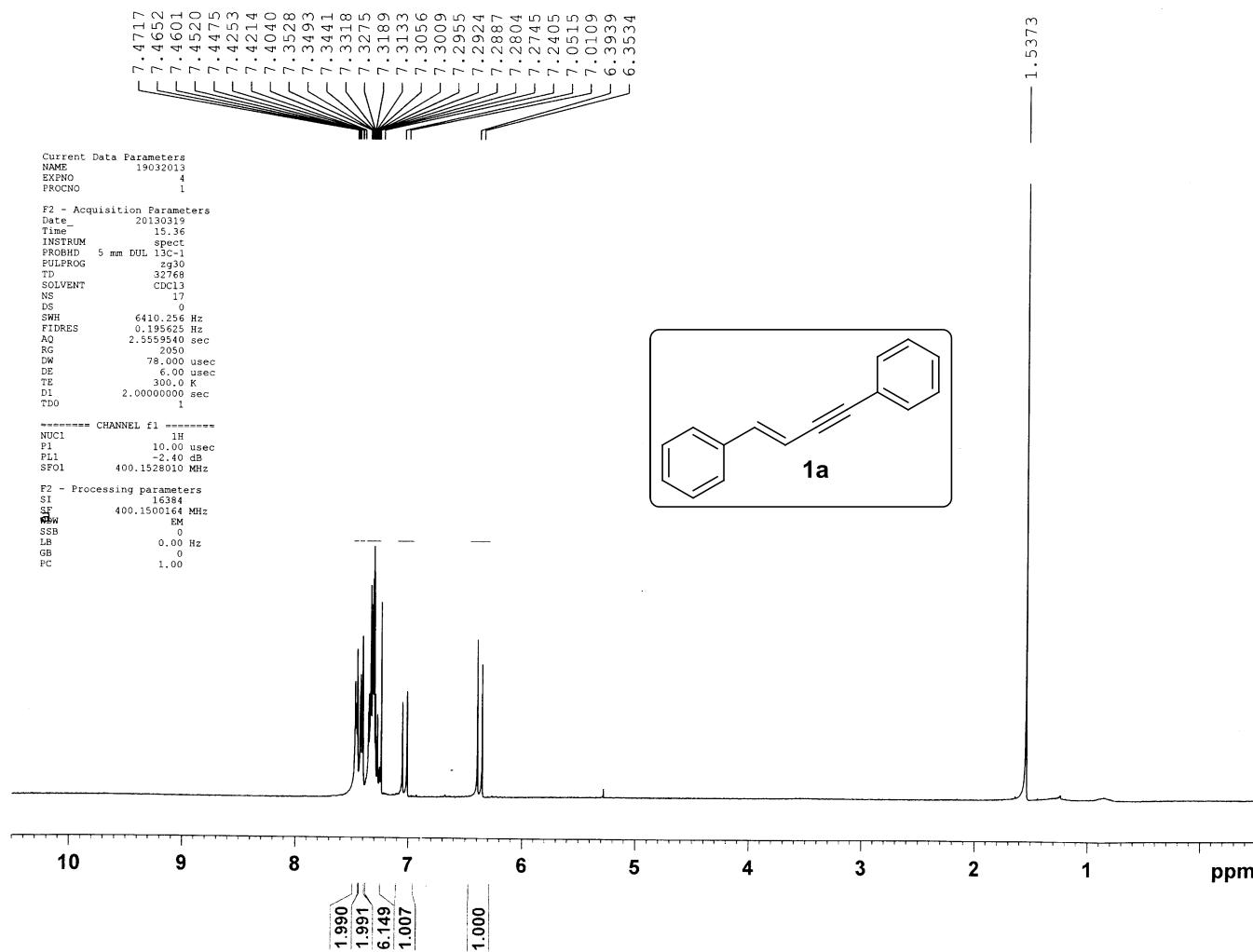
Yellow Oil; ^1H NMR (600 MHz, CDCl_3): δ 7.31 ~ 7.23 (m, 6 H), 6.41 (dd, J = 1.2, 1.2 Hz, 1 H), 6.35 (dd, J = 1.8, 1.8 Hz, 1 H), 6.25 ~ 6.21 (m, 2 H), 4.48 ~ 4.46 (m, 1 H), 2.95 ~ 2.81 (m, 2 H), 1.97 (bs, 1 H); ^{13}C NMR (150 MHz, CDCl_3): δ 152.3, 141.9, 137.6, 130.1, 129.5, 128.5, 126.6, 118.6, 111.2, 108.0, 73.0, 44.1; HRMS: calcd. for $\text{C}_{14}\text{H}_{14}\text{O}_2$: 214.0994; Found: 214.0991.

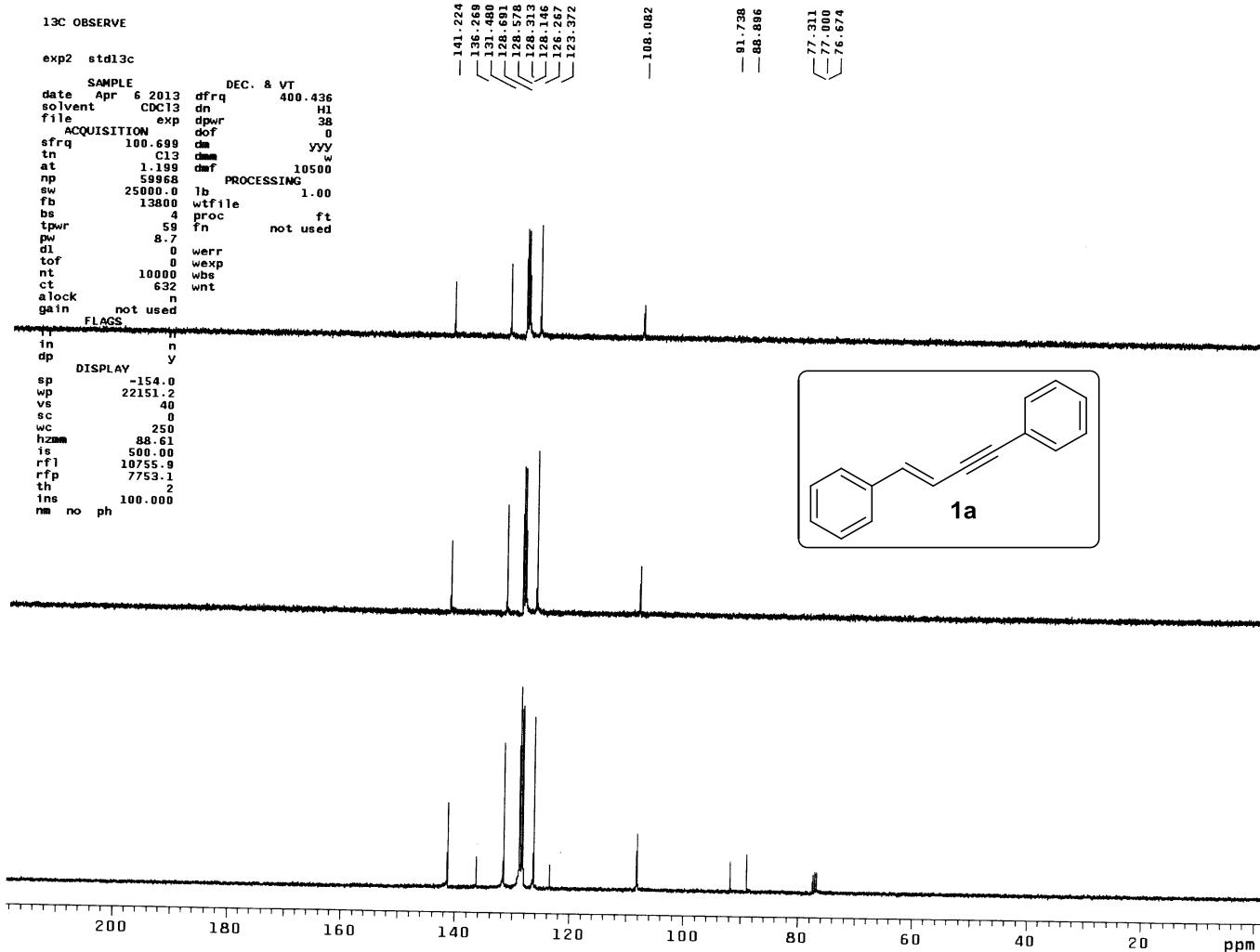
Spectral data for (E)-1-phenyl-4-(thiophen-2-yl)but-3-en-2-ol (5l)

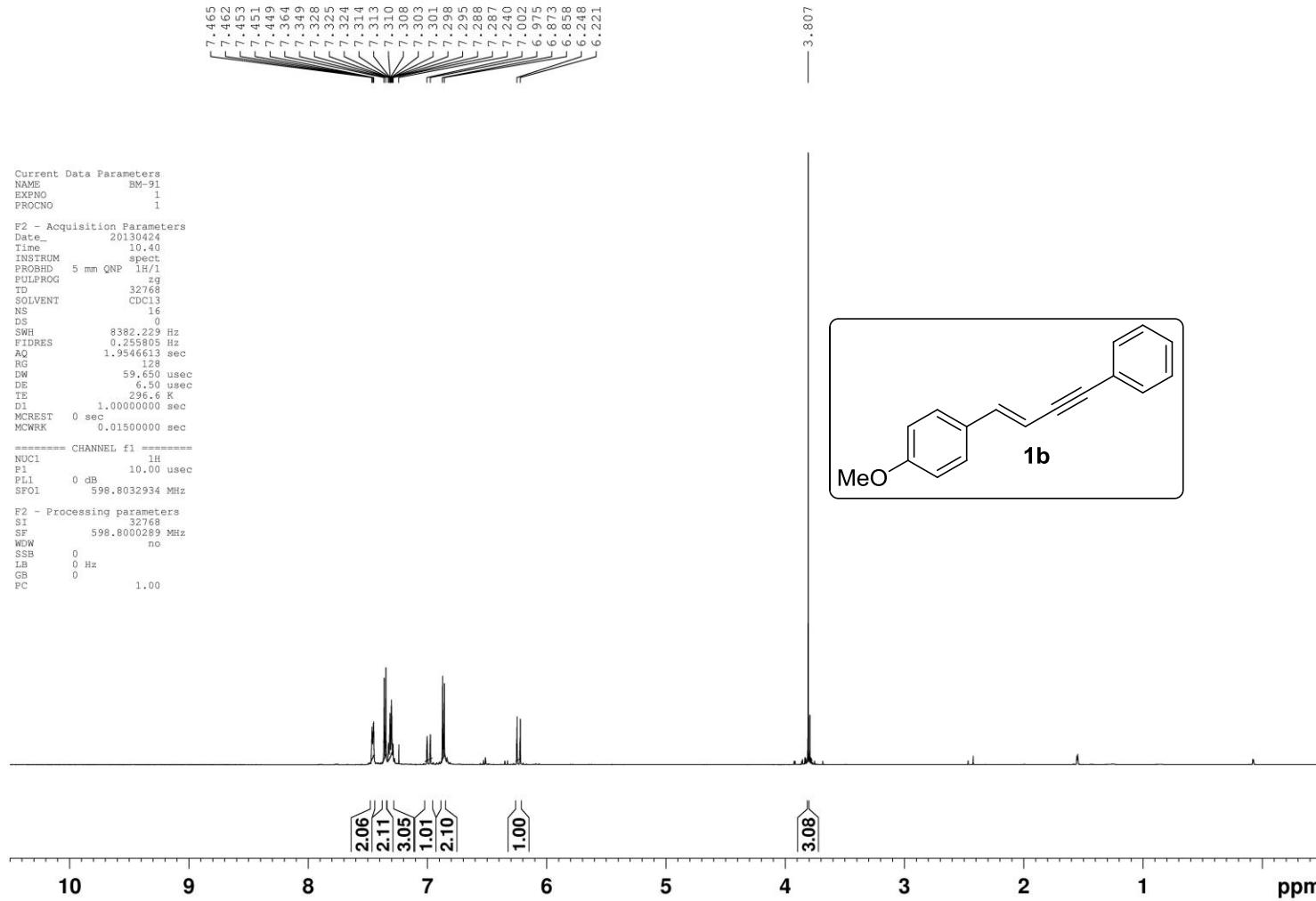


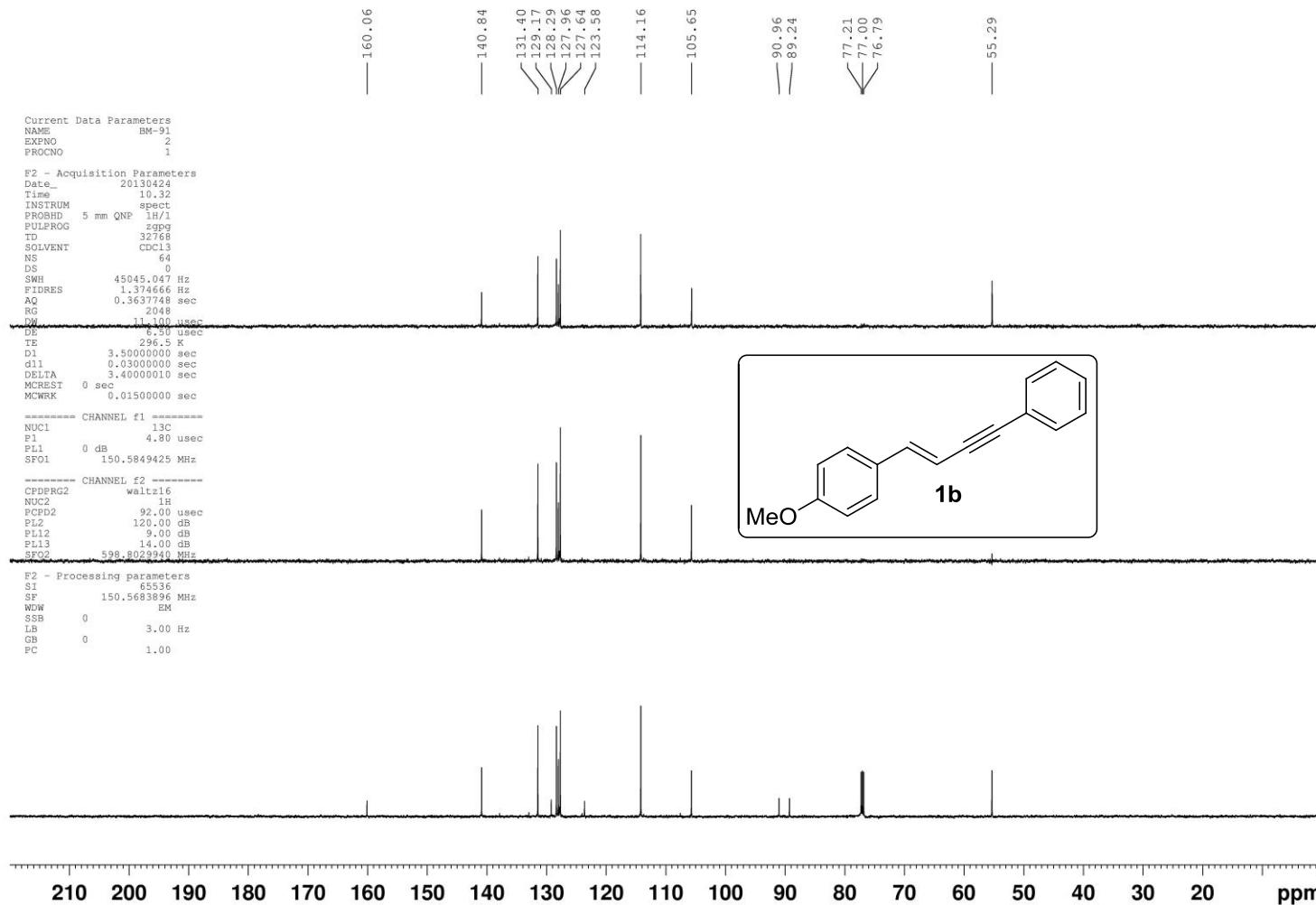
Brown Solid; ^1H NMR (600 MHz, CDCl_3): δ 7.33 ~ 7.31 (m, 2 H), 7.25 ~ 7.23 (m, 3 H), 7.15 (d, J = 5.1 Hz, 1 H), 6.96 ~ 6.93 (m, 2 H), 6.73 (d, J = 15.6 Hz, 1 H), 6.14 (dd, J = 15.6, 6.2 Hz, 1 H), 4.49 ~ 4.46 (m, 1 H), 2.96 ~ 2.83 (m, 2 H), 1.81 (bs, 1 H); ^{13}C NMR

(150 MHz, CDCl₃): δ 141.8, 137.5, 131.1, 129.5, 128.5, 127.3, 126.6, 125.8, 124.3, 123.5, 73.1, 44.1; HRMS: calcd. for C₁₄H₁₄OS: 230.0765; Found: 230.9852.









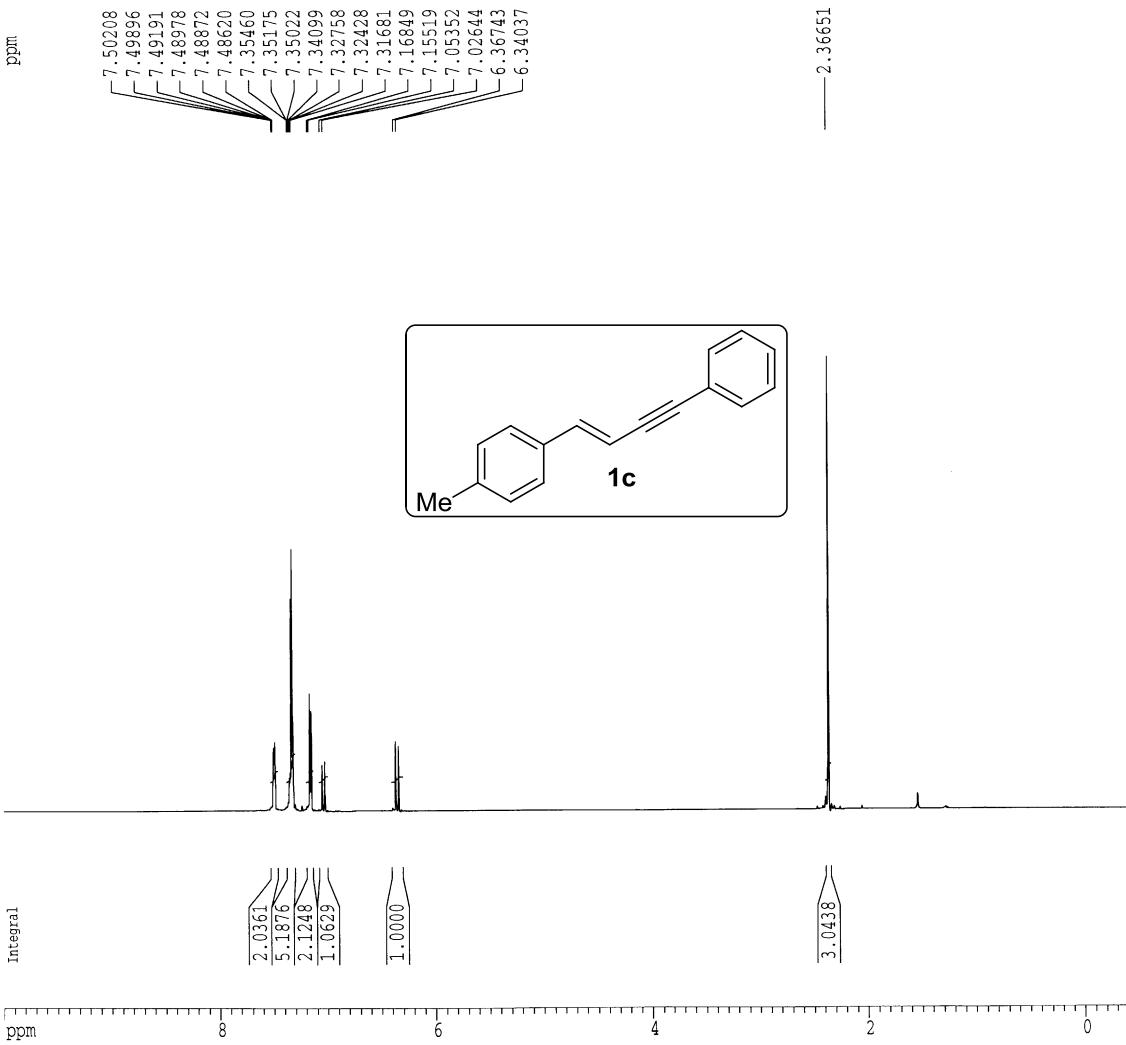
Current Data Parameters
NAME BM-106
EXPNO 1
PROCNO 1

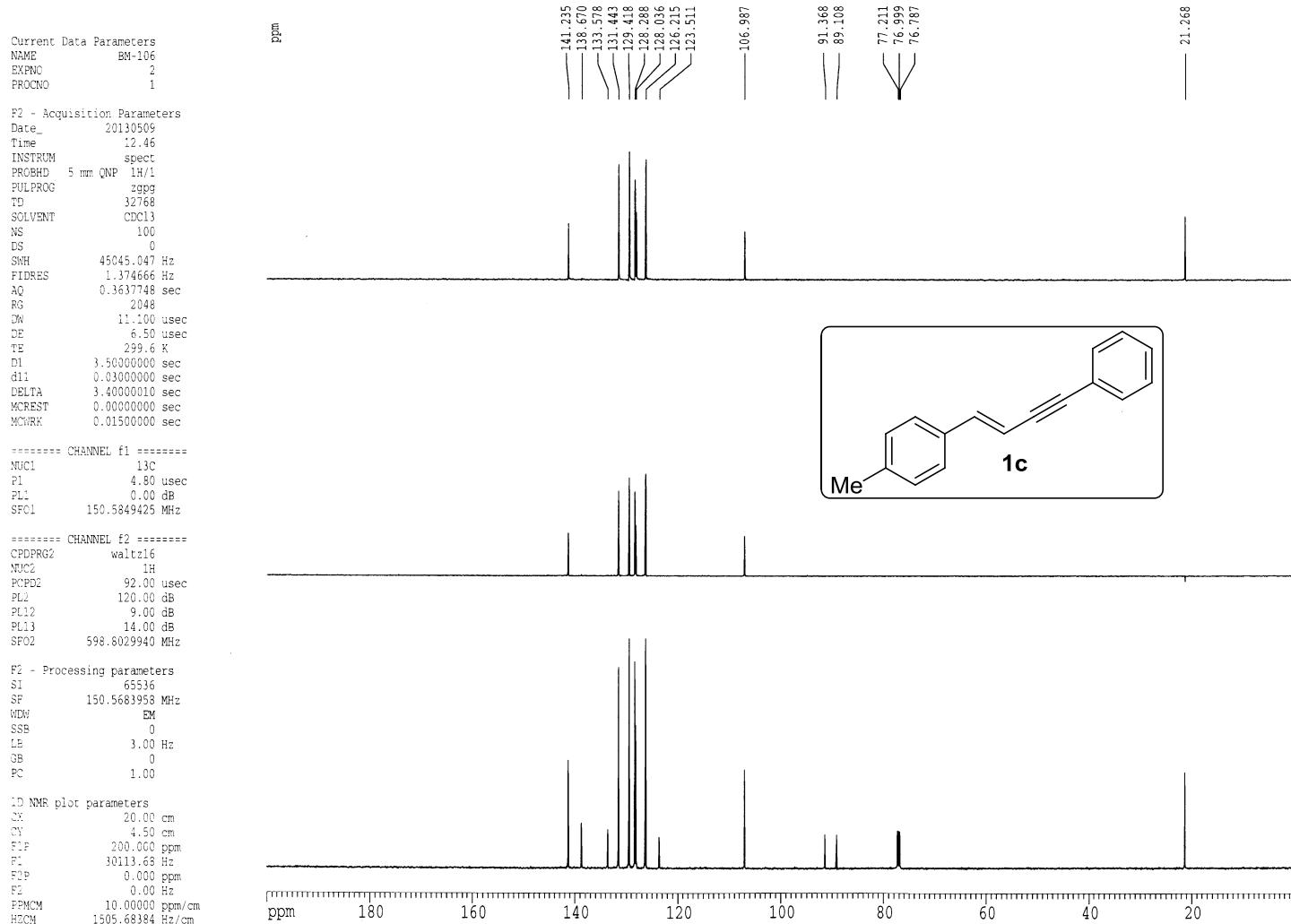
F2 - Acquisition Parameters
Date_ 20130509
Time 16.06
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 32768
SOLVENT CDCl3
NS 16
DS 0
SWH 8382.229 Hz
FIDRES 0.255805 Hz
AQ 1.9546613 sec
RG 128
DM 59.650 usec
DE 6.50 usec
TE 298.7 K
D1 2.0000000 sec
MCREST 0.0000000 sec
MCURK 0.0150000 sec

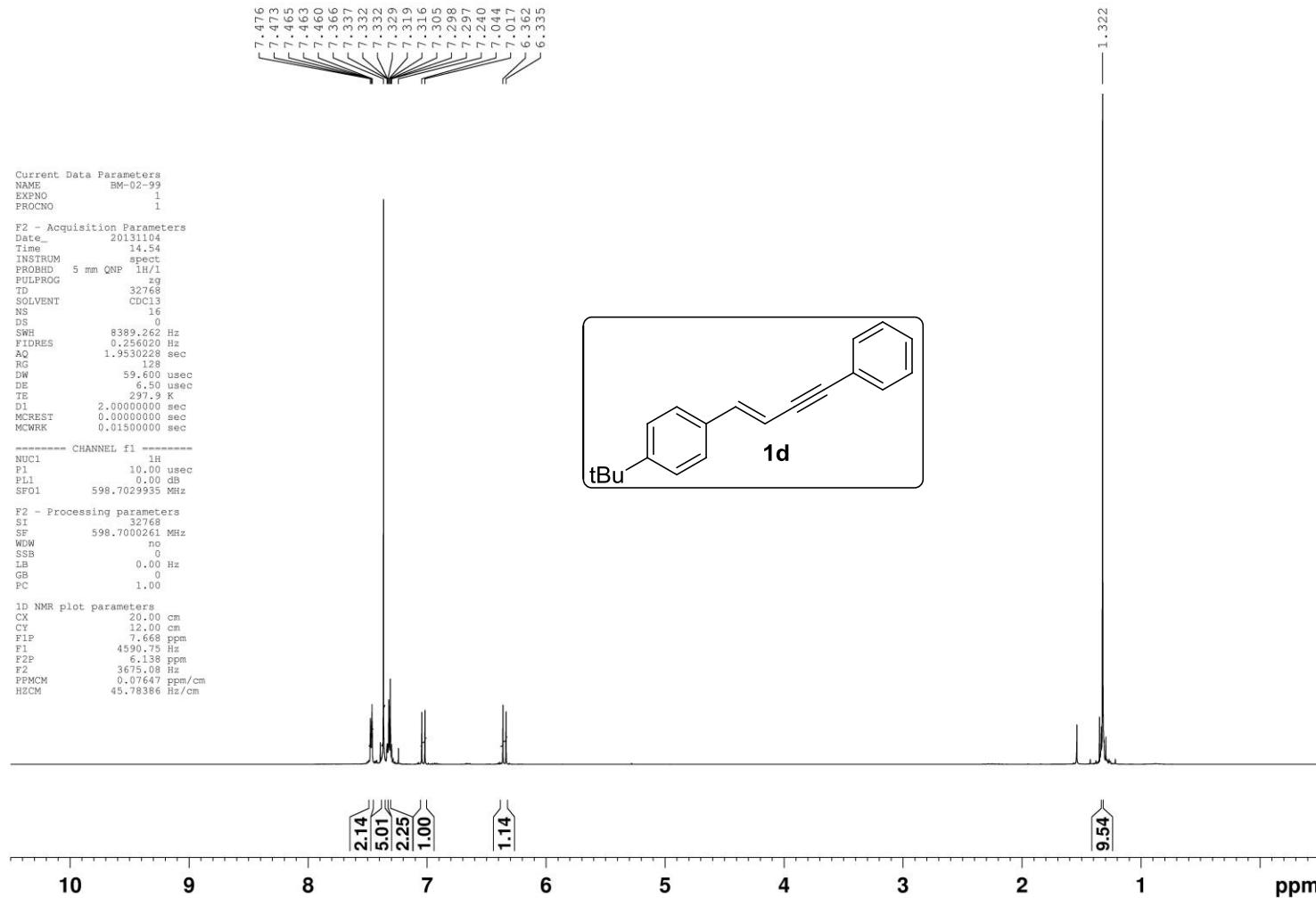
***** CHANNEL f1 *****
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SF01 598.8029940 MHz

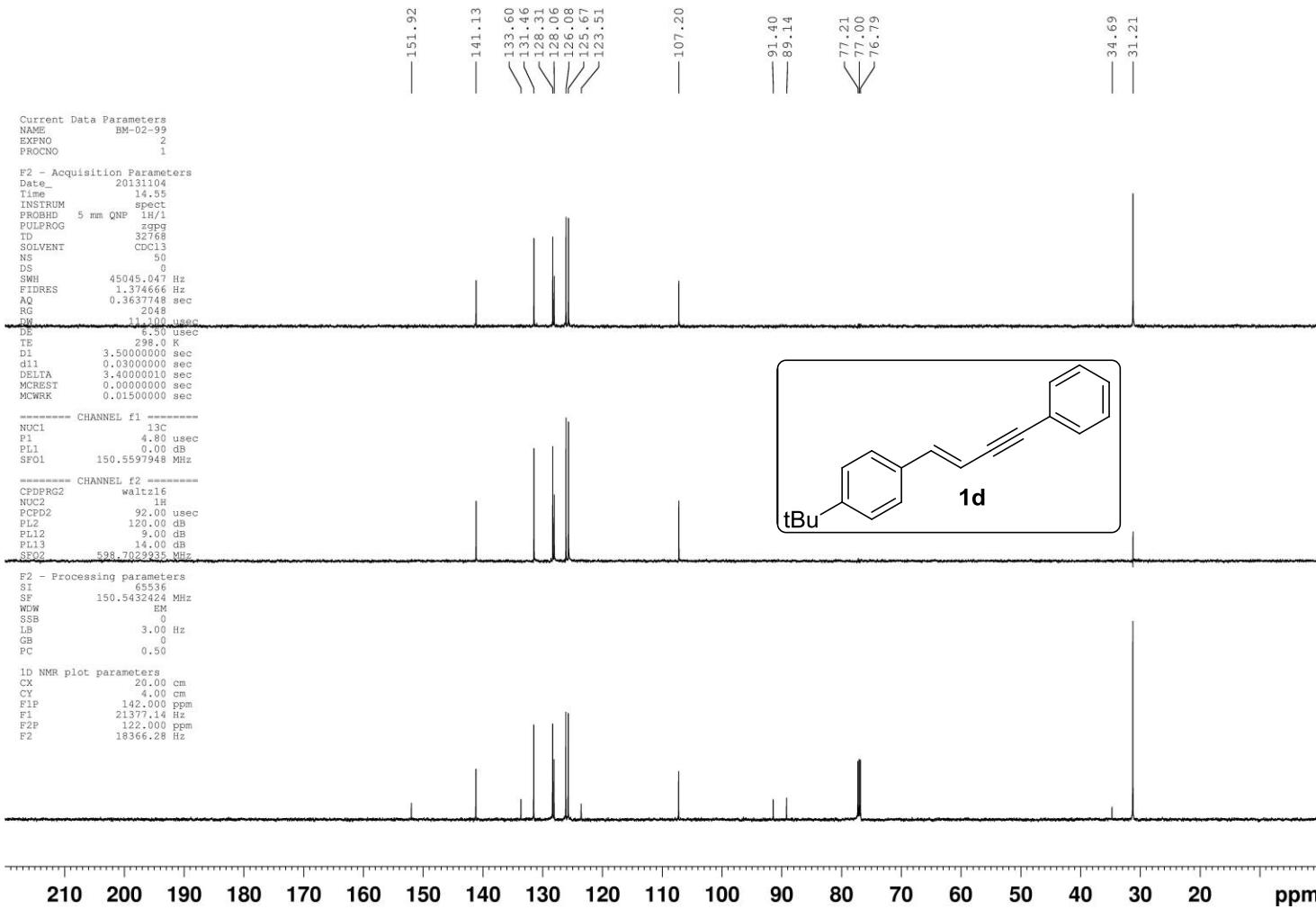
F2 - Processing parameters
SI 32768
SP 598.8000296 MHz
WDW no
SSB 0
LB 0.00 Hz
GB 0
PC 1.00

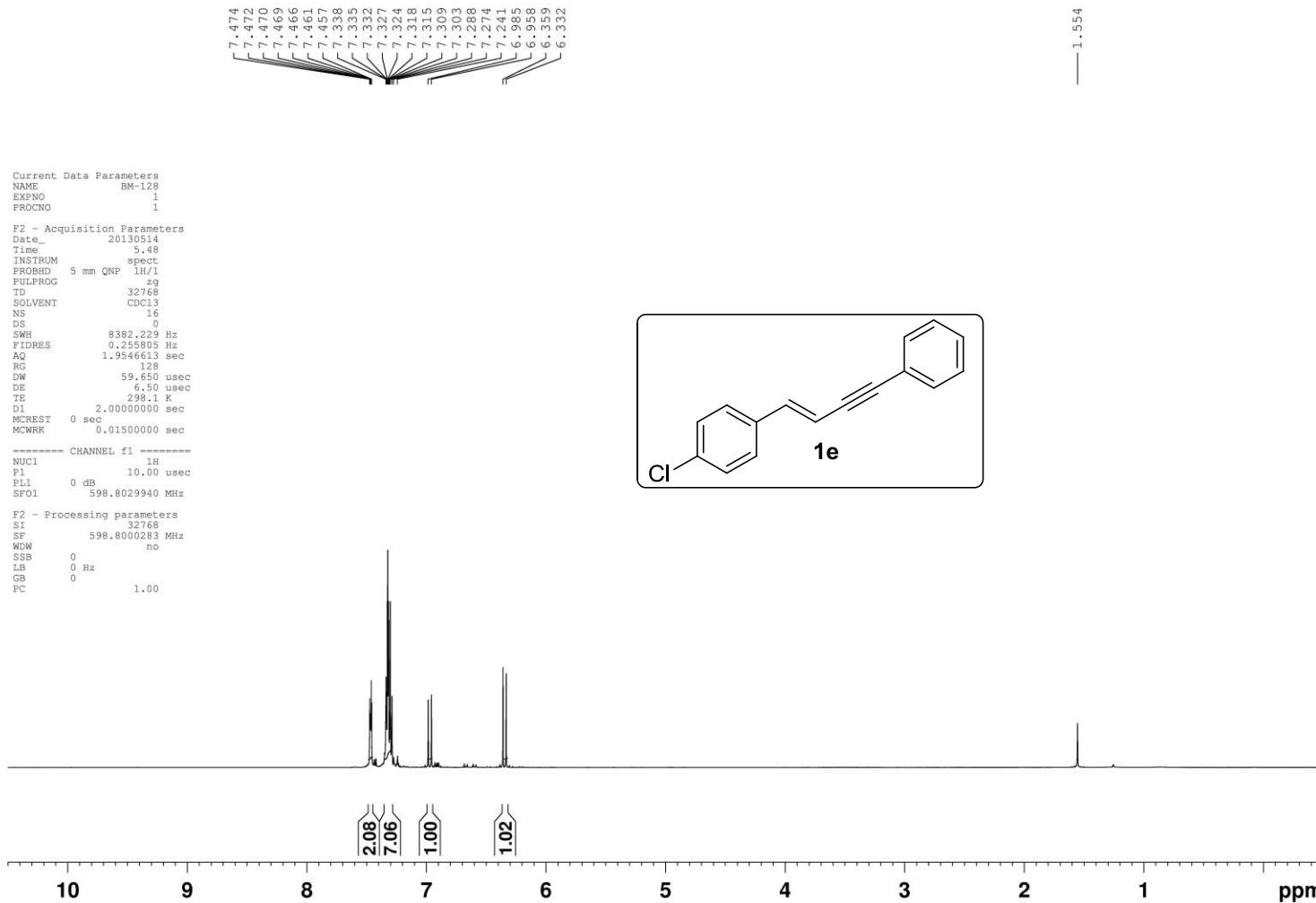
1D NMR plot parameters
CX 20.00 cm
CY 8.00 cm
F1P 10.000 ppm
F1 5988.00 Hz
F2P -0.500 ppm
F2 -299.40 Hz
PPMCM 0.52500 ppm/cm
HEGM 314.37003 Hz/cm

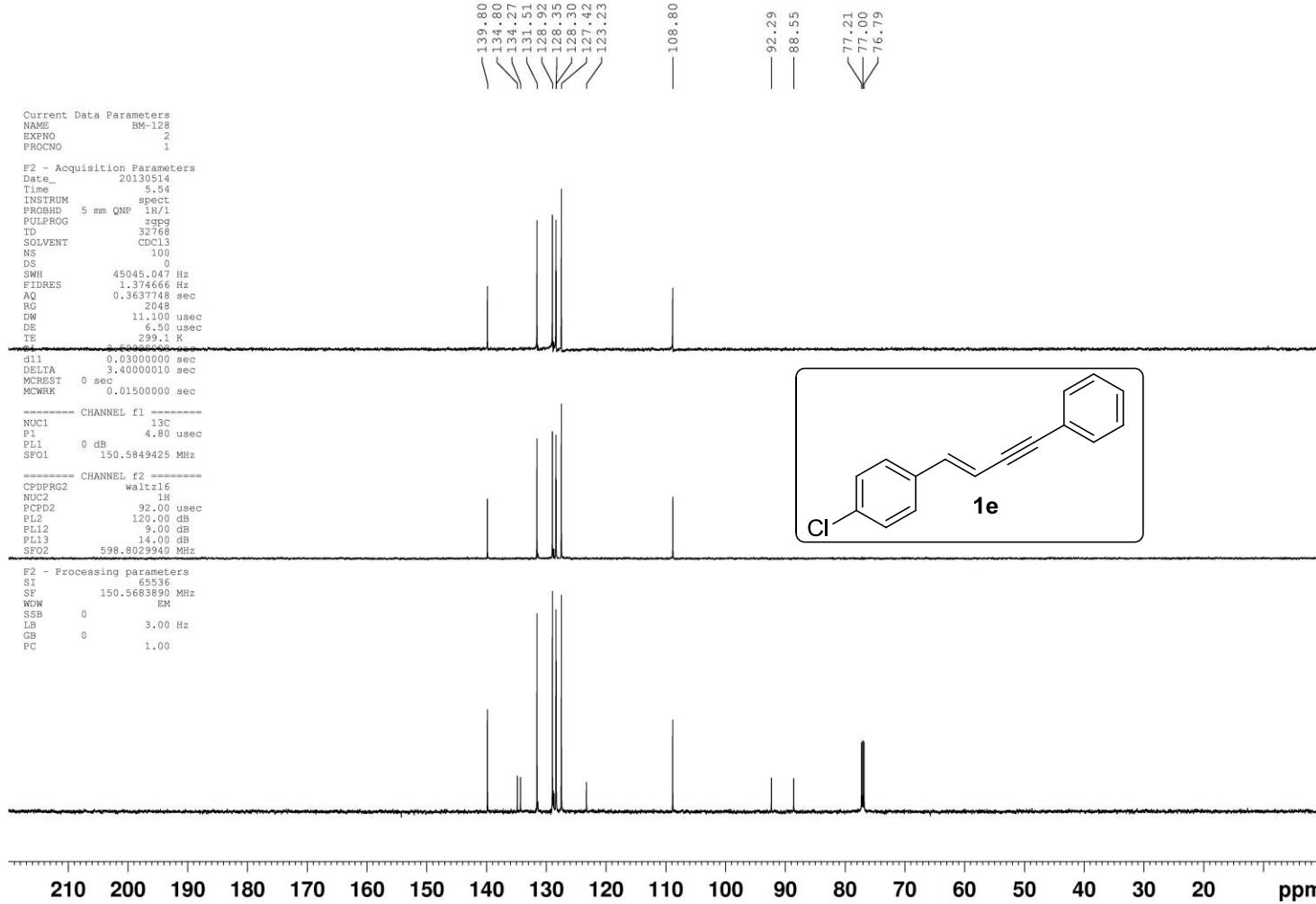












Current Data Parameters
NAME BM-116
EXPN: 1
PRCNO 1

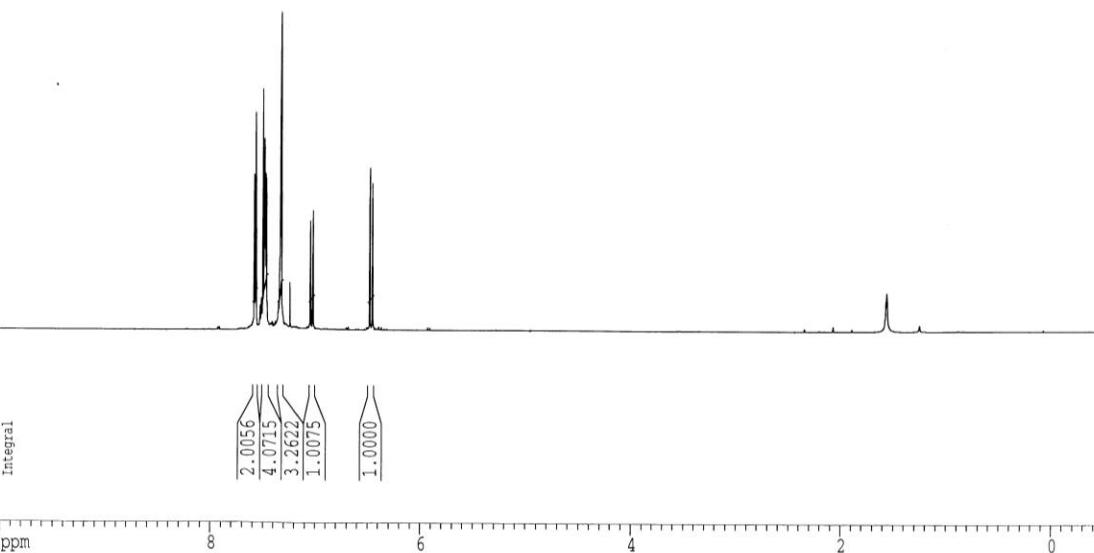
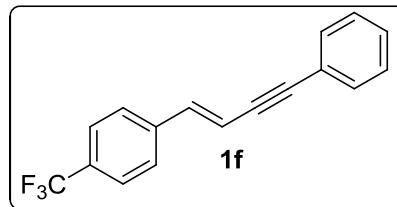
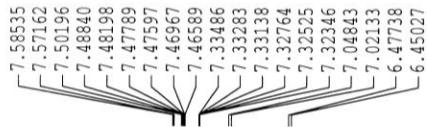
F2 - Acquisition Parameters
Date_ 20130509
Time 16.12
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 32768
SOLVENT CDCl3
NS 16
DS 0
SWH 8382.229 Hz
FIDRES 0.255805 Hz
AQ 1.954613 sec
RG 128
DW 59.650 usec
DE 6.50 usec
TE 298.7 K
D1 2.0000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

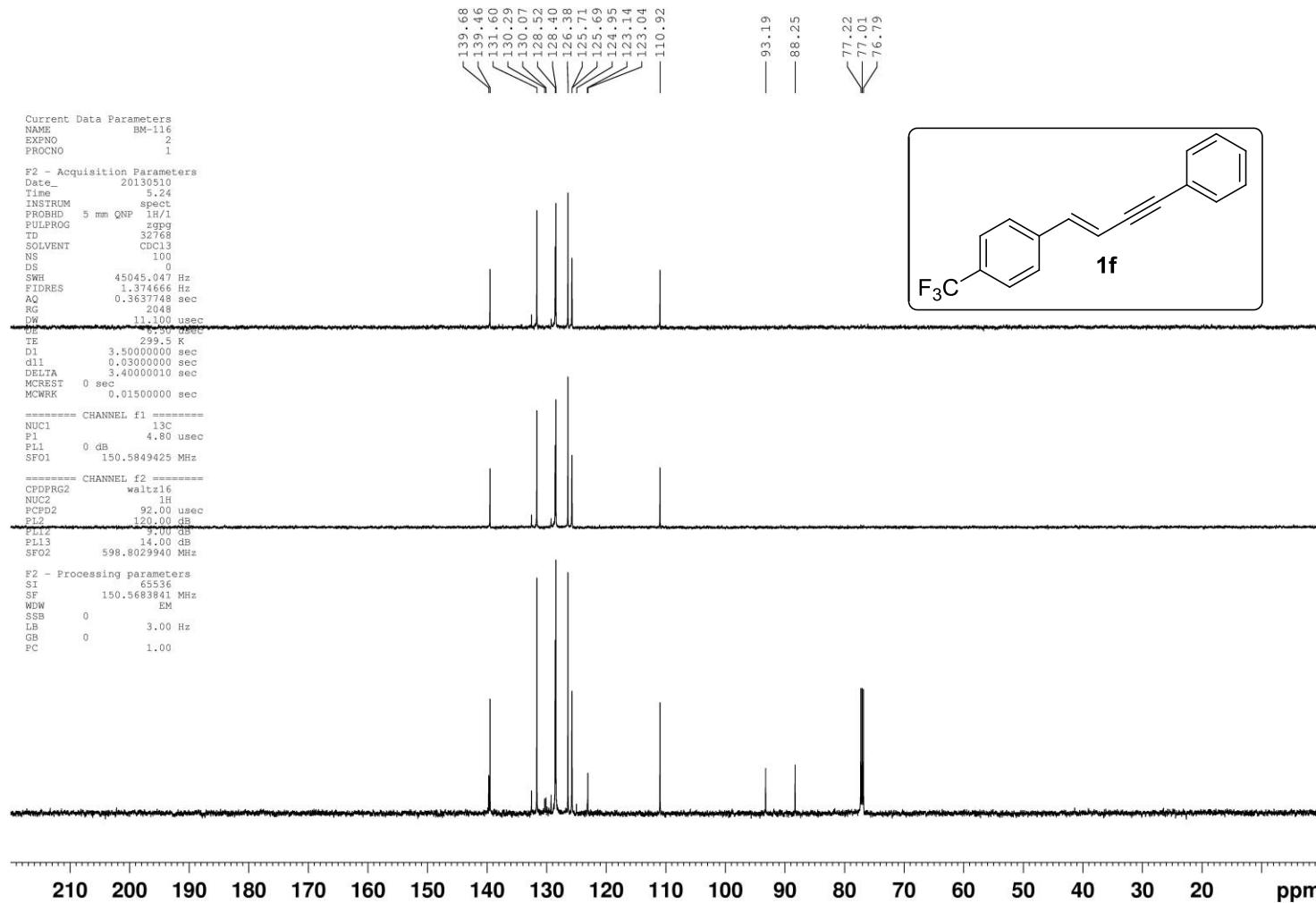
***** CHANNEL f1 *****
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SF01 598.8029940 MHz

F2 - Processing parameters
SI 32768
SF 598.8000285 MHz
NDW no
SSB 0
LB 0.00 Hz
GB 0
PC 1.00

1D NMR plot parameters
CX 20.00 cm
CY 6.00 cm
F1P 10.000 ppm
F1 5988.00 Hz
F2P -0.500 ppm
F2 -299.40 Hz
PPCM 0.52500 ppm/cm
HCMM 314.37003 Hz/cm

ppm





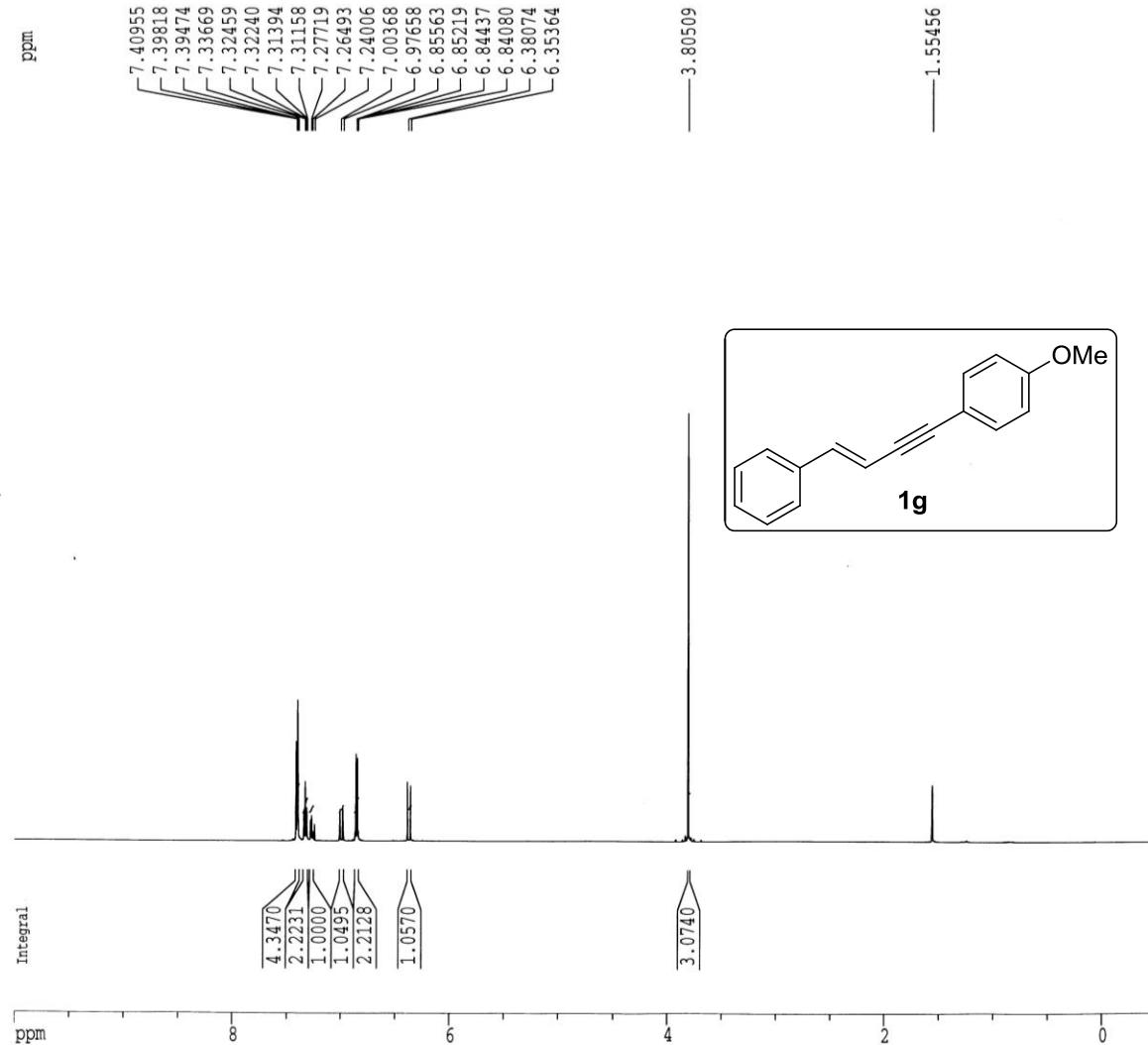
Current Data Parameters
 NAME BM-02-201
 EXPNO 1
 PROCNO 1

F2 - Acquisition Parameters
 Date_ 20140316
 Time 18.50
 INSTRUM spect
 PROBHD 5 mm QNP 1H/1
 PULPROG zg
 TD 33556
 SOLVENT CDCl3
 NS 16
 DS 0
 SWH 8389.262 Hz
 FIDRES 0.250008 Hz
 AQ 1.9999876 sec
 RG 512
 DW 59.600 usec
 DE 6.50 usec
 TE 294.7 K
 D1 2.0000000 sec
 MCREST 0.0000000 sec
 MCWRK 0.0150000 sec

===== CHANNEL f1 =====
 NUC1 1H
 P1 10.00 usec
 PL1 0.00 dB
 SFO1 598.7029935 MHz

F2 - Processing parameters
 SI 32768
 SF 598.7000263 MHz
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 LB 0.00 Hz
 GB 0
 PC 1.00

1D NMR plot parameters
 CX 20.00 cm
 CY 8.00 cm
 F1P 10.000 ppm
 F1 5987.00 Hz
 F2P -0.500 ppm
 F2 -299.35 Hz
 PPMCM 0.52500 ppm/cm
 HZCM 314.31750 Hz/cm



Current Data Parameters
NAME BM-02-201
EXPNO 2
PROCNO 1

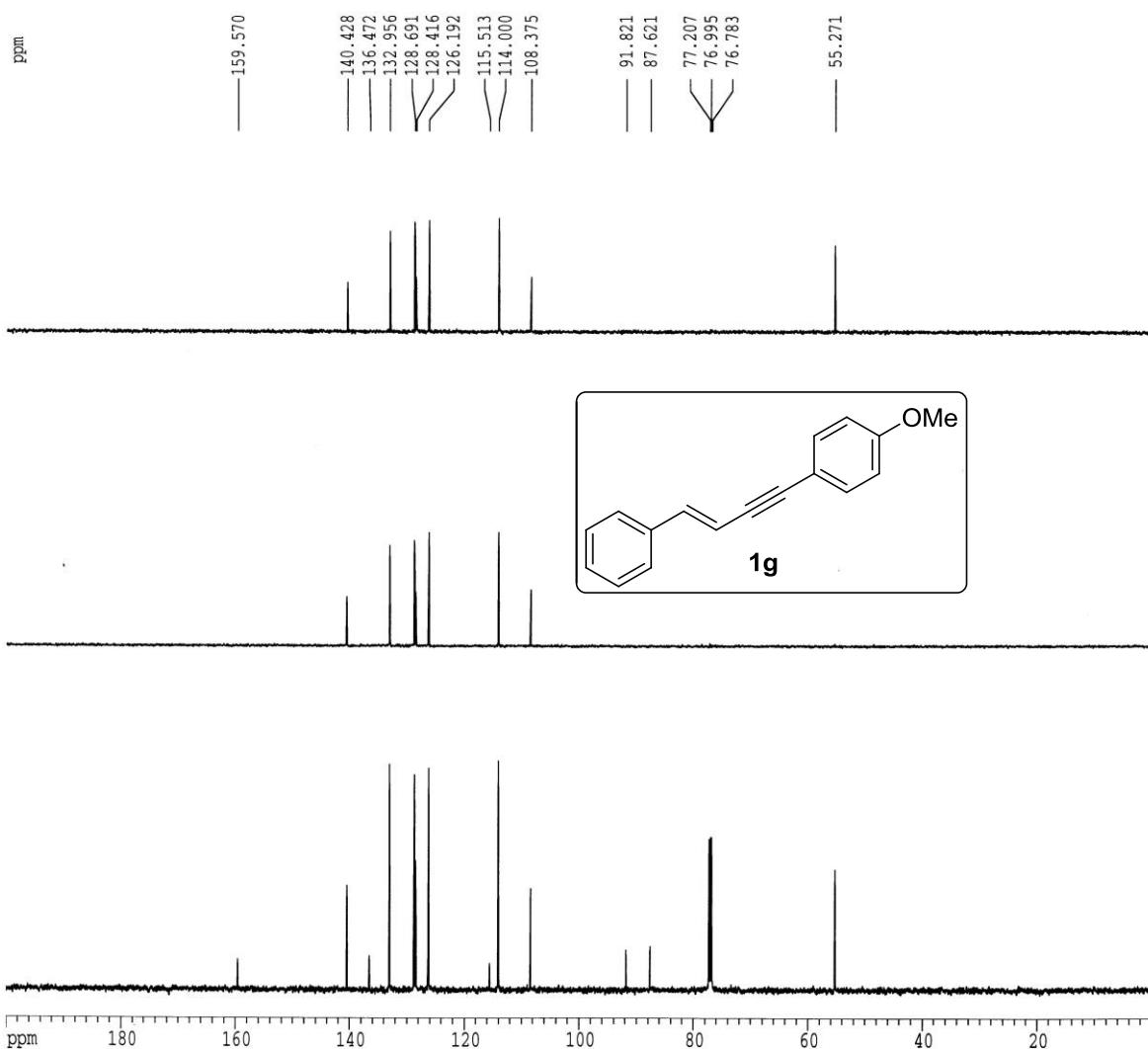
F2 - Acquisition Parameters
Date 20140316
Time 18.57
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT CDCl3
NS 100
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 296.2 K
D1 3.5000000 sec
d11 0.0300000 sec
DELTA 3.40000010 sec
MCREST 0.0000000 sec
MCWRK 0.01500000 sec

===== CHANNEL f1 =====
NUC1 13C
PL 4.80 usec
PL1 0.00 dB
SFO1 150.5597948 MHz

===== CHANNEL f2 =====
CPDPGR2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SFO2 598.7029935 MHz

F2 - Processing parameters
SI 65536
SF 150.5432418 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 0.50

1D NMR plot parameters
CX 20.00 cm
CY 4.00 cm
F1P 200.000 ppm
F1 30108.65 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPMCM 10.00000 ppm/cm
HZCM 1505.43237 Hz/cm



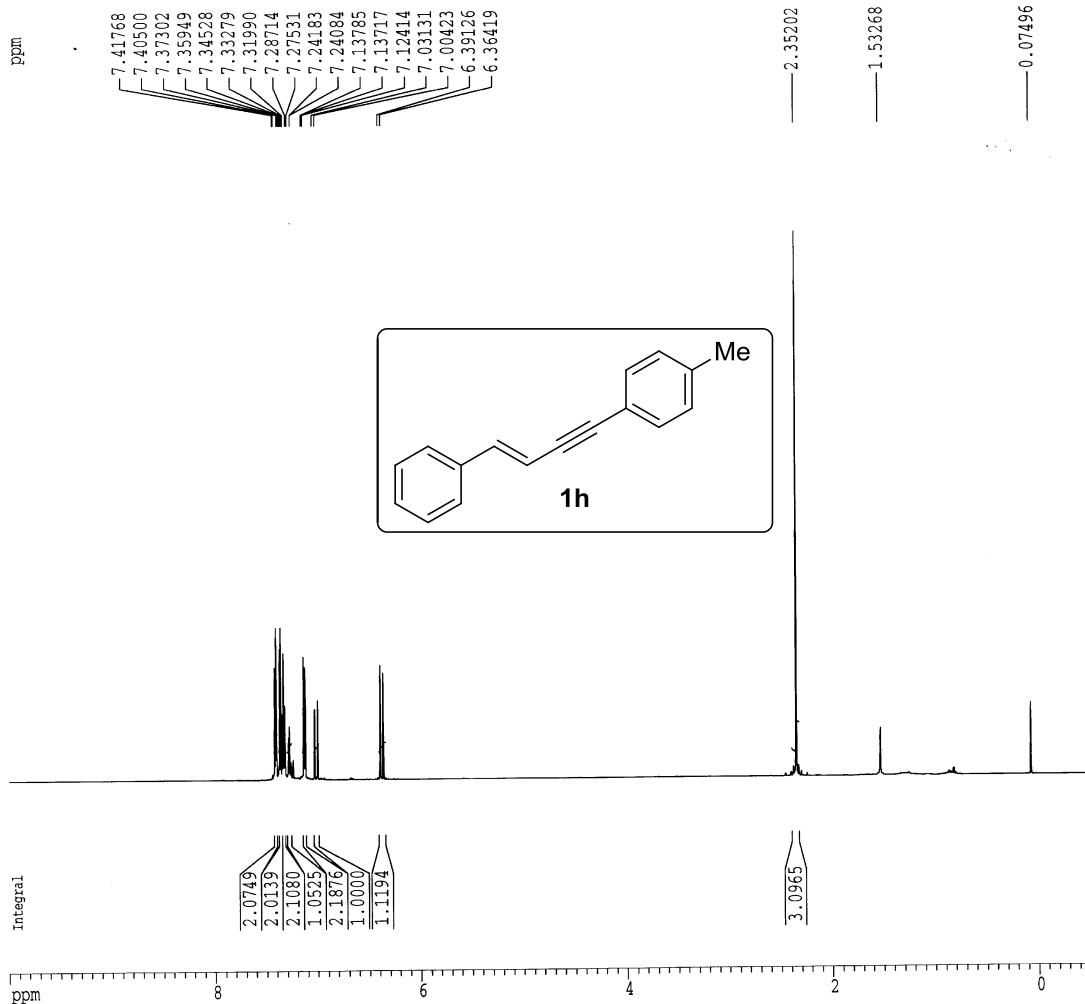
Current Data Parameters
NAME RD-1298
EXPNO 1
PROCNO 1

F2 - Acquisition Parameters
Date 20130703
Time 17.14
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 32768
SOLVENT CDCl3
NS 16
DS 0
SWH 8382.229 Hz
FIDRES 0.255805 Hz
AQ 1.9546613 sec
RG 256
DW 59.650 usec
DE 6.50 usec
TE 300.7 K
D1 1.0000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

===== CHANNEL f1 =====
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SF01 598.8026946 MHz

F2 - Processing parameters
SI 32768
SF 598.8000282 MHz
WWD no
SSB 0
LB 0.00 Hz
GB 0
PC 1.00

1D NMR plot parameters
CX 20.00 cm
CY 10.00 cm
F1P 10.000 ppm
F1 5988.00 Hz
F2P -0.500 ppm
F2 -299.40 Hz
PPMCM 0.52500 ppm/cm
HZCM 314.37003 Hz/cm



Current Data Parameters
NAME RD-1298
EXPNO 2
PROCNO 1

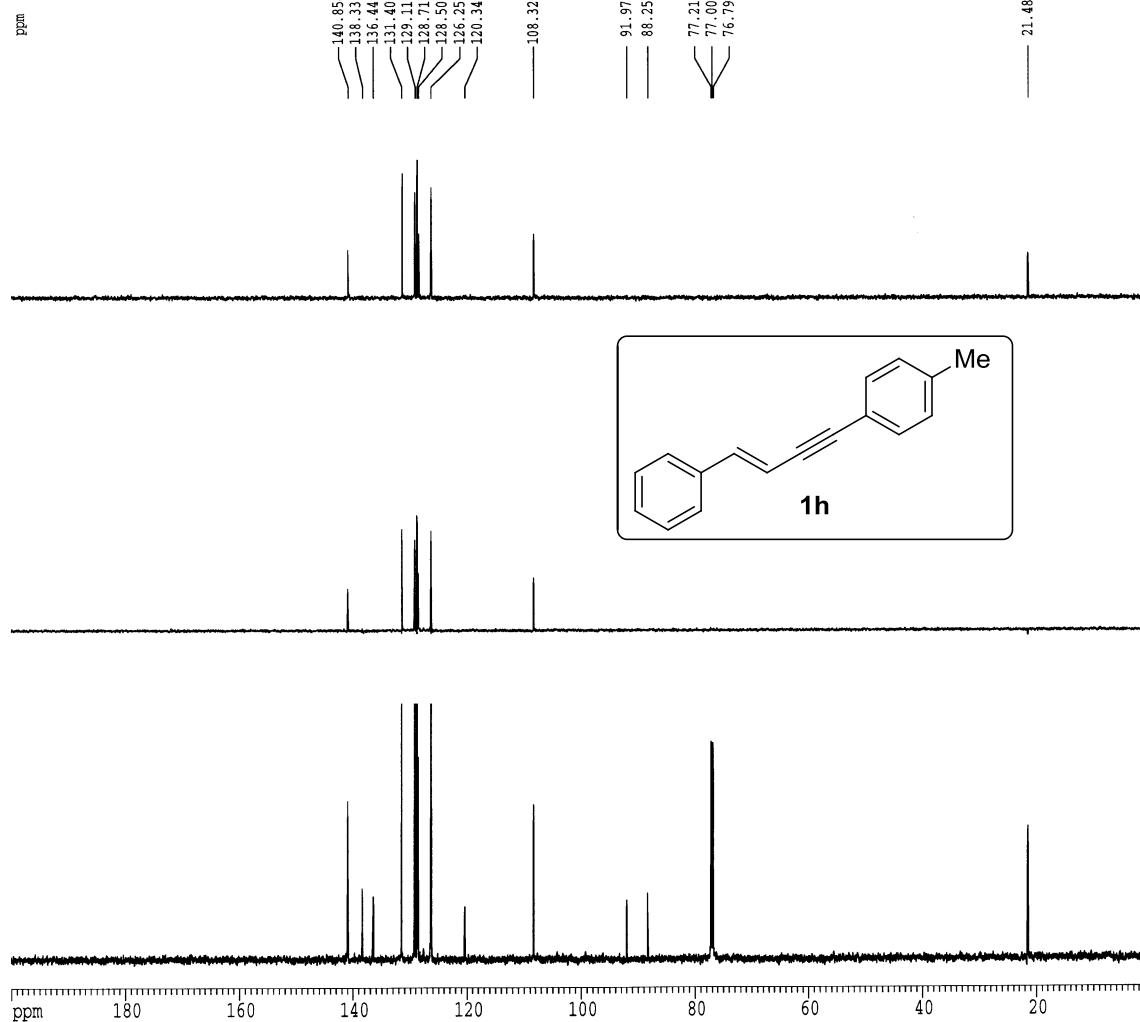
F2 - Acquisition Parameters
Date 20130703
Time 17.15
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT CDCl3
NS 100
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 300.8 K
D1 3.5000000 sec
d11 0.0300000 sec
DELTA 3.4000010 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SF01 150.5849425 MHz

===== CHANNEL f2 =====
CPDPRG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SF02 598.8029940 MHz

F2 - Processing parameters
SI 65536
SF 150.5683855 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 0.50

1D NMR plot parameters
CX 20.00 cm
CY 6.00 cm
F1P 200.00 ppm
F1 30113.68 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPCM 10.00000 ppm/cm
HzCM 1505.68384 Hz/cm



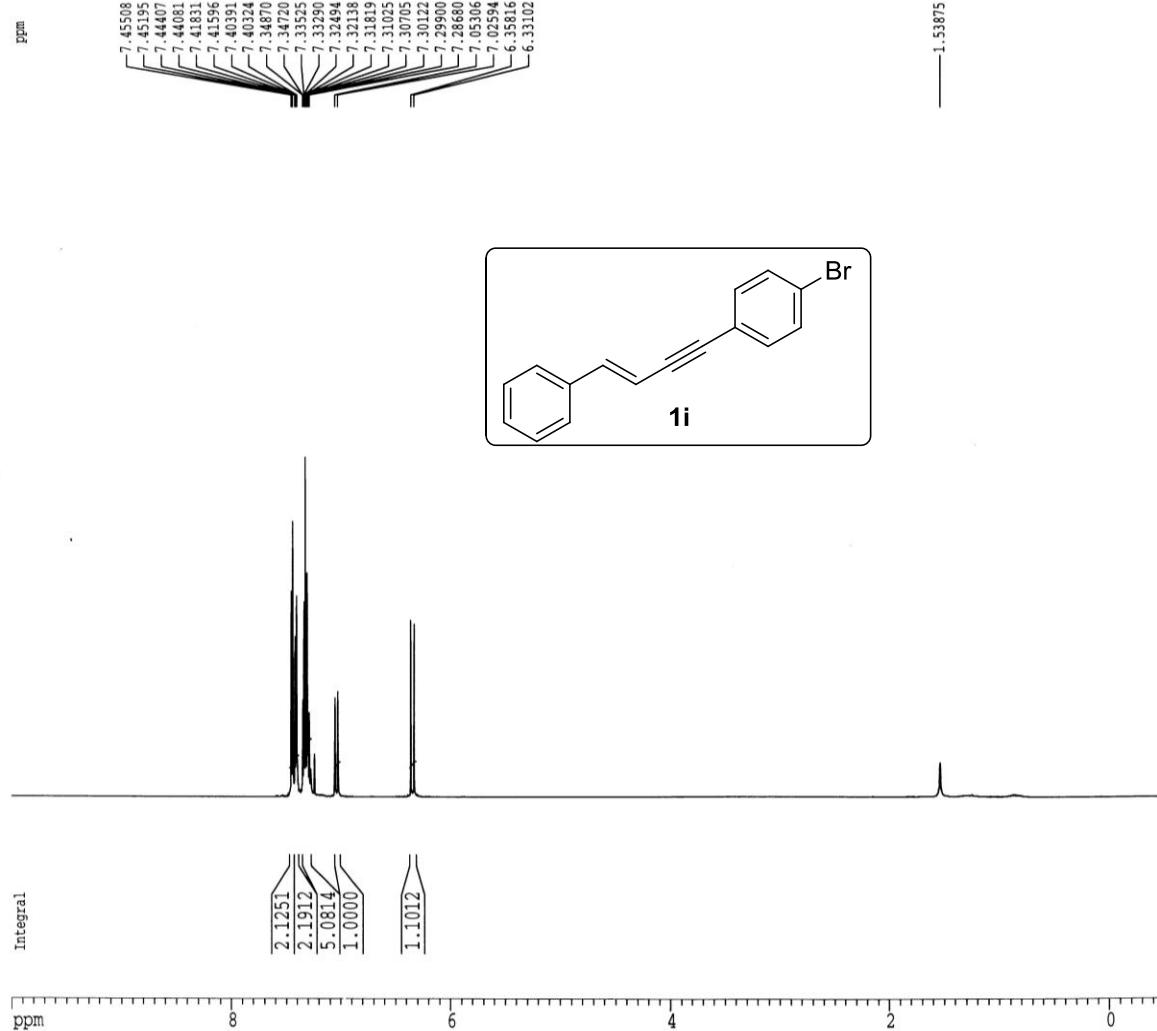
Current Data Parameters
NAME BM-02-35
EXPNO 1
PROCNO 1

F2 - Acquisition Parameters
Date_ 20130903
Time 16.16
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 32768
SOLVENT CDCl3
NS 16
DS 0
SWH 9578.544 Hz
FIDRES 0.292314 Hz
AQ 1.7105396 sec
RG 128
DW 52.200 usec
DE 6.50 usec
TE 301.8 K
D1 1.0000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

===== CHANNEL f1 =====
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SF01 598.8041916 MHz

F2 - Processing parameters
SI 32768
SF 598.8000293 MHz
WDW no
SSB 0
LB 0.00 Hz
GB 0
PC 1.00

1D NMR plot parameters
CX 20.00 cm
CY 6.00 cm
F1P 10.000 ppm
F1 5988.00 Hz
F2P -0.500 ppm
F2 -299.40 Hz
PPCM 0.52500 ppm/cm
HZCM 314.37003 Hz/cm



Current Data Parameters
NAME BM-02-35
EXPNO 2
PROCNO 1

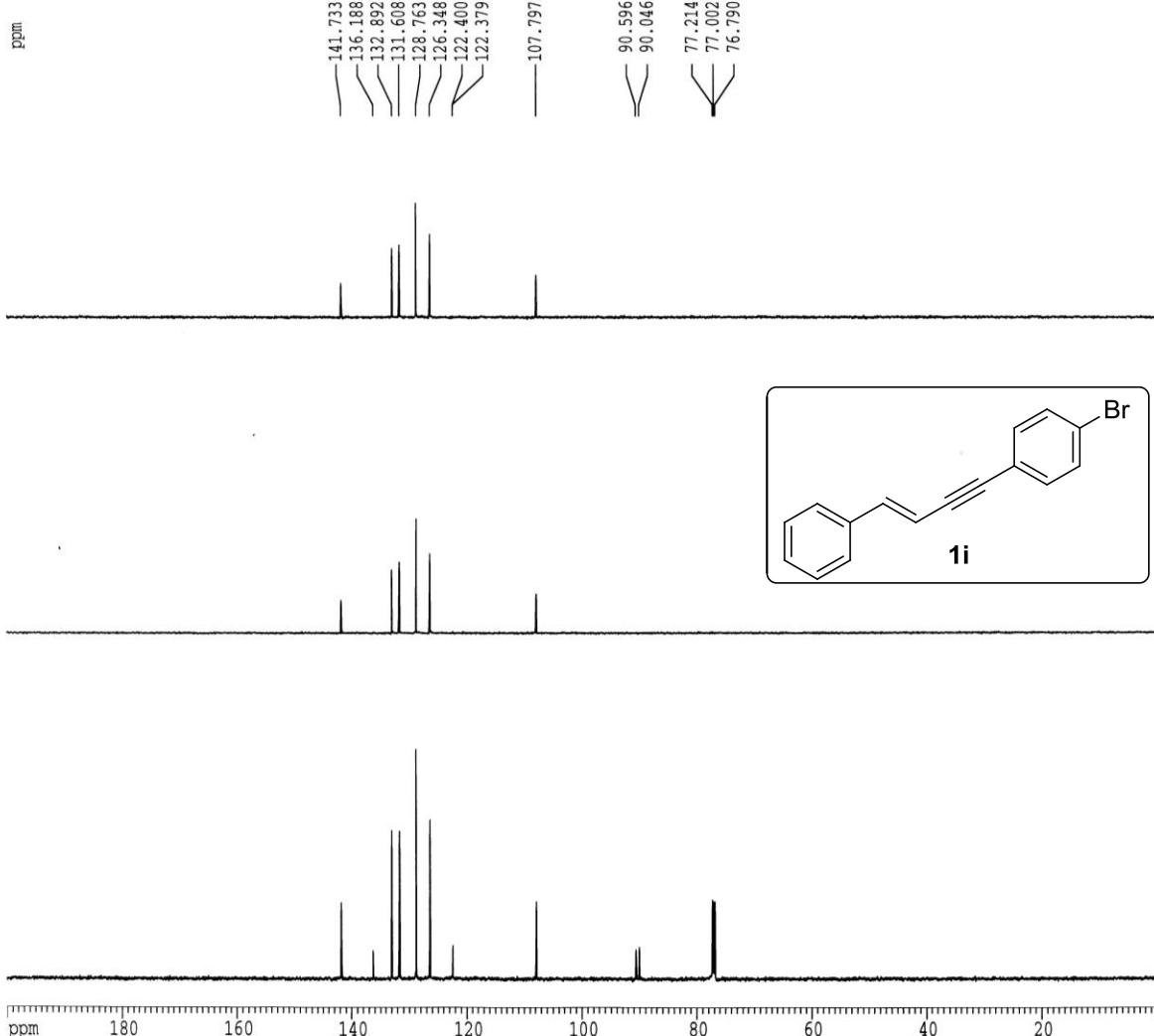
F2 - Acquisition Parameters
Date_ 20130903
Time 16.23
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT CDCl3
NS 100
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 302.8 K
D1 3.5000000 sec
d11 0.0300000 sec
DELTA 3.4000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SPO1 150.5849425 MHz

===== CHANNEL f2 =====
CPDPG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SFQ2 598.8029940 MHz

F2 - Processing parameters
SI 65536
SF 150.5683855 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 0.50

1D NMR plot parameters
CX 20.00 cm
CY 4.00 cm
F1P 200.000 ppm
F1 30113.68 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPCM 10.00000 ppm/cm
HZCM 1505.68384 Hz/cm



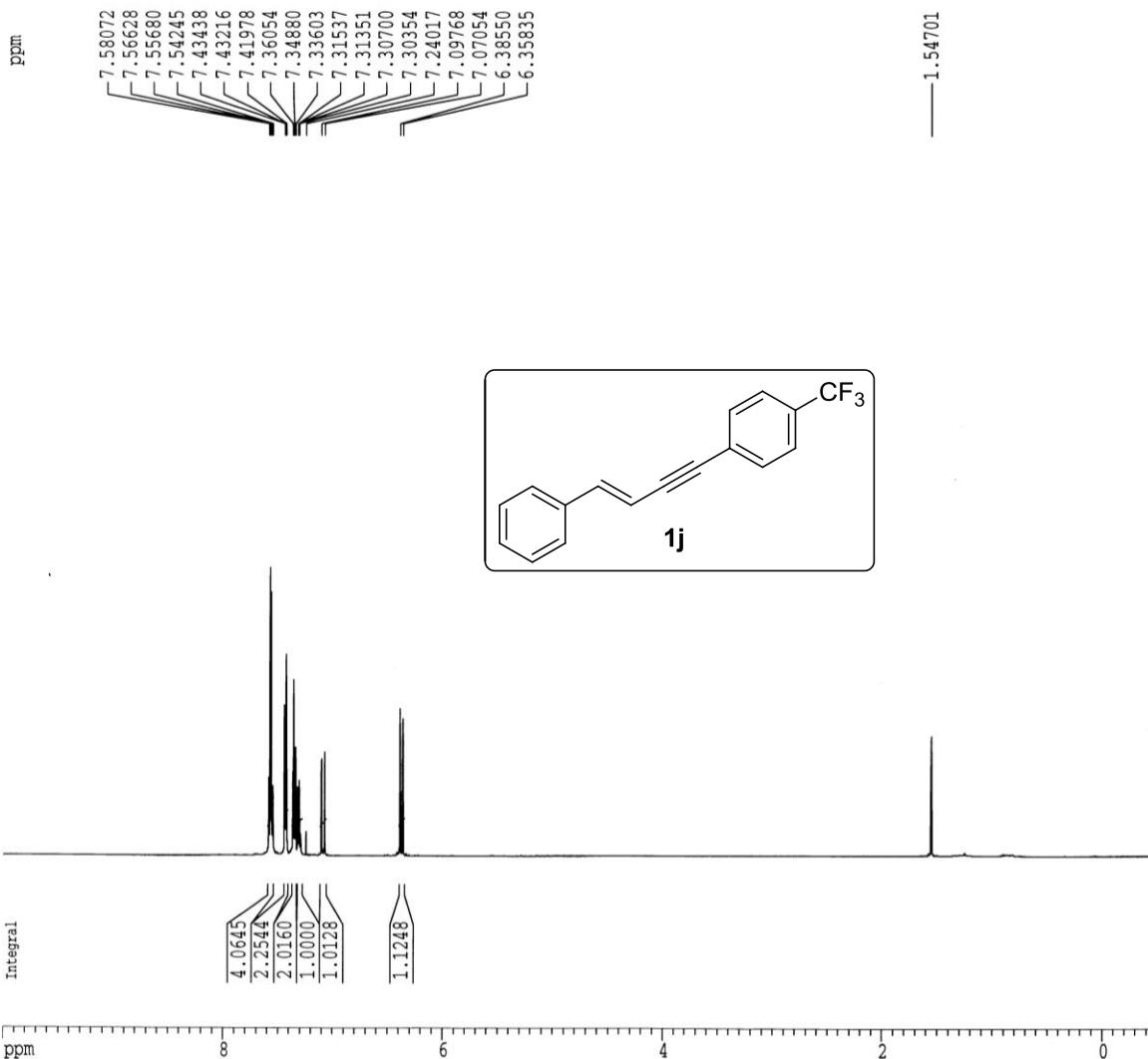
Current Data Parameters
NAME BM-02-202
EXPNO 1
PROCNO 1

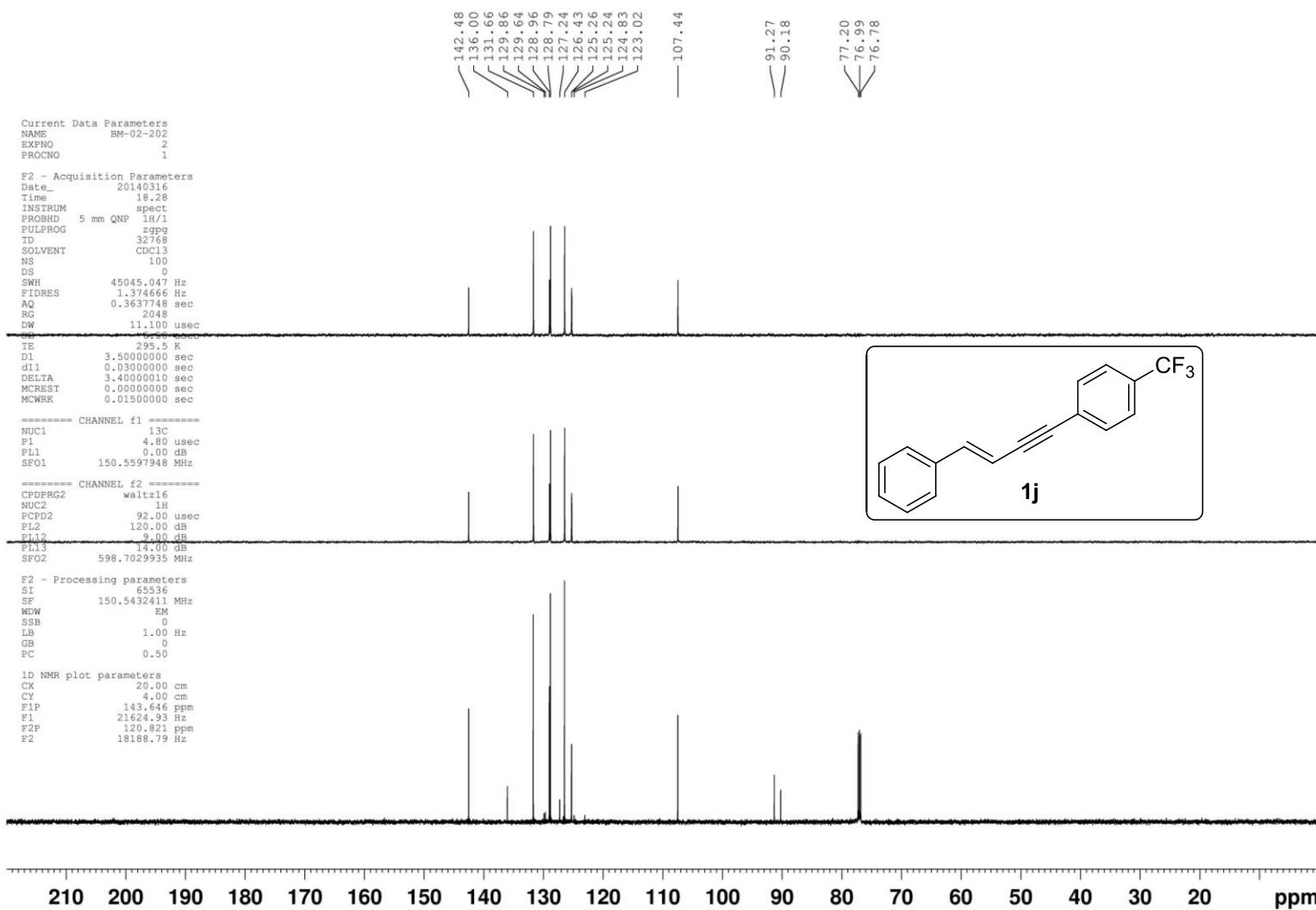
F2 - Acquisition Parameters
Date 20140316
Time 18.25
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 33556
SOLVENT CDCl3
NS 16
DS 0
SWH 8389.262 Hz
FIDRES 0.250008 Hz
AQ 1.9999876 sec
RG 512
DW 59.600 usec
DE 6.50 usec
TE 294.6 K
D1 2.0000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

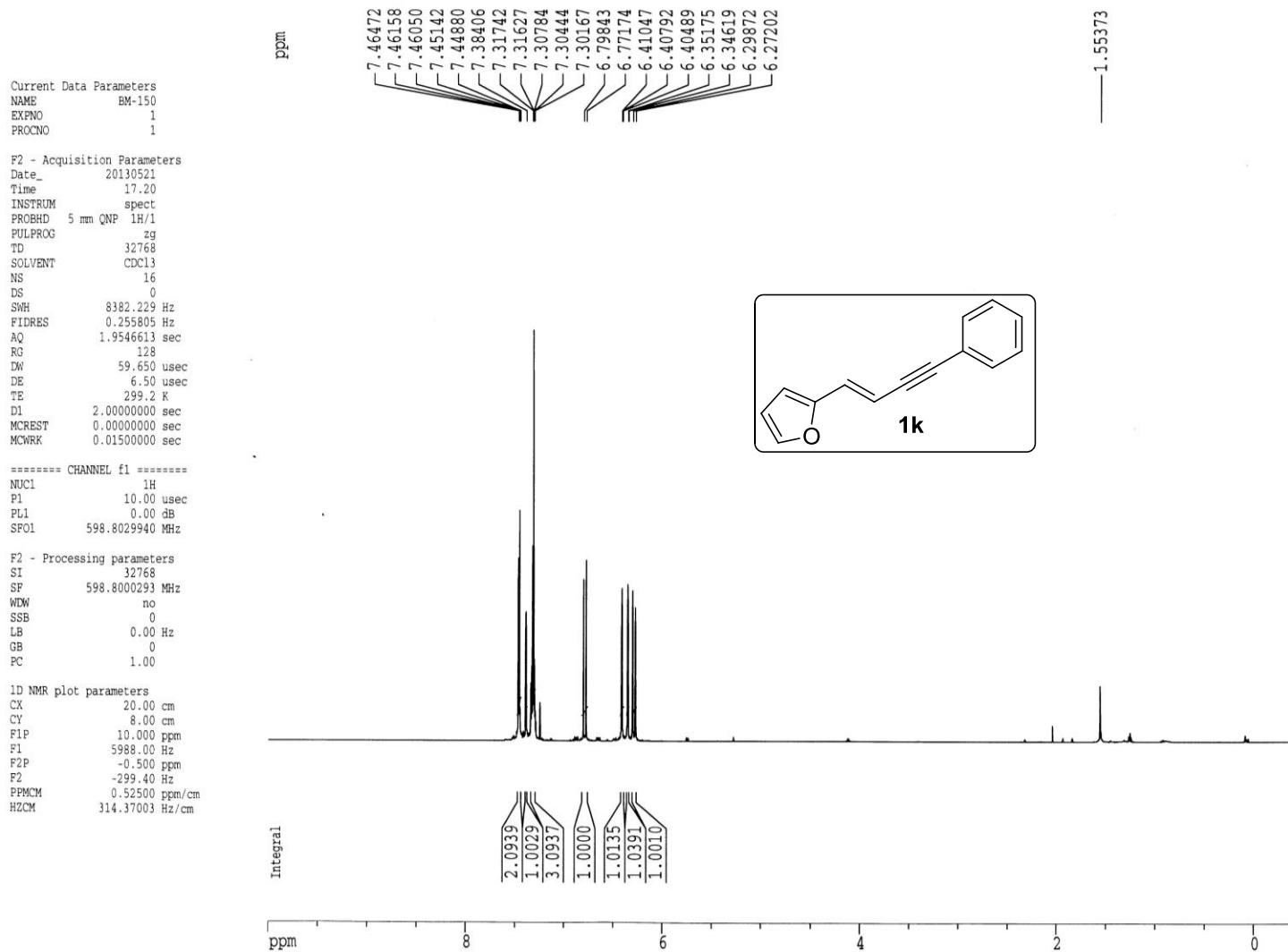
===== CHANNEL f1 =====
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SF01 598.7029935 MHz

F2 - Processing parameters
SI 32768
SF 598.7000261 MHz
WDW no
SSB 0
LB 0.00 Hz
GB 0
PC 1.00

1D NMR plot parameters
CX 20.00 cm
CY 5.00 cm
F1P 10.000 ppm
F1 5987.00 Hz
F2P -0.500 ppm
F2 -299.35 Hz
PPMOM 0.52500 ppm/cm
HZCM 314.31750 Hz/cm







Current Data Parameters
NAME BM-150
EXPNO 2
PROCNO 1

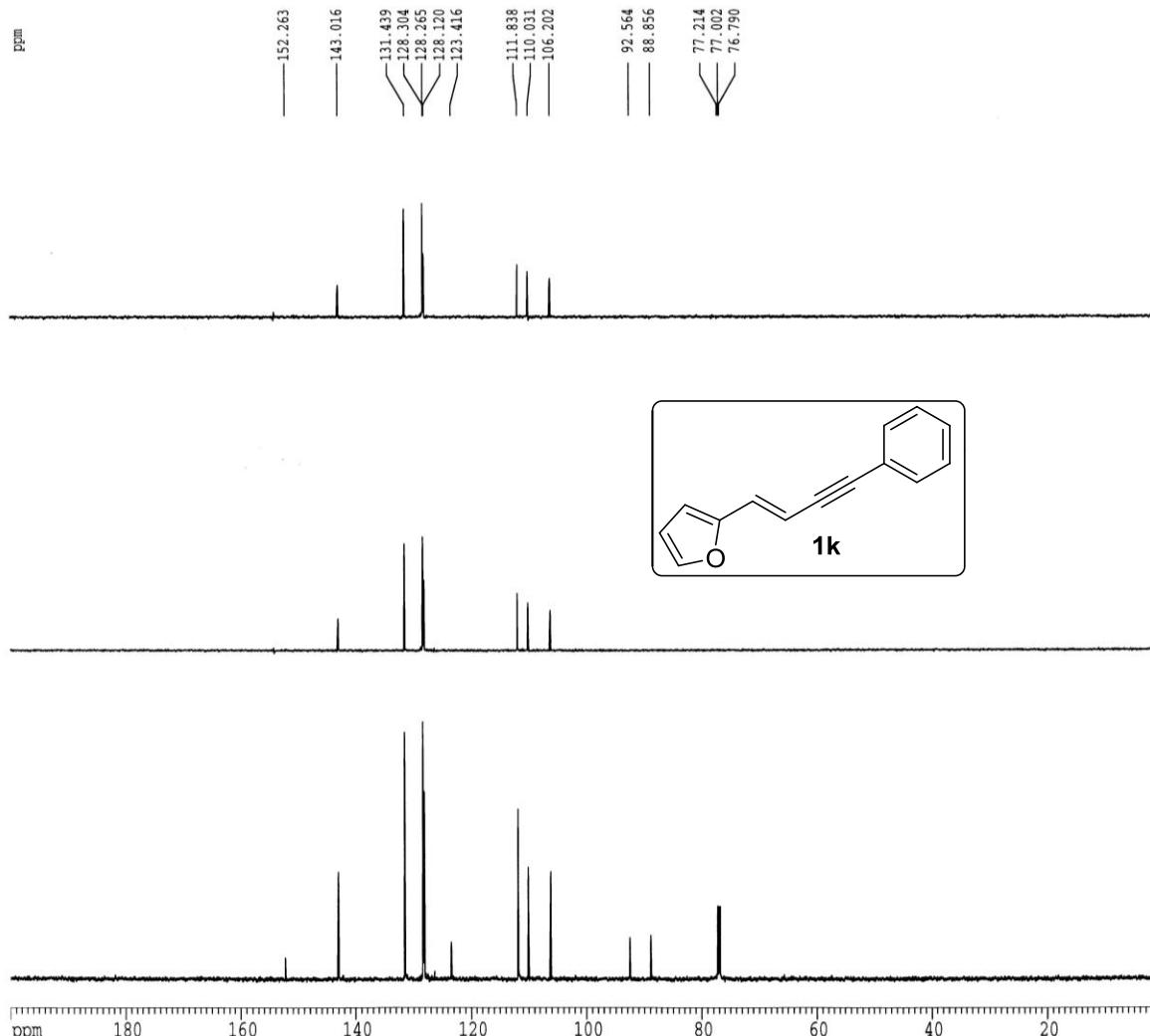
F2 - Acquisition Parameters
Date_ 20130521
Time 17.20
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT CDCl3
NS 51
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 299.2 K
D1 3.5000000 sec
d11 0.0300000 sec
DELTA 3.40000010 sec
MCREST 0.00000000 sec
MCWRK 0.01500000 sec

===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SFO1 150.5849425 MHz

===== CHANNEL f2 =====
CPDPG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SFO2 598.8029940 MHz

F2 - Processing parameters
SI 65536
SF 150.5683897 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 1.00

1D NMR plot parameters
CX 20.00 cm
CY 4.50 cm
F1P 200.000 ppm
F1 30113.68 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPCM 10.00000 ppm/cm
HZCM 1505.68384 Hz/cm



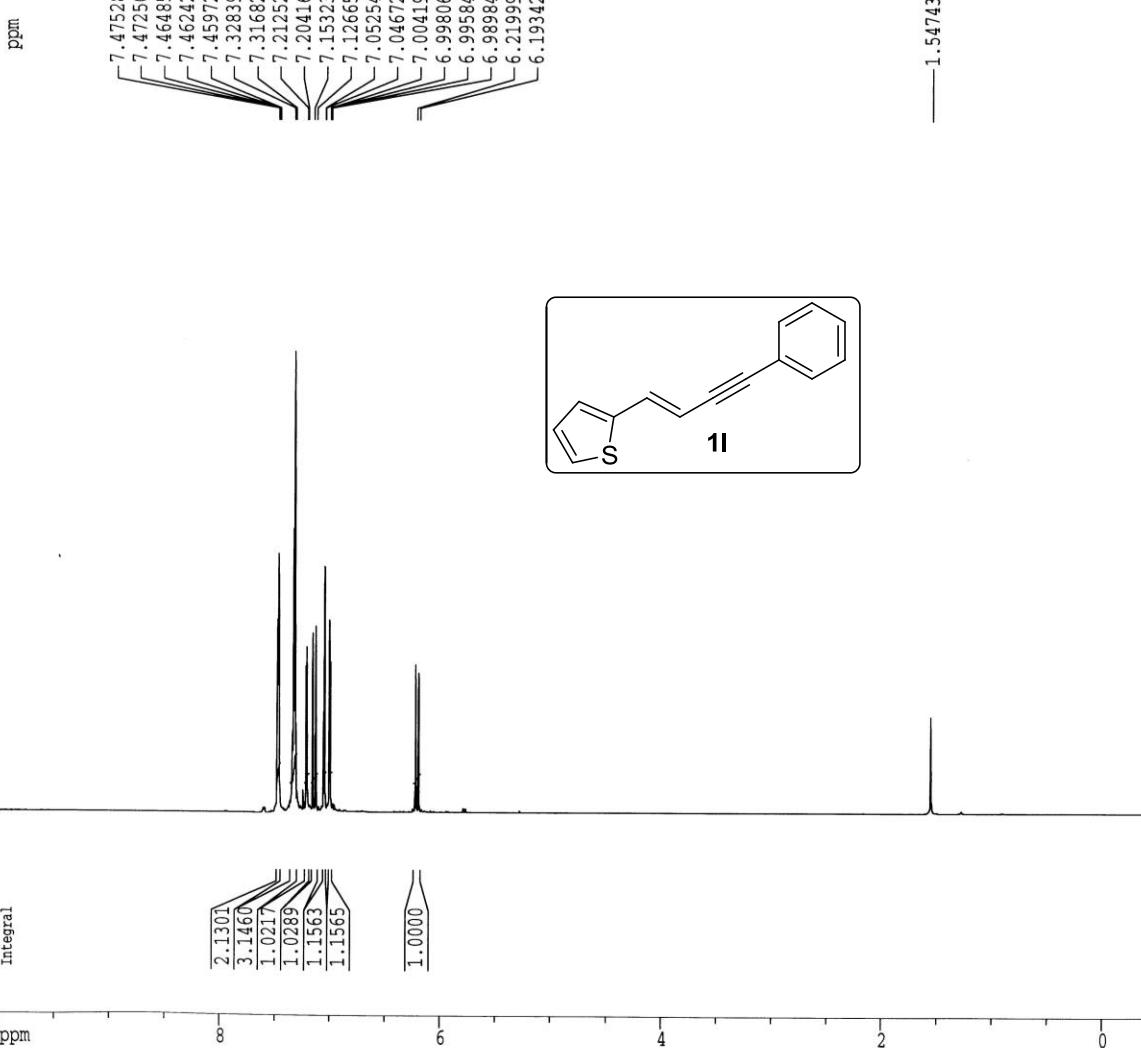
Current Data Parameters
NAME BM-158
EXPNO 1
PROCNO 1

F2 - Acquisition Parameters
Date 20130526
Time 20.29
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 32768
SOLVENT CDCl3
NS 16
DS 0
SWH 8382.229 Hz
FIDRES 0.255805 Hz
AQ 1.9546613 sec
RG 128
DW 59.650 usec
DE 6.50 usec
TE 300.0 K
D1 2.0000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

===== CHANNEL f1 =====
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SFO1 598.8029940 MHz

F2 - Processing parameters
SI 32768
SF 598.8000285 MHz
WDW no
SSB 0
LB 0.00 Hz
GB 0
PC 2.00

1D NMR plot parameters
CX 20.00 cm
CY 8.00 cm
F1P 10.000 ppm
F1 5988.00 Hz
F2P -0.500 ppm
F2 -299.40 Hz
PPCM 0.52500 ppm/cm
HZCM 314.37003 Hz/cm



Current Data Parameters
 NAME BM-158
 EXPNO 2
 PROCN0 1

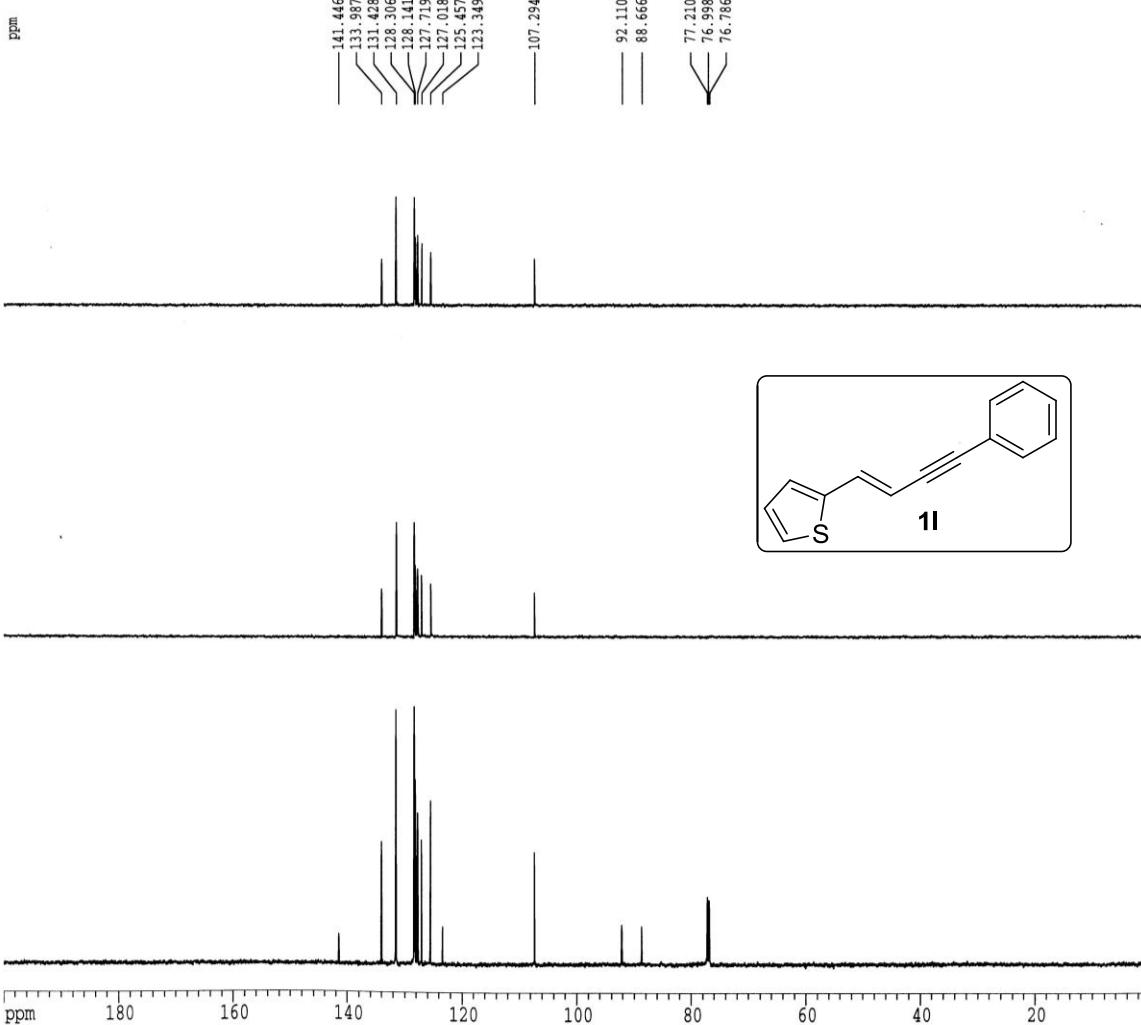
F2 - Acquisition Parameters
 Date_ 20130526
 Time 20.30
 INSTRUM spect
 PROBHD 5 mm QNP 1H/1
 PULPROG zpgd
 TD 32768
 SOLVENT CDCl3
 NS 35
 DS 0
 SWH 45045.047 Hz
 FIDRES 1.374666 Hz
 AQ 0.3637748 sec
 RG 2048
 DW 11.100 usec
 DE 6.50 usec
 TE 300.2 K
 D1 3.5000000 sec
 d11 0.0300000 sec
 DELTA 3.40000010 sec
 MCREST 0.00000000 sec
 MCWRK 0.01500000 sec

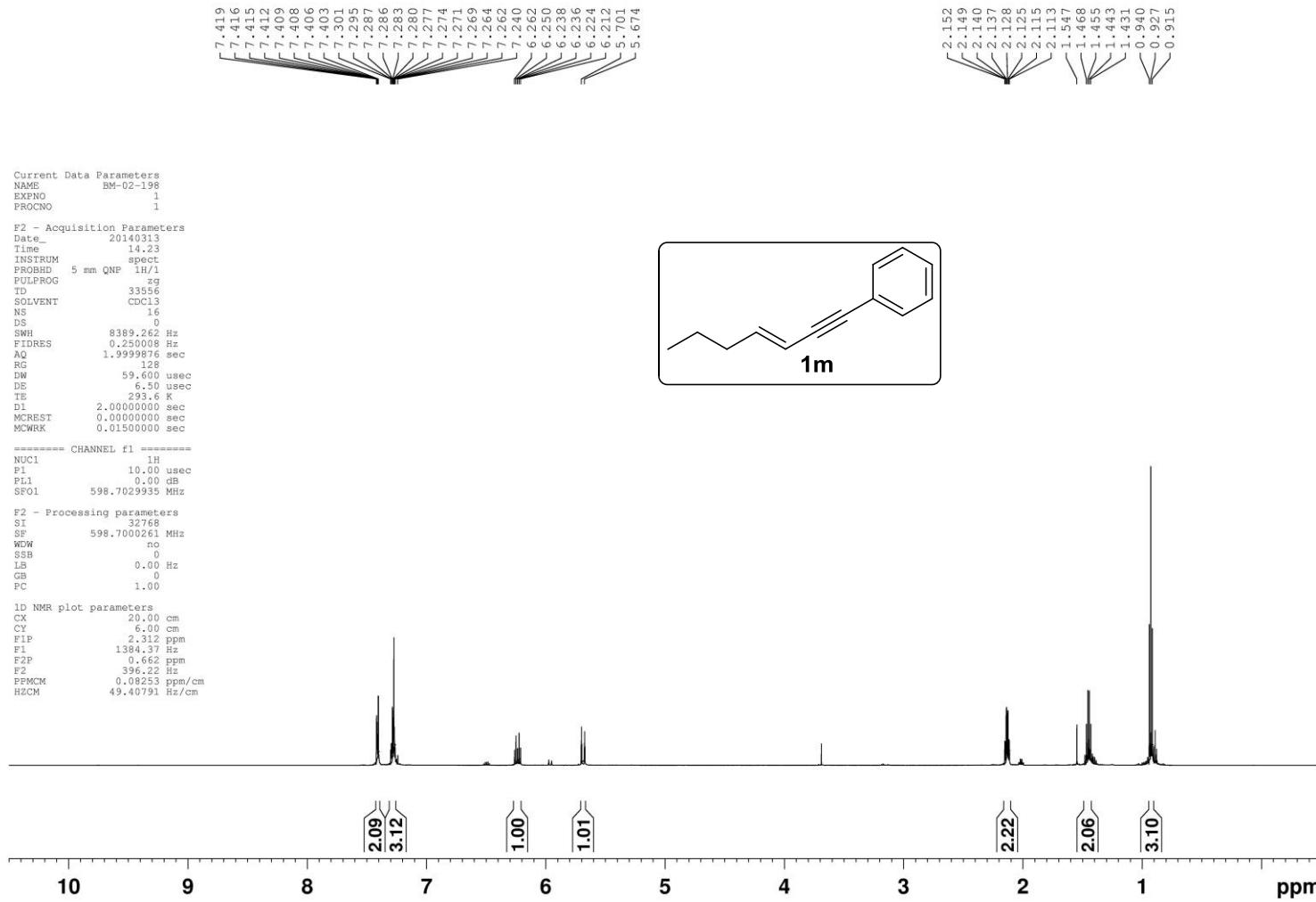
===== CHANNEL f1 =====
 NUC1 13C
 P1 4.80 usec
 PL1 0.00 dB
 SF01 150.5849425 MHz

===== CHANNEL f2 =====
 CPDPRG2 waltz16
 NUC2 1H
 PCPD2 92.00 usec
 PL2 120.00 dB
 PL12 9.00 dB
 PL13 14.00 dB
 SF02 598.8029940 MHz

F2 - Processing parameters
 SI 65536
 SF 150.5683945 MHz
 WDW EM
 SSB 0
 LB 3.00 Hz
 GB 0
 PC 1.00

1D NMR plot parameters
 CX 20.00 cm
 CY 4.50 cm
 F1P 200.000 ppm
 F1 30113.68 Hz
 F2P 0.000 ppm
 F2 0.00 Hz
 PPMCM 10.00000 ppm/cm
 HZCM 1505.68384 Hz/cm





Current Data Parameters
NAME BM-02-198
EXPNO 2
PROCNO 1

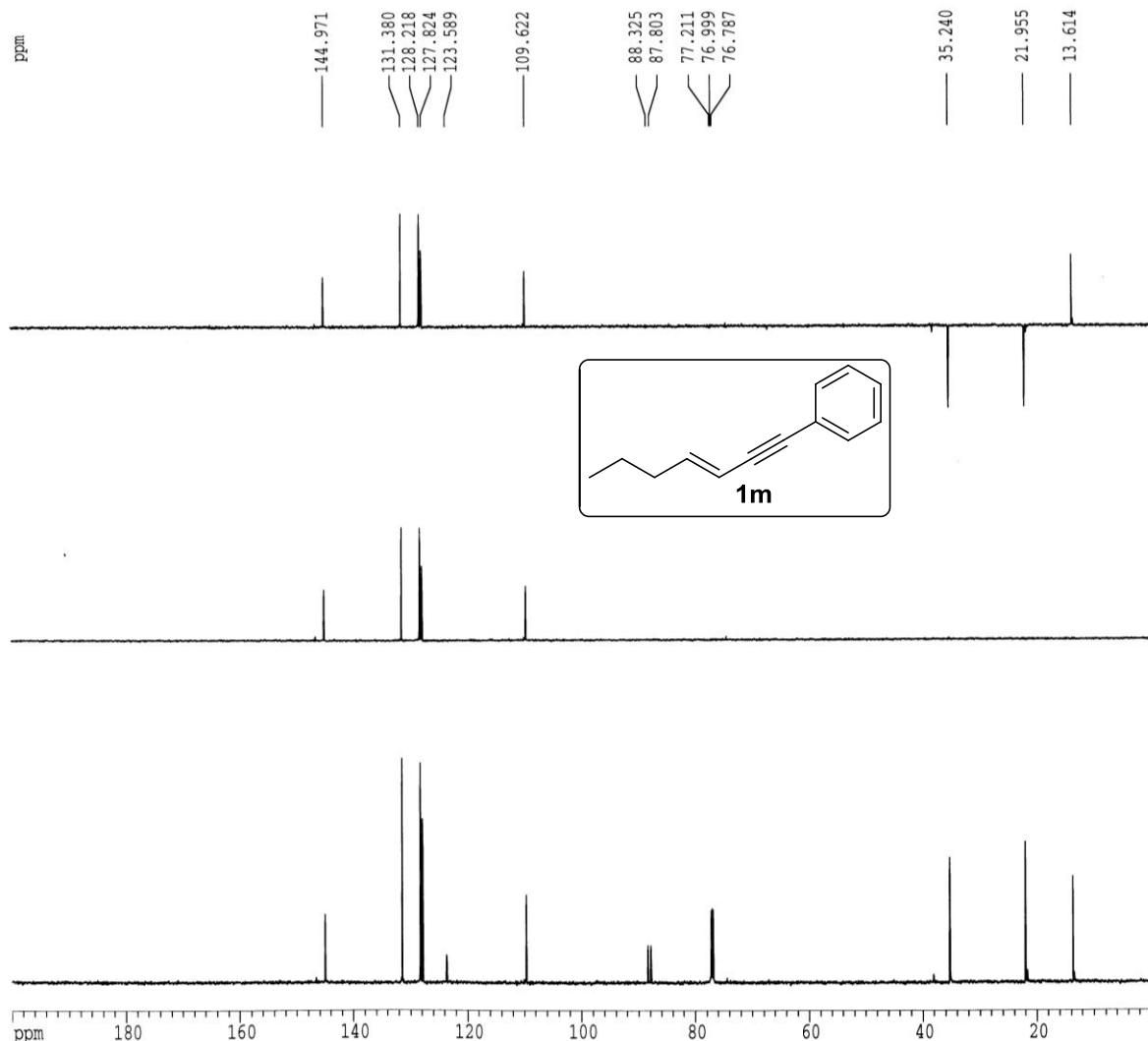
F2 - Acquisition Parameters
Date 20140313
Time 14.25
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT CDCl3
NS 100
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 294.3 K
D1 3.5000000 sec
d1 0.0300000 sec
DELTA 3.40000010 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

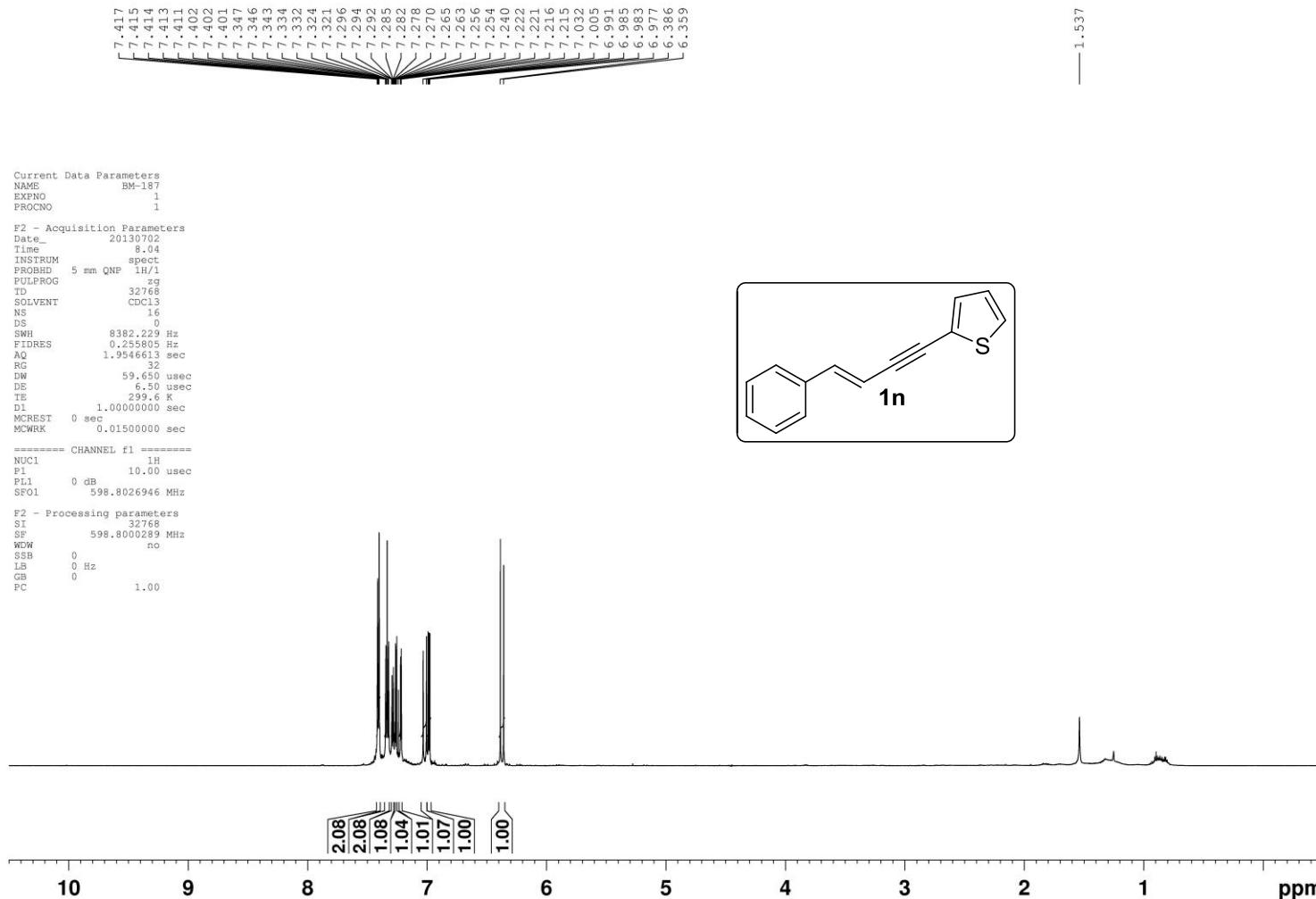
===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SF01 150.5597948 MHz

===== CHANNEL f2 =====
CPDPFRG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SF02 598.7029935 MHz

F2 - Processing parameters
S1 65536
SF 150.5432431 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 0.50

1D NMR plot parameters
CX 20.00 cm
CY 4.00 cm
F1P 200.000 ppm
F1 30108.65 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPMCM 10.00000 ppm/cm
HZCM 1505.43237 Hz/cm





Current Data Parameters
NAME BM-187
EXPNO 2
PROCNO 1

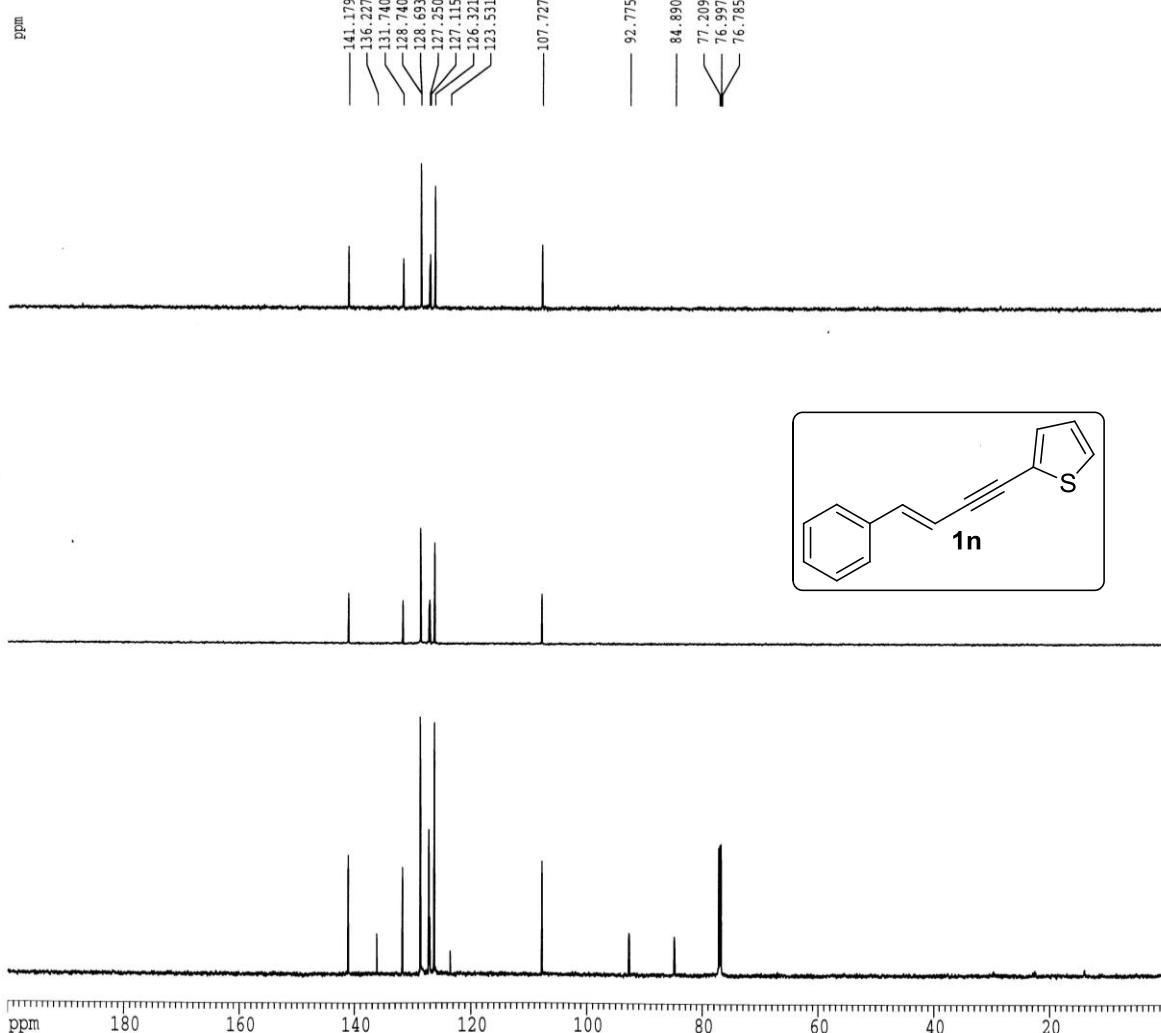
F2 - Acquisition Parameters
Date_ 20130701
Time 17:17
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT CDCl3
NS 205
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 300.6 K
D1 3.5000000 sec
d11 0.0300000 sec
DELTA 3.4000010 sec
MCREST 0.0000000 sec
MCVRK 0.0150000 sec

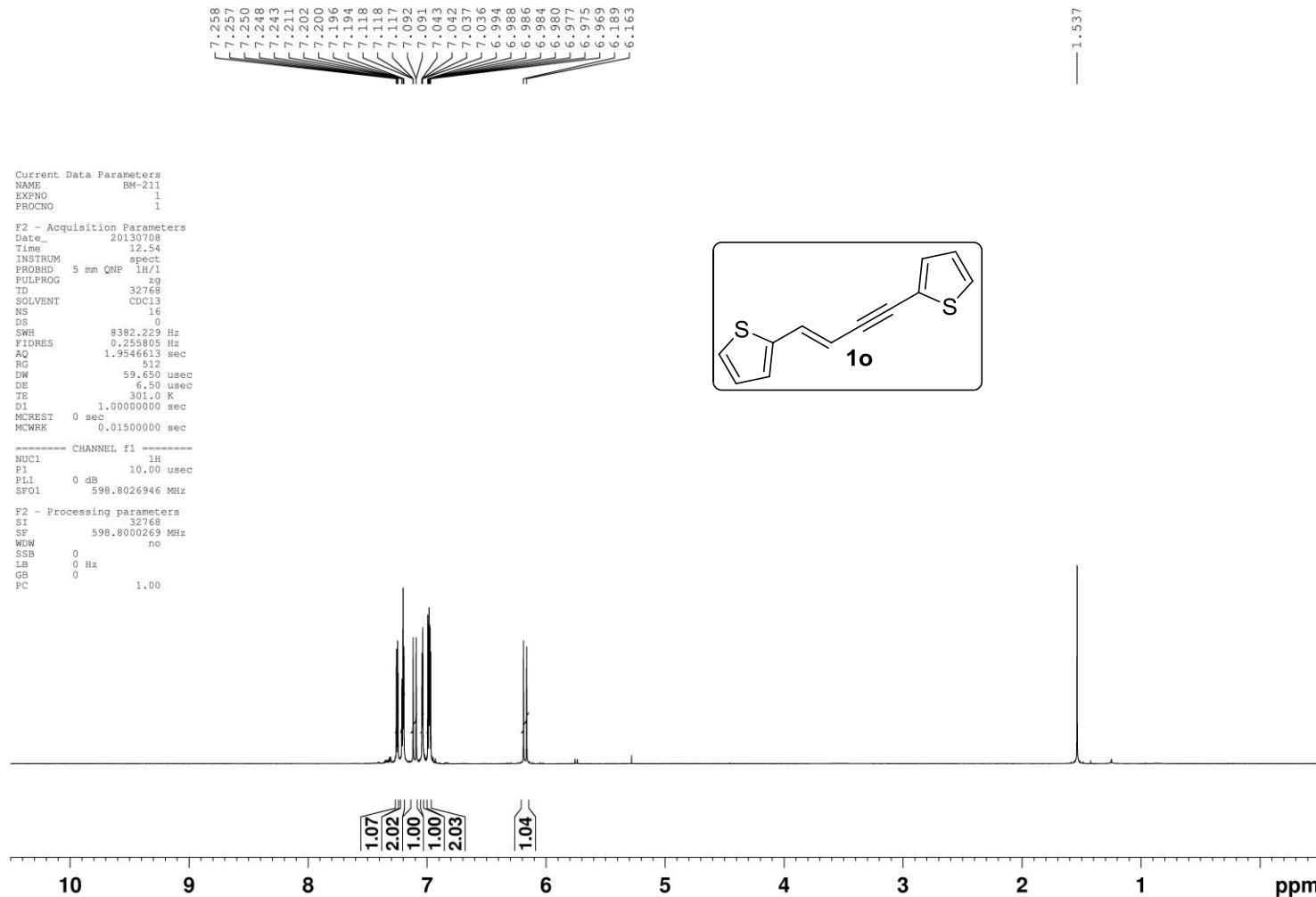
===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SFO1 150.5849425 MHz

===== CHANNEL f2 =====
CPDPRG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SFO2 598.8029940 MHz

F2 - Processing parameters
SI 65536
SF 150.5683869 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 0.50

1D NMR plot parameters
CX 20.00 cm
CY 4.50 cm
PIP 200.000 ppm
F1 30113.68 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPMCM 10.00000 ppm/cm
HZCM 1505.68384 Hz/cm





Current Data Parameters
NAME BM-211
EXPNO 2
PROCNO 1

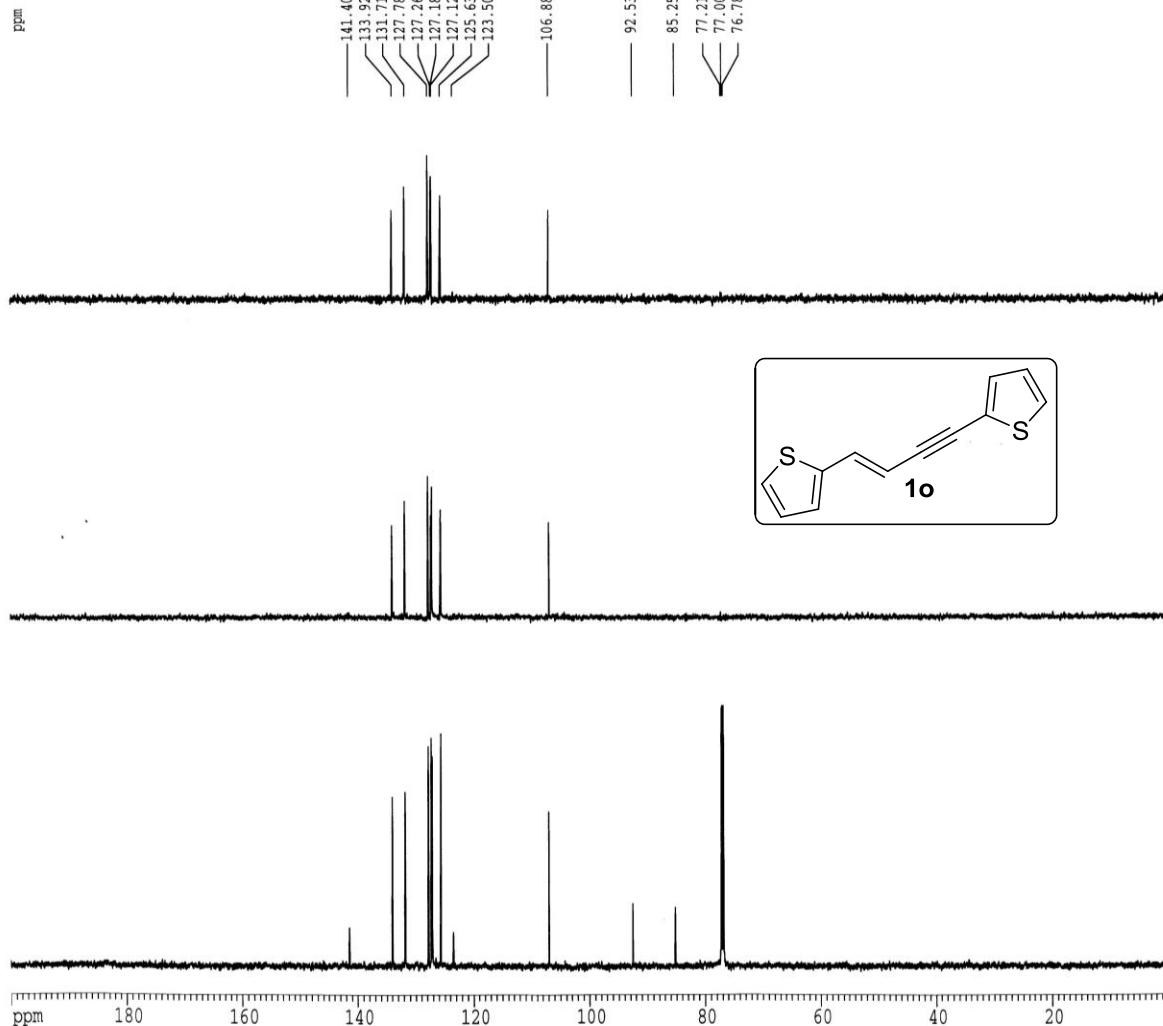
F2 - Acquisition Parameters
Date_ 20130707
Time 16.08
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT CDCl3
NS 325
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 300.6 K
D1 3.5000000 sec
d11 0.0300000 sec
DELTA 3.40000010 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SF01 150.5849425 MHz

===== CHANNEL f2 =====
CPDPRG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SF02 598.8029940 MHz

F2 - Processing parameters
SI 65536
SF 150.5683848 MHz
WDW EM
SSB 0
LB 3.00 Hz
GE 0
PC 0.50

1D NMR plot parameters
CX 20.00 cm
CY 4.50 cm
F1P 200.000 ppm
F1 30113.68 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPMCM 10.00000 ppm/cm
HZCM 1505.68384 Hz/cm



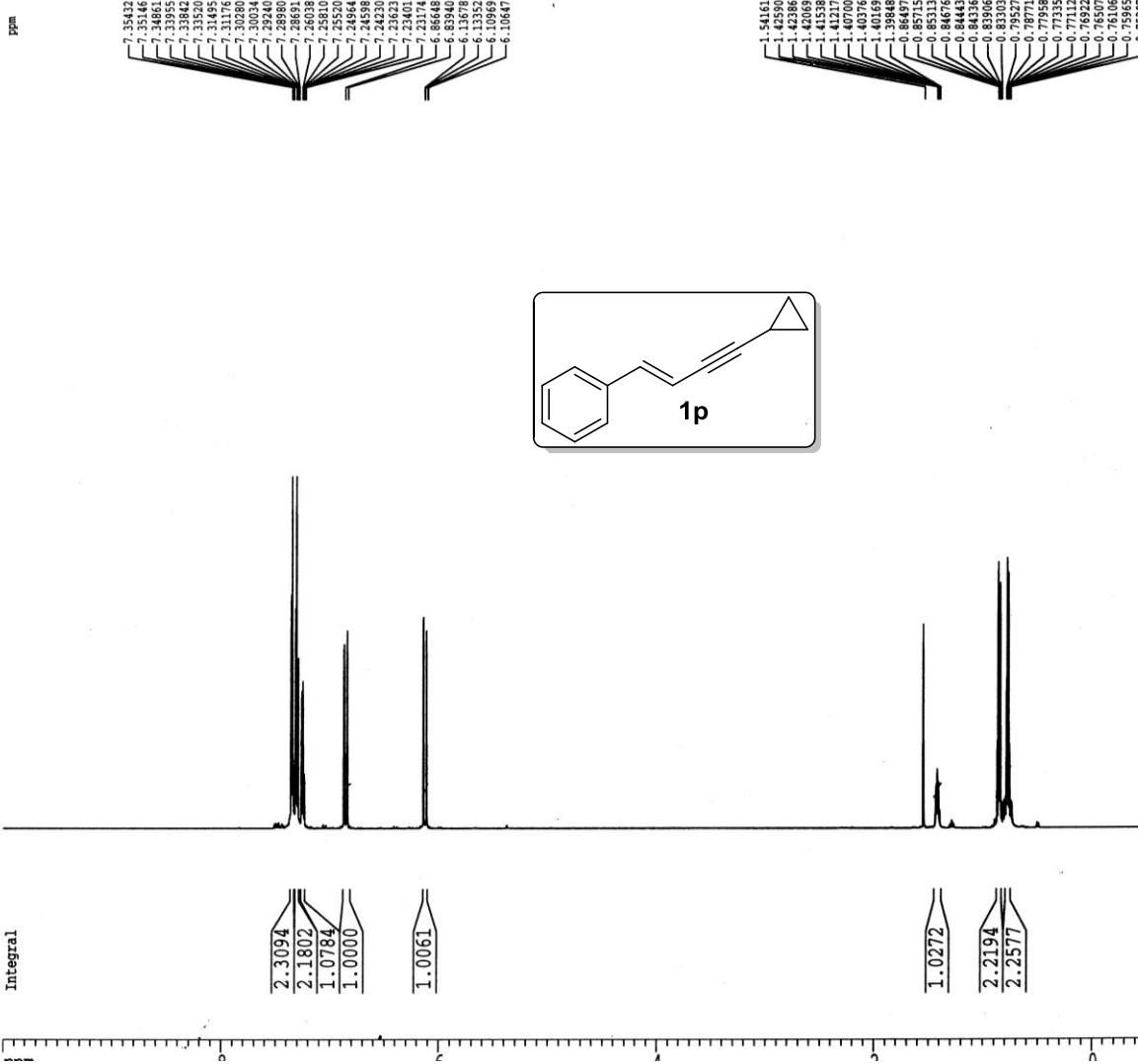
Current Data Parameters
 NAME BM-03-08
 EXPNO 1
 PROCNO 1

F2 - Acquisition Parameters
 Date 20140521
 Time 20.40
 INSTRUM spect
 PROBHD 5 mm QNP 1H/1
 PULPROG zg
 TD 32768
 SOLVENT CDCl3
 NS 16
 DS 0
 SWH 8389.262 Hz
 FIDRES 0.256020 Hz
 AQ 1.9530228 sec
 RG 256
 DW 59.600 usec
 DE 6.50 usec
 TE 298.5 K
 D1 2.0000000 sec
 MCREST 0.0000000 sec
 MCWRK 0.01500000 sec

===== CHANNEL f1 =====
 NUC1 1H
 P1 10.00 usec
 PL1 3.00 dB
 SFO1 598.7029935 MHz

F2 - Processing parameters
 SI 32768
 SF 598.7000173 MHz
 WDW no
 SSB 0
 LB 0.00 Hz
 GB 0
 PC 1.00

1D NMR plot parameters
 CX 20.00 cm
 CY 6.00 cm
 F1P 10.000 ppm
 F1 5987.00 Hz
 F2P -0.500 ppm
 F2 -299.35 Hz
 PPMCM 0.52500 ppm/cm
 HZCM 314.31750 Hz/cm



Current Data Parameters
NAME EM-03-08
EXPNO 2
PROCNO 1

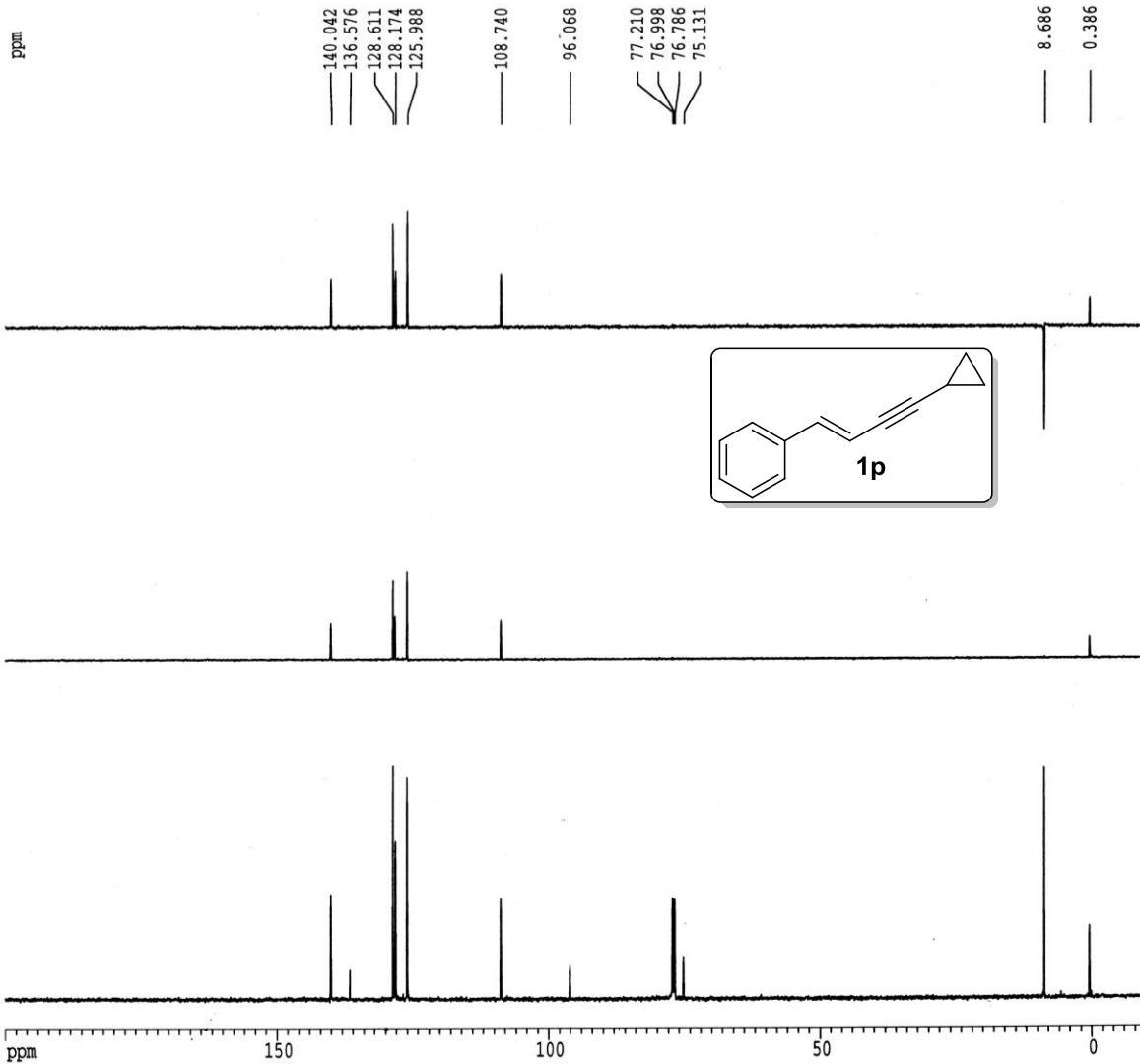
F2 - Acquisition Parameters
Date_ 20140521
Time 20.51
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT CDCl3
NS 166
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 299.7 K
D1 3.5000000 sec
d11 0.03000000 sec
DELTA 3.40000010 sec
MCREST 0.0000000 sec
MCWRK 0.01500000 sec

===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SF01 150.5597948 MHz

===== CHANNEL f2 =====
CPDPRG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SF02 598.7029935 MHz

F2 - Processing parameters
SI 65536
SF 150.5432397 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 0.50

1D NMR plot parameters
CX 20.00 cm
CY 4.00 cm
F1P 200.000 ppm
F1 30108.65 Hz
F2P -10.000 ppm
F2 -1505.43 Hz
PPMCM 10.50000 ppm/cm
HZCM 1580.70410 Hz/cm



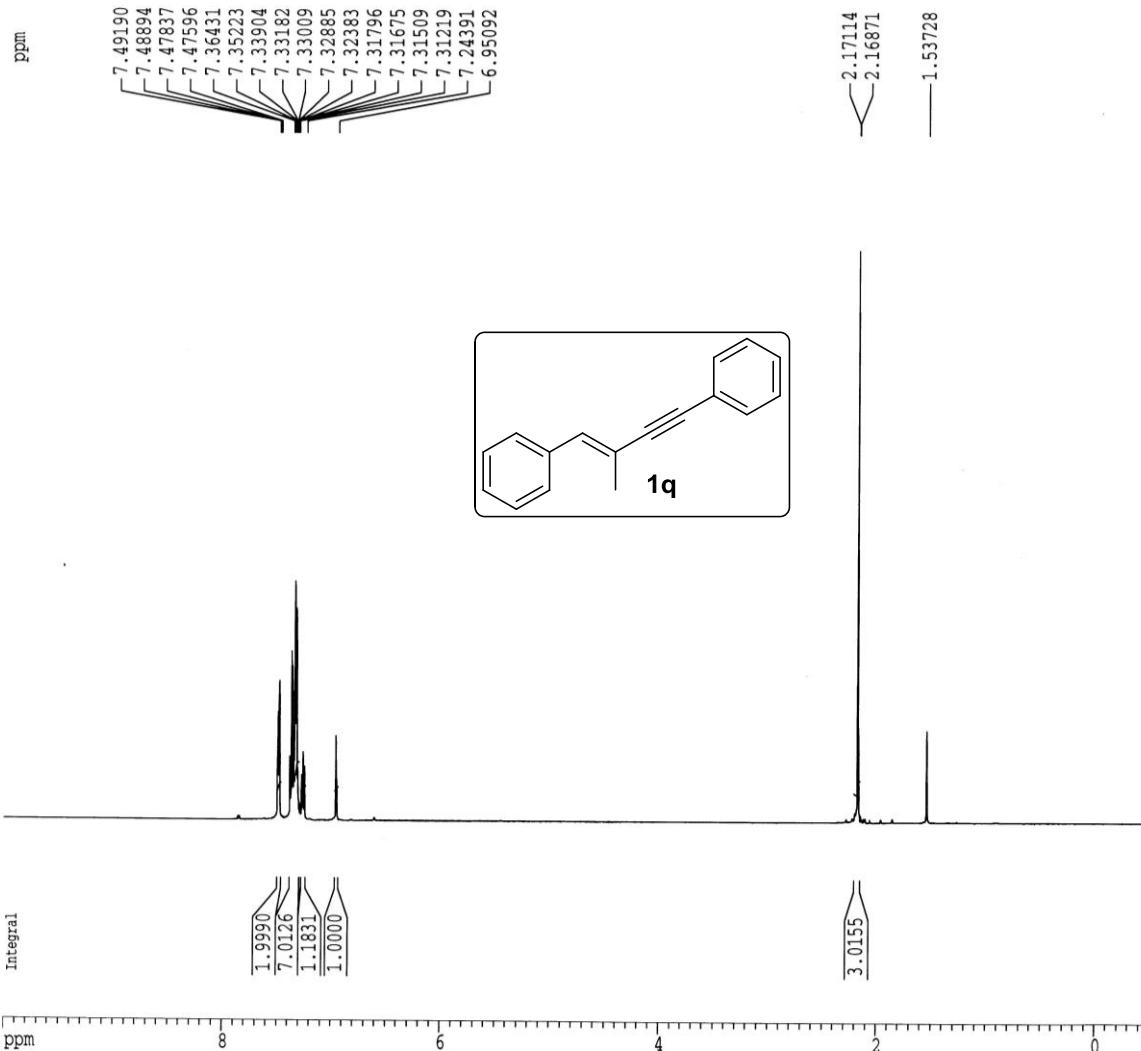
Current Data Parameters
NAME BM-191
EXPNO 1
PROCNO 1

F2 - Acquisition Parameters
Date_ 20130701
Time 17.46
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 32768
SOLVENT CDCl3
NS 16
DS 0
SWH 8382.229 Hz
FIDRES 0.255805 Hz
AQ 1.9546613 sec
RG 32
DW 59.650 usec
DE 6.50 usec
TE 300.2 K
D1 1.0000000 sec
MCREST 0.0000000 sec
MCWKR 0.0150000 sec

===== CHANNEL f1 =====
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SF01 598.8026946 MHz

F2 - Processing parameters
SI 32768
SF 598.8000269 MHz
WDW no
SSB 0
LB 0.00 Hz
GB 0
PC 1.00

1D NMR plot parameters
CX 20.00 cm
CY 10.00 cm
F1P 10.000 ppm
F1 5988.00 Hz
F2P -0.500 ppm
F2 -299.40 Hz
PPCM 0.52500 ppm/cm
HZCM 314.37003 Hz/cm



Current Data Parameters
NAME BM-191
EXPNO 2
PROCNO 1

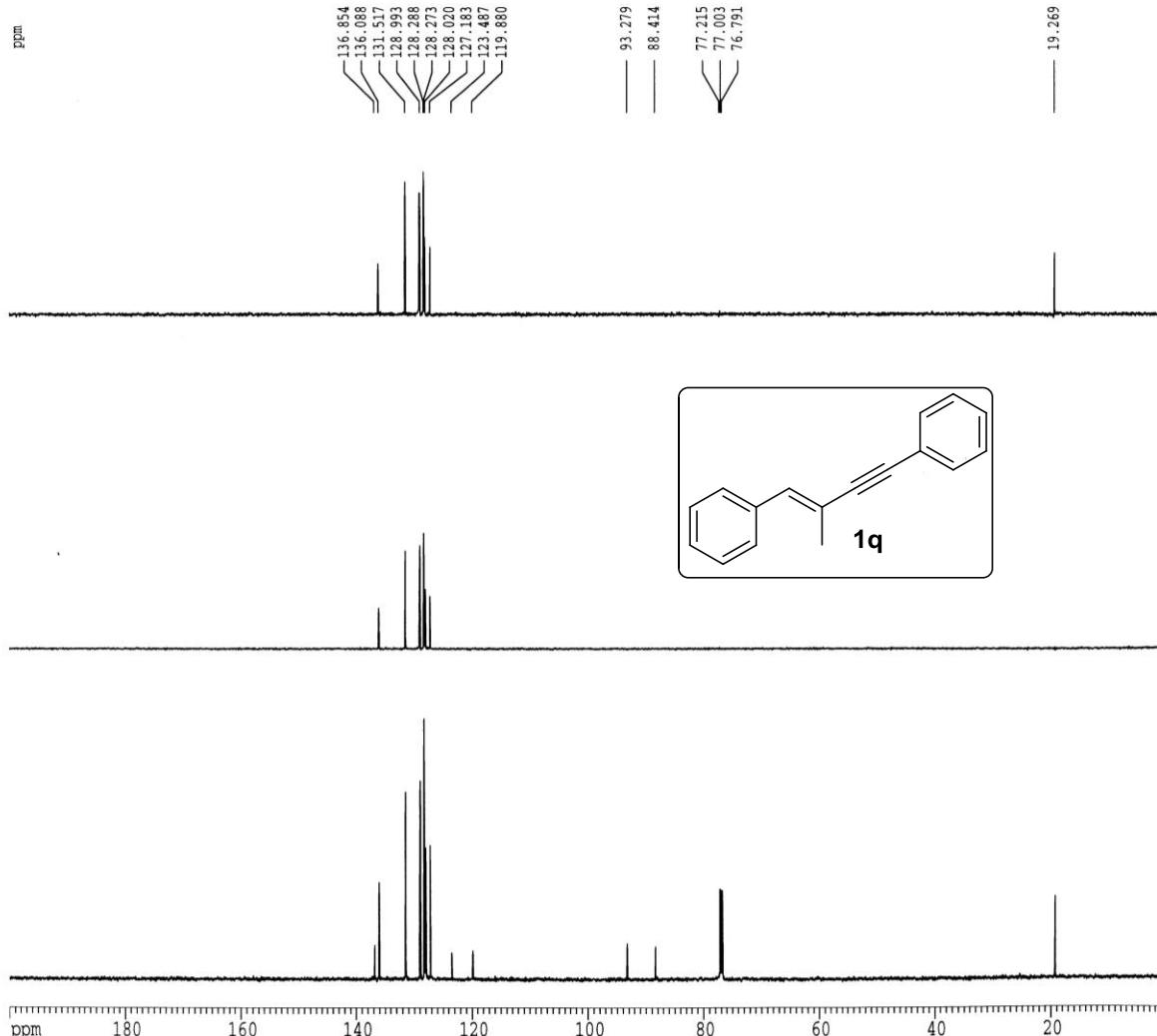
F2 - Acquisition Parameters
Date 20130701
Time 17:53
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT CDCl3
NS 100
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 301.0 K
DI 3.5000000 sec
d1 0.0300000 sec
DELT1 3.4000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

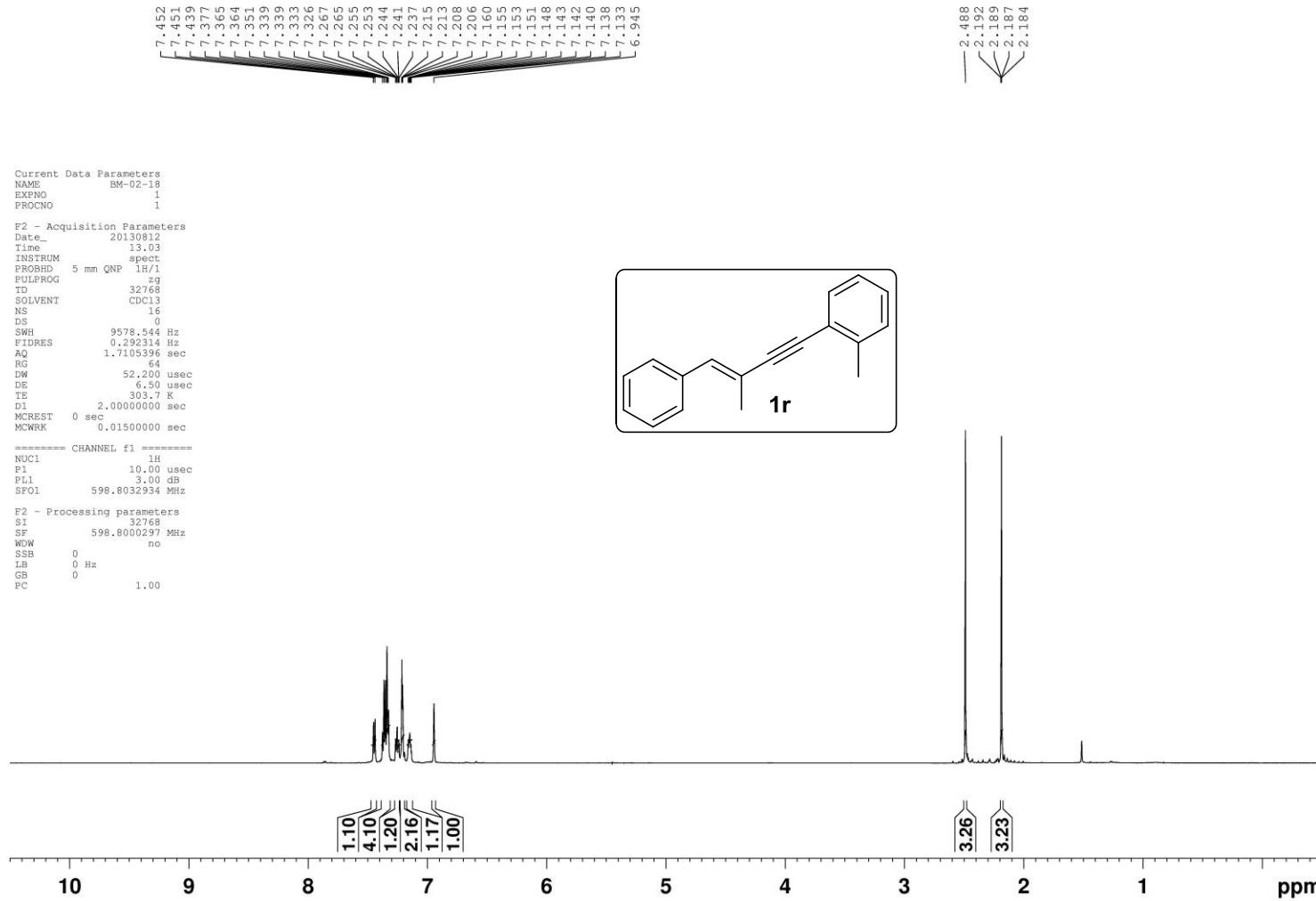
===== CHANNEL f1 =====
NUC1 13C
PI 4.80 usec
PL1 0.00 dB
SFO1 150.5849425 MHz

===== CHANNEL f2 =====
CPDPG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SFO2 598.8029940 MHz

F2 - Processing parameters
SI 65536
SF 150.5681869 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 0.50

1D NMR plot parameters
CX 20.00 cm
CY 4.50 cm
FLP 200.000 ppm
F1 30113.68 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPCM 10.00000 ppm/cm
HZCM 1505.68384 Hz/cm





Current Data Parameters
NAME BM-02-18
EXPNO 2
PROCNO 1

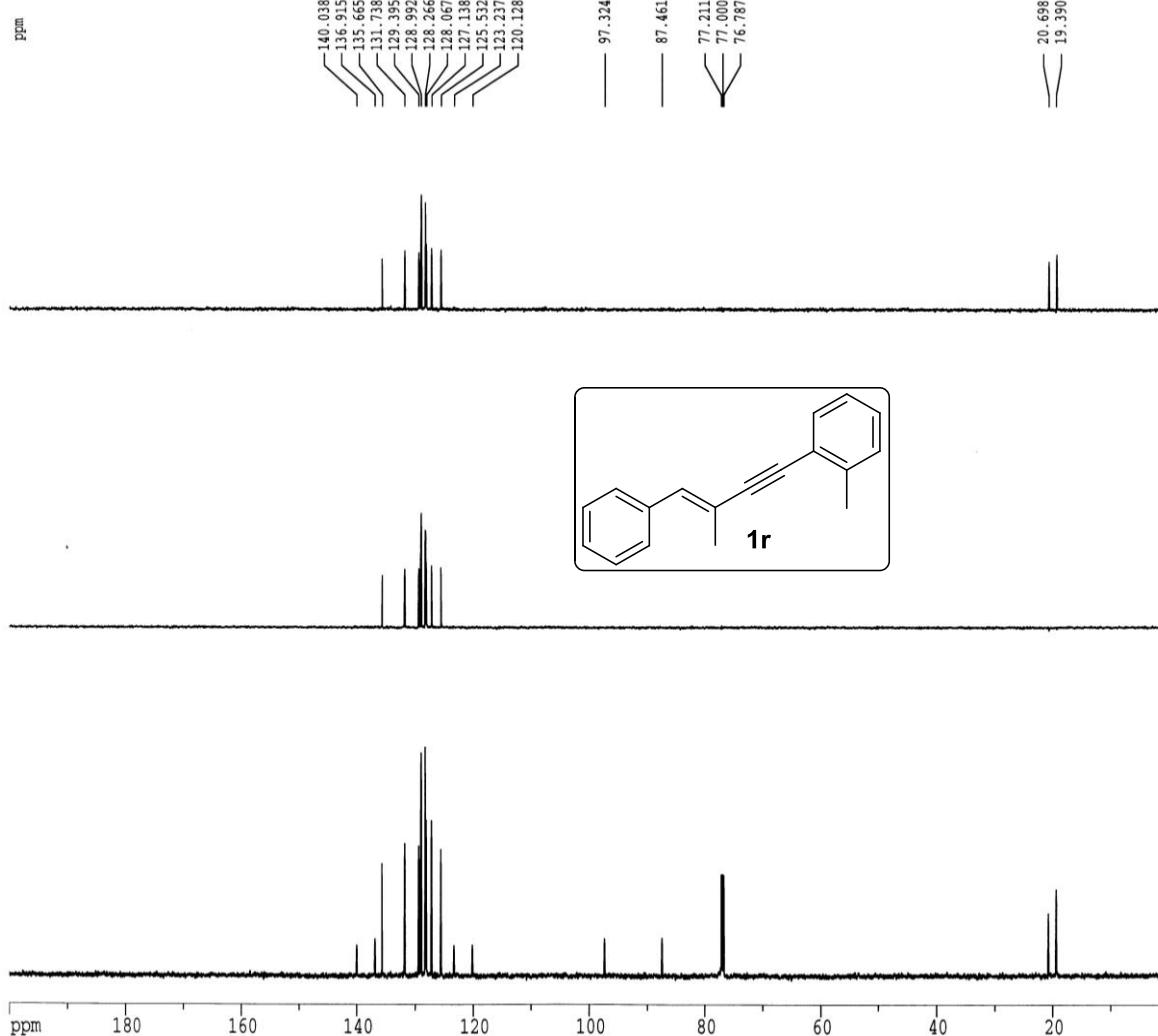
F2 - Acquisition Parameters
Date_ 20130811
Time 22.07
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT CDCl3
NS 66
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 304.5 K
D1 3.5000000 sec
d11 0.0300000 sec
DELTA 3.4000010 sec
MCREST 0.0000000 sec
MCWRK 0.01500000 sec

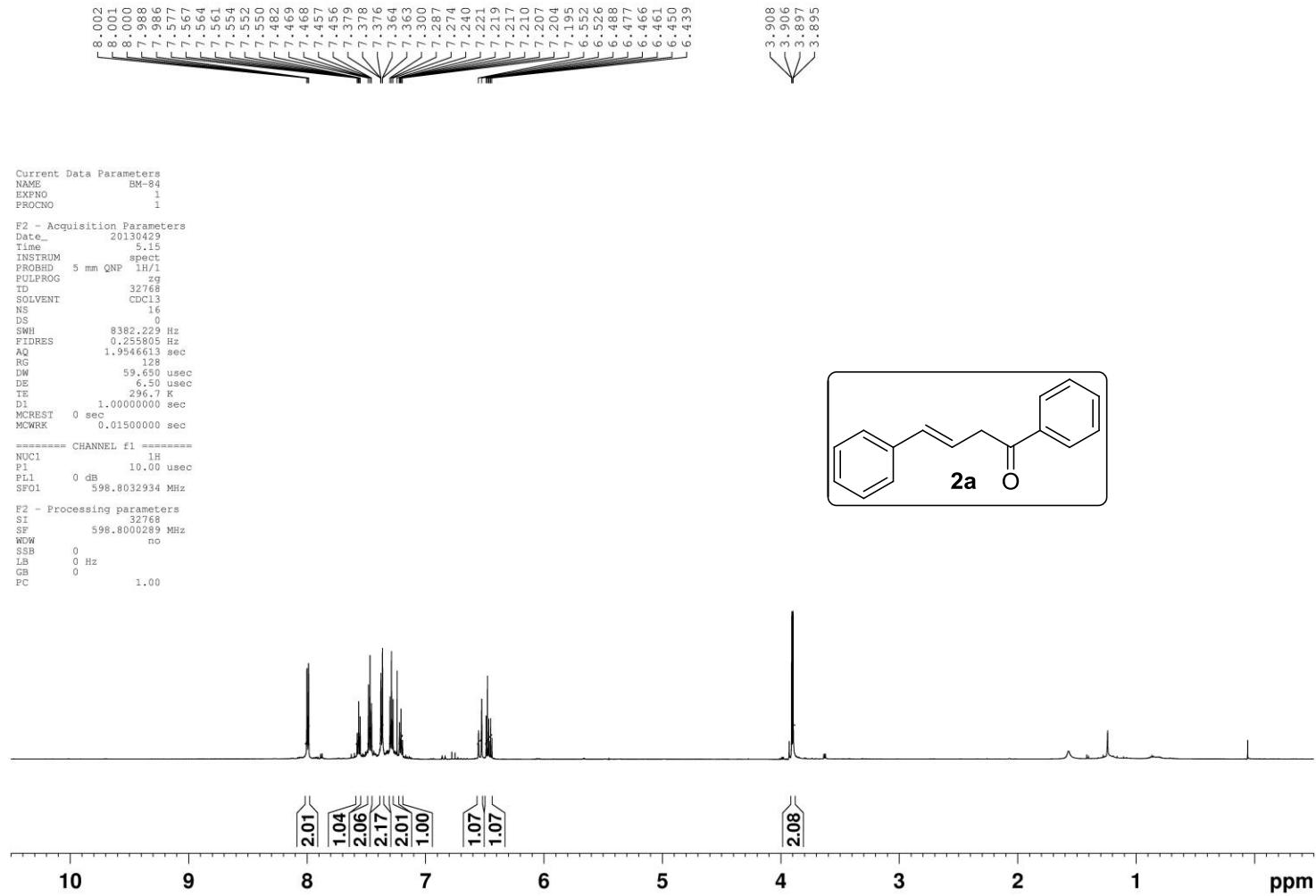
===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SPO1 150.5849425 MHz

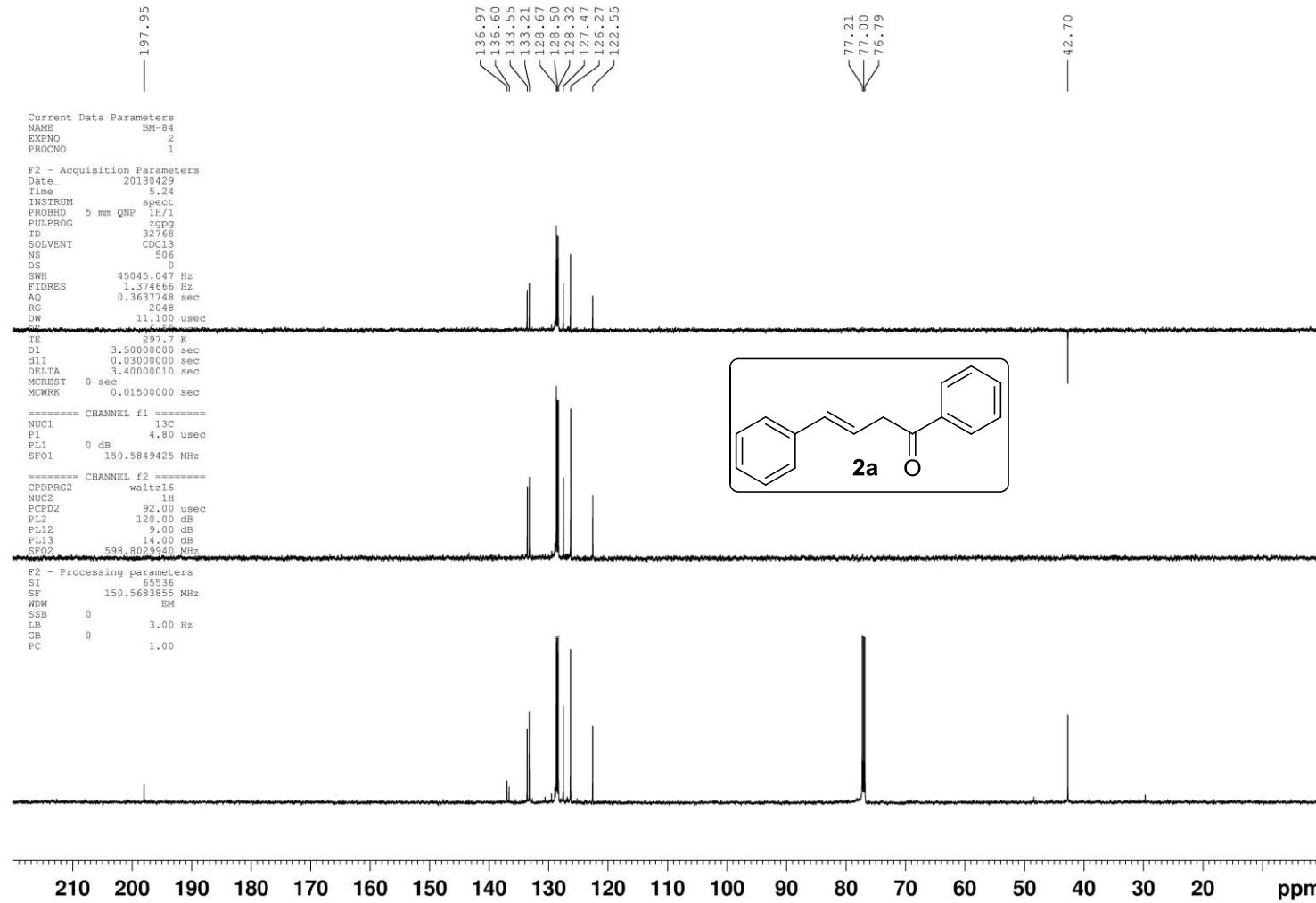
===== CHANNEL f2 =====
CPDPG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SPO2 598.8029940 MHz

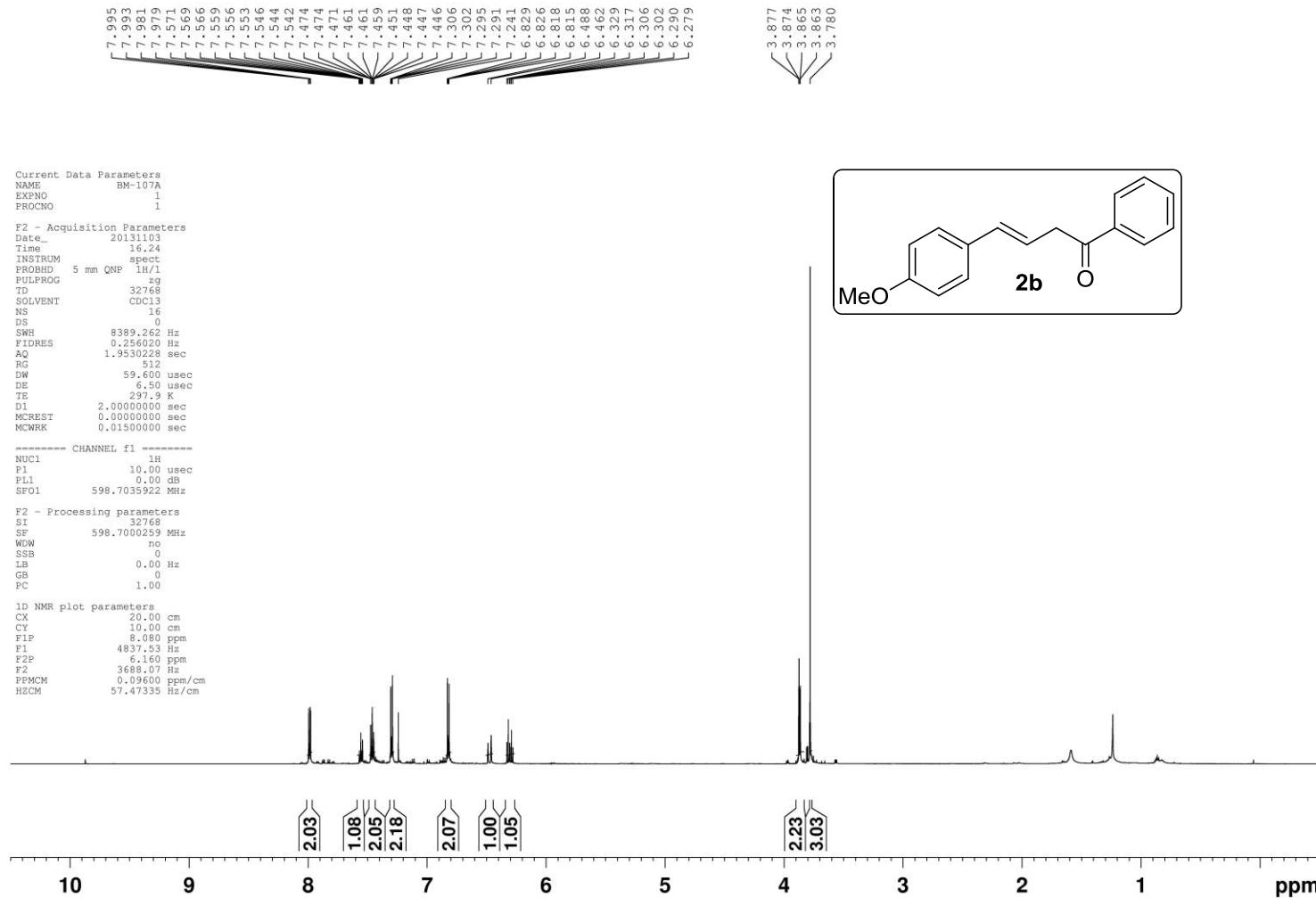
F2 - Processing parameters
SI 65536
SF 150.5683869 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 0.50

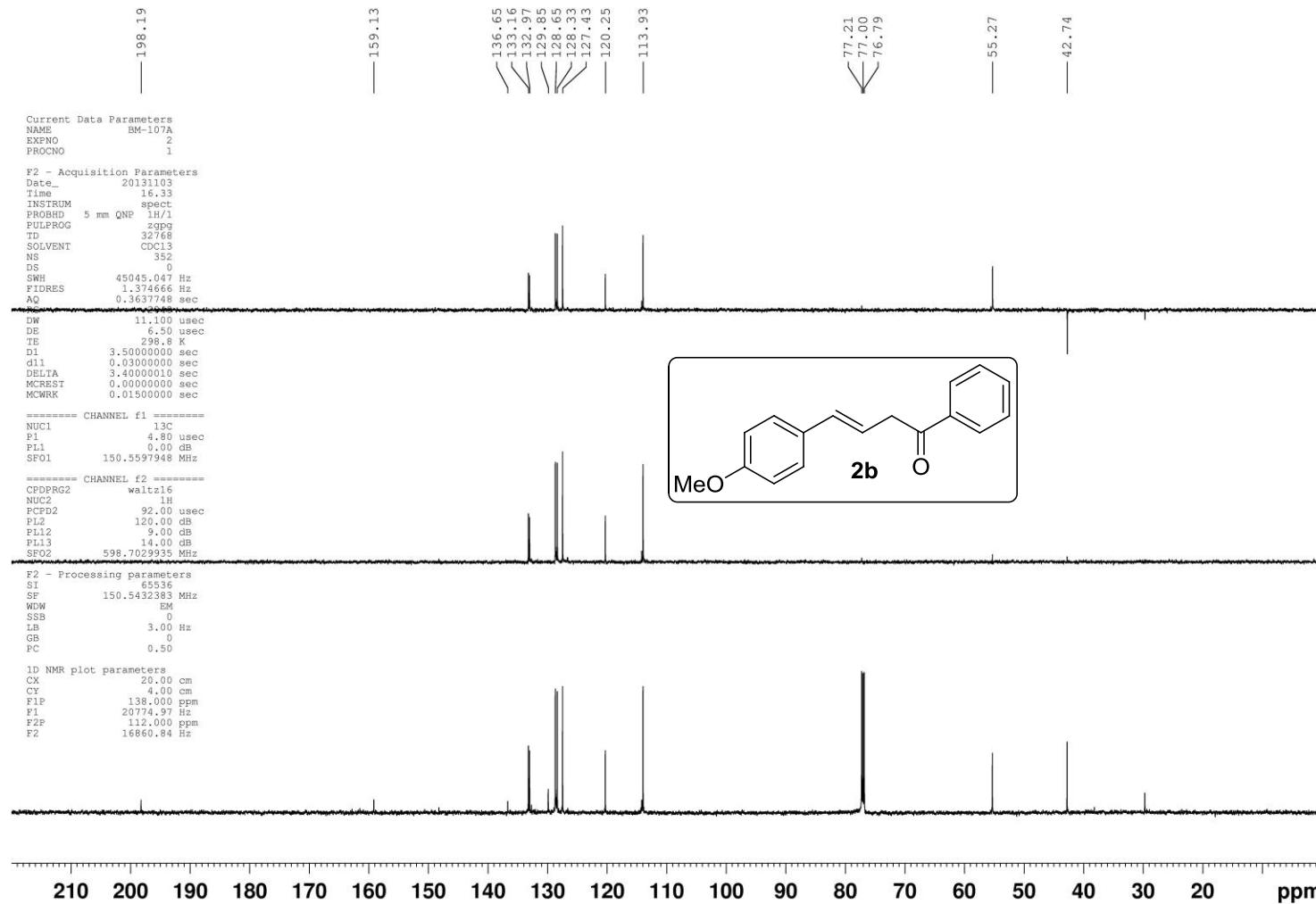
1D NMR plot parameters
CX 20.00 cm
CY 4.00 cm
F1P 200.000 ppm
F1 30113.68 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPCM 10.00000 ppm/cm
HZCM 1505.68384 Hz/cm

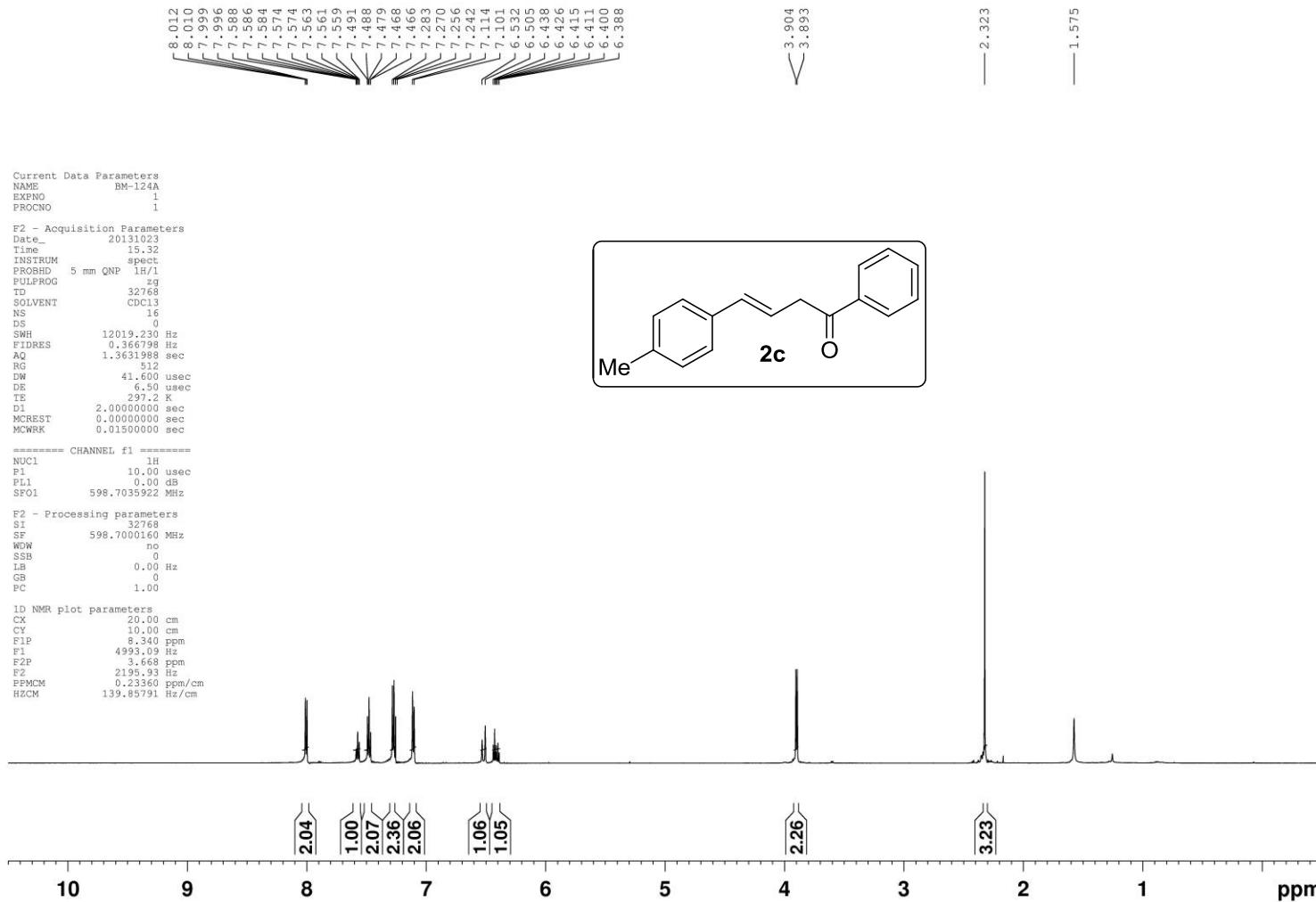












Current Data Parameters
 NAME BM-124A
 EXPNO 2
 PROCNO 1

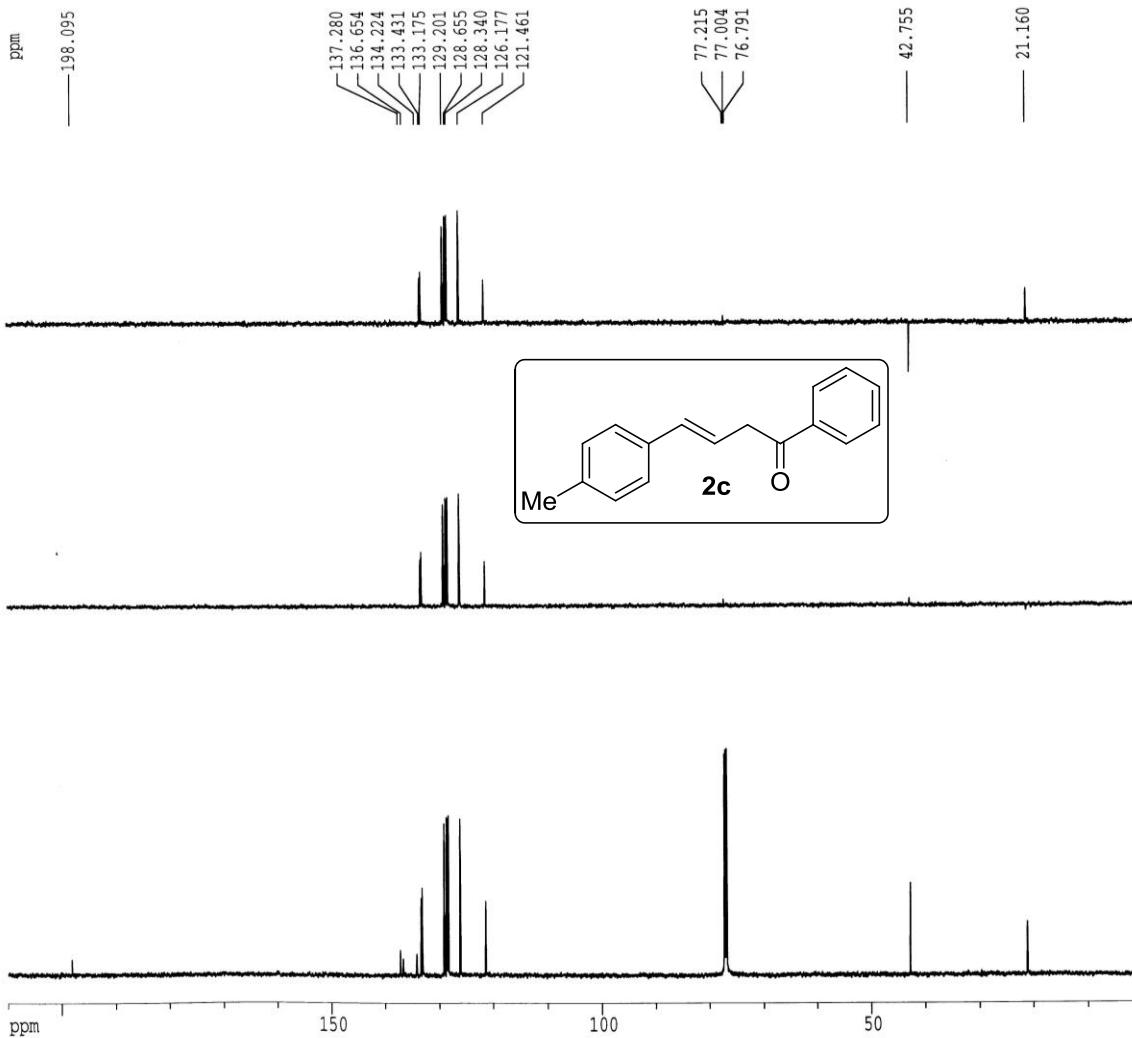
P2 - Acquisition Parameters
 Date_ 20131023
 Time 15.34
 INSTRUM spect
 PROBHD 5 mm QNP 1H/1
 PULPROG zgpg
 TD 32768
 SOLVENT CDCl3
 NS 500
 DS 0
 SWH 45045.047 Hz
 FIDRES 1.374666 Hz
 AQ 0.3637748 sec
 RG 2048
 DW 11.100 usec
 DE 6.50 usec
 TE 297.5 K
 D1 3.5000000 sec
 d11 0.0300000 sec
 DELTA 3.40000010 sec
 MCREST 0.0000000 sec
 MWKR 0.01500000 sec

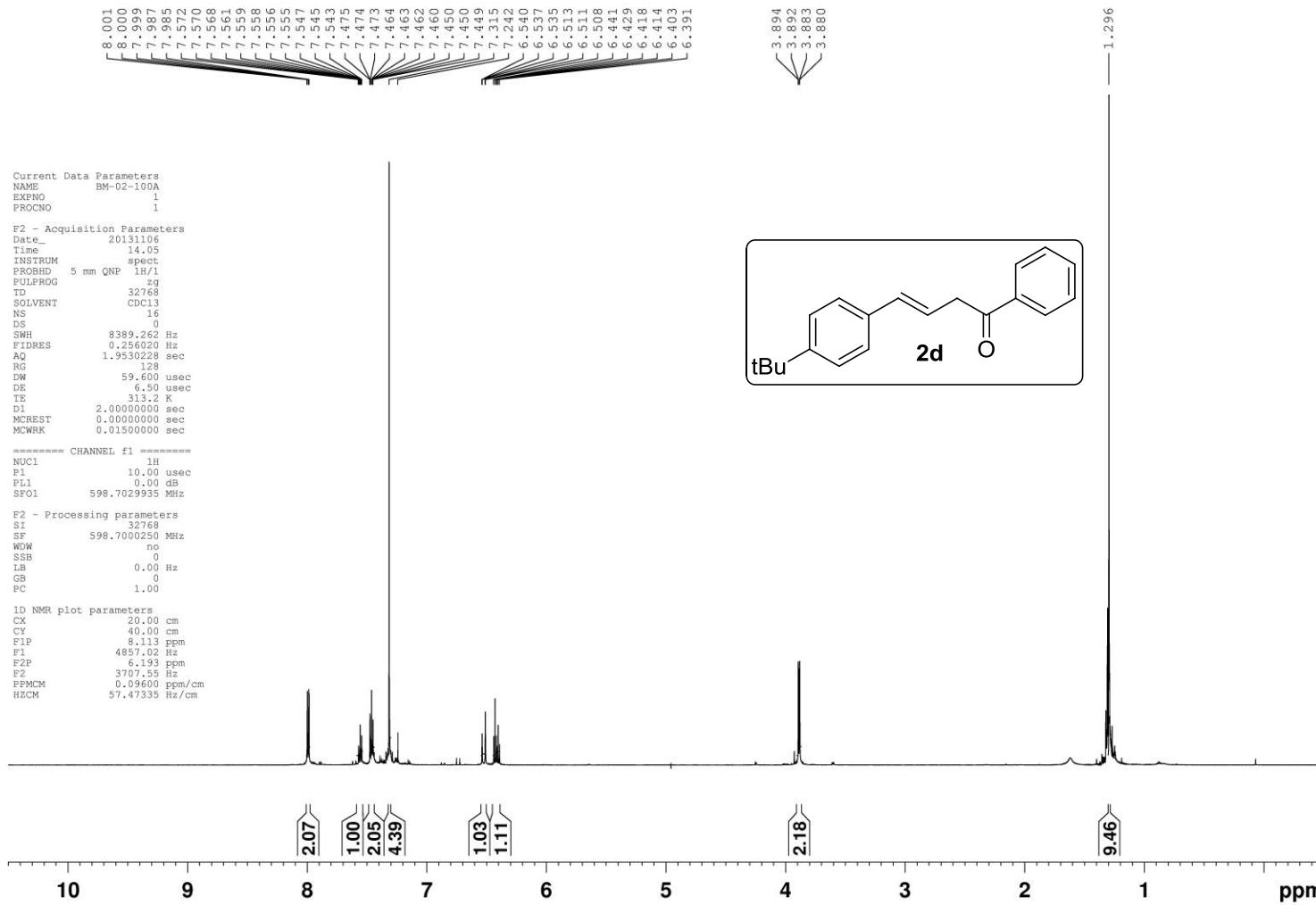
===== CHANNEL f1 =====
 NUC1 13C
 P1 4.80 usec
 PL1 0.00 dB
 SF01 150.5597948 MHz

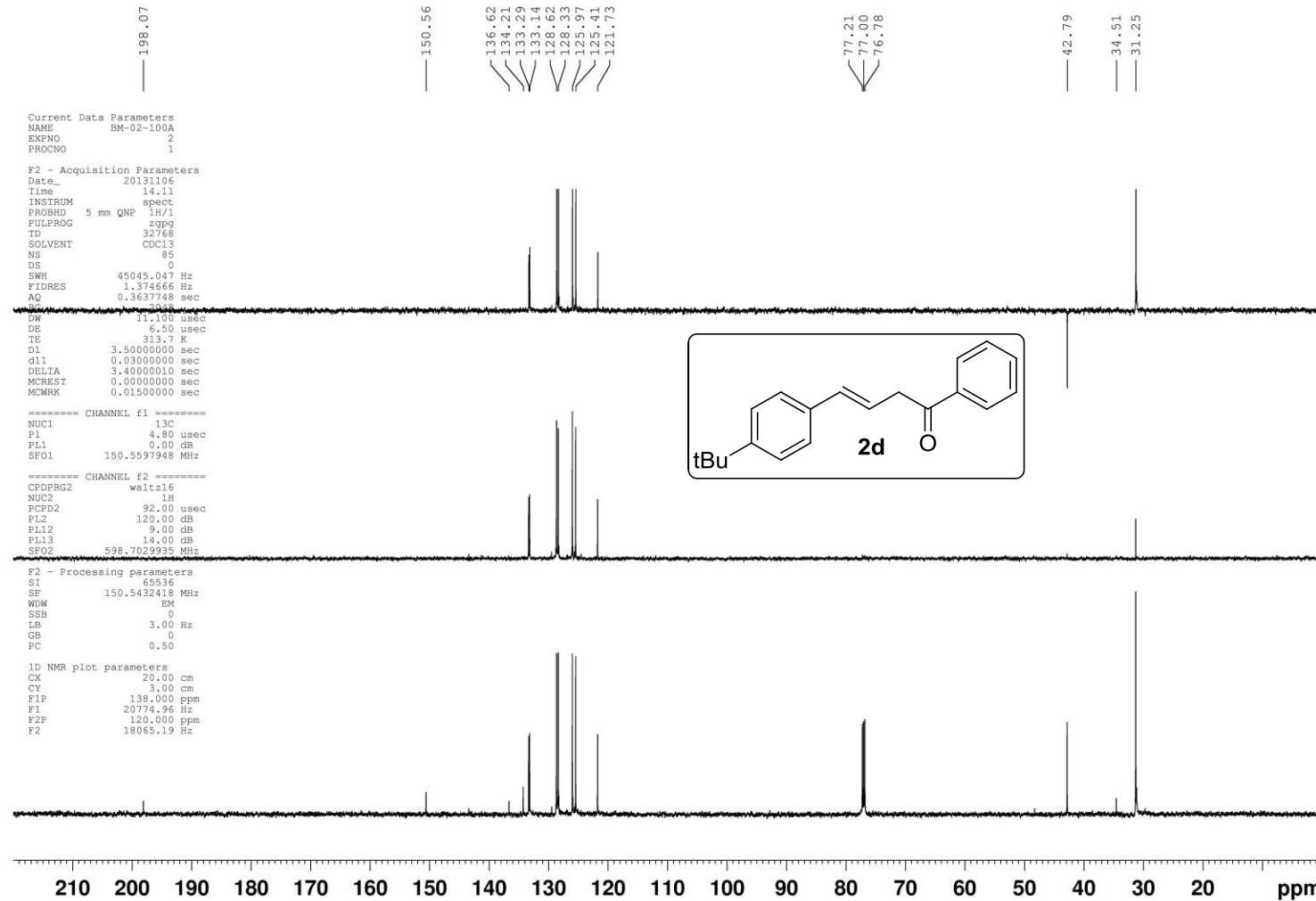
===== CHANNEL f2 =====
 CPDPRG2 waltz16
 NUC2 1H
 PCPD2 92.00 usec
 PL2 120.00 dB
 PL12 9.00 dB
 PL13 14.00 dB
 SF02 598.7029935 MHz

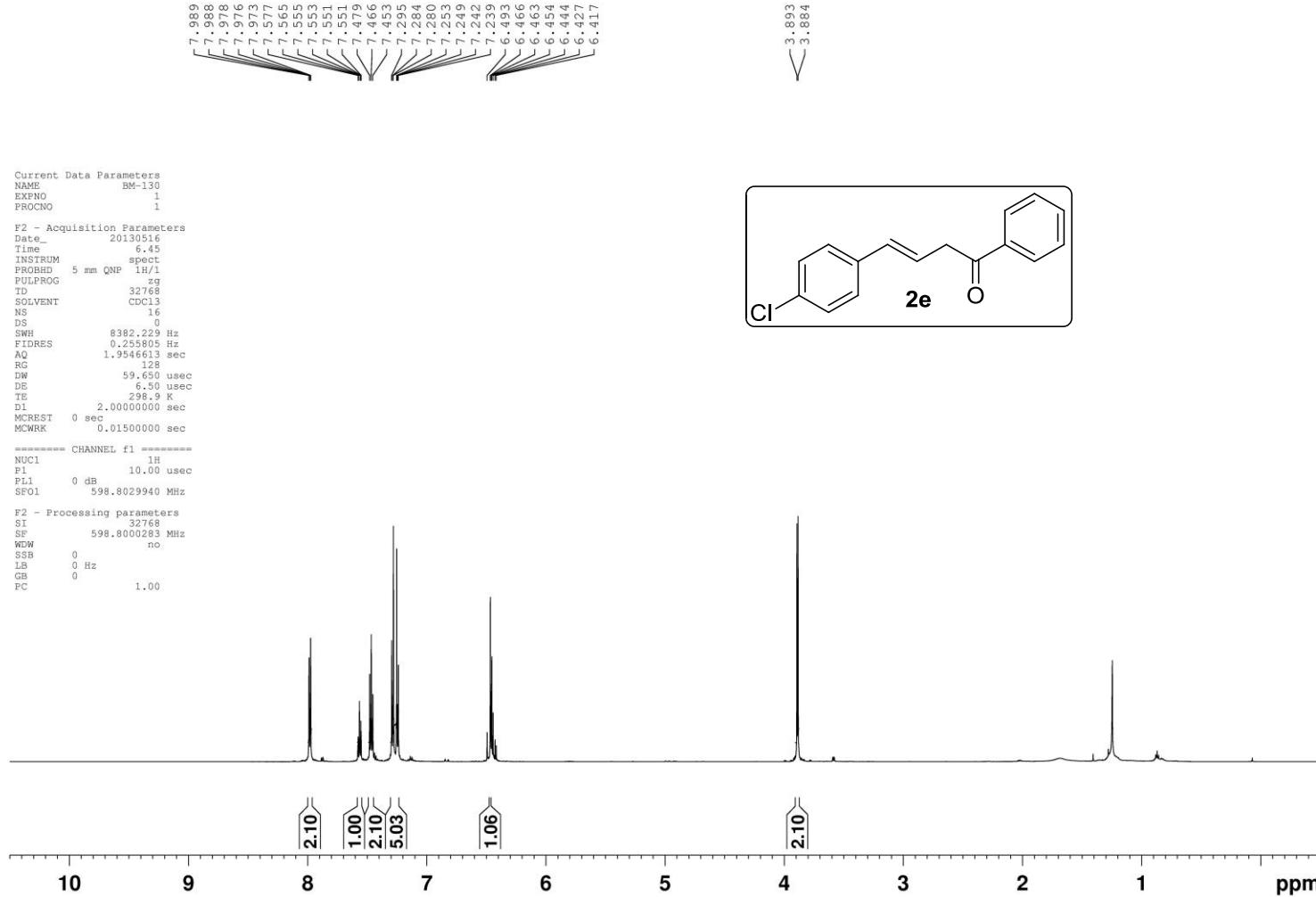
F2 - Processing parameters
 SI 65536
 SF 150.5432376 MHz
 WDW EM
 SSB 0
 LB 3.00 Hz
 GB 0
 PC 0.50

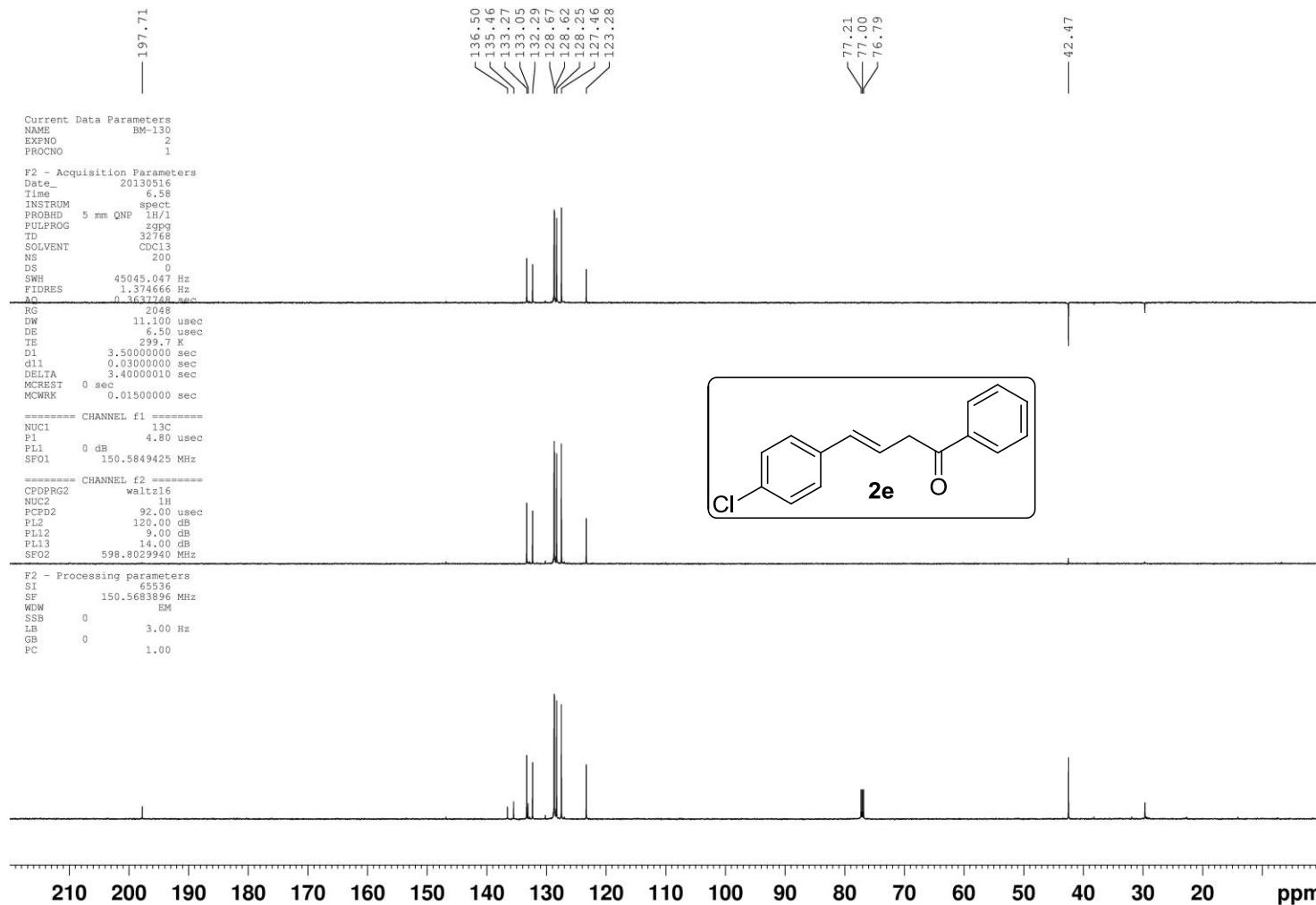
1D NMR plot parameters
 CX 20.00 cm
 CY 4.00 cm
 F1P 210.000 ppm
 F1 31614.08 Hz
 F2P 0.000 ppm
 F2 0.00 Hz
 PPMCM 10.50000 ppm/cm
 HZCM 1580.70410 Hz/cm











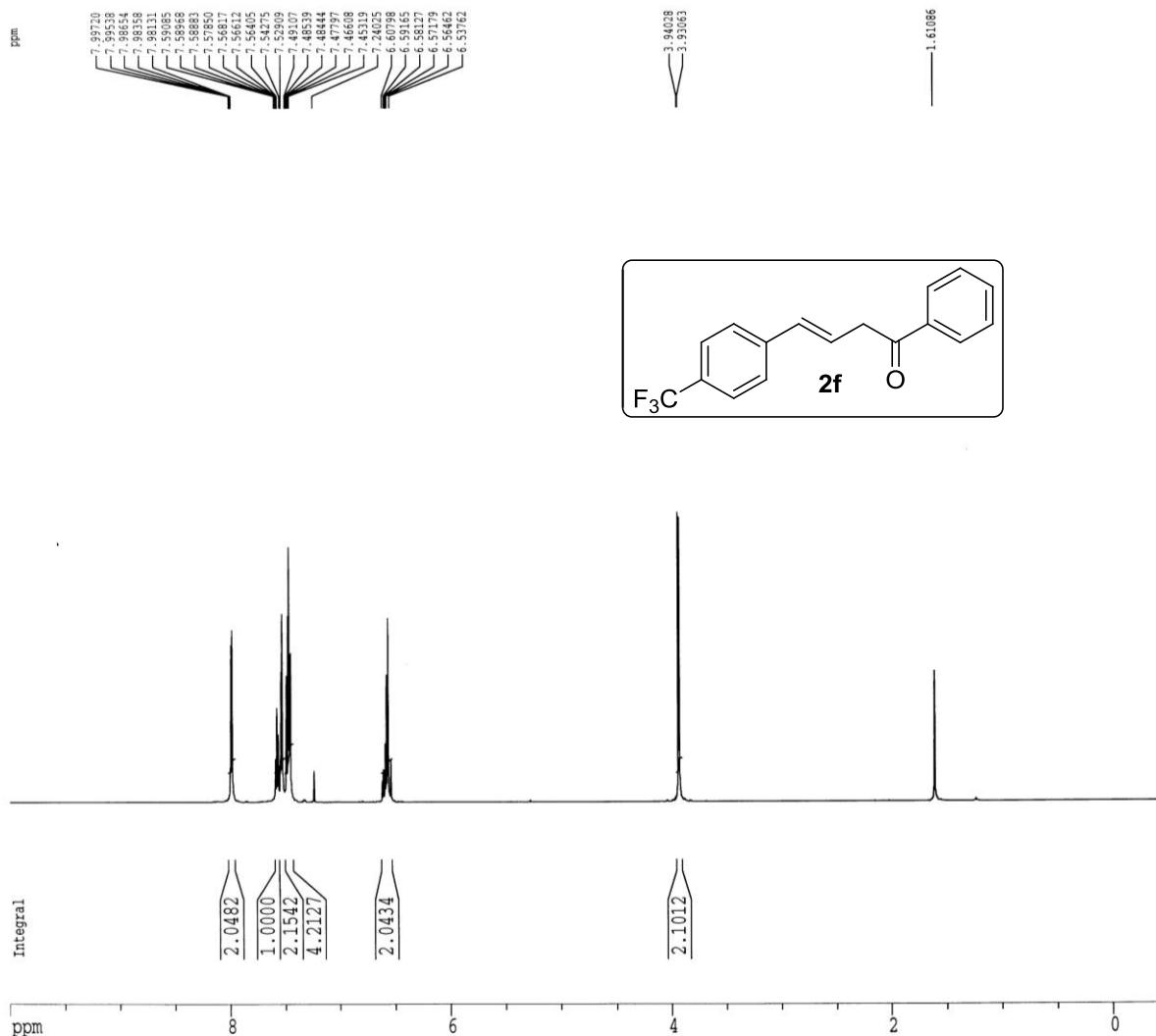
Current Data Parameters
NAME BM-126B
EXPNO 1
PROCNO 1

F2 - Acquisition Parameters
Date 20140311
Time 19.52
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 32768
SOLVENT CDCl3
NS 16
DS 0
SWH 8382.229 Hz
FIDRES 0.255805 Hz
AQ 1.9546613 sec
RG 128
DW 59.650 usec
DE 6.50 usec
TE 293.5 K
D1 2.0000000 sec
MCREST 0.0000000 sec
MCWRK 0.01500000 sec

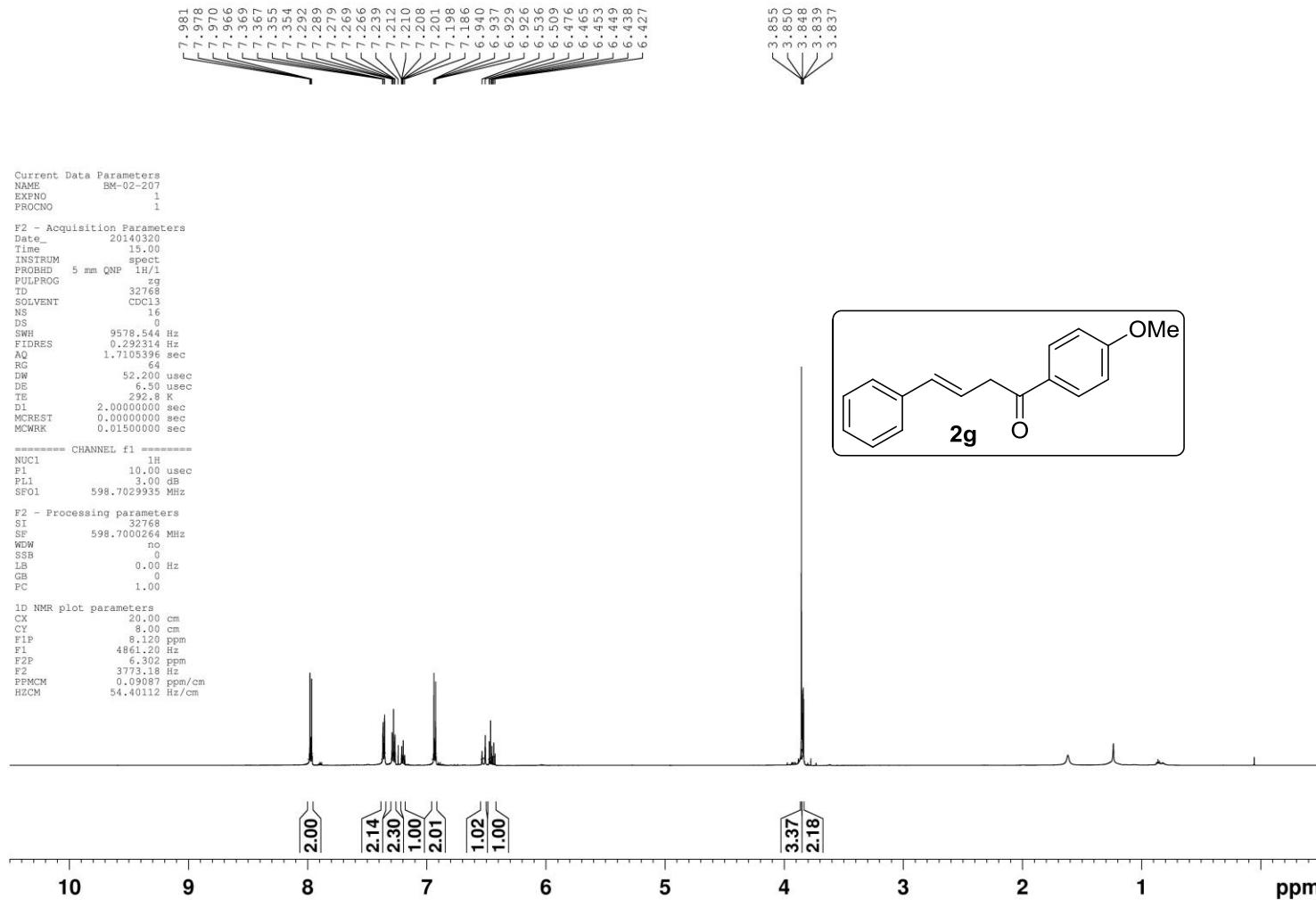
===== CHANNEL f1 =====
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SP01 598.7032929 MHz

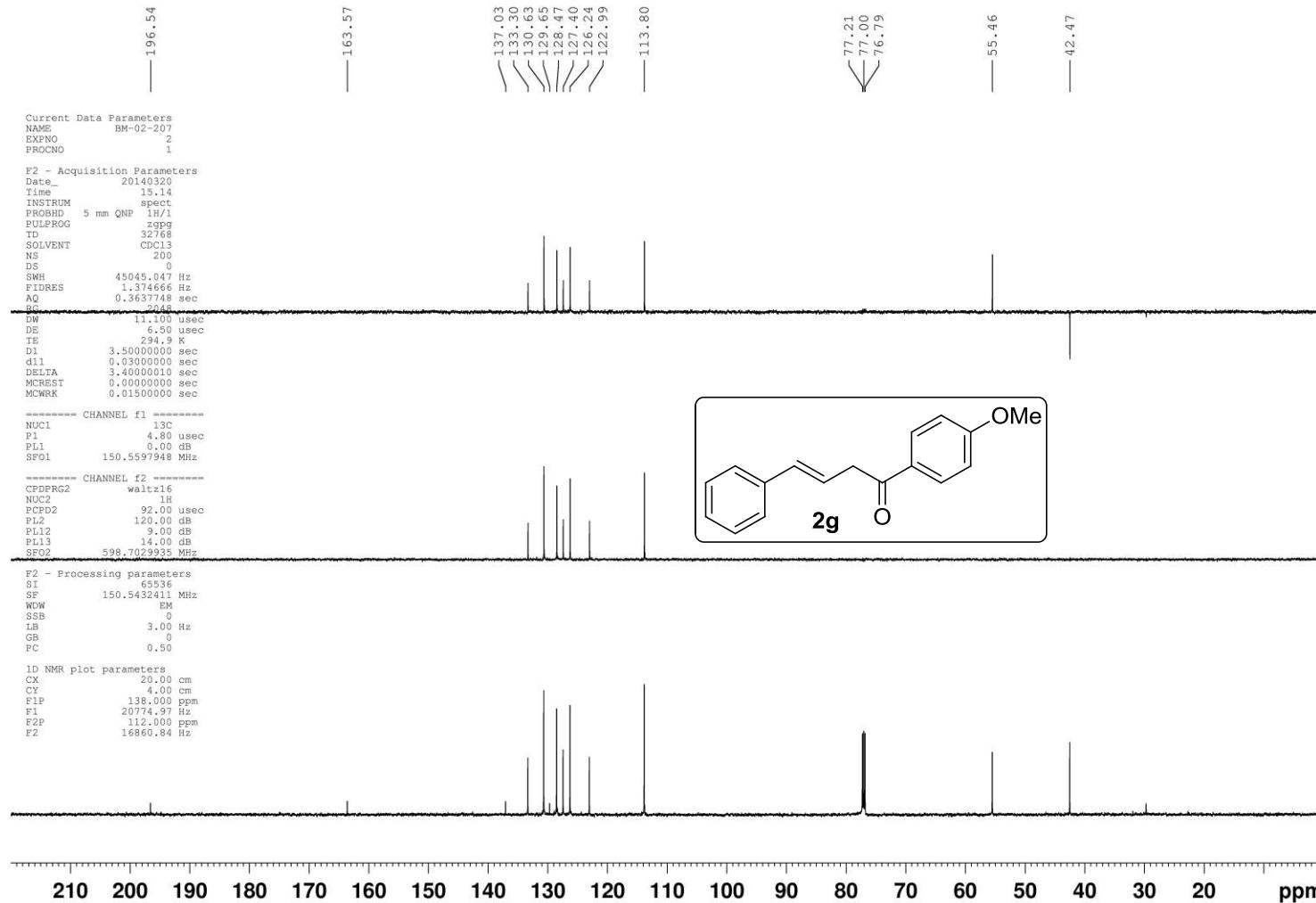
F2 - Processing parameters
SI 32768
SF 598.7000259 MHz
WDW no
SSB 0
LB 0.00 Hz
GB 0
PC 1.00

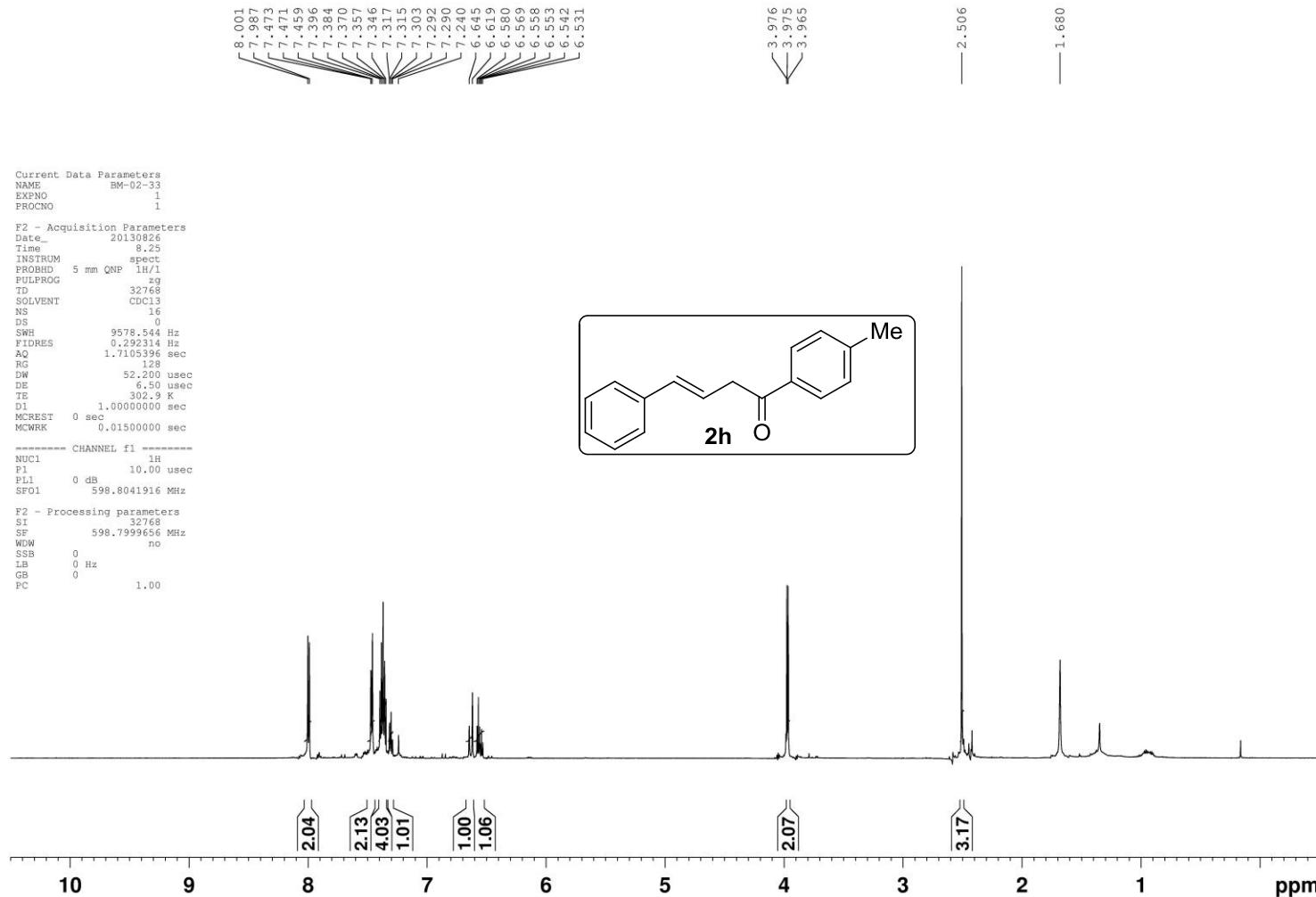
1D NMR plot parameters
CX 20.00 cm
CY 5.00 cm
F1P 10.000 ppm
F1 5987.00 Hz
F2P -0.500 ppm
F2 -299.35 Hz
PPCM 0.52500 ppm/cm
HZCM 314.31750 Hz/cm











Current Data Parameters
NAME BM-02-33
EXPNO 2
PROCNO 1

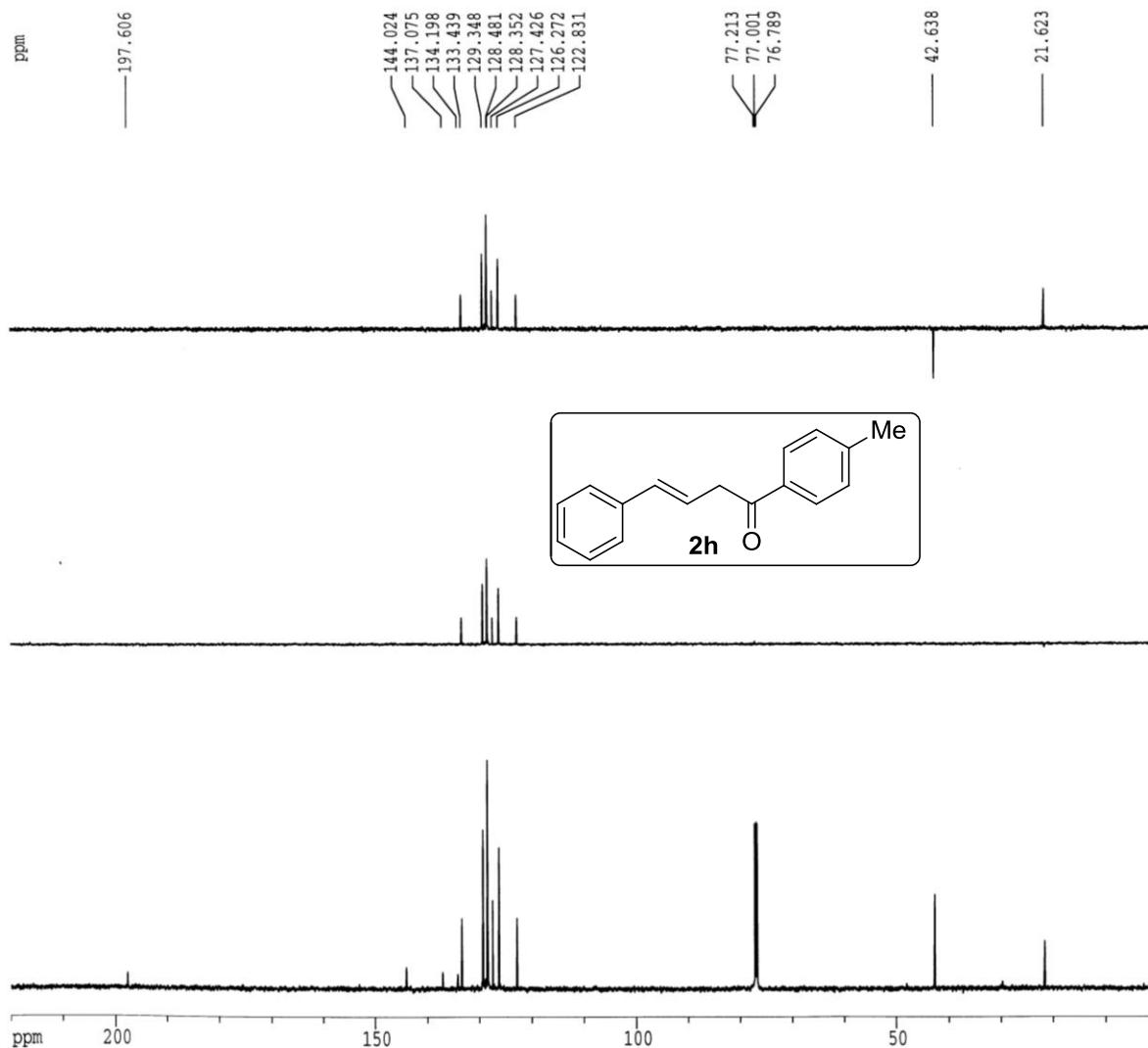
F2 - Acquisition Parameters
Date_ 20130825
Time 17.46
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT DMSO
NS 300
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 304.2 K
D1 3.5000000 sec
d11 0.03000000 sec
DELTA 3.40000010 sec
MCREST 0.0000000 sec
MCWRK 0.01500000 sec

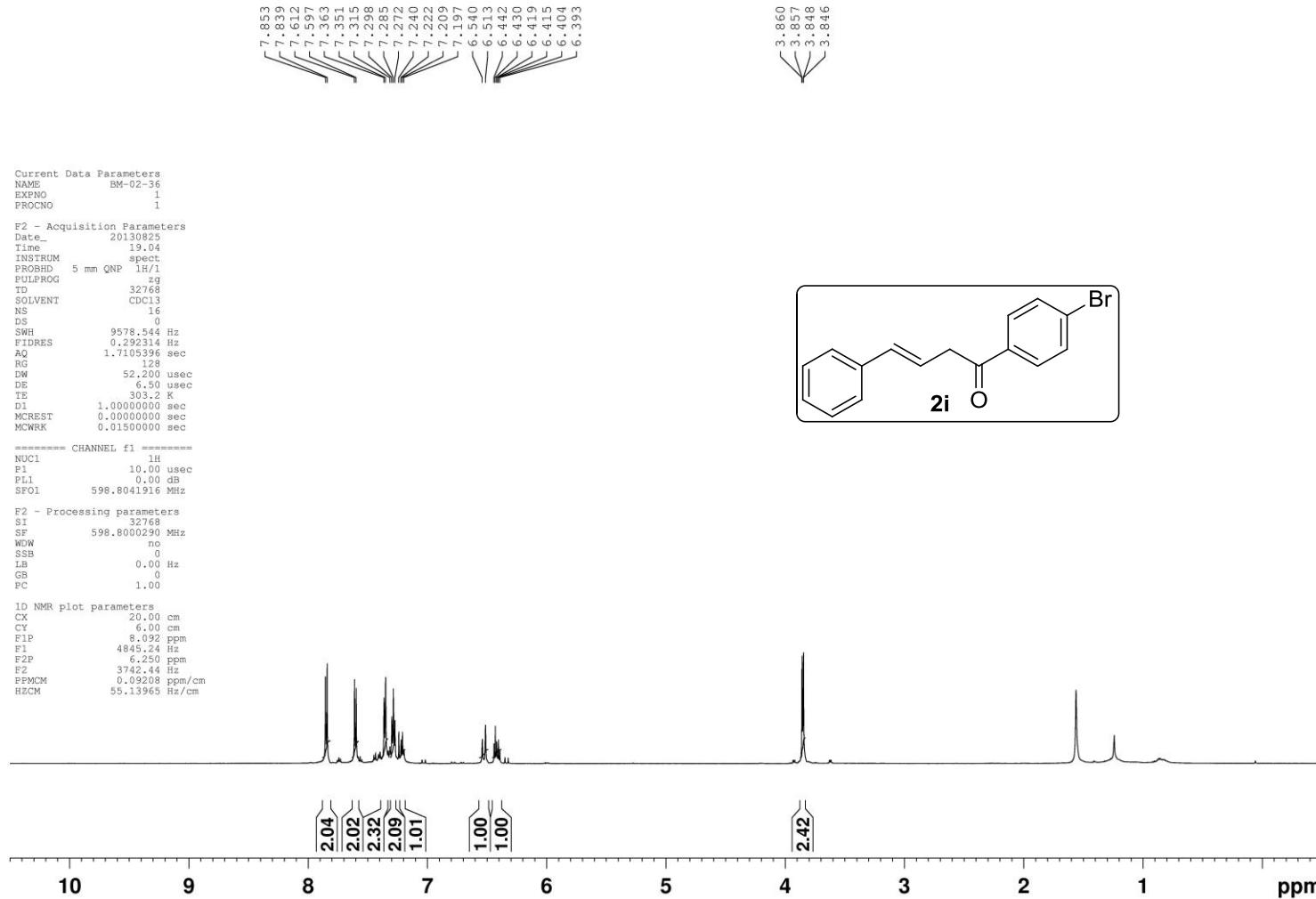
===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SF01 150.5849425 MHz

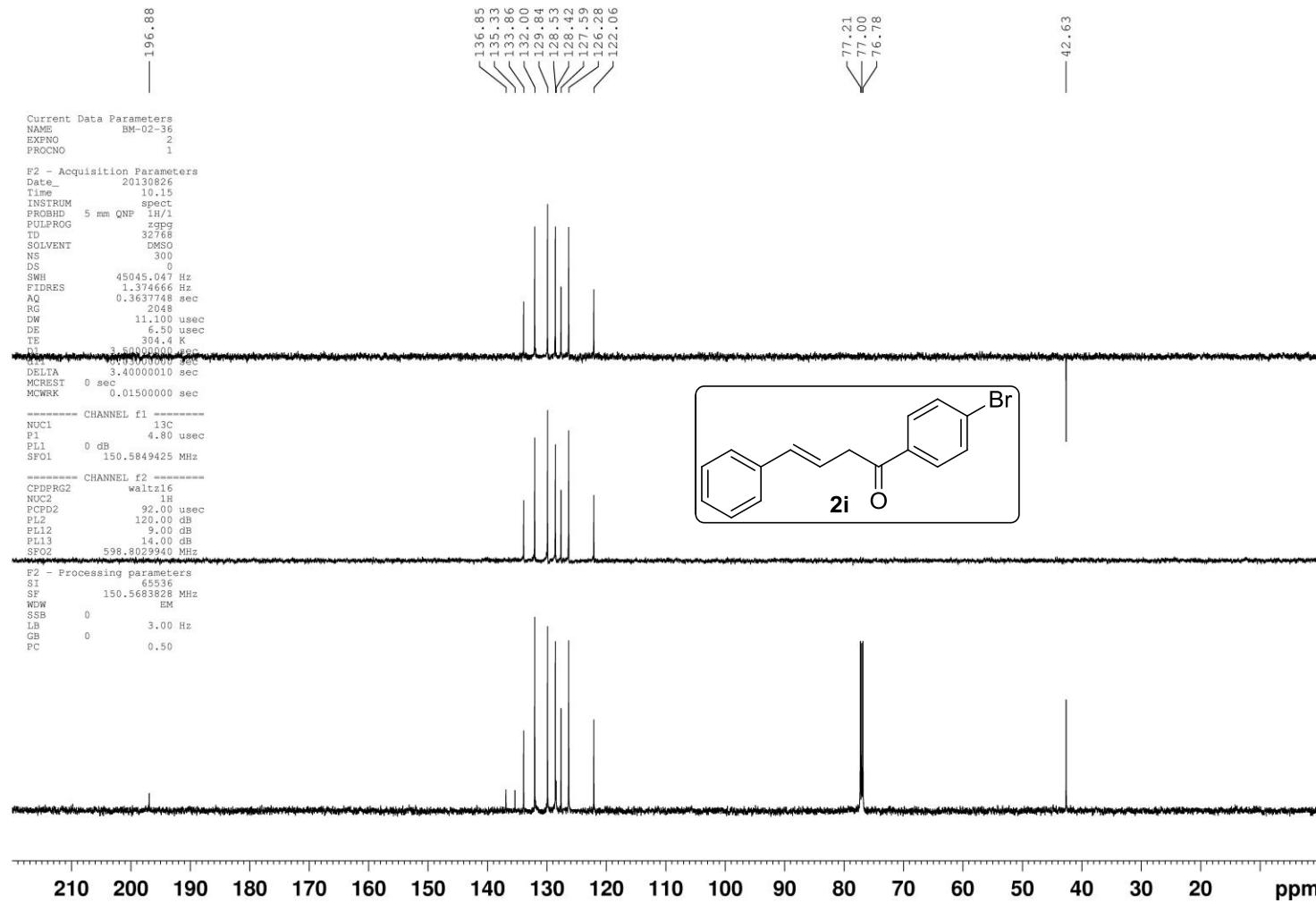
===== CHANNEL F2 =====
CPDPRG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SF02 598.8029940 MHz

F2 - Processing parameters
SI 65536
SF 150.5683821 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 0.50

1D NMR plot parameters
CX 20.00 cm
CY 4.00 cm
F1P 220.000 ppm
F1 33125.04 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPCM 11.00000 ppm/cm
HZCM 1656.25208 Hz/cm







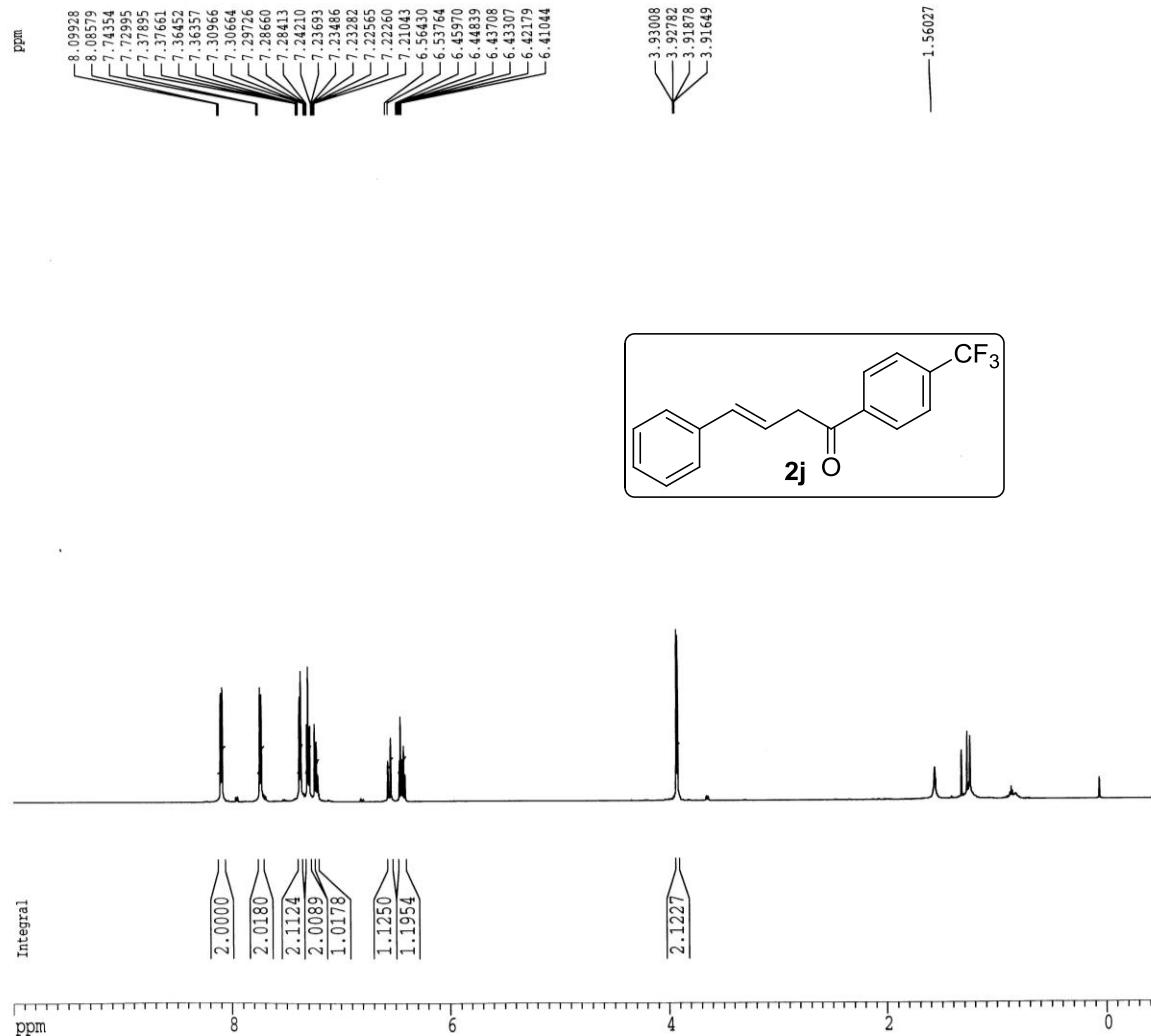
Current Data Parameters
 NAME BM-02-204A
 EXPNO 11
 PROCNO 1

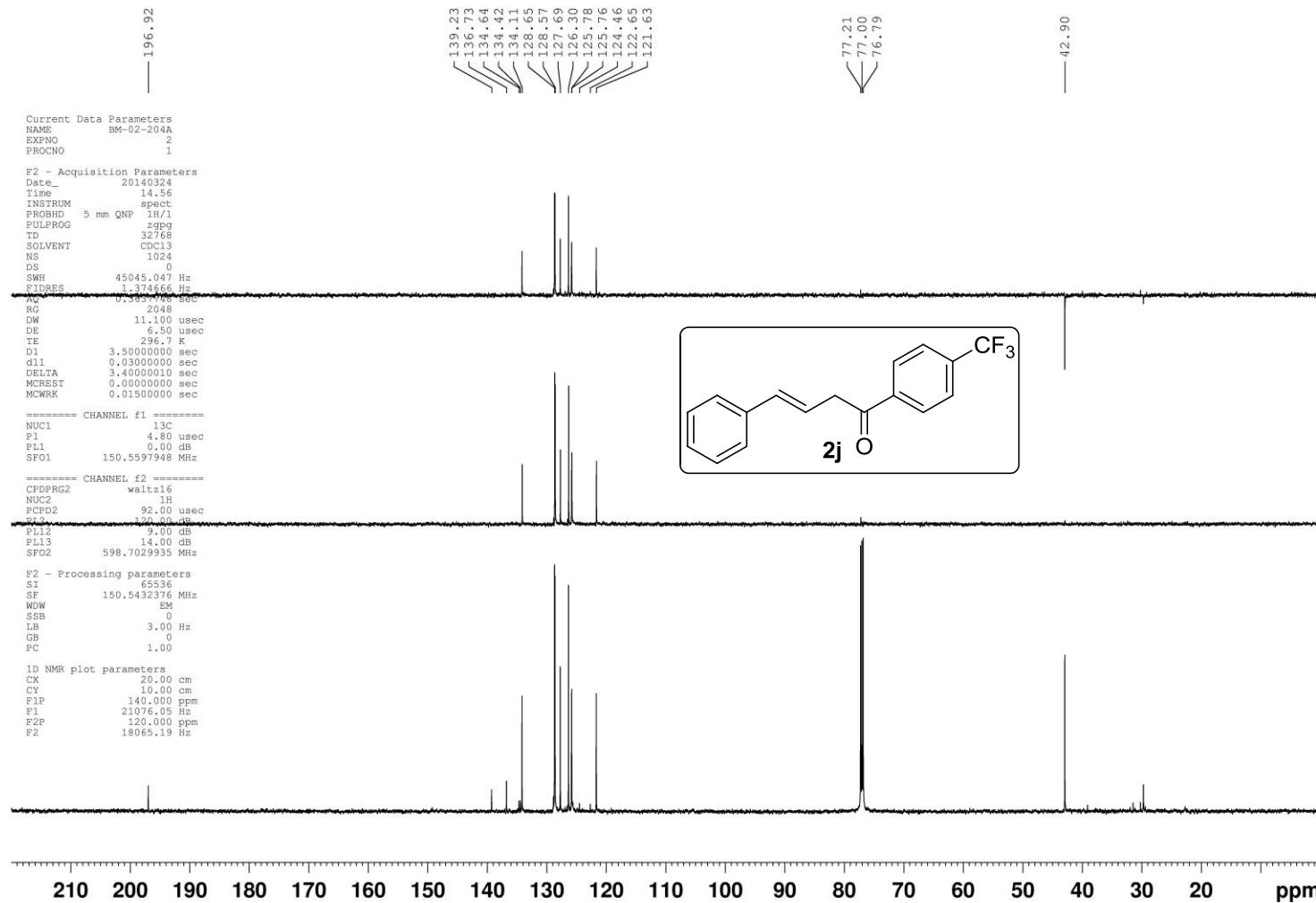
F2 - Acquisition Parameters
 Date_ 20140327
 Time_ 19.47
 INSTRUM spect
 PROBHD 5 mm QNP 1H/1
 PULPROG zg
 TD 32768
 SOLVENT CDCl3
 NS 16
 DS 0
 SWH 8382.229 Hz
 FIDRES 0.255805 Hz
 AQ 1.9546613 sec
 RG 128
 DW 59.650 usec
 DE 6.50 usec
 TE 295.6 K
 D1 2.0000000 sec
 MCREST 0.0000000 sec
 MCWRK 0.0150000 sec

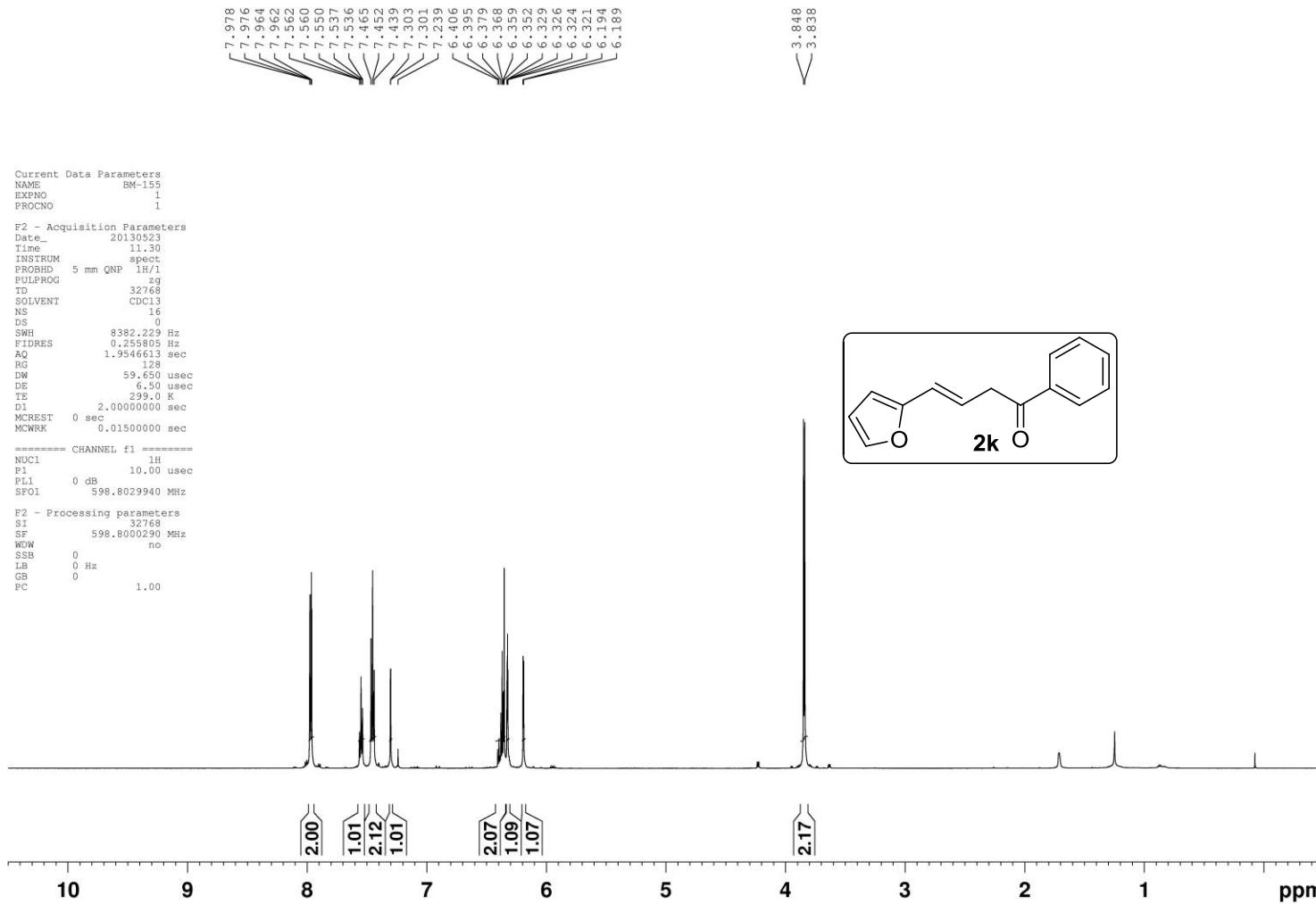
===== CHANNEL f1 =====
 NUC1 1H
 P1 10.00 usec
 PL1 0.00 dB
 SFO1 598.7035922 MHz

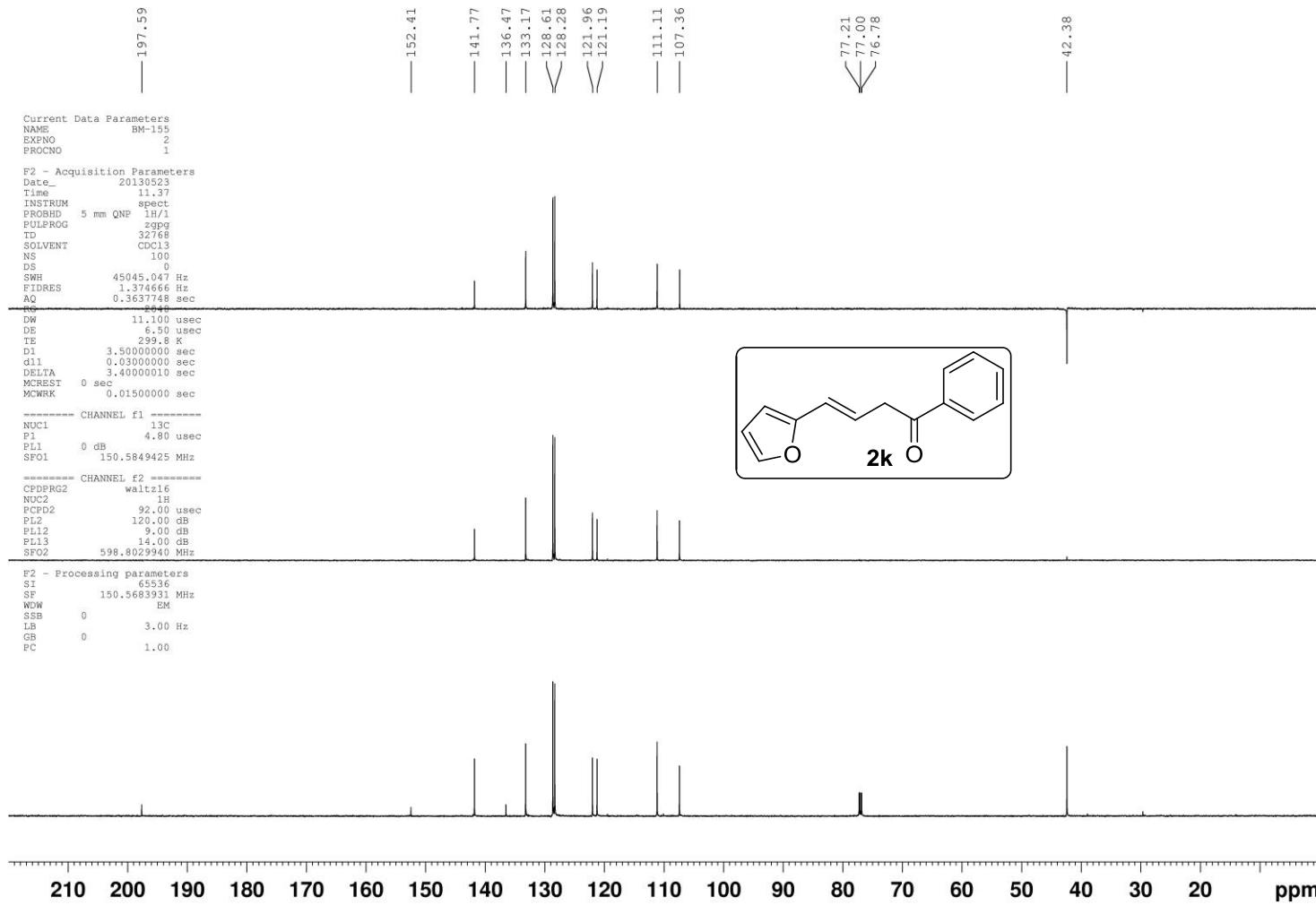
F2 - Processing parameters
 SI 32768
 SF 598.7000250 MHz
 WDW no
 SSB 0
 LB 0.00 Hz
 GB 0
 PC 1.00

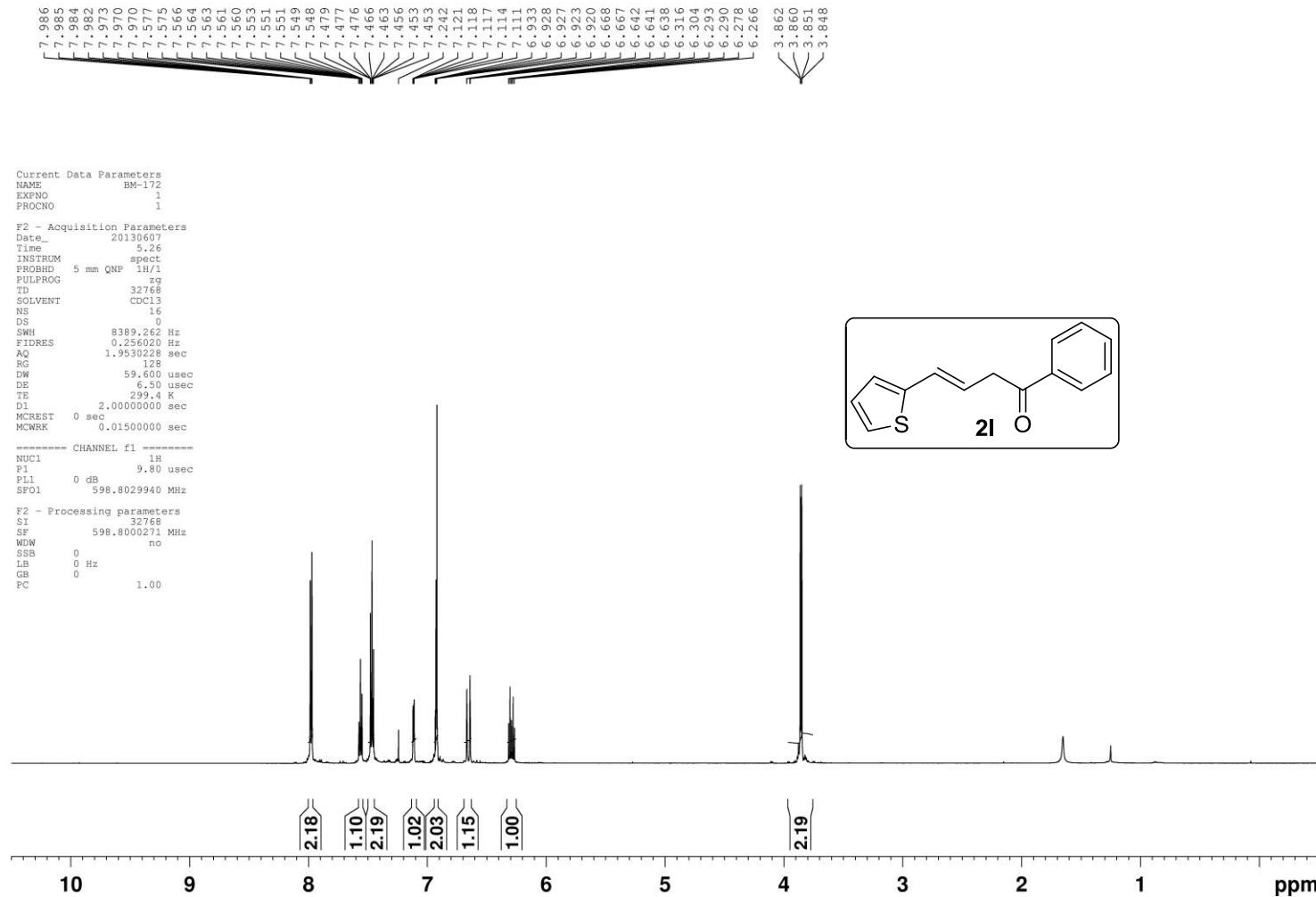
1D NMR plot parameters
 CX 20.00 cm
 CY 3.00 cm
 F1P 10.000 ppm
 F1 5987.00 Hz
 F2P -0.500 ppm
 F2 -299.35 Hz
 PPMCM 0.52500 ppm/cm
 HZCM 314.31750 Hz/cm











Current Data Parameters
NAME BM-172
EXPNO 2
PROCNO 1

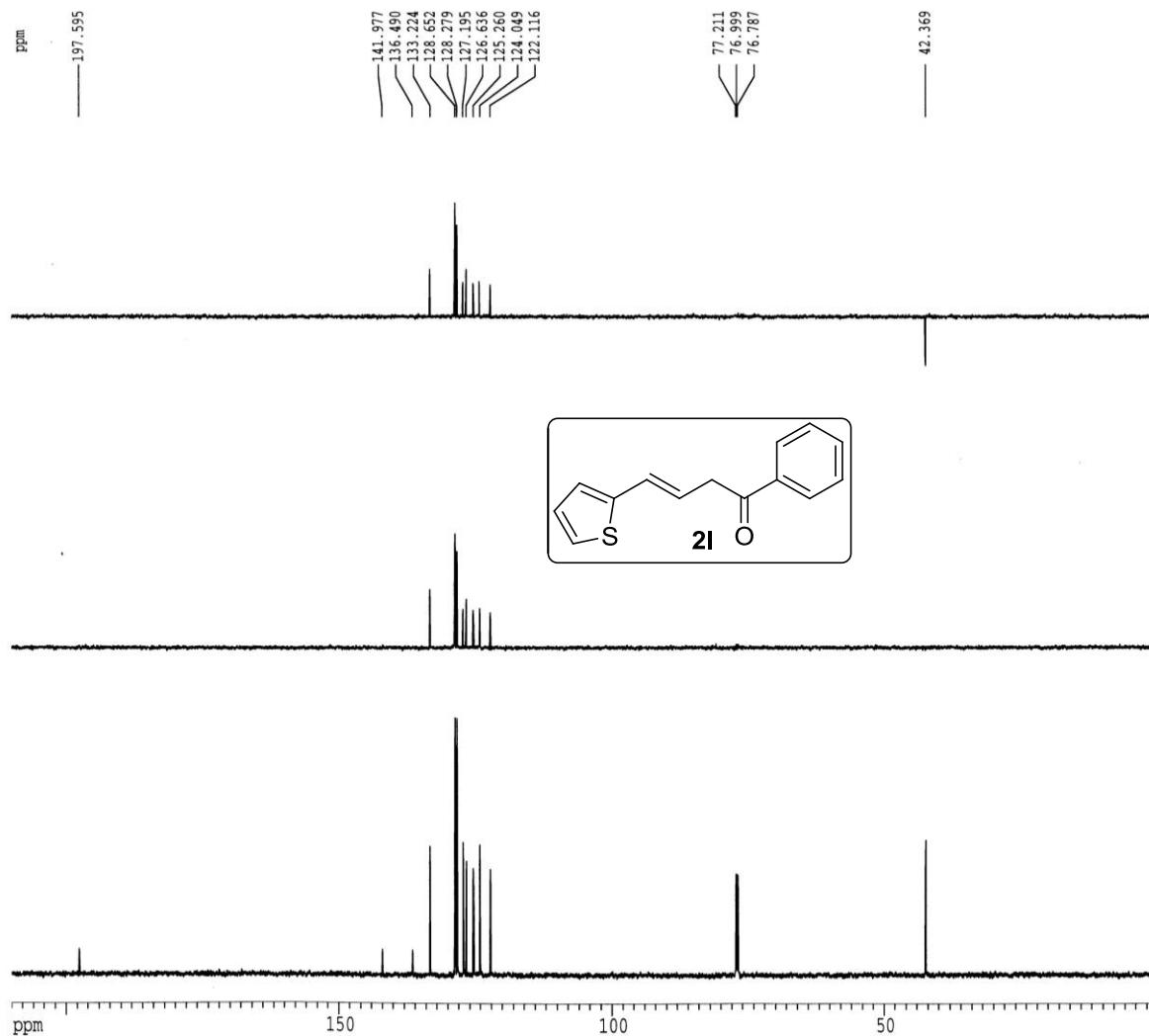
F2 - Acquisition Parameters
Date_ 20130606
Time 14.18
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT CDCl3
NS 46
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 299.7 K
D1 3.5000000 sec
d11 0.0300000 sec
DELTA 3.4000000 sec
MCREST 0.0000000 sec
MCWRK 0.01500000 sec

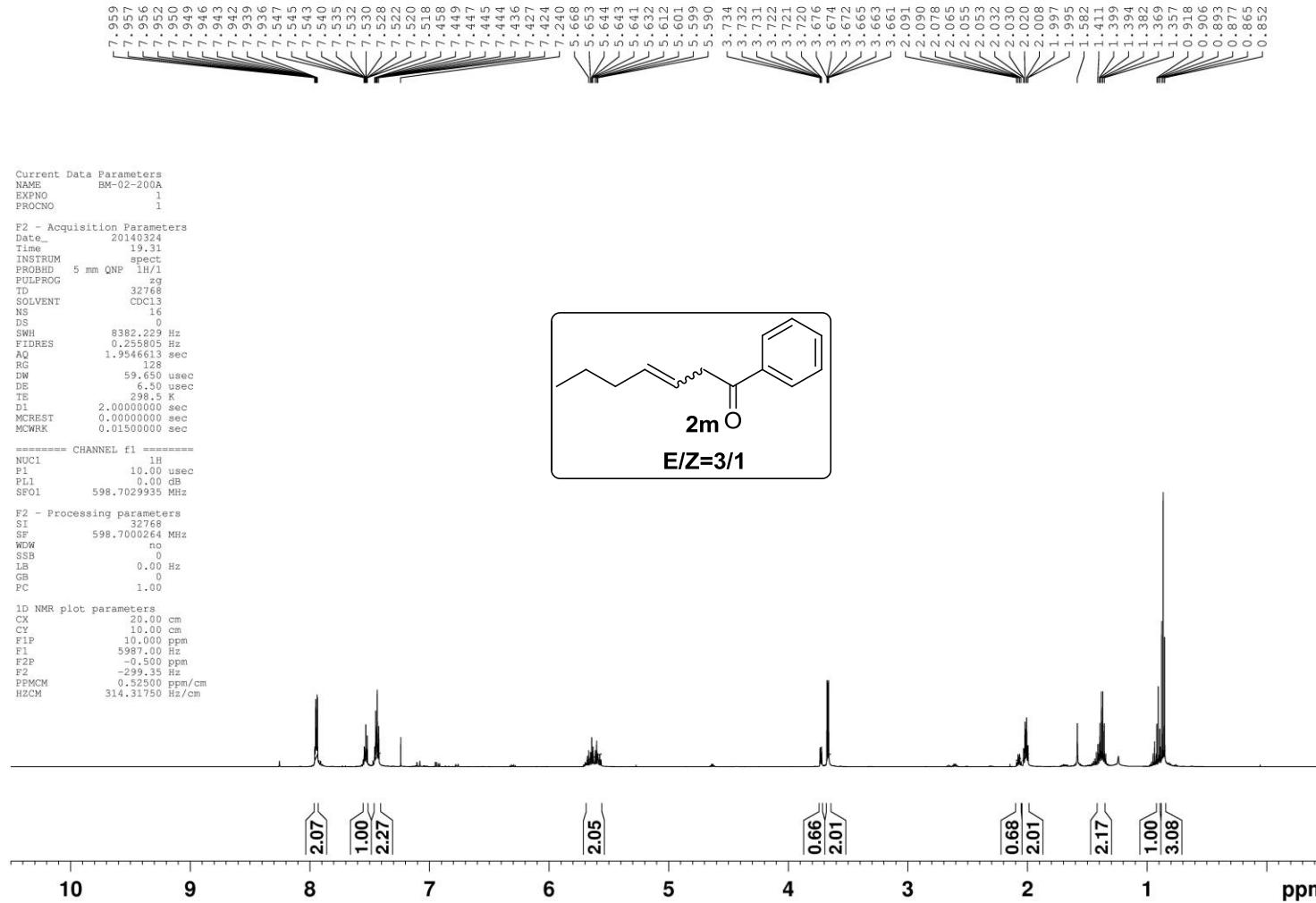
===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SF01 150.5849425 MHz

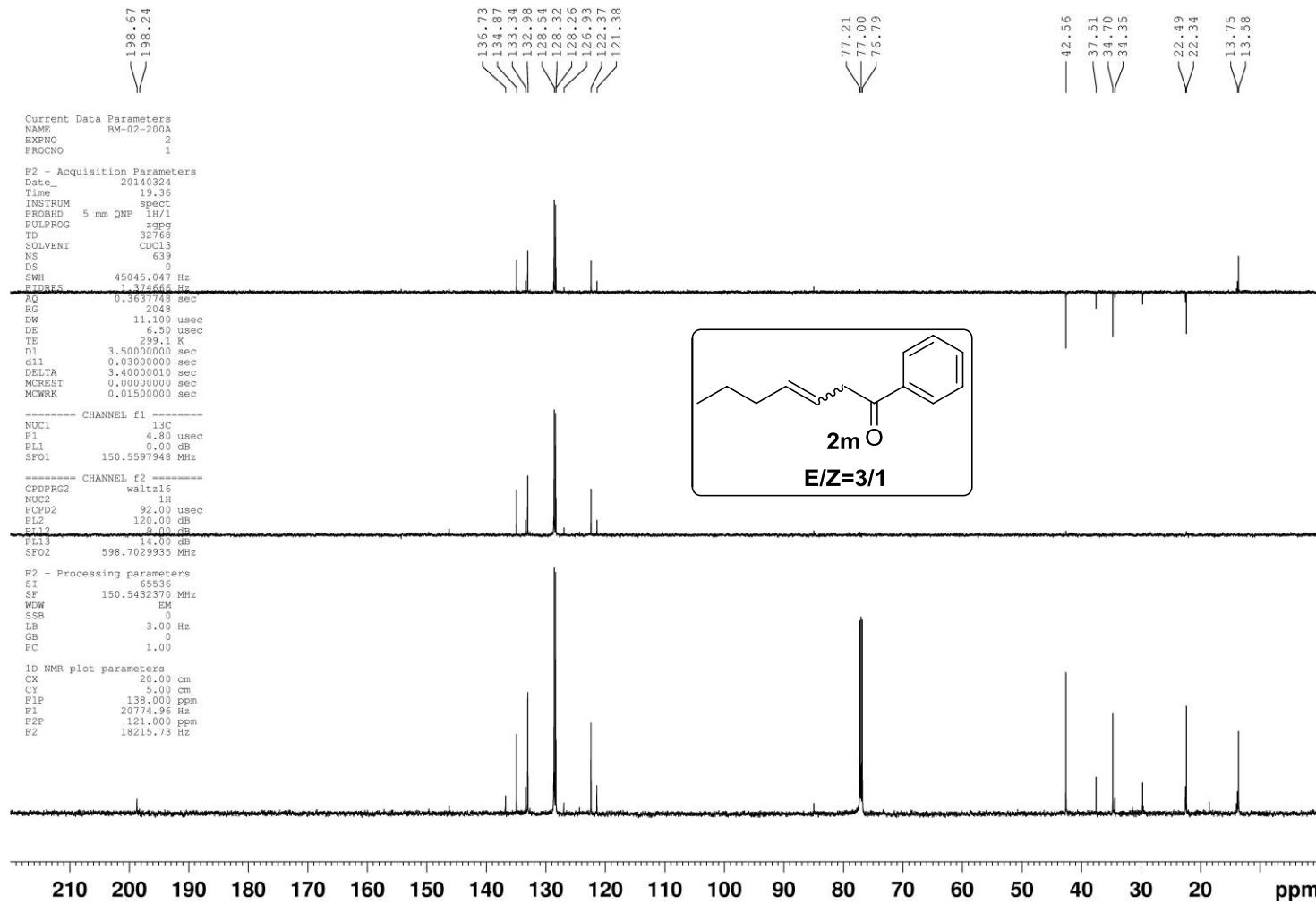
===== CHANNEL f2 =====
CPDPRG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SF02 598.8029940 MHz

F2 - Processing parameters
SI 65536
SF 150.5683910 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 1.00

1D NMR plot parameters
CX 20.00 cm
CY 4.50 cm
F1P 210.000 ppm
F1 31619.36 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPCM 10.50000 ppm/cm
HZCM 1580.96814 Hz/cm







Current Data Parameters
NAME BM-195
EXPNO 1
PROCNO 1

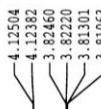
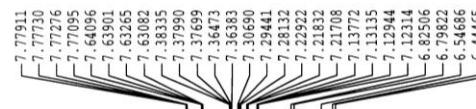
F1 - Acquisition Parameters
Date 20140323
Time 18.45
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 32768
SOLVENT CDCl3
NS 16
DS 0
SWH 8382.229 Hz
FIDRES 0.255805 Hz
AQ 1.9546613 sec
RG 128
DW 59.650 usec
DE 6.50 usec
TE 295.3 K
D1 2.0000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

===== CHANNEL f1 =====
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SF01 598.7029935 MHz

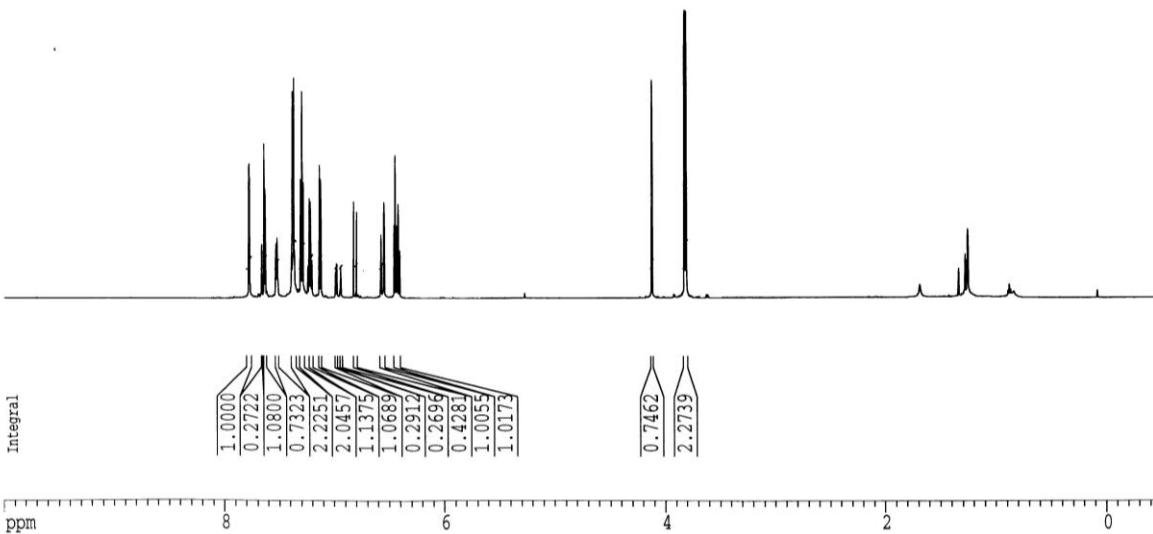
F2 - Processing parameters
SI 32768
SF 598.7000250 MHz
WDW no
SSB 0
LB 0.00 Hz
GB 0
PC 1.00

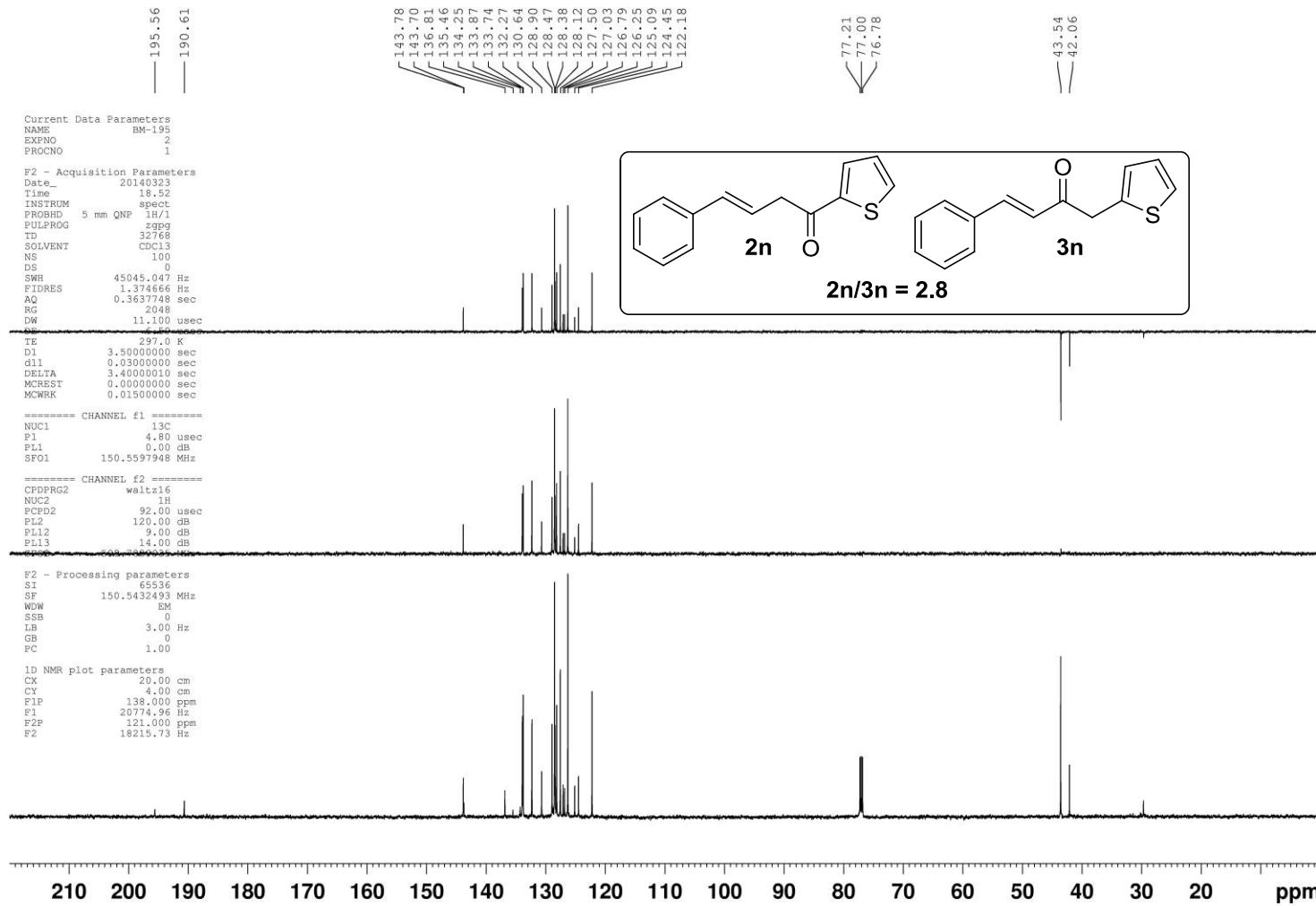
1D NMR plot parameters
CX 20.00 cm
CY 5.00 cm
F1P 10.000 ppm
F1 5987.00 Hz
F2P -0.500 ppm
F2 -299.35 Hz
PPCM 0.52500 ppm/cm
HZCM 314.31750 Hz/cm

ppm



$$2n/3n = 2.8$$





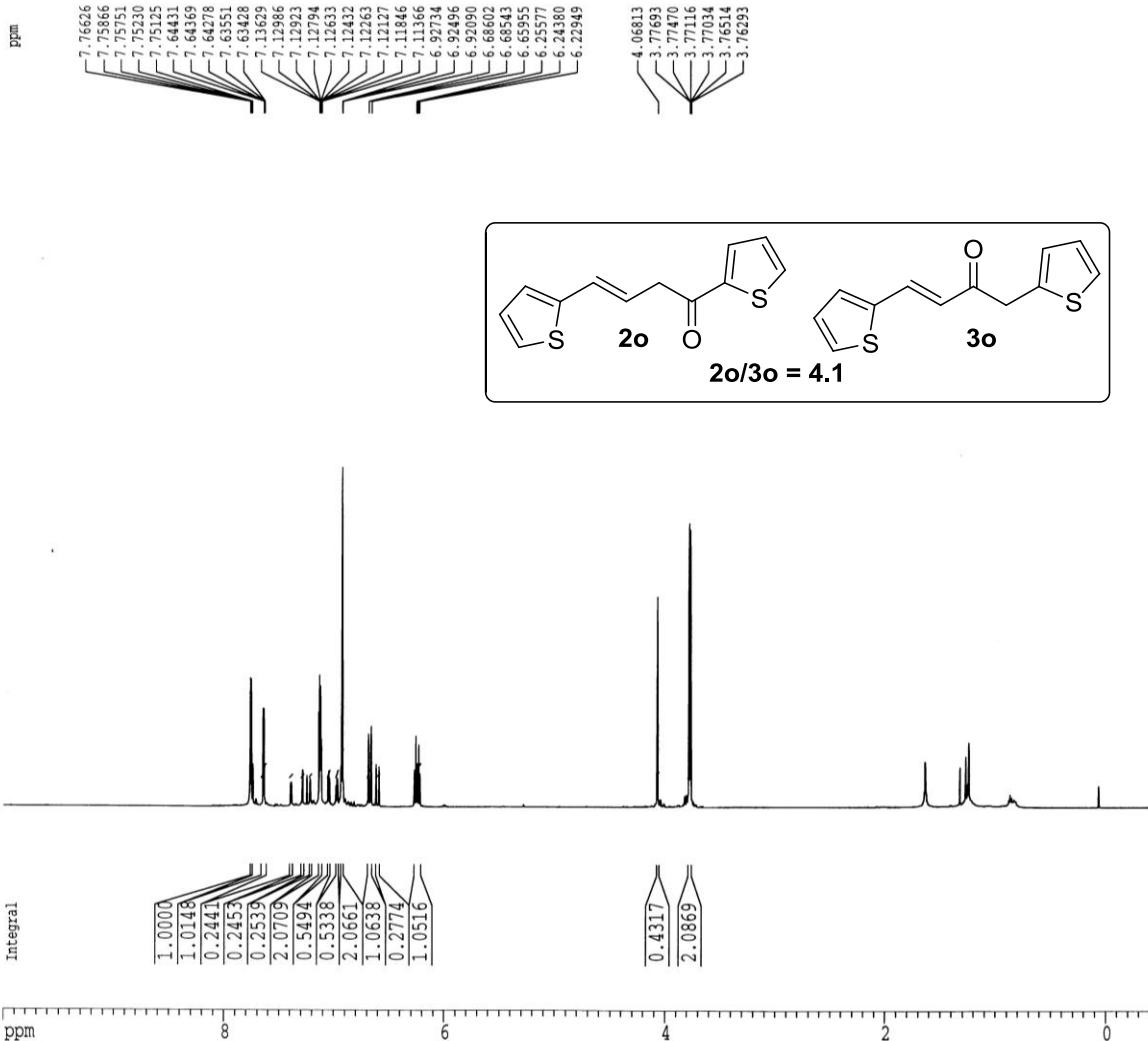
Current Data Parameters
NAME BM-216
EXPNO 1
PROCNO 1

F2 - Acquisition Parameters
Date_ 20140323
Time 19.47
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 32768
SOLVENT CDCl3
NS 16
DS 0
SWH 8382.229 Hz
FIDRES 0.255805 Hz
AQ 1.9546613 sec
RG 128
DM 59.650 usec
DE 6.50 usec
TE 295.5 K
D1 2.0000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

***** CHANNEL f1 *****
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SFO1 598.7029935 MHz

F2 - Processing parameters
SI 32768
SF 598.7000250 MHz
WDW no
SSB 0
LB 0.00 Hz
GB 0
PC 1.00

1D NMR plot parameters
CX 20.00 cm
CY 6.00 cm
F1P 10.000 ppm
F1 5987.00 Hz
F2P -0.500 ppm
F2 -299.35 Hz
PPCM 0.52500 ppm/cm
HZCM 314.31750 Hz/cm



Current Data Parameters
 NAME EM-216
 EXPNO 2
 PROCNO 1

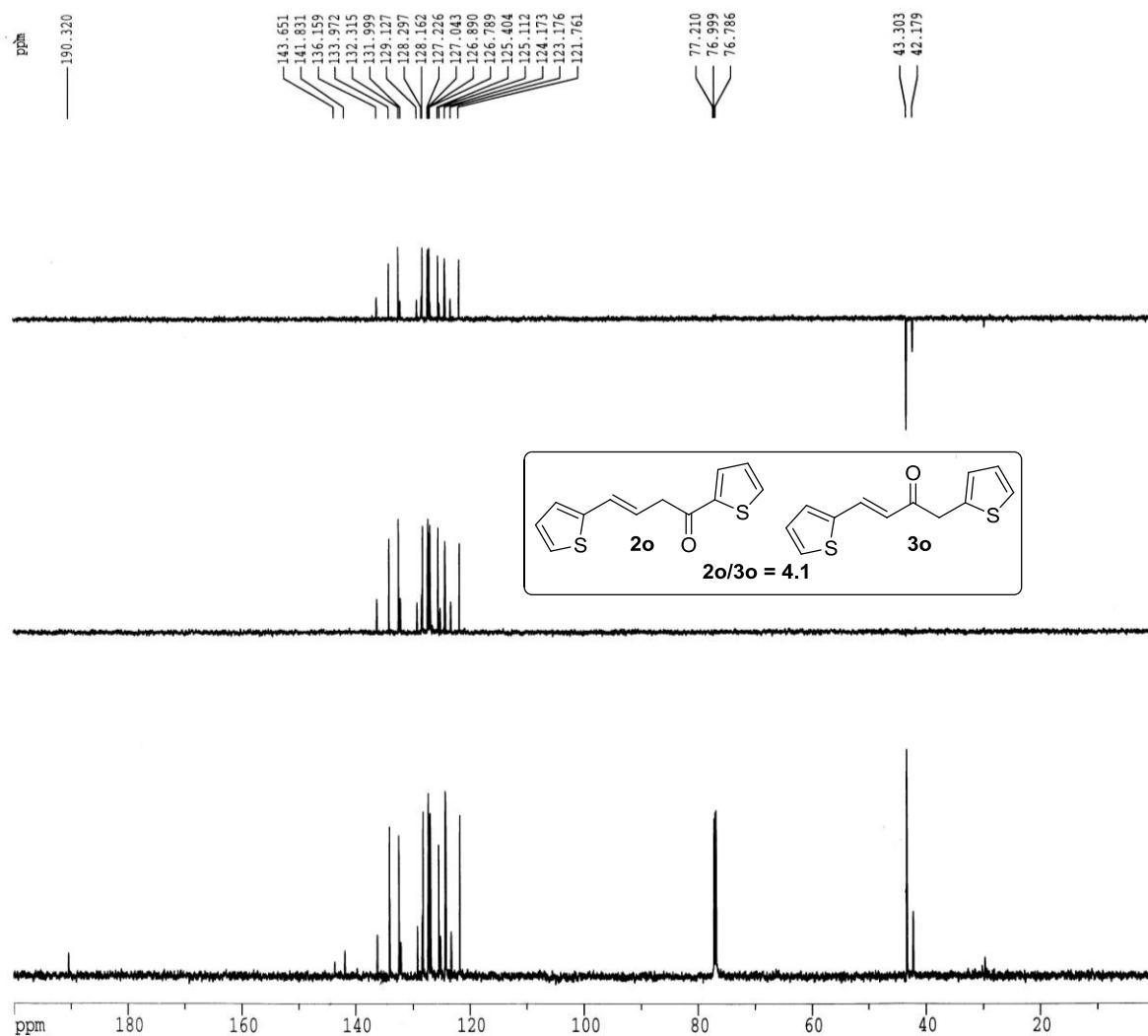
F2 - Acquisition Parameters
 Date_ 20140323
 Time 19.54
 INSTRUM spect
 PROBHD 5 mm QNP 1H/1
 PULPROG zgpg
 TD 32768
 SOLVENT CDCl3
 NS 100
 DS 0
 SWH 45045.047 Hz
 FIDRES 1.374666 Hz
 AQ 0.3637748 sec
 RG 2048
 DW 11.100 usec
 DE 6.50 usec
 TE 297.0 K
 D1 3.5000000 sec
 d11 0.0300000 sec
 DELTA 3.40000010 sec
 MCREST 0.0000000 sec
 MCWRK 0.01500000 sec

===== CHANNEL f1 =====
 NUC1 13C
 P1 4.80 usec
 PL1 0.00 dB
 SFO1 150.5597948 MHz

===== CHANNEL f2 =====
 CDPGR2 waltz16
 NUC2 1H
 PCPD2 92.00 usec
 PL2 120.00 dB
 PL12 9.00 dB
 PL13 14.00 dB
 SFO2 598.7029935 MHz

F2 - Processing parameters
 SI 65536
 SF 150.5432431 MHz
 WDW EM
 SSB 0
 LB 3.00 Hz
 GB 0
 PC 1.00

1D NMR plot parameters
 CX 20.00 cm
 CY 4.00 cm
 F1P 200.000 ppm
 F1 30108.65 Hz
 F2P 0.000 ppm
 F2 0.00 Hz
 PPMCM 10.00000 ppm/cm
 HZCM 1505.43237 Hz/cm



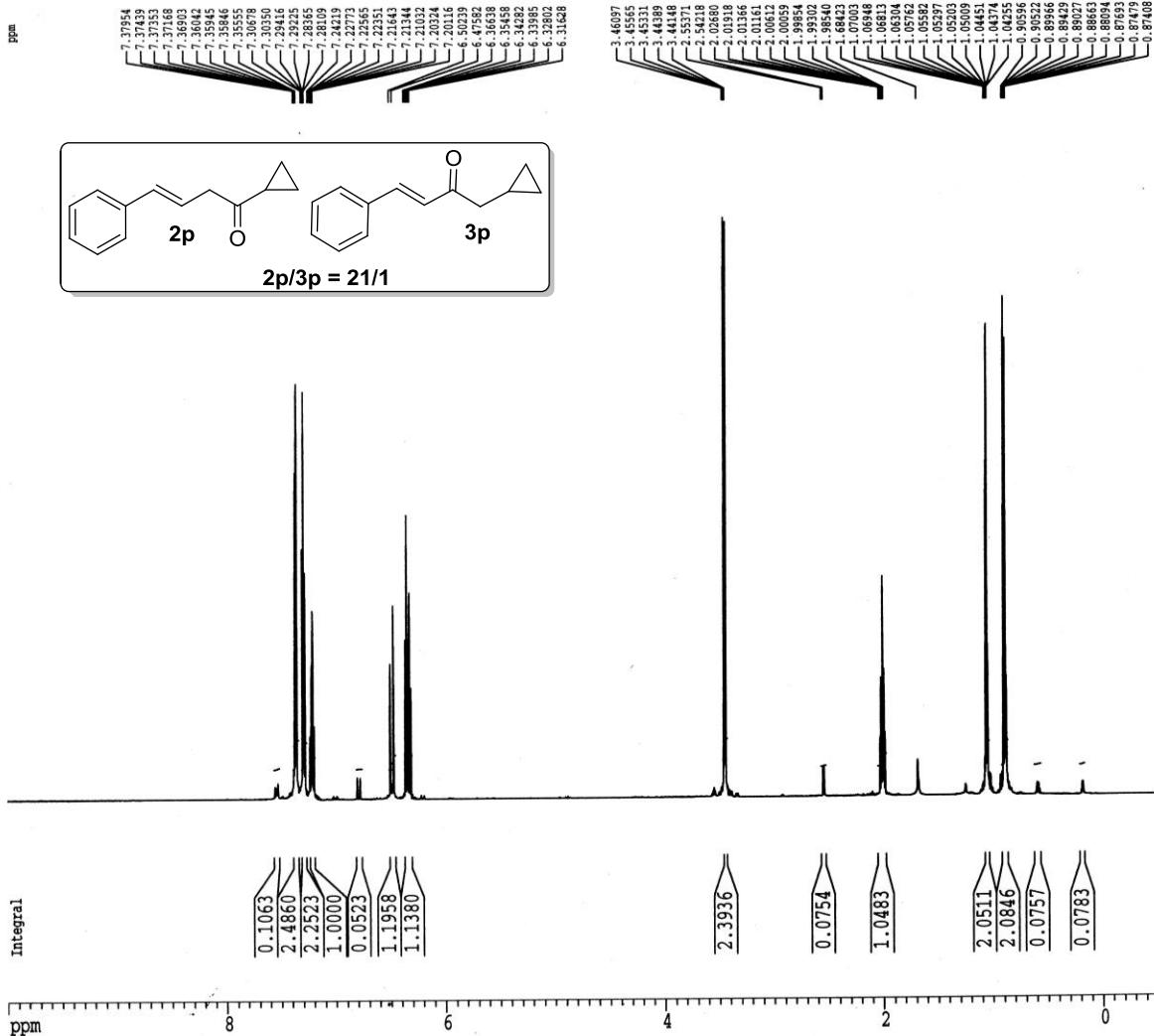
Current Data Parameters
 NAME BM-03-11
 EXPNO 1
 PROCNO 1

F2 - Acquisition Parameters
 Date_ 20140522
 Time 20.01
 INSTRUM spect
 PROBHD 5 mm QNP 1H/1
 PULPROG zg
 TD 32768
 SOLVENT CDCl3
 NS 16
 DS 0
 SWH 8389.262 Hz
 FIDRES 0.256020 Hz
 AQ 1.9530228 sec
 RG 128
 DW 59.600 usec
 DE 6.50 usec
 TE 298.2 K
 D1 2.0000000 sec
 MCREST 0.0000000 sec
 MCWRK 0.0150000 sec

===== CHANNEL f1 =====
 NUC1 1H
 PI 10.00 usec
 PLL 3.00 dB
 SFO1 598.7029935 MHz

F2 - Processing parameters
 SI 32768
 SF 598.7000248 MHz
 WDW no
 SSB 0
 LB 0.00 Hz
 GB 0
 PC 1.00

1D NMR plot parameters
 CX 20.00 cm
 CY 10.00 cm
 F1P 10.000 ppm
 F1 5987.00 Hz
 F2P -0.500 ppm
 F2 -299.35 Hz
 PPMCM 0.52500 ppm/cm
 HZCM 314.31750 Hz/cm



Current Data Parameters
NAME BM-03-11
EXPNO 2
PROCNO 1

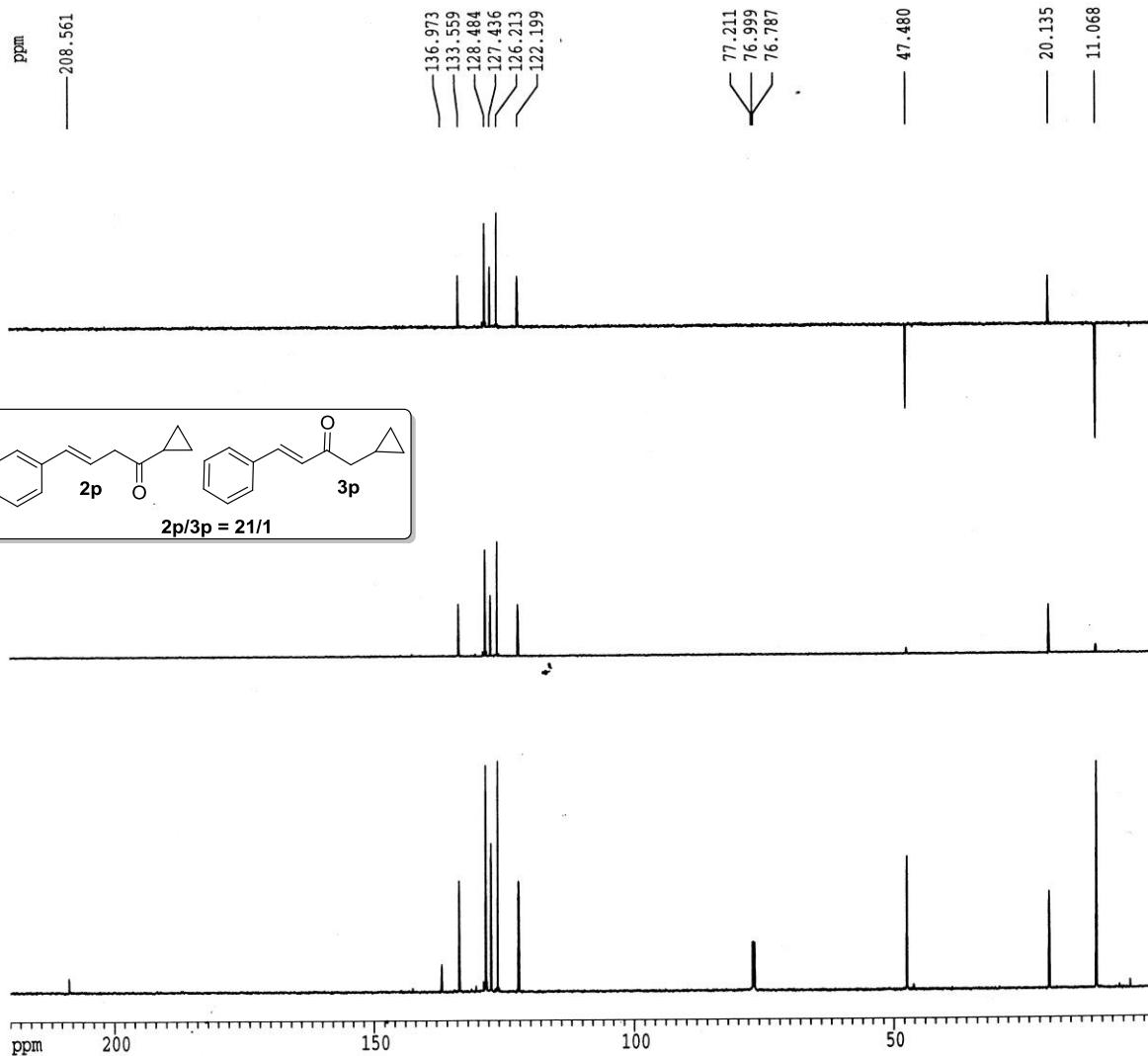
F2 - Acquisition Parameters
Date_ 20140522
Time 20.15
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zpgc
TD 32768
SOLVENT CDCl3
NS 200
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 299.4 K
D1 3.5000000 sec
d11 0.0300000 sec
DELTA 3.4000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

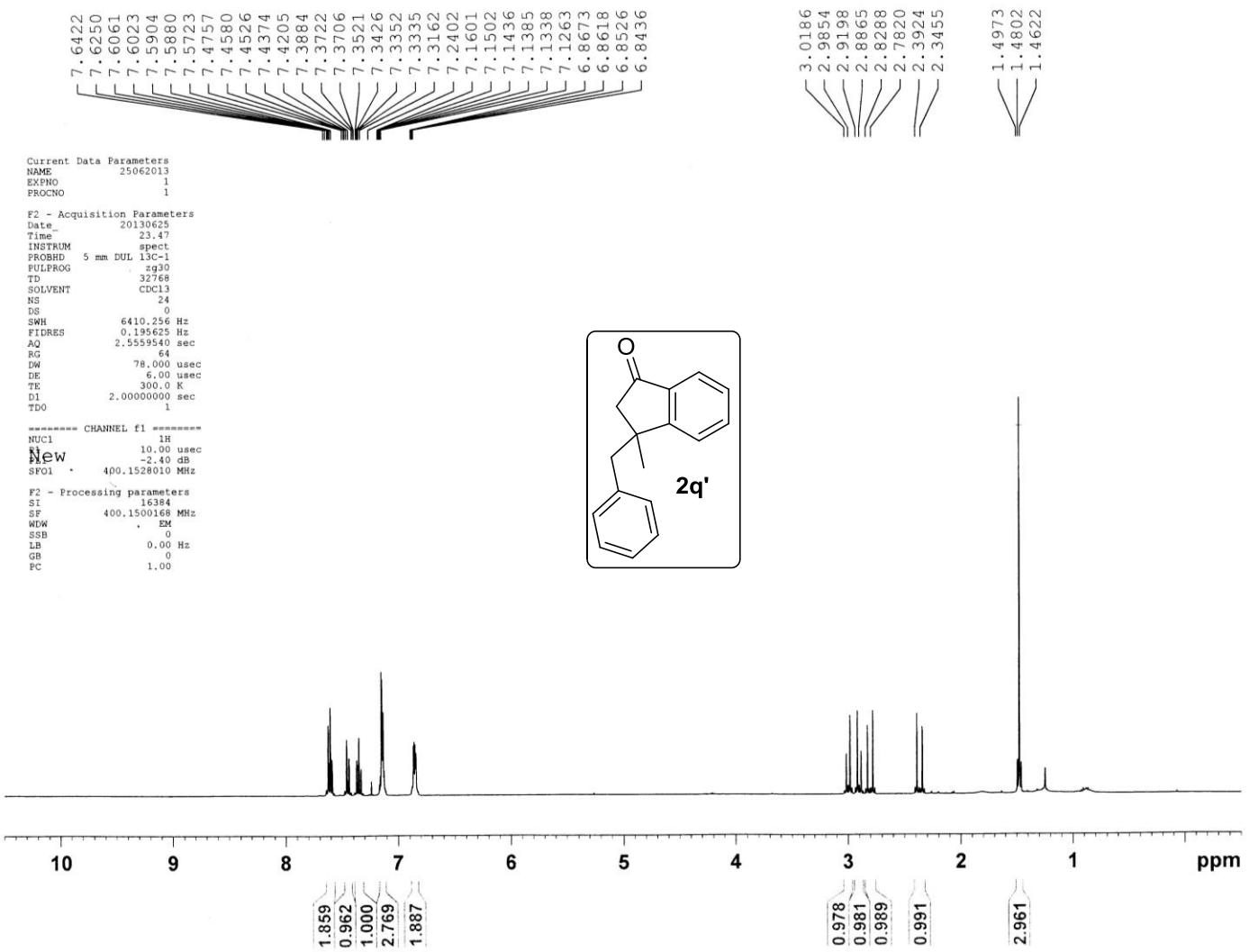
===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SF01 150.5597948 MHz

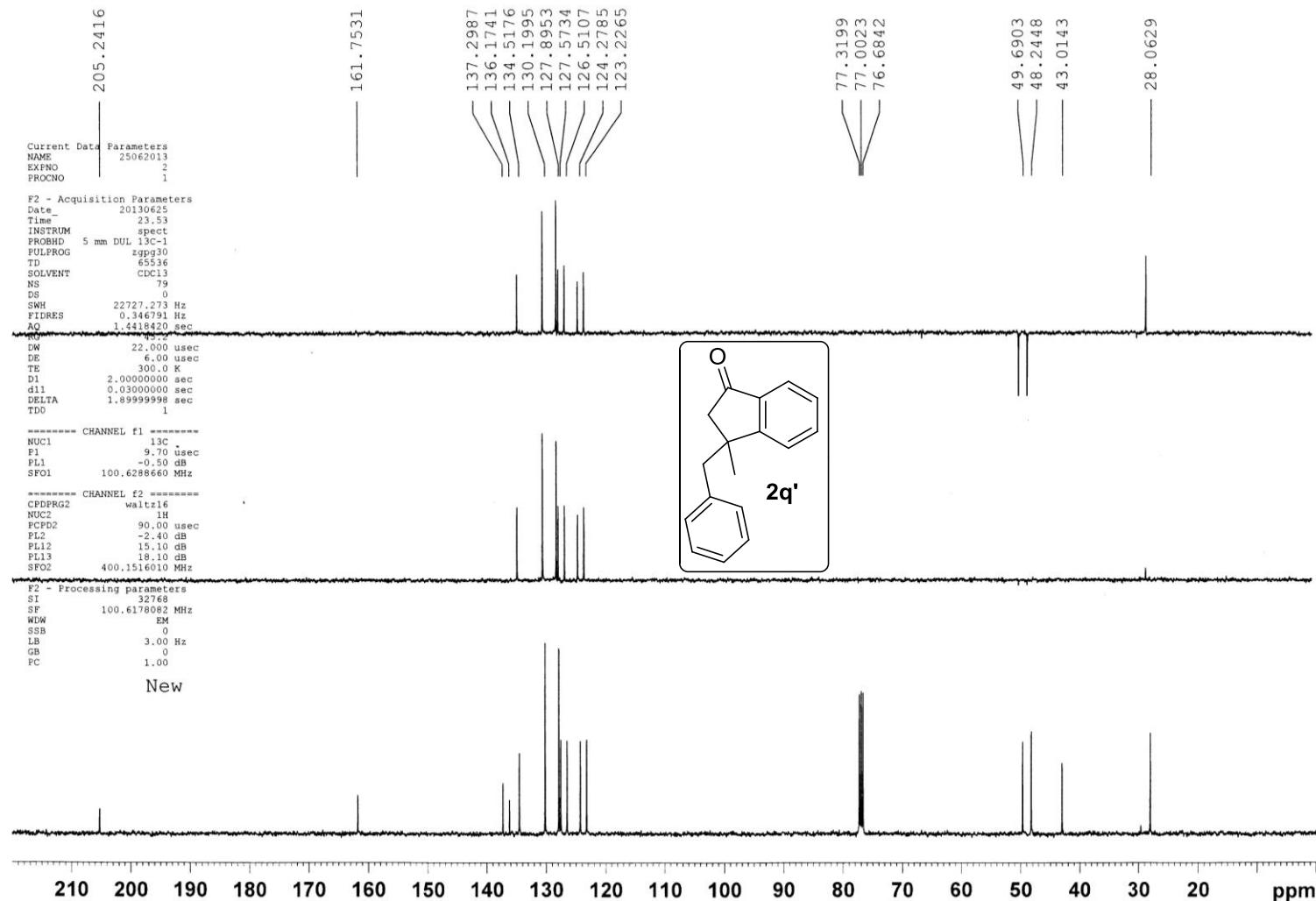
===== CHANNEL f2 =====
CPDPGR2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SF02 598.7029935 MHz

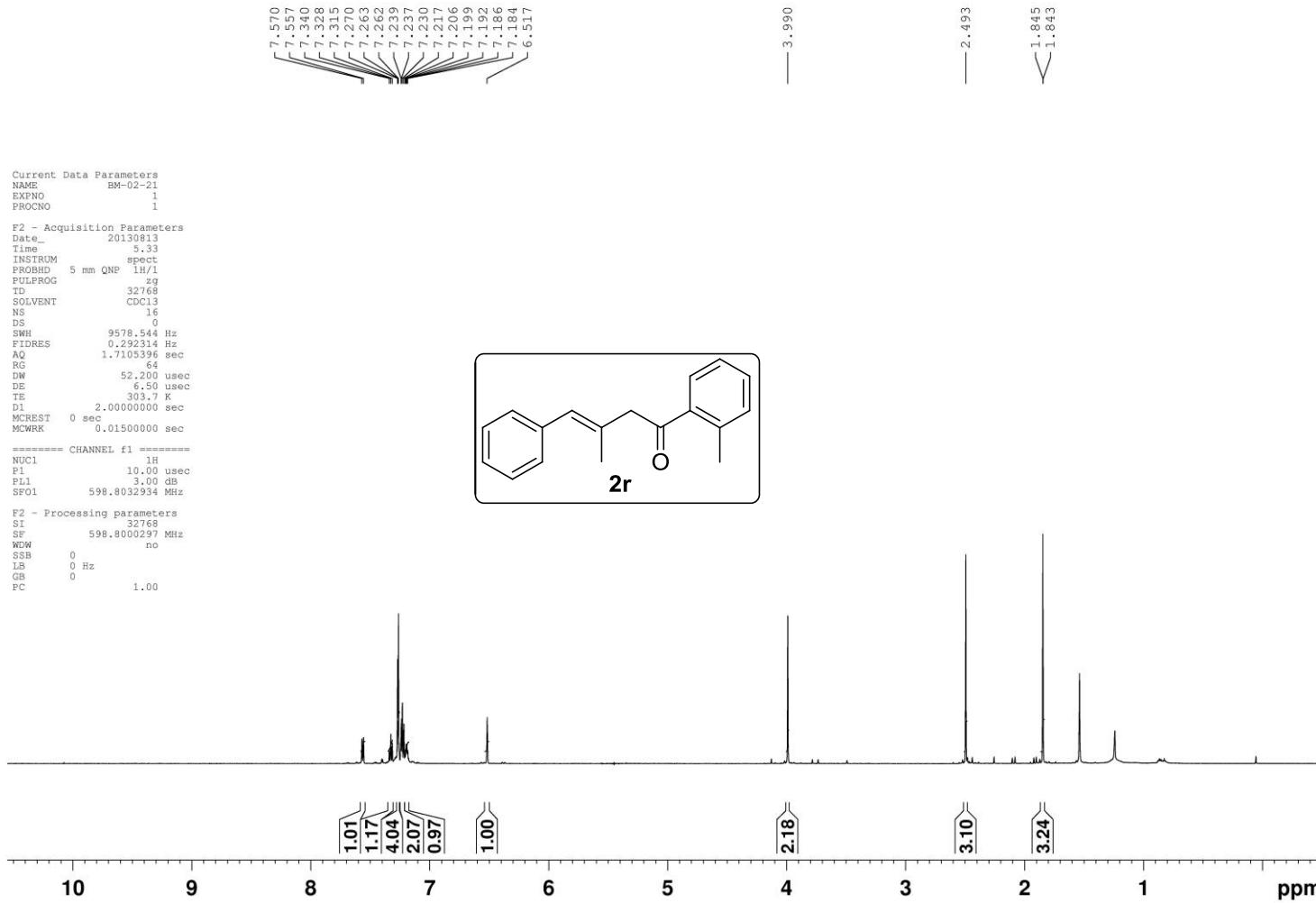
F2 - Processing parameters
SI 65536
SF 150.5432445 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 0.50

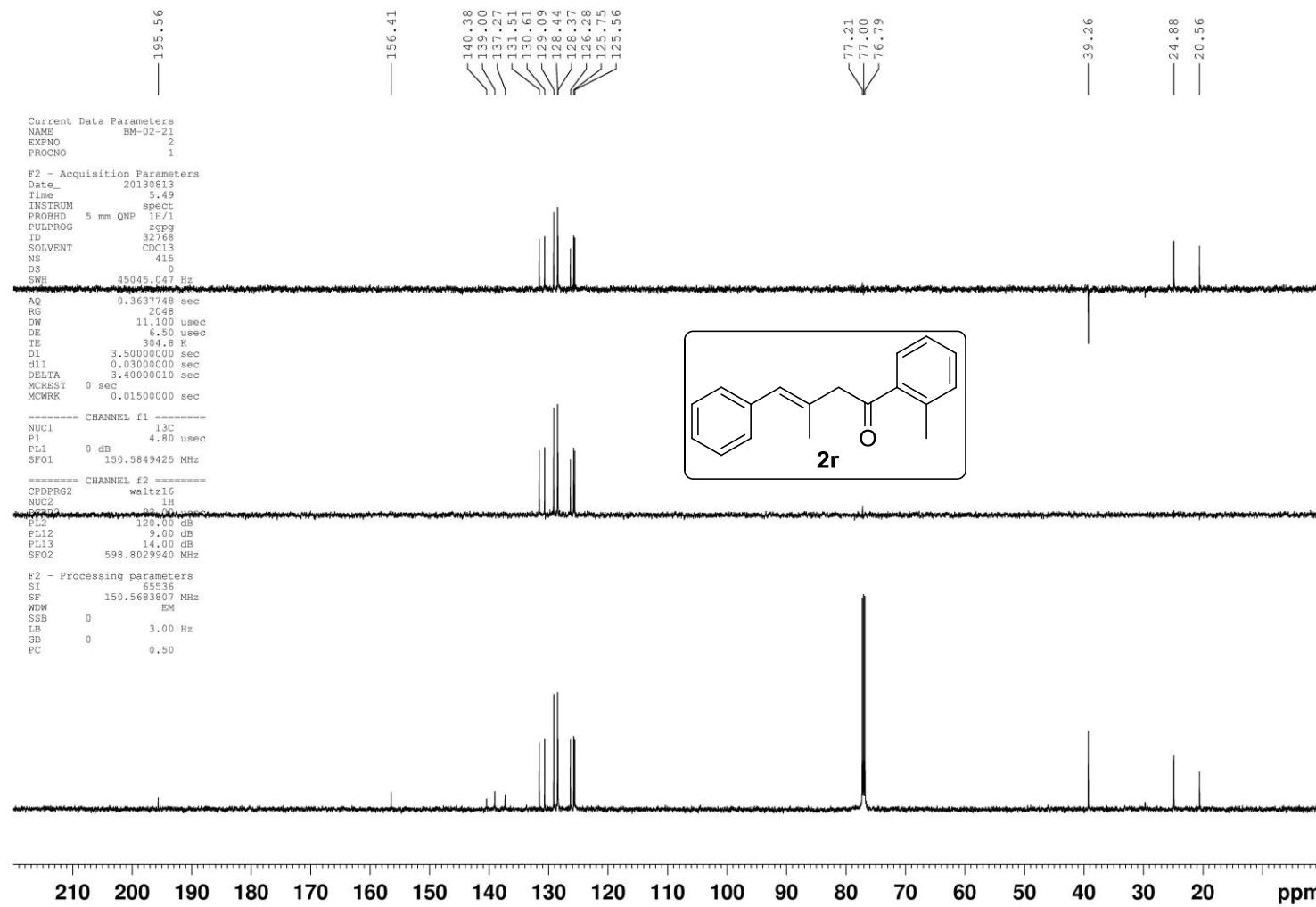
1D NMR plot parameters
CX 20.00 cm
CY 4.00 cm
F1P 220.000 ppm
F1 33119.51 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPCM 11.00000 ppm/cm
HZCM 1655.97571 Hz/cm











Current Data Parameters
NAME BM-86
EXPNO 1
PROCNO 1

```

F2 - Acquisition Parameters
Date_      20130501
Time       14.14
INSTRUM   spect
PROBHD   5 mm QNP 1H/1
PULPROG  zg
TD        32768
SOLVENT    CDCl3
NS         16
DS          0
SWH       8382.229 Hz
FIDRES   0.255805 Hz
AQ        1.9546611 sec
RG          128
DW        59.650 usec
DE        6.50 usec
TE        296.5 K
D1    1.0000000 sec
MCREST   0.0000000 sec
MCRWKR  0.01500000 sec

```

===== CHANNEL f1 =====
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SF01 598.8032934 MHz

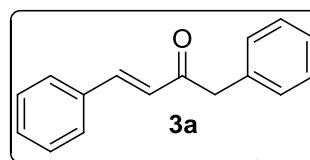
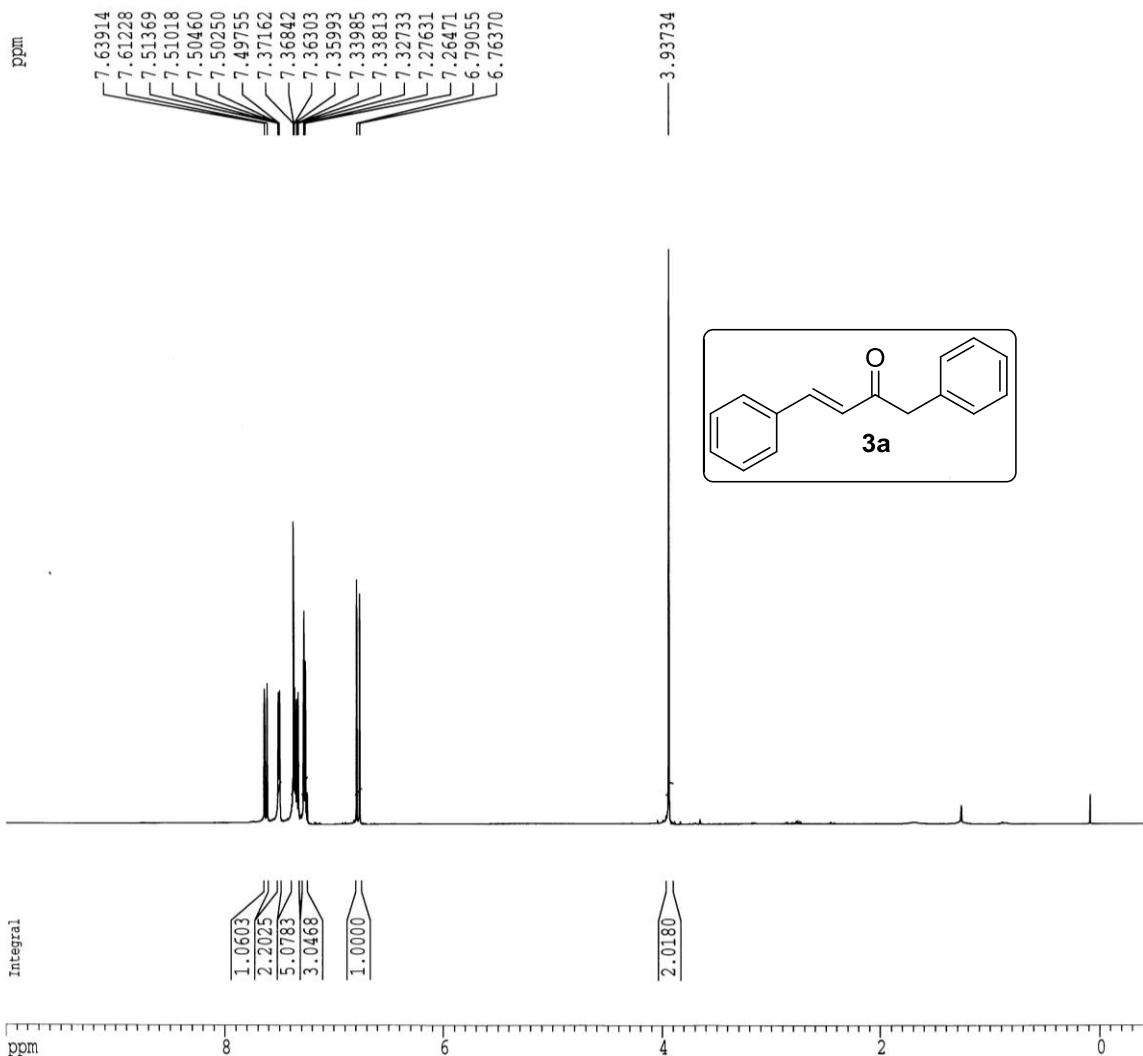
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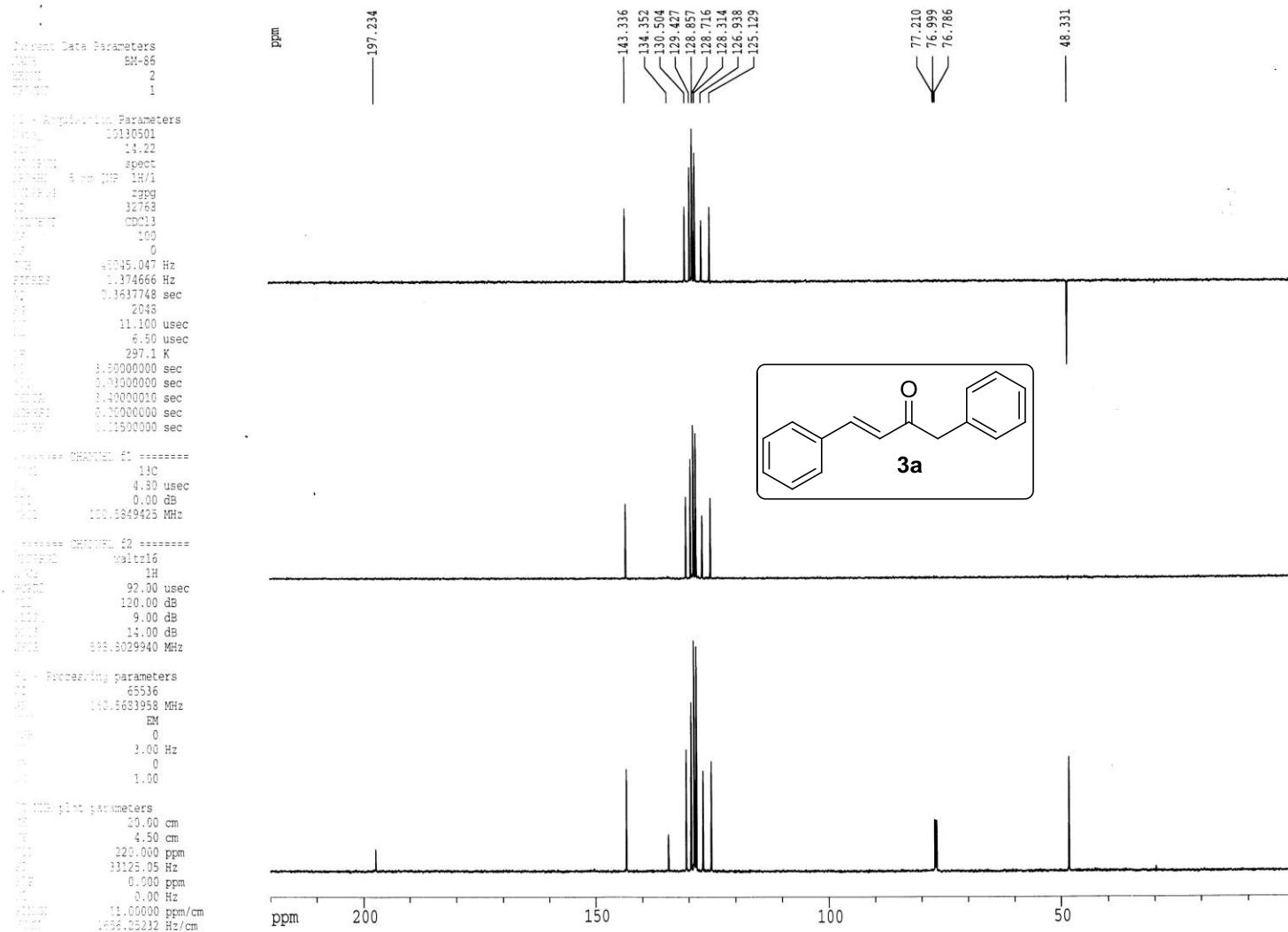
F2 - Processing parameters
SI          32768
SF          598.8000286 MHz
WDW         no
SSB         0
LB          0.00 Hz
GB         0
PC          1.00

```

1D NMR plot parameters

CX	20.00	cm
CY	10.00	cm
F1P	10.000	ppm
F1	5988.00	Hz
F2P	-0.500	ppm
F2	-299.40	Hz
PPCMC	0.52500	ppm/cm
HZCM	314.37003	Hz/cm





Current Data Parameters
NAME BM-105
EXPNO 1
PROCNO 1

```

F2 - Acquisition Parameters
Date_   20130501
Time    14.34
INSTRUM spect
PROBHD  5 mm QNP 1H/1
PULPROG zg
TD      32768
SOLVENT CDC13
NS      16
DS      0
SWH    8382.229 Hz
FIDRES 0.255805 Hz
AQ     1.954661 sec
RG      128
DW      59.650 usec
DE      6.50 usec
TE      296.3 K
D1      1.0000000 sec
MCREST 0.0000000 sec
MCWRK  0.0150000 sec

```

===== CHANNEL f1 =====
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SF01 598.8032934 MHz

```

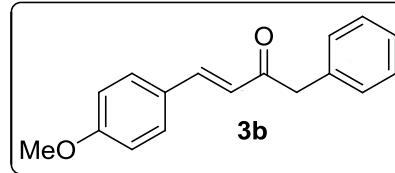
F2 - Processing parameters
SI          32768
SF         598.8000284 MHz
WDW        no
SSB        0
LB          0.00 Hz
GB        0
PC        1.00

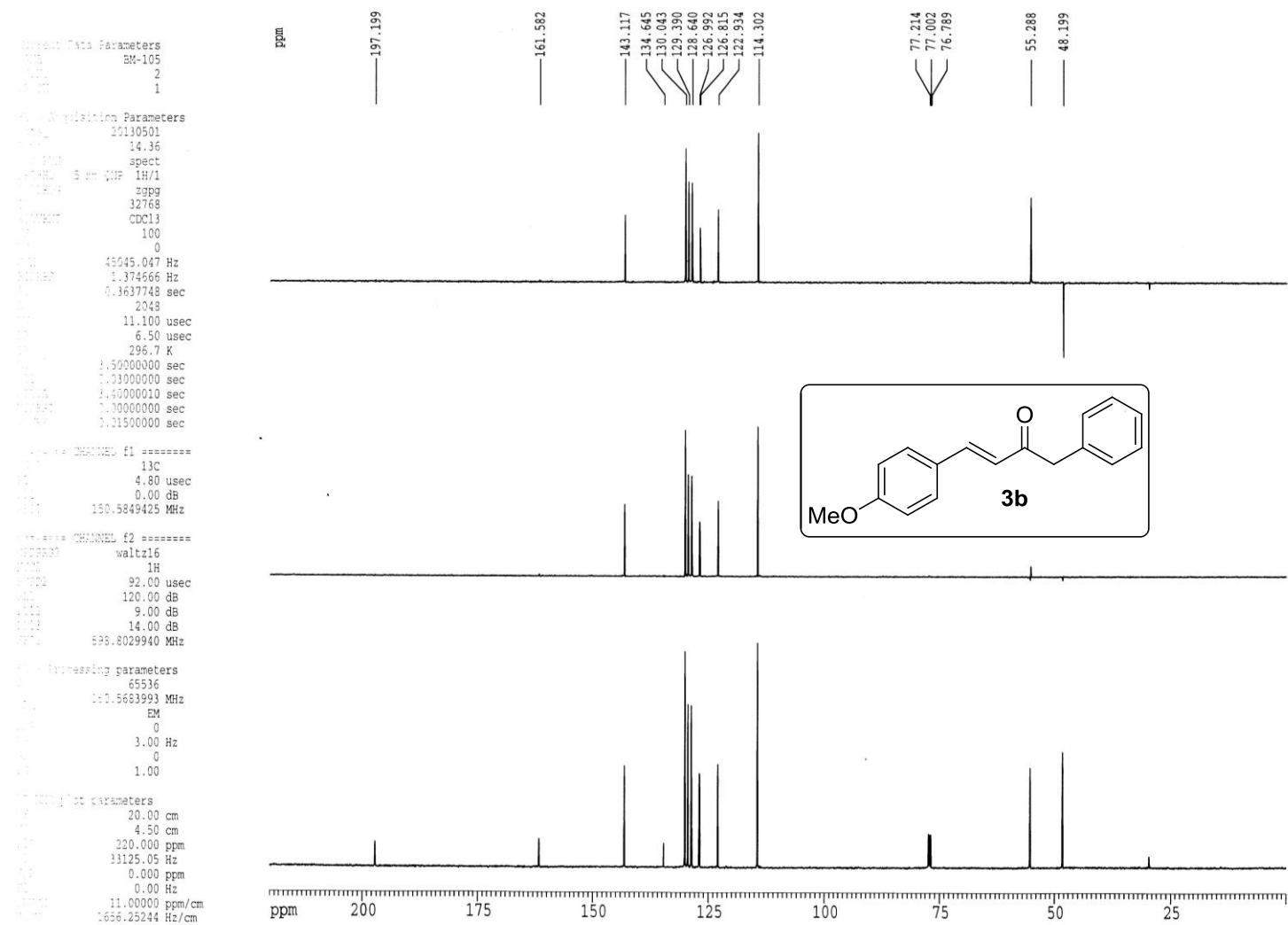
```

```

1D NMR plot parameters
CX           20.00 cm
CY           15.00 cm
F1P          10.000 ppm
F1           5988.00 Hz
F2P          -0.500 ppm
F2           -299.40 Hz
PPMCM        0.52500 ppm/cm
HzCM         314.37003 Hz/cm

```





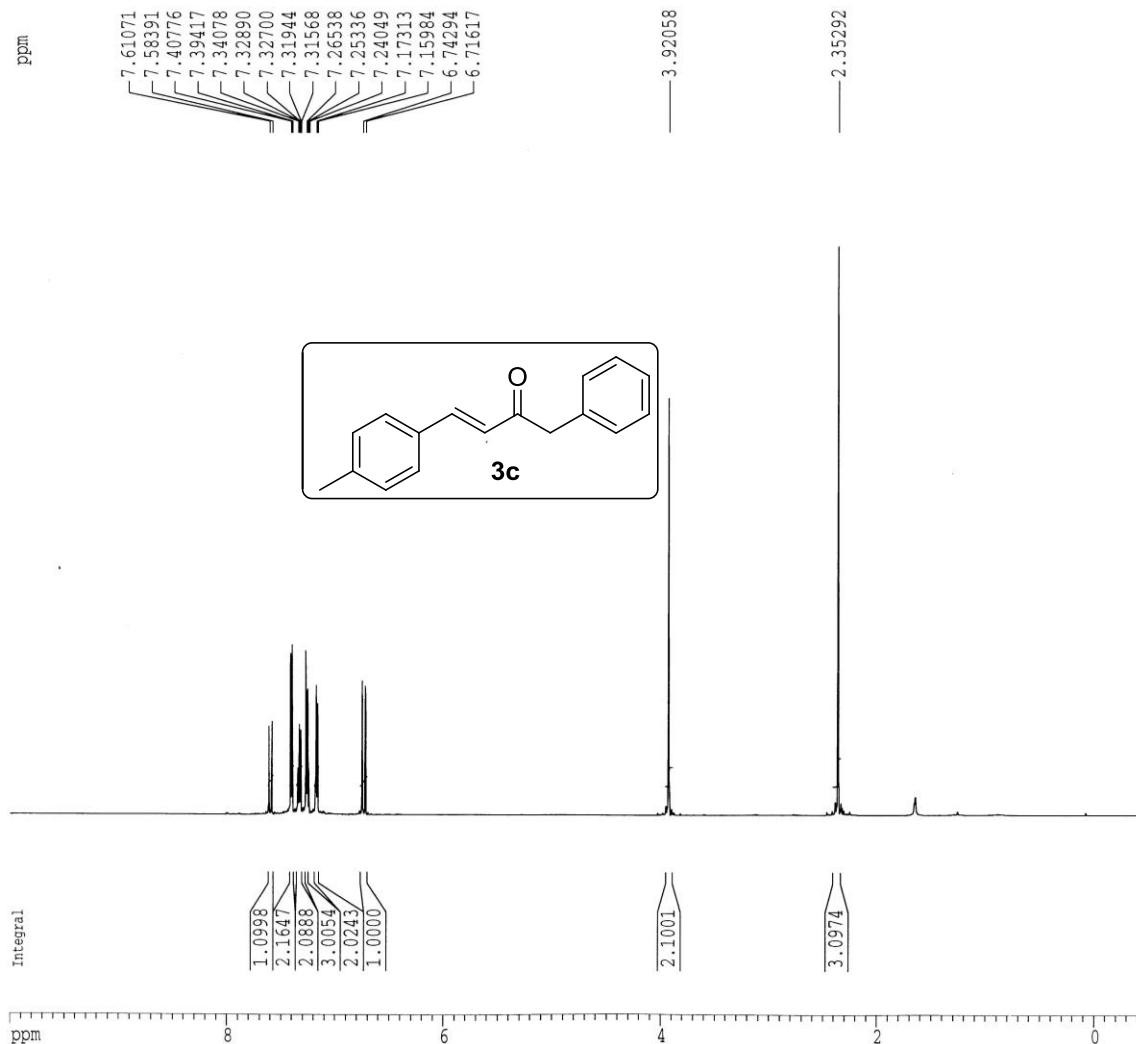
Current Data Parameters
NAME BM-123
EXPNO 1
PROCNO 1

F2 - Acquisition Parameters
Date_ 20110509
Time 15.28
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 32768
SOLVENT CDCl3
NS 16
DS 0
SWH 8382.229 Hz
FIDRES 0.255805 Hz
AQ 1.9546613 sec
RG 128
DM 59.650 usec
DE 6.50 usec
TE 298.7 K
D1 2.0000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

===== CHANNEL f1 =====
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SFO1 598.8029940 MHz

F2 - Processing parameters
SI 32768
SF 598.8000283 MHz
WDW no
SSB 0
LB 0.00 Hz
GB 0
PC 1.00

1D NMR plot parameters
CX 20.00 cm
CY 10.00 cm
F1P 10.000 ppm
F1 5988.00 Hz
F2P -0.500 ppm
F2 -299.40 Hz
PPCM 0.52500 ppm/cm
HZCM 314.37003 Hz/cm



Current Data Parameters
NAME EM-123
EXPNO 2
PROCNO 1

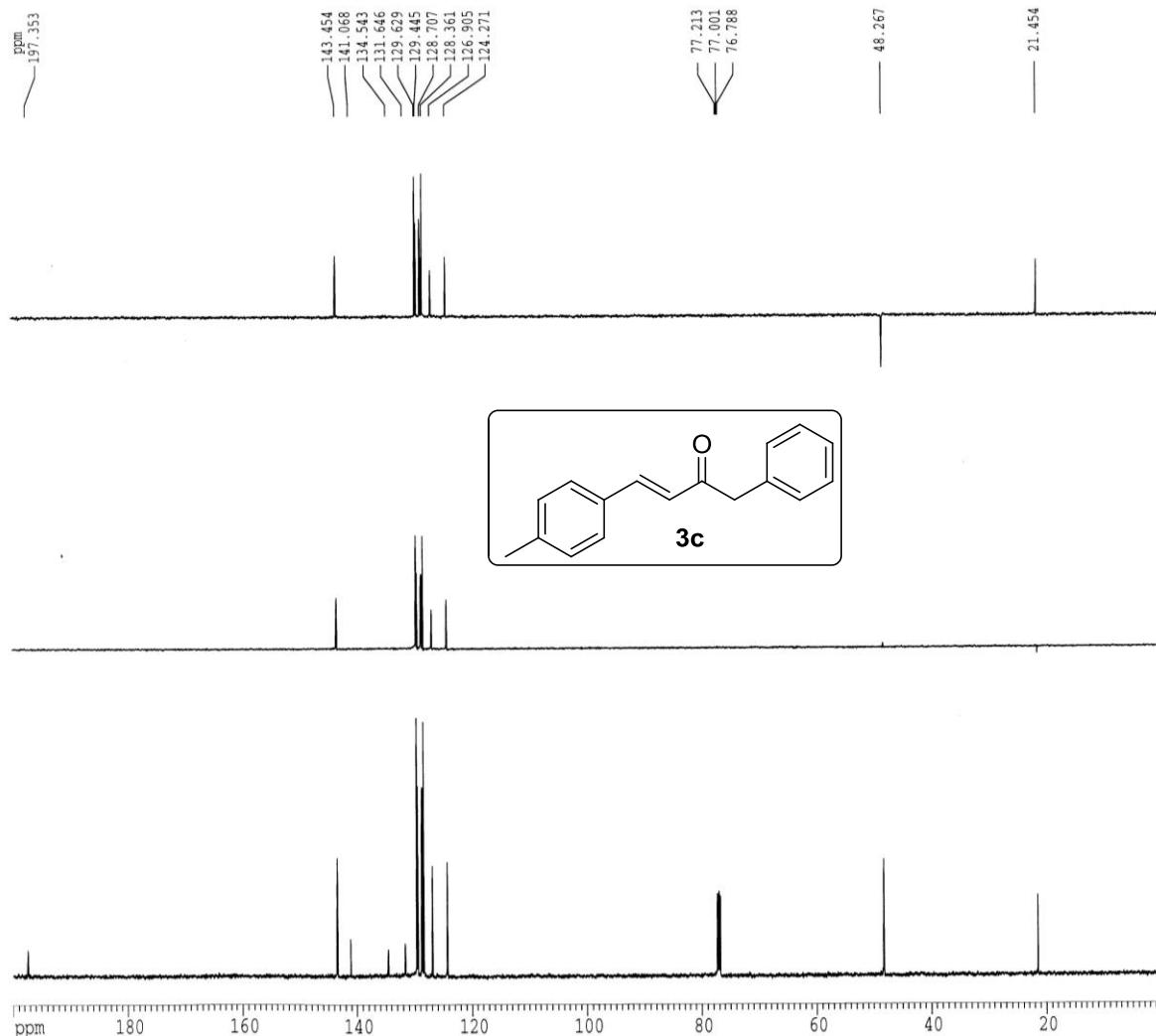
F2 - Acquisition Parameters
Date_ 20110509
Time 15:34
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT CDCl3
NS 100
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 299.5 K
D1 3.5000000 sec
d11 0.0300000 sec
DELTA 3.40000010 sec
MCREST 0.0000000 sec
MCRWKR 0.0150000 sec

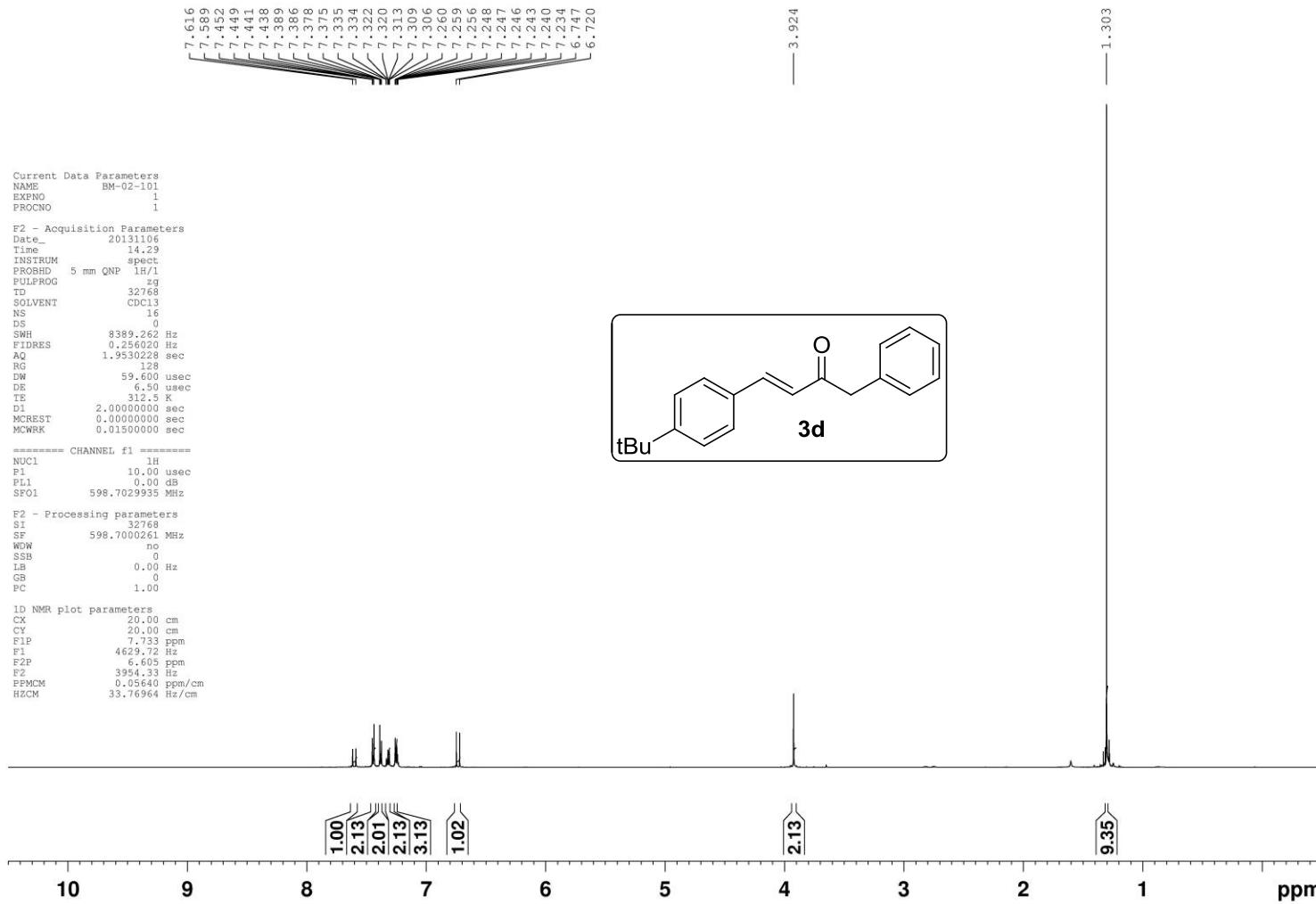
***** CHANNEL f1 *****
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SF01 150.5849425 MHz

***** CHANNEL f2 *****
CPDPRG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SF02 598.8029940 MHz

F2 - Processing parameters
SI 65536
SF 150.5683896 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 1.00

1D NMR plot parameters
CX 20.00 cm
CY 4.50 cm
FLP 300.000 ppm
F1 30113.68 Hz
P2P 0.00 ppm
F2 0.00 Hz
PPMCM 10.0000 ppm/cm
HZCM 1505.68384 Hz/cm





Current Data Parameters
NAME BM-02-101
EXPNO 2
PROCNO 1

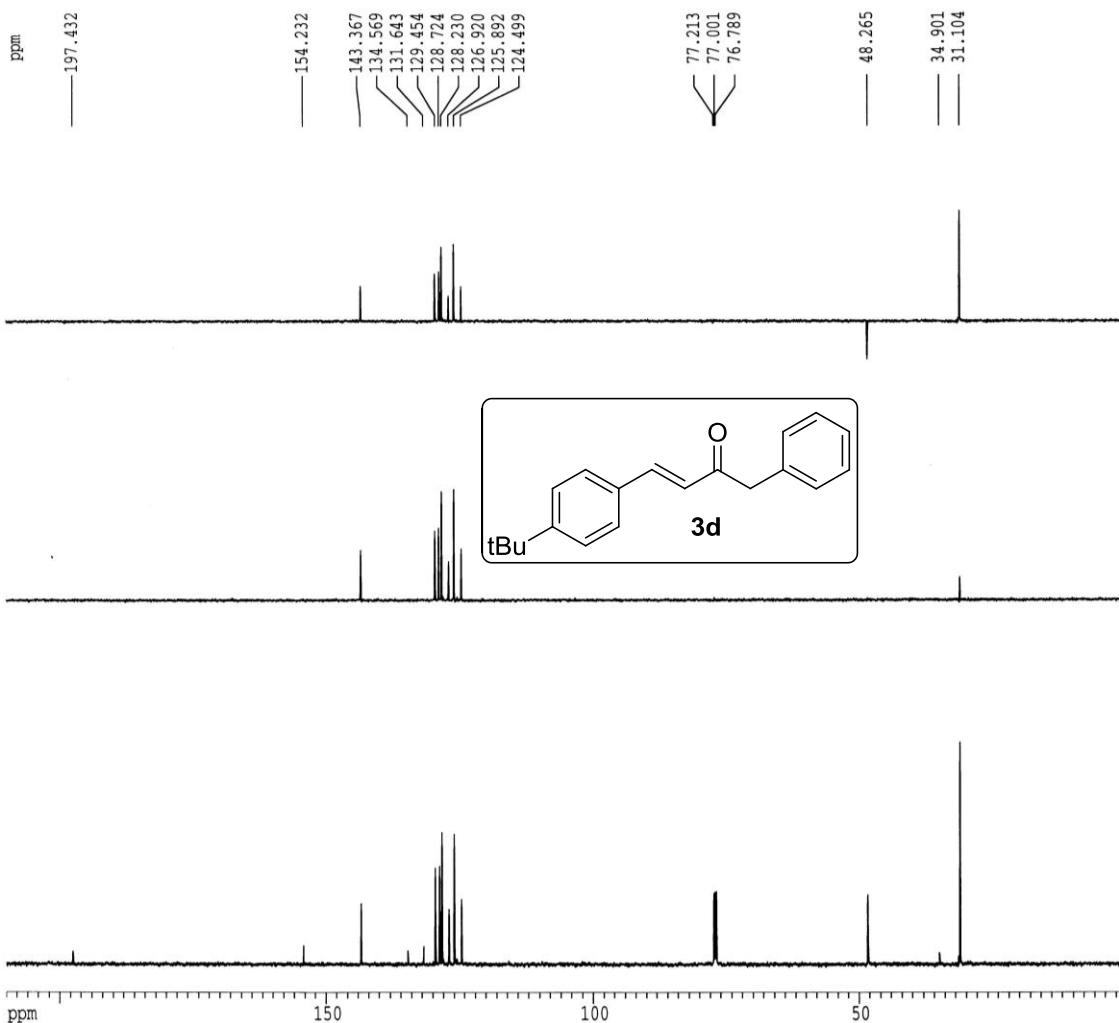
F2 - Acquisition Parameters
Date_ 20131106
Time 14.36
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT CDCl3
NS 100
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 313.7 K
D1 3.5000000 sec
d11 0.0300000 sec
DELTA 3.40000010 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

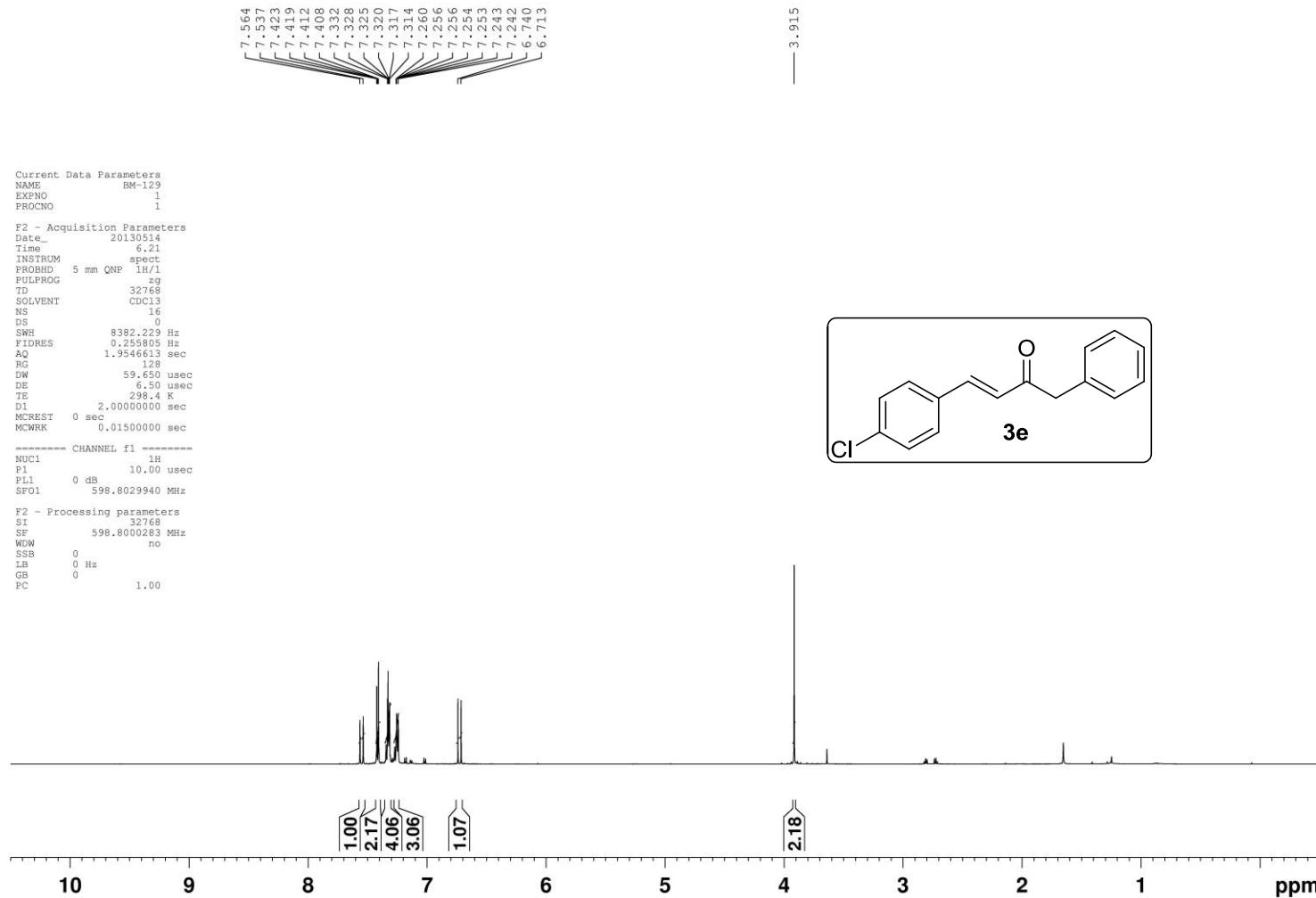
===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SFO1 150.5597948 MHz

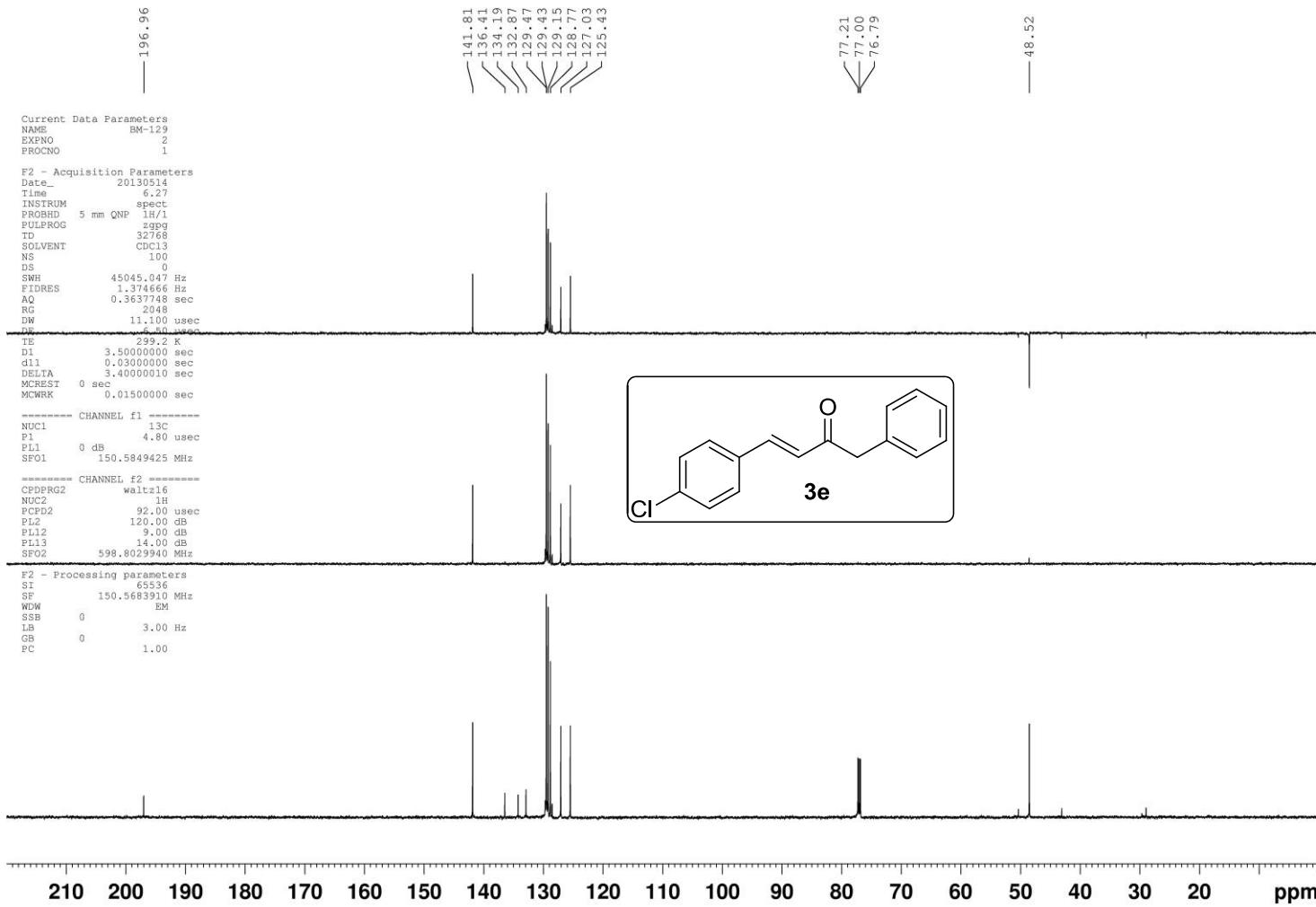
===== CHANNEL f2 =====
CPDPRG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SFO2 598.7029935 MHz

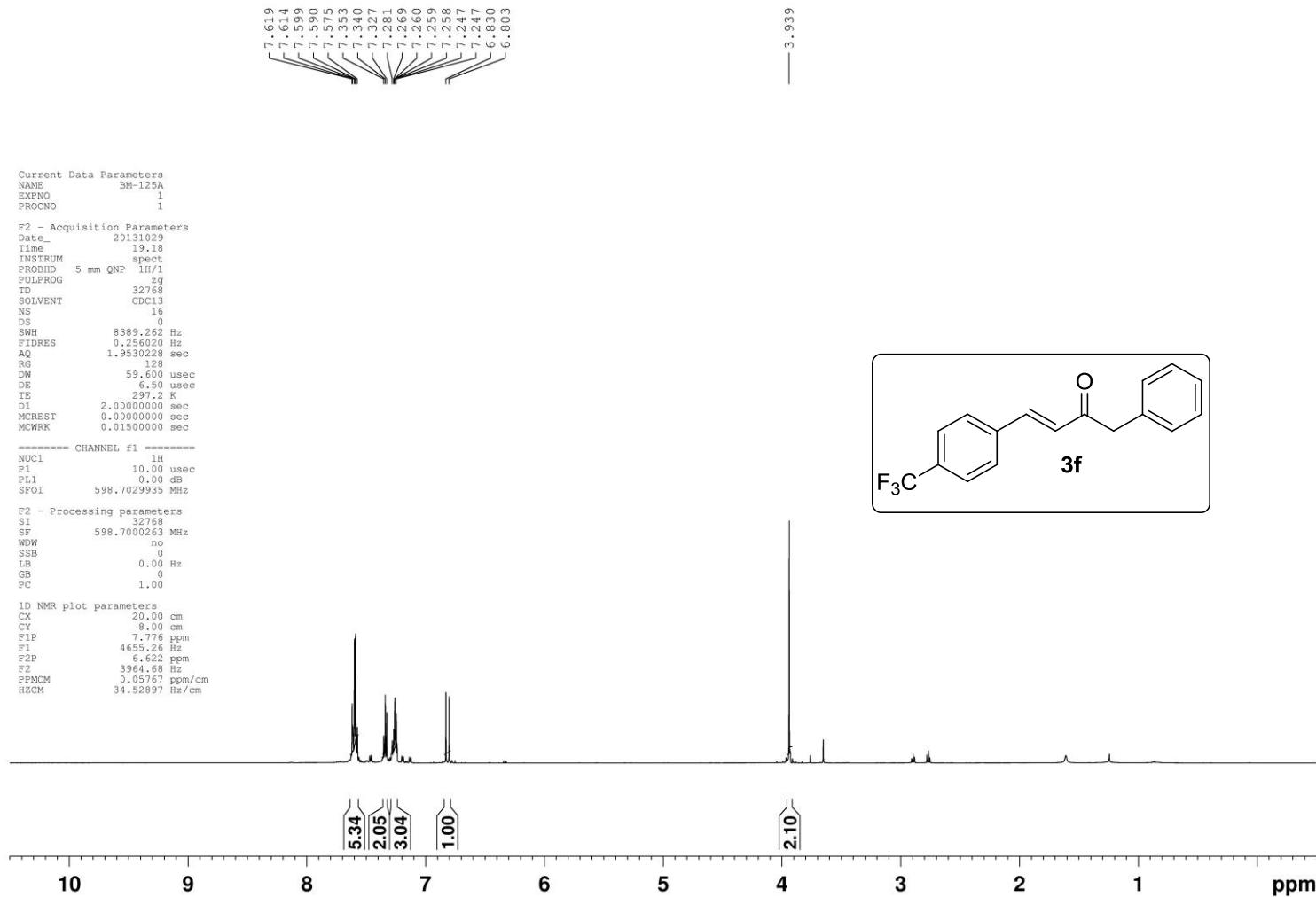
F2 - Processing parameters
SI 65536
SF 150.5432404 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 0.50

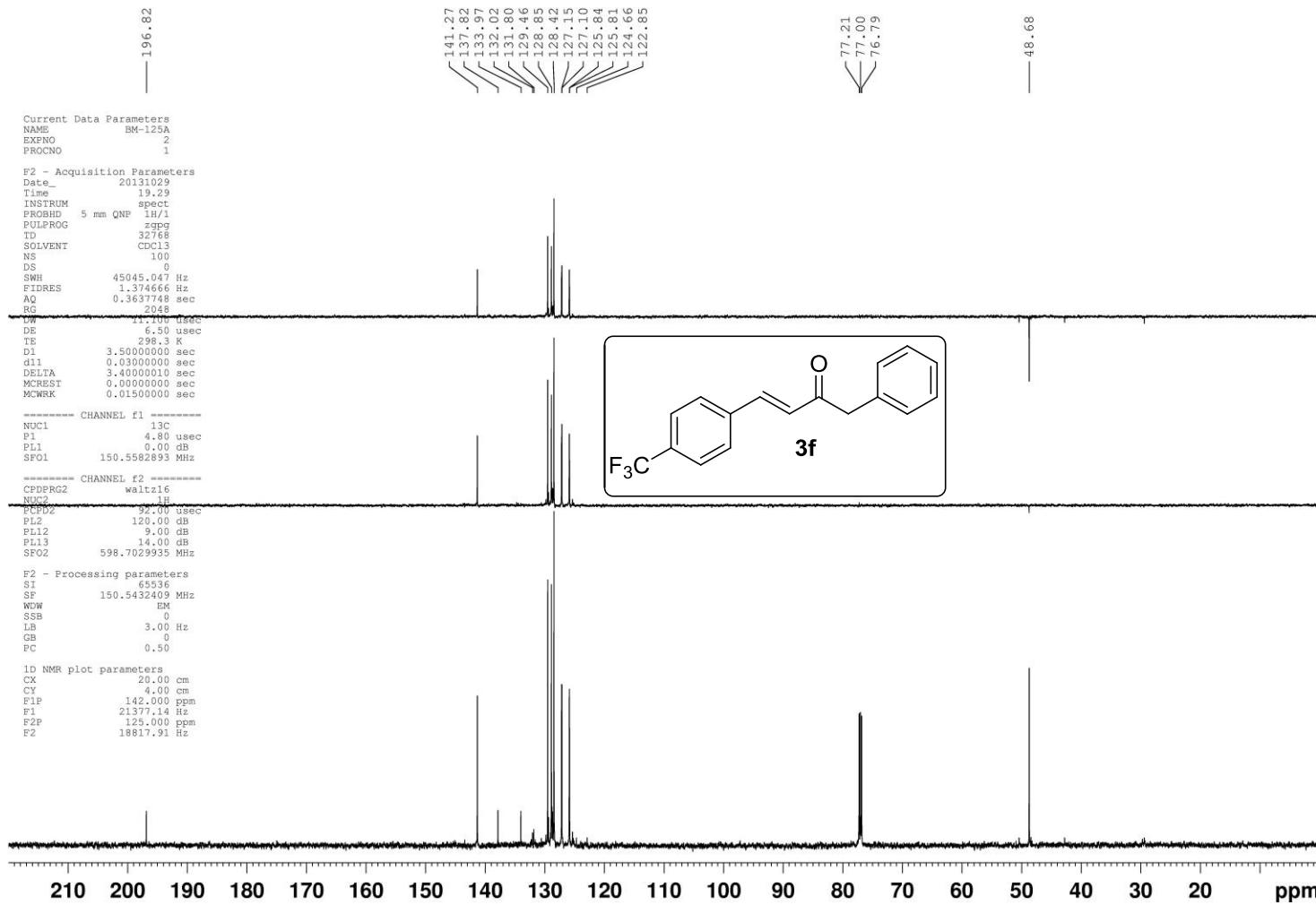
1D NMR plot parameters
CX 20.00 cm
CY 4.00 cm
F1P 210.000 ppm
F1 31614.08 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPMCM 10.50000 ppm/cm
HZCM 1580.70410 Hz/cm

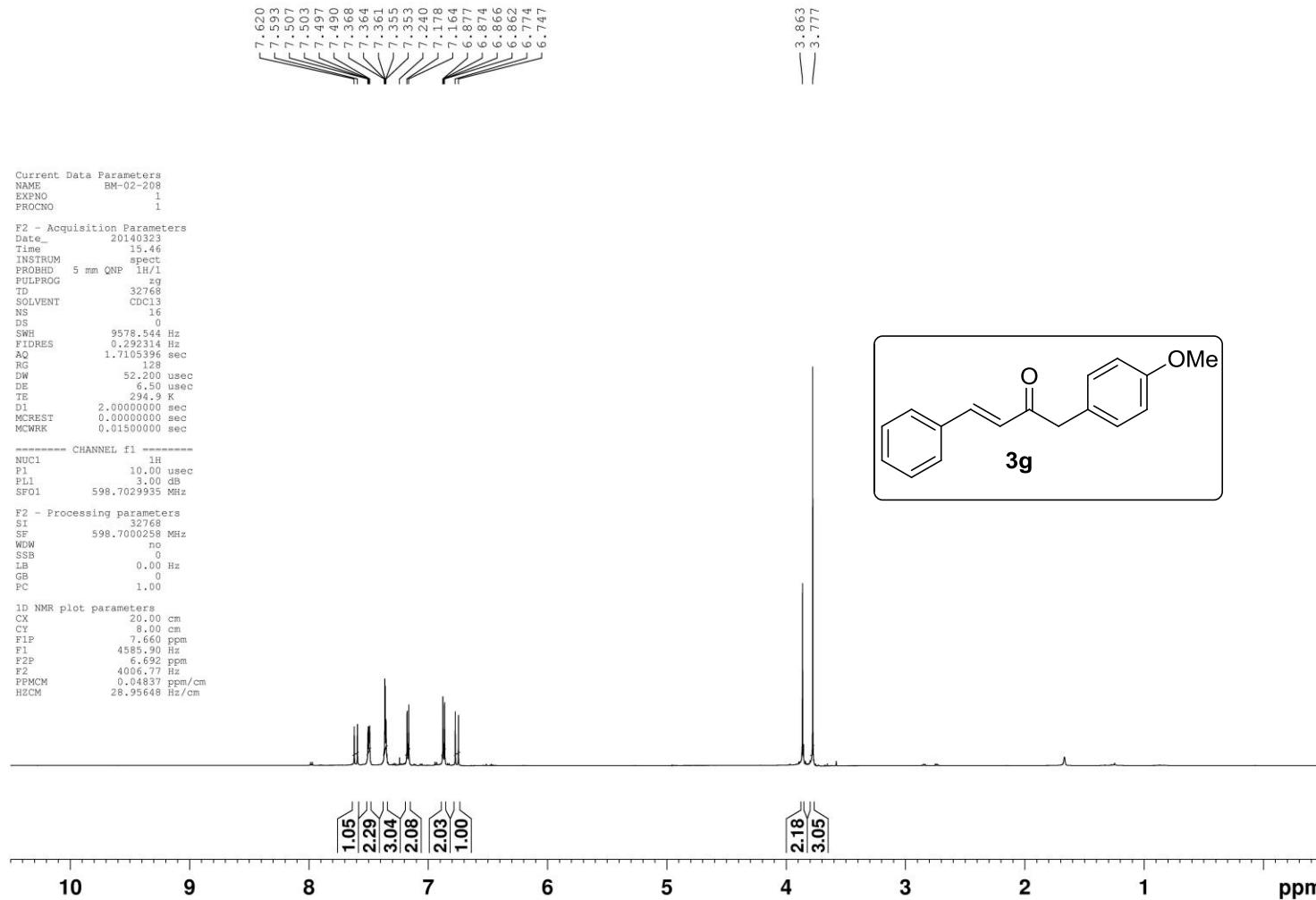


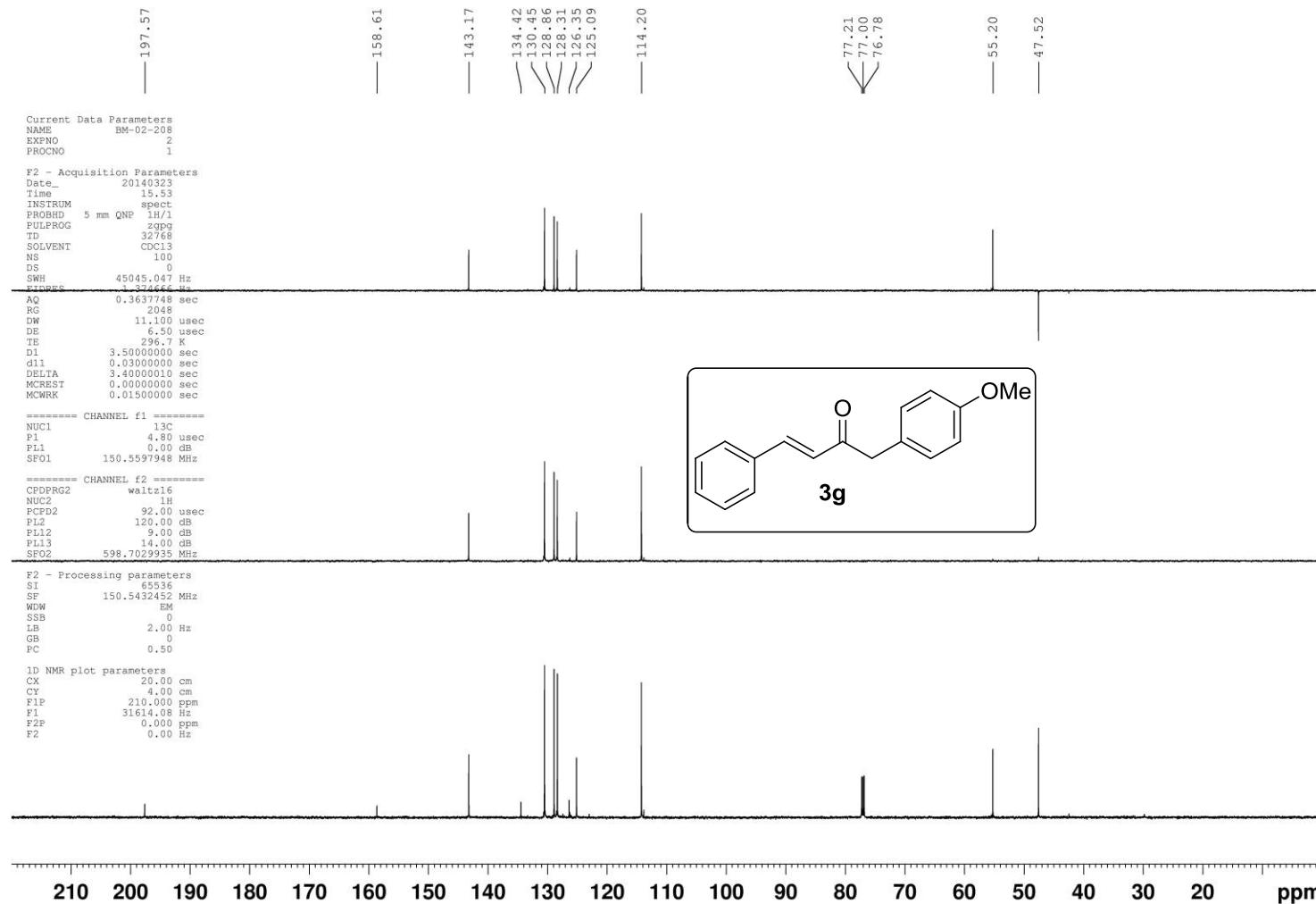


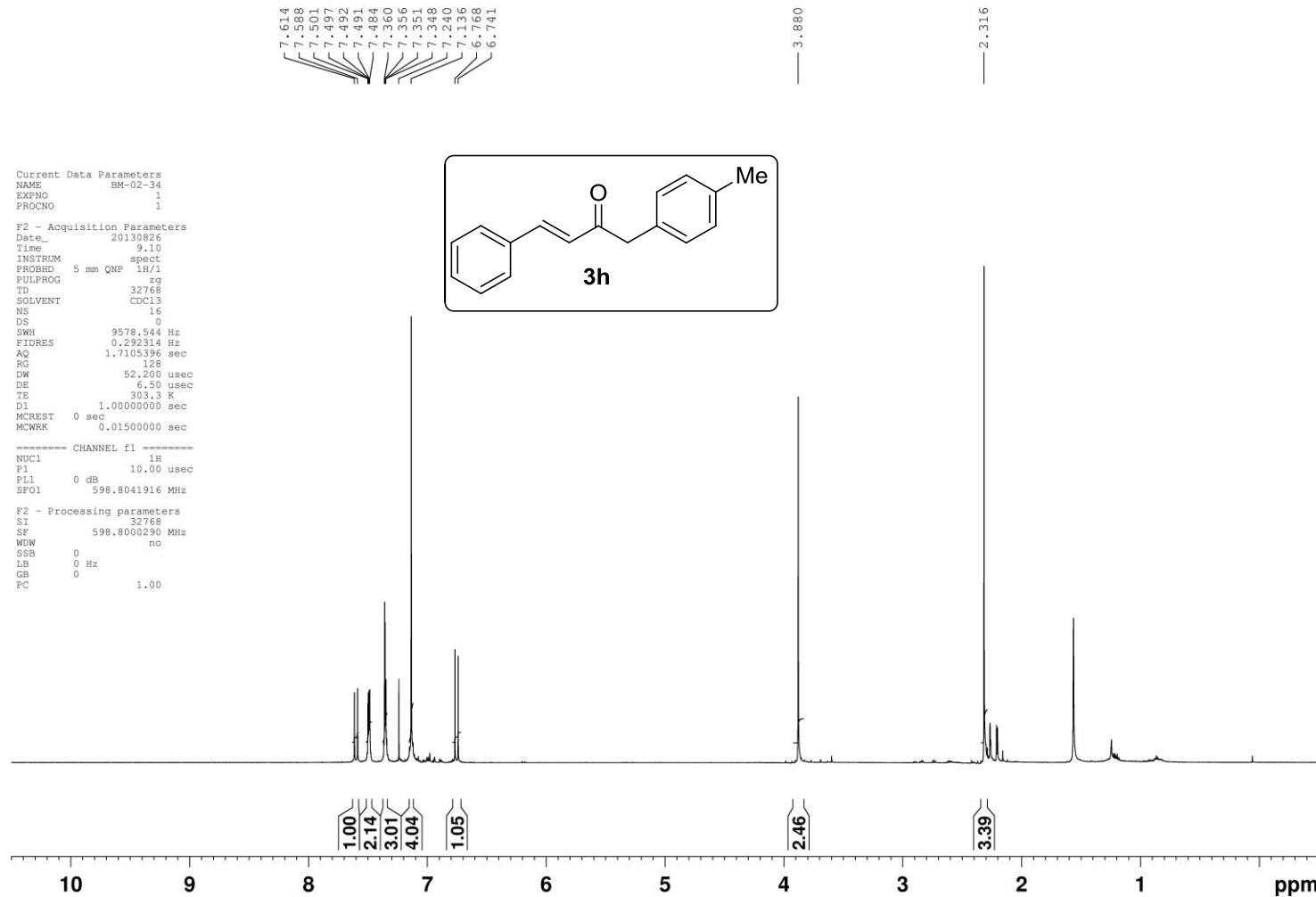


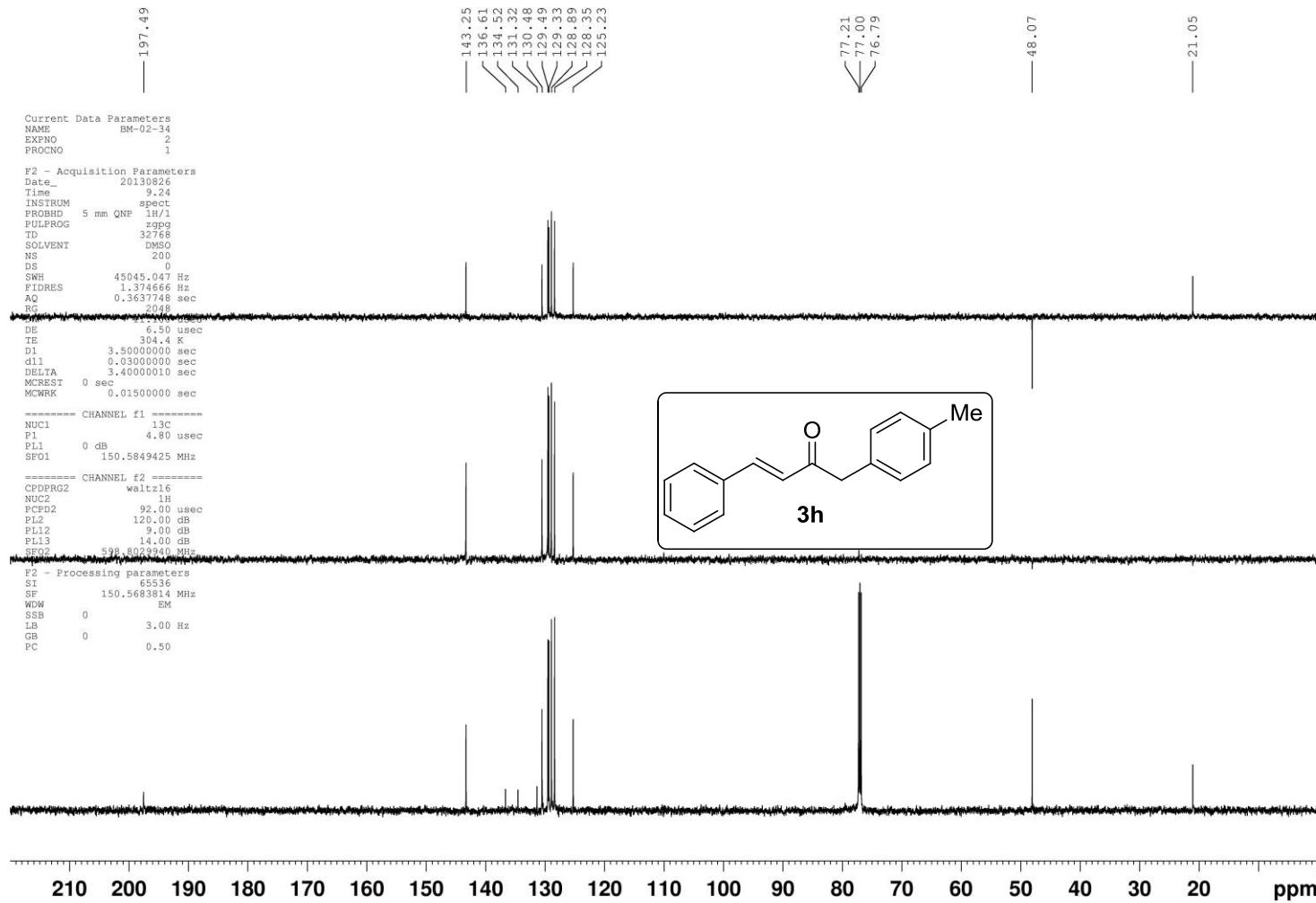












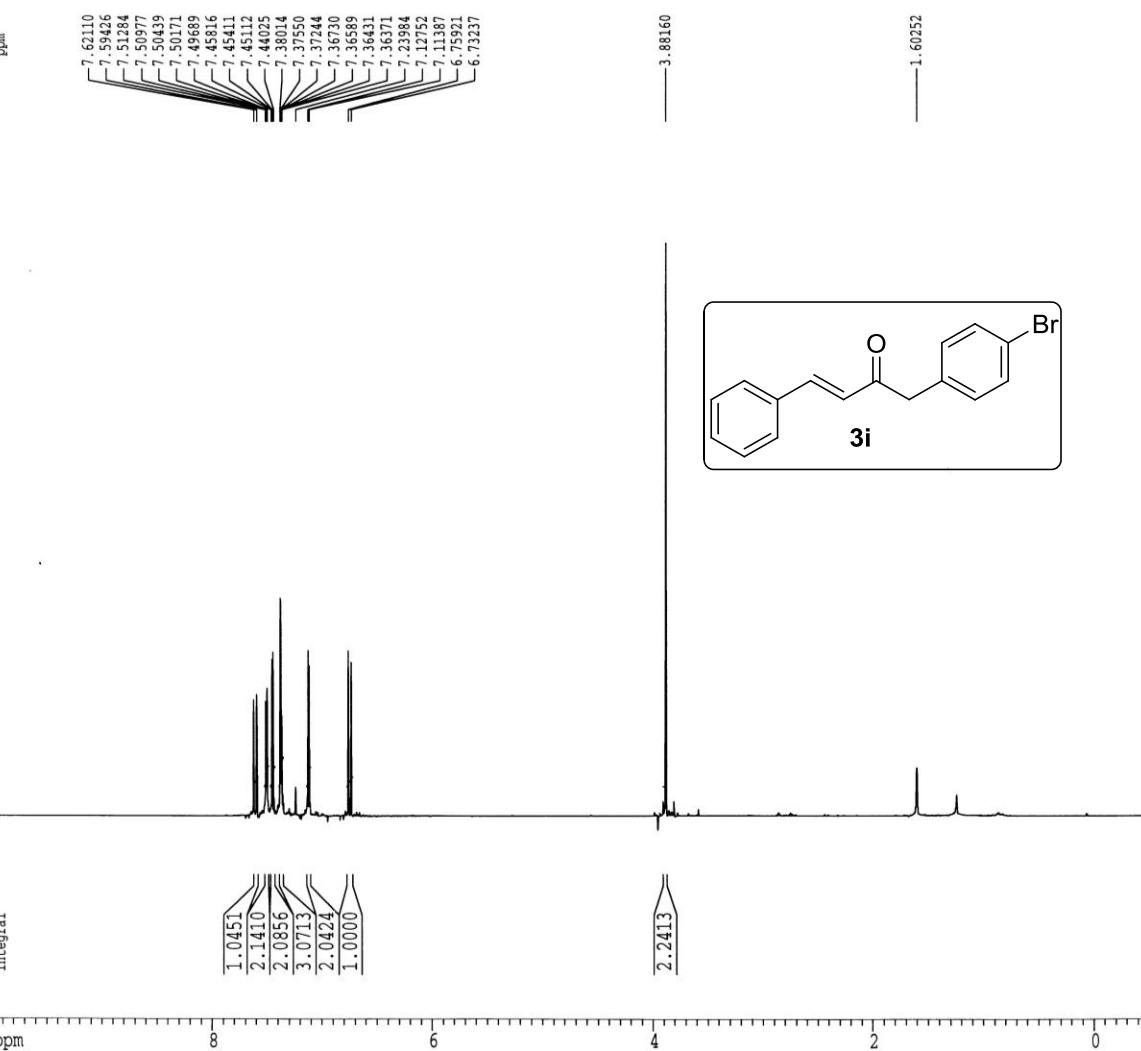
Current Data Parameters
 NAME BM-02-37
 EXPNO 1
 PROCNO 1

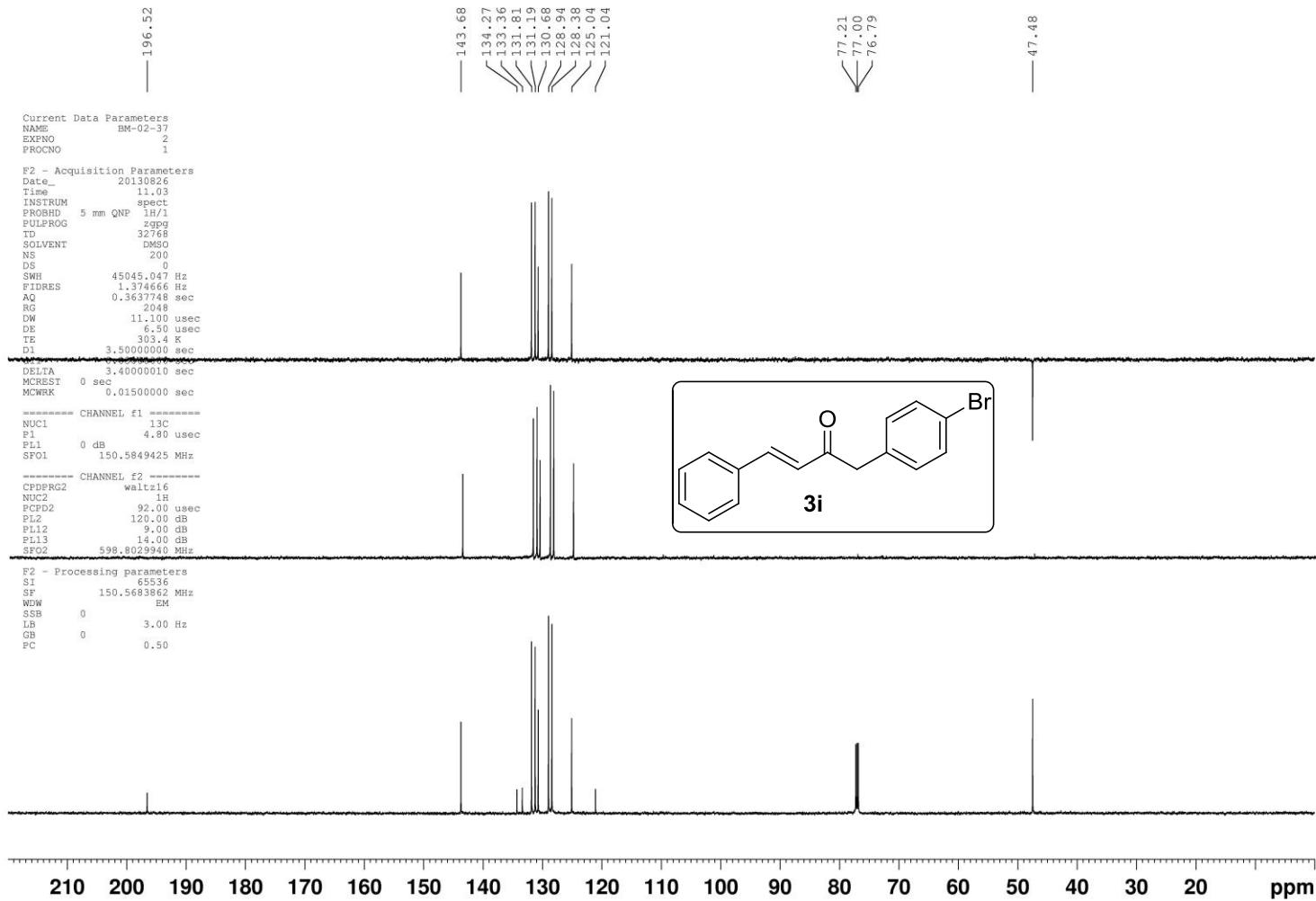
F2 - Acquisition Parameters
 Date_ 20130825
 Time 20.31
 INSTRUM spect
 PROBHD 5 mm QNP 1H/1
 PULPROG zg
 TD 32768
 SOLVENT CDCl3
 NS 16
 DS 0
 SWH 9578.544 Hz
 FIDRES 0.292314 Hz
 AQ 1.7105396 sec
 RG 64
 DW 52.200 usec
 DE 6.50 usec
 TE 303.8 K
 D1 1.0000000 sec
 MCREST 0.0000000 sec
 MCWRK 0.0150000 sec

===== CHANNEL f1 =====
 NUC1 1H
 P1 10.00 usec
 PL1 0.00 dB
 SF01 598.8041916 MHz

F2 - Processing parameters
 SI 32768
 SF 598.8000288 MHz
 WDW no
 SSB 0
 LB 0.00 Hz
 GB 0
 PC 1.00

1D NMR plot parameters
 CX 20.00 cm
 CY 10.00 cm
 F1P 10.000 ppm
 F1 5988.00 Hz
 F2P -0.500 ppm
 F2 -299.40 Hz
 PPMCM 0.52500 ppm/cm
 HZCM 314.37003 Hz/cm





Current Data Parameters
NAME BM-02-203
EXPNO 1
PROCNO 1

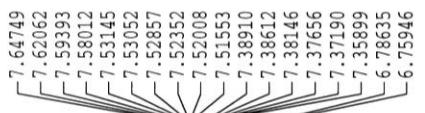
F1 - Acquisition Parameters
Date_ 20140318
Time 19.19
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 33556
SOLVENT CDCl3
NS 16
DS 0
SWH 8389.262 Hz
FIDRES 0.250008 Hz
AQ 1.9999876 sec
RG 128
DW 59.600 usec
DE 6.50 usec
TE 294.8 K
D1 2.0000000 sec
MCREST 0.0000000 sec
MCNMRK 0.0150000 sec

***** CHANNEL f1 *****
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SPO1 598.7029935 MHz

F1 - Processing parameters
SI 32768
SF 598.7000261 MHz
WDW no
SSB 0
LB 0.00 Hz
GB 0
PC 1.00

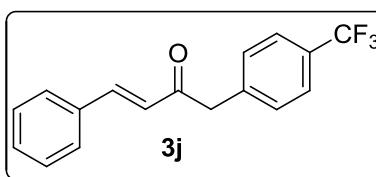
1D NMR plot parameters
CX 20.00 cm
CY 6.00 cm
F1P 10.000 ppm
F1 5987.00 Hz
F2P -0.500 ppm
F2 -299.35 Hz
PPMCM 0.52500 ppm/cm
HZCM 314.31750 Hz/cm

ppm



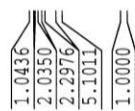
4.00783

-1.60692



Integral

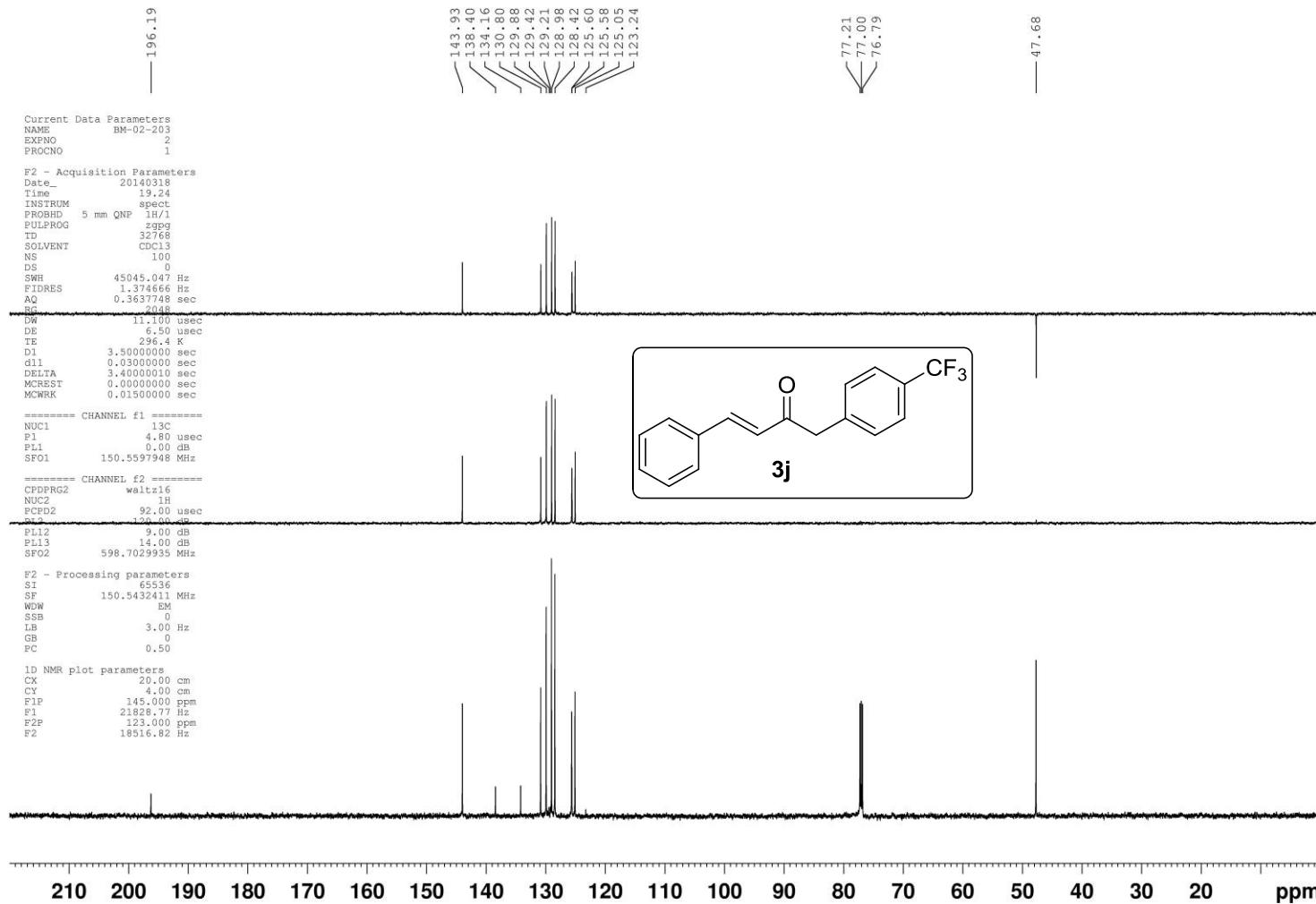
ppm

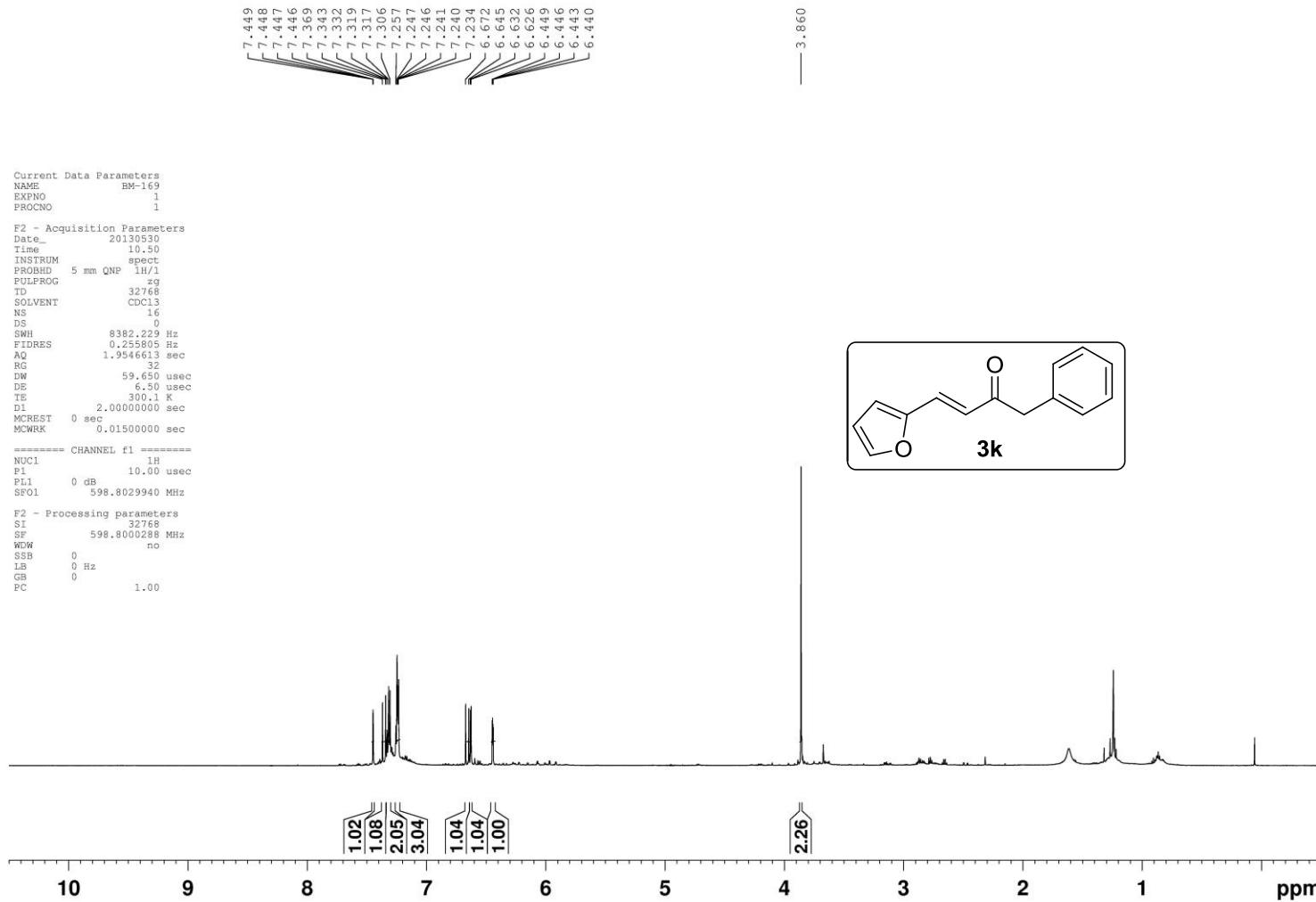


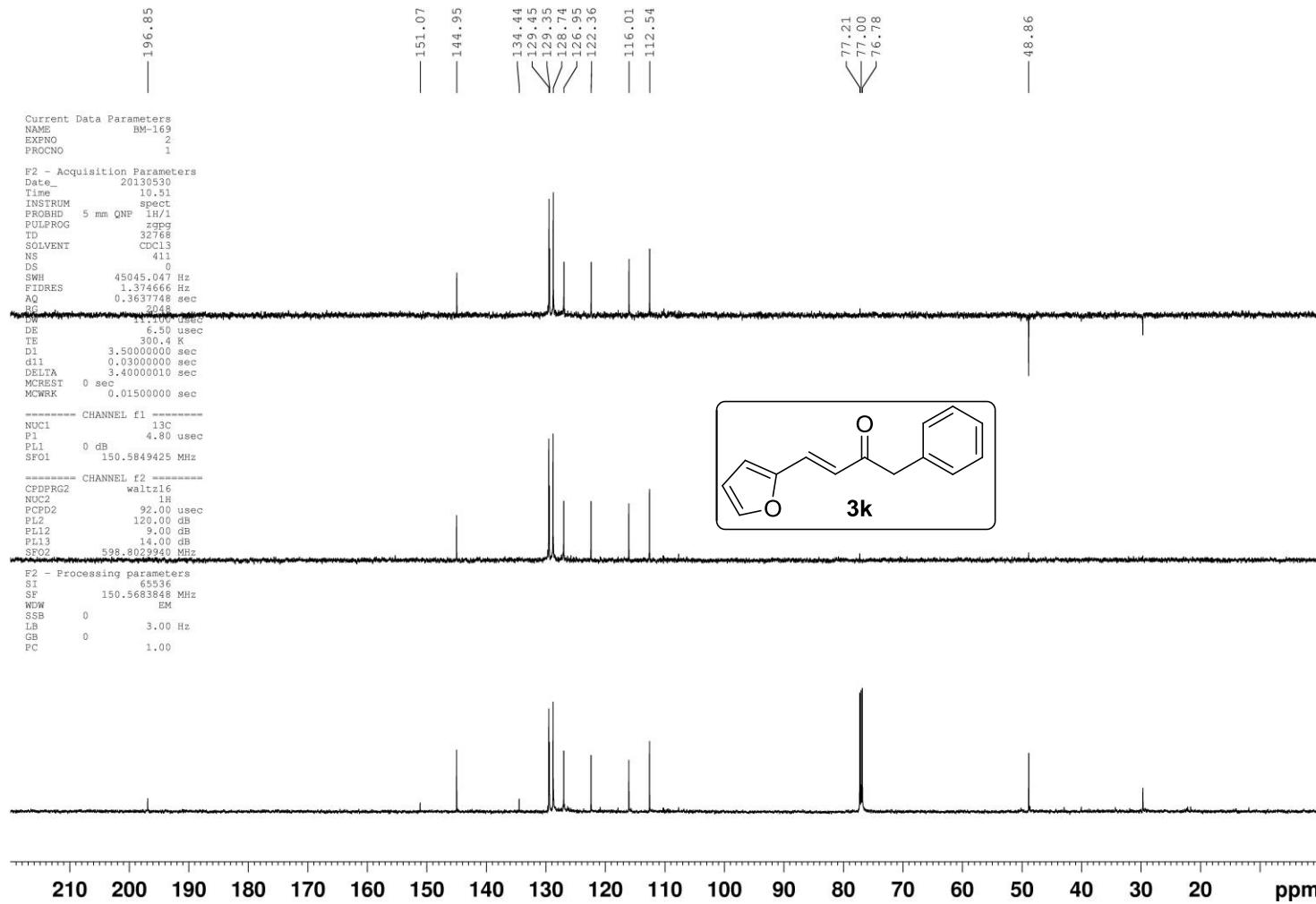
4

2

0







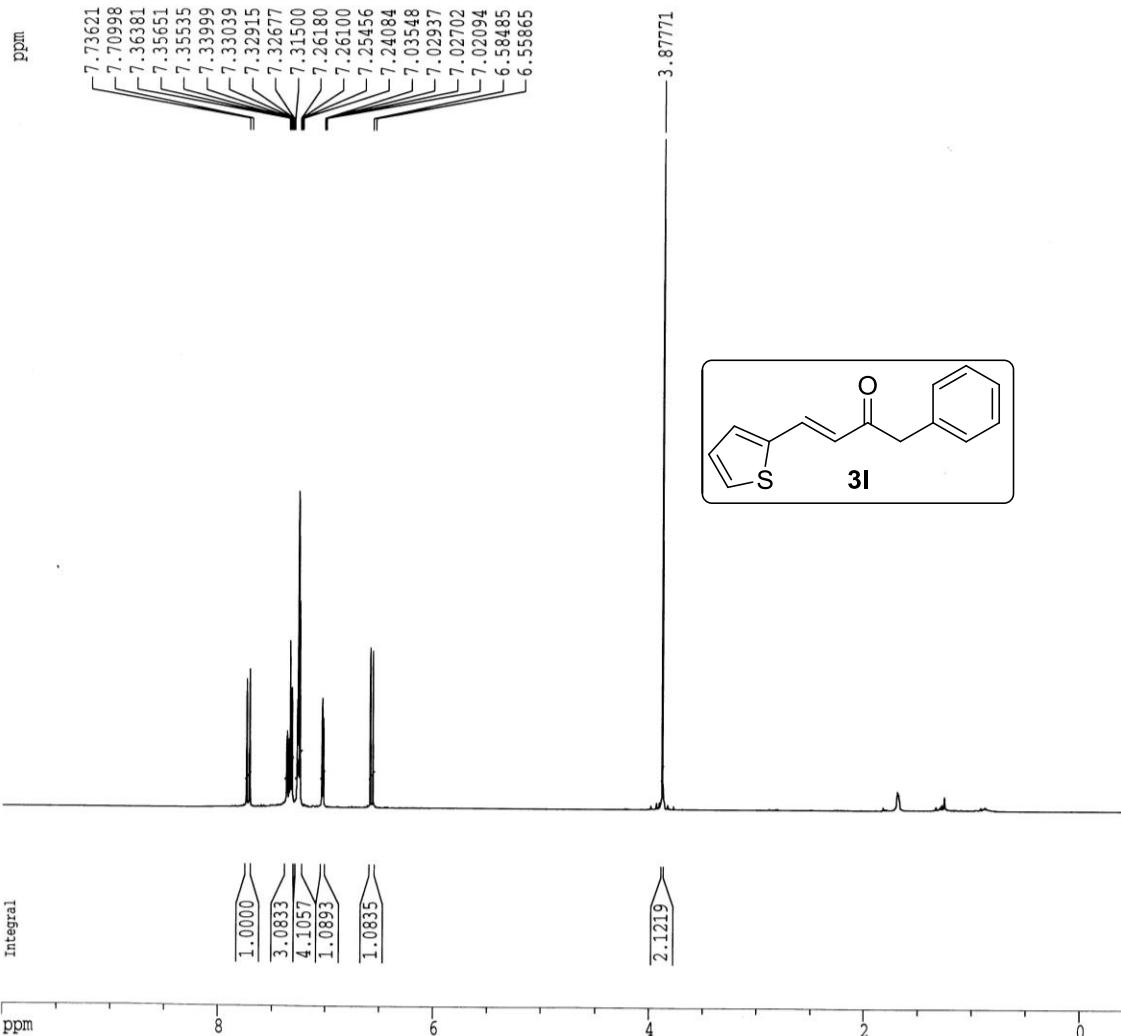
Current Data Parameters
NAME BM-159
EXPNO 1
PROCNO 1

F2 - Acquisition Parameters
Date_ 20130526
Time 20.40
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 32768
SOLVENT CDCl3
NS 16
DS 0
SWH 8382.229 Hz
FIDRES 0.255805 Hz
AQ 1.9546613 sec
RG 128
DW 59.650 usec
DE 6.50 usec
TE 300.0 K
D1 2.0000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

===== CHANNEL f1 =====
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SFO1 598.802940 MHz

F2 - Processing parameters
SI 32768
SF 598.8000285 MHz
WDW no
SSB 0
LB 0.00 Hz
GB 0
PC 2.00

1D NMR plot parameters
CX 20.00 cm
CY 12.00 cm
F1P 10.000 ppm
F1 5988.00 Hz
F2P -0.500 ppm
F2 -299.40 Hz
PPCM 0.52500 ppm/cm
HZCM 314.37003 Hz/cm



Current Data Parameters
NAME BM-159
EXPNO 2
PROCNO 1

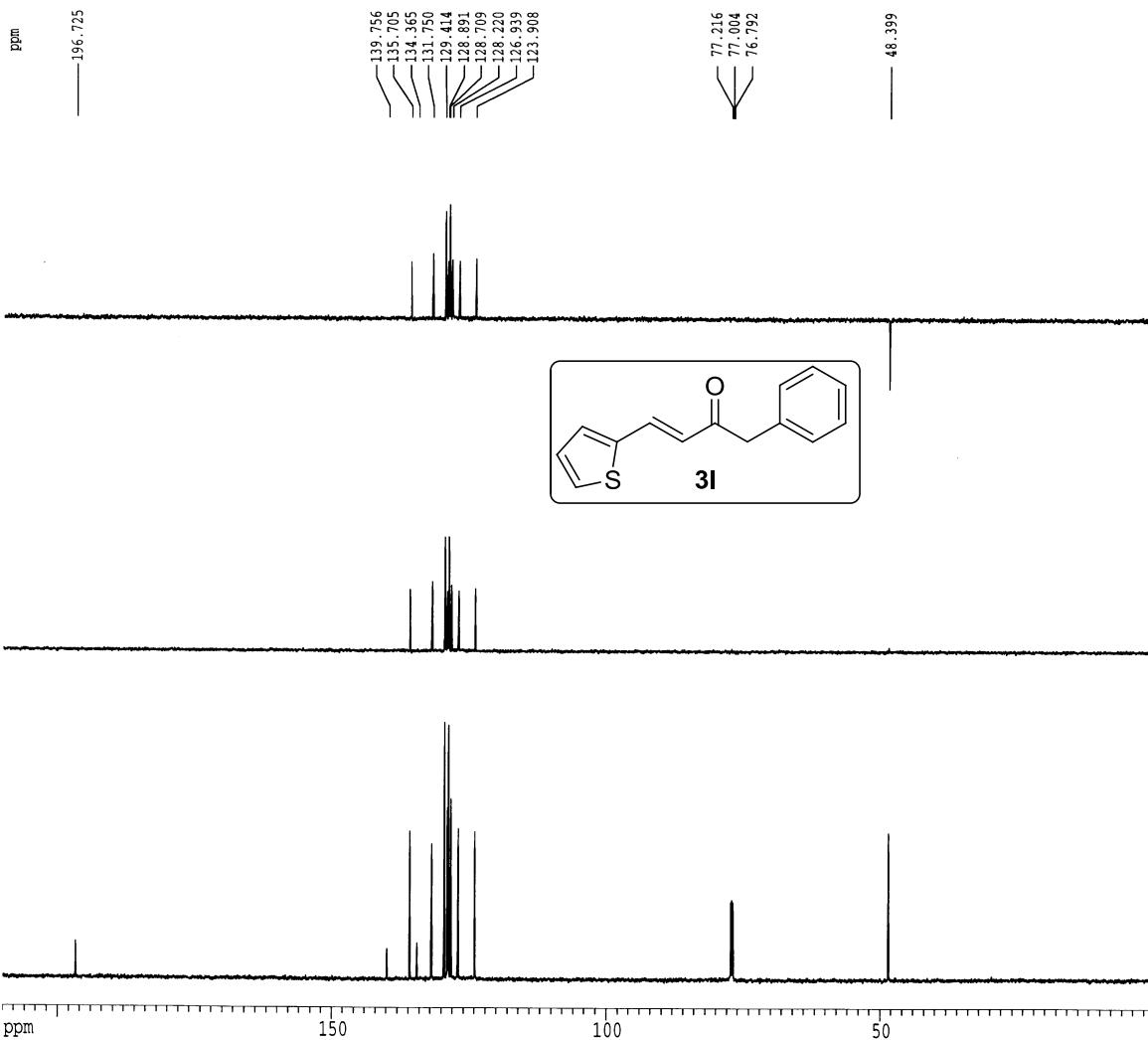
F2 - Acquisition Parameters
Date_ 20130526
Time 20.41
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zpgq
TD 32768
SOLVENT CDCl3
NS 100
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.00 usec
DE 6.50 usec
TE 300.1 K
D1 3.5000000 sec
d11 0.0300000 sec
DELTA 3.4000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SF01 150.5849425 MHz

===== CHANNEL f2 =====
CPDPRG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SF02 598.8029940 MHz

F2 - Processing parameters
SI 65536
SF 150.5683924 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 1.00

1D NMR plot parameters
CX 20.00 cm
CY 4.50 cm
F1P 210.000 ppm
F1 31619.36 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPCM 10.50000 ppm/cm
HZCM 1580.96814 Hz/cm



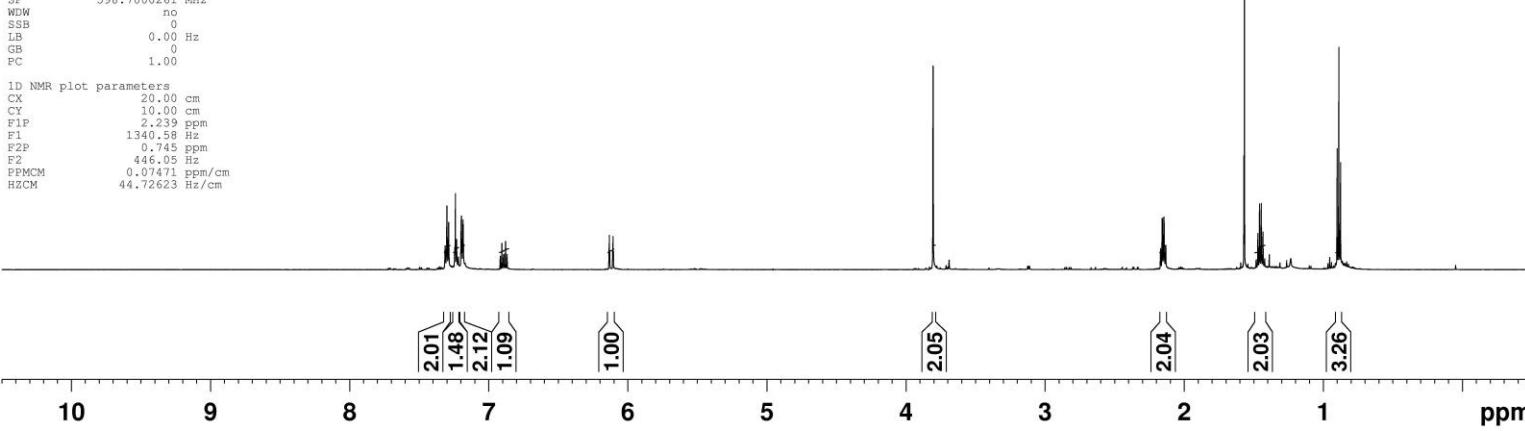
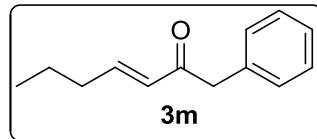
Current Data Parameters
NAME BM-02-199A
EXPNO 1
PROCNO 1

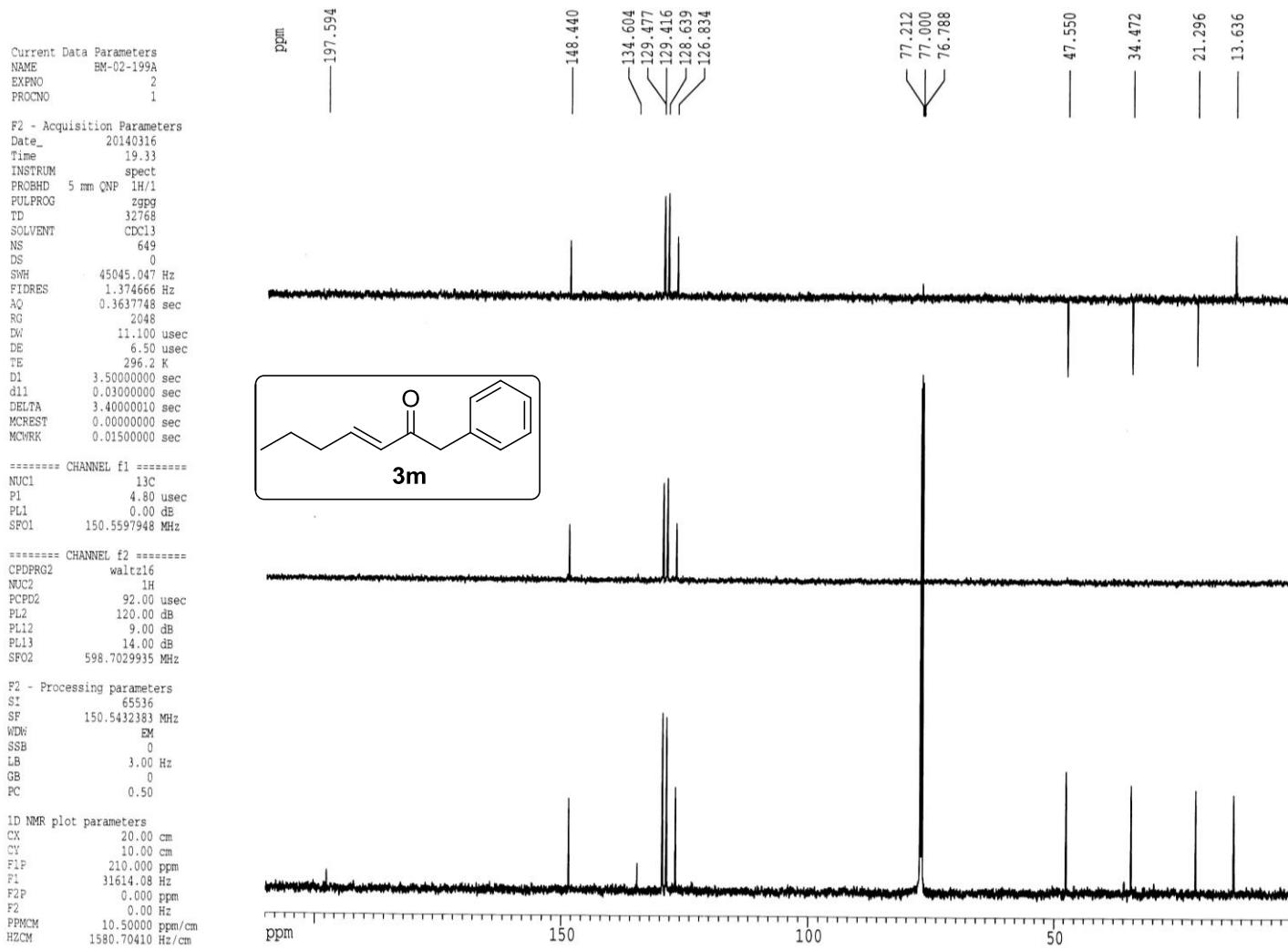
F2 - Acquisition Parameters
Date 20140316
Time 19.22
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 3356
SOLVENT CDCl3
NS 16
DS 0
SWH 8389.262 Hz
FIDRES 0.250008 Hz
AQ 1.9999876 sec
RG 512
DW 59.00 usec
DE 6.50 usec
TE 294.5 °
D1 2.0000000 sec
MCREST 0.0000000 sec
MCWKR 0.0150000 sec

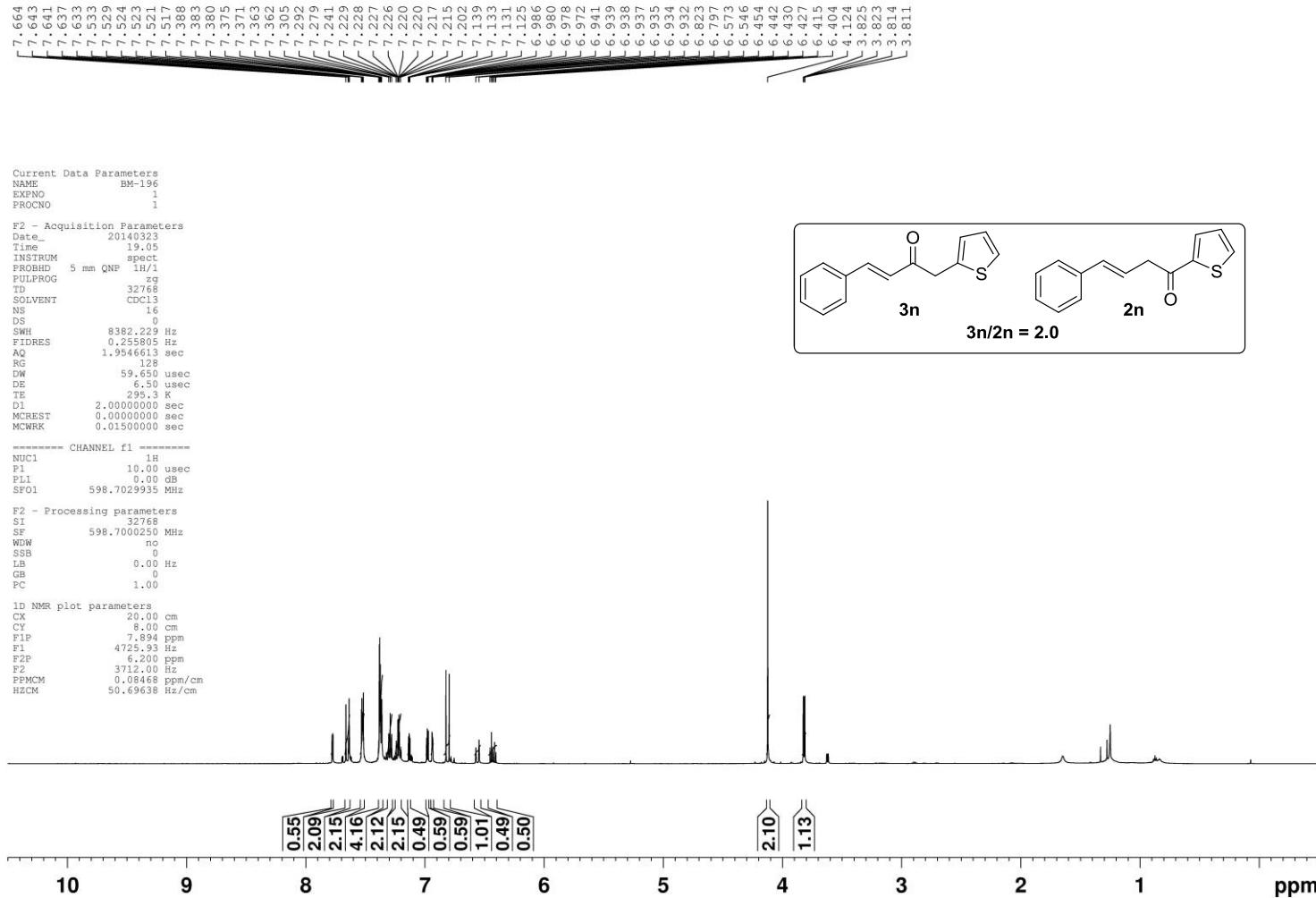
CHANNEL f1 -----
NUC1 1H
F1 10.00 usec
PL1 0.00 dB
SF01 598.7029935 MHz

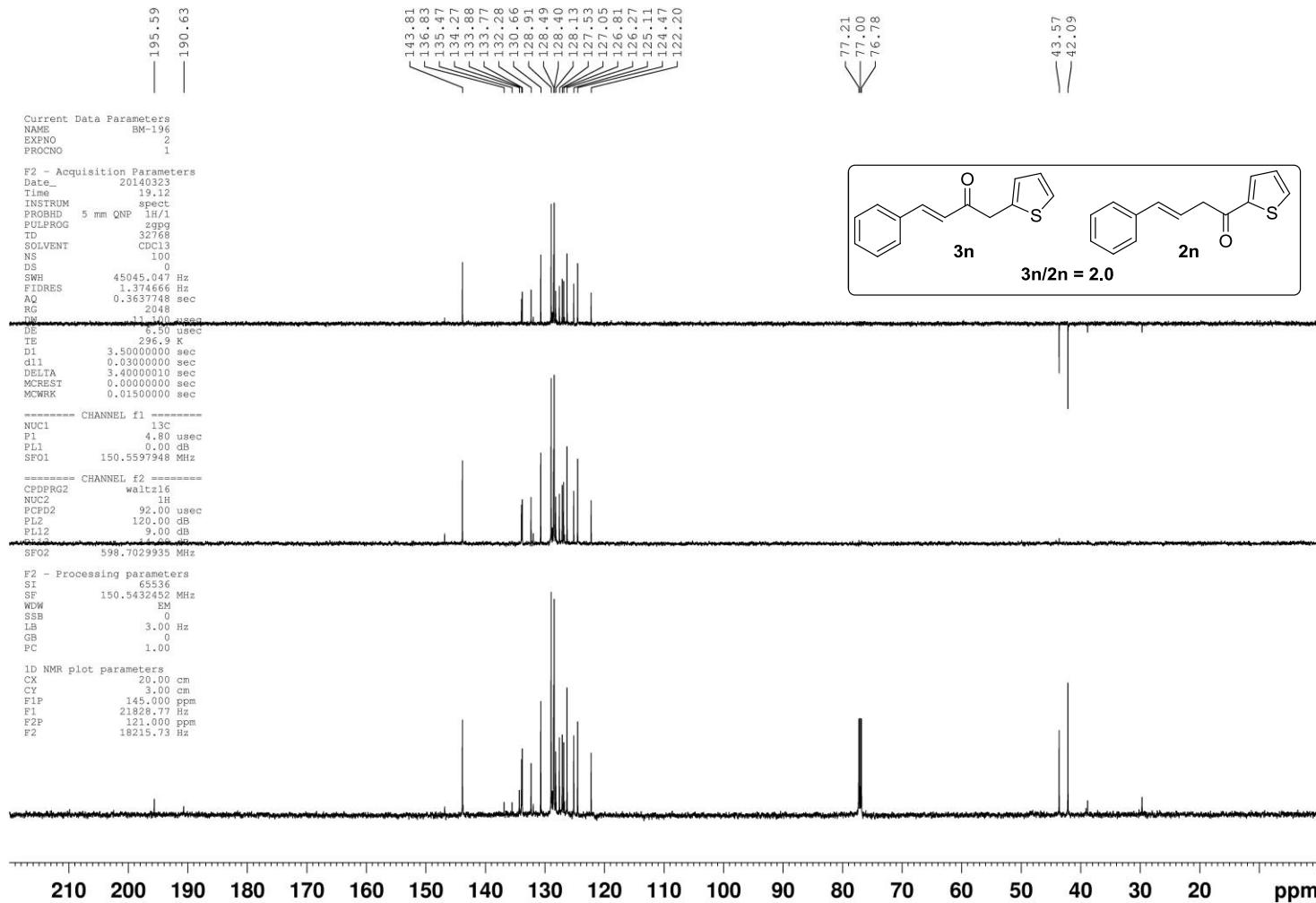
F2 - Processing parameters
SI 32768
SF 598.7000261 MHz
NDW 1
SSB 0
LB 0.00 Hz
GB 0
PC 1.00

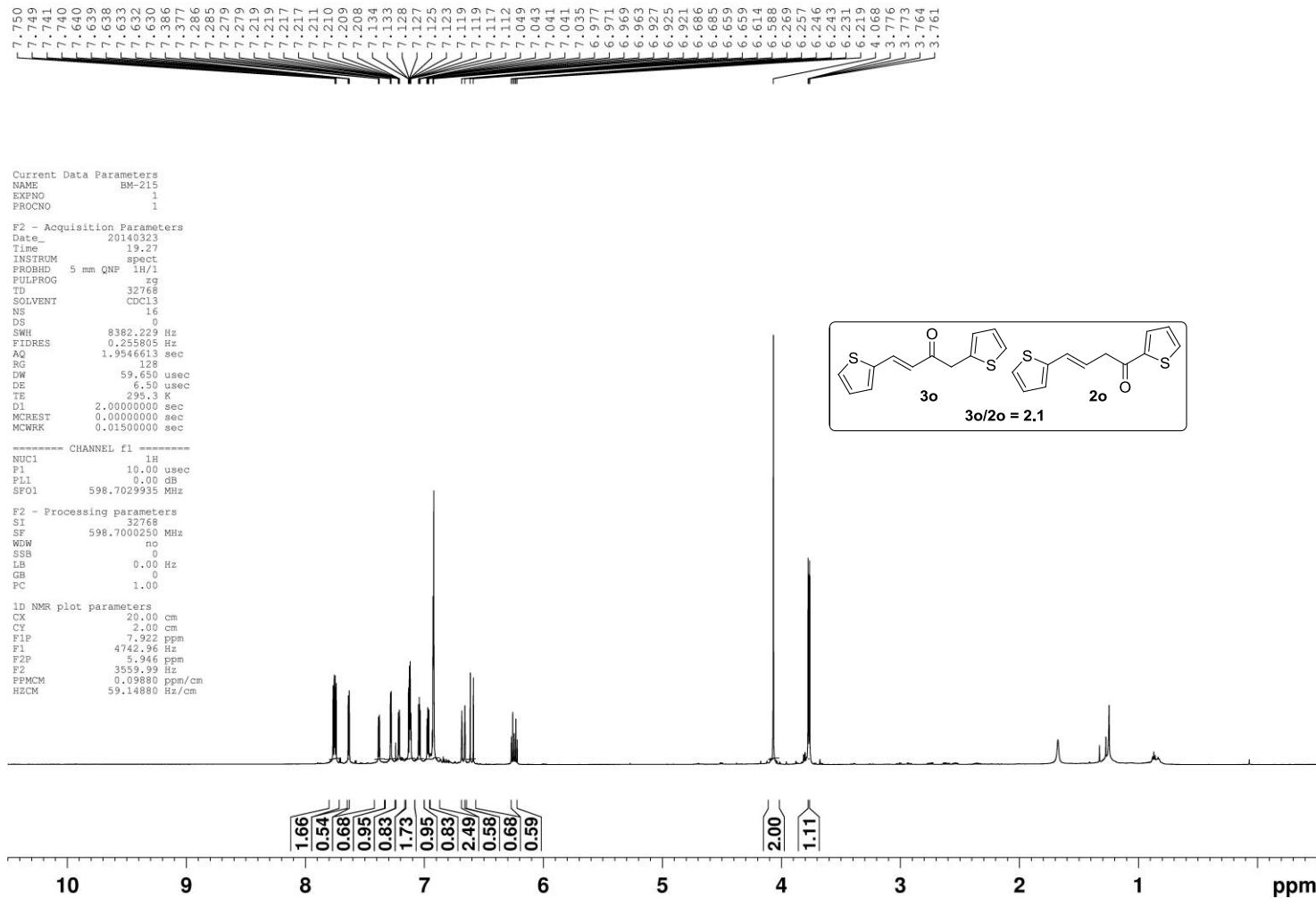
1D NMR plot parameters
CX 20.00 cm
CY 1.00 cm
F1P 2.239 ppm
F1 1340.58 Hz
F2P 0.745 ppm
F2 446.05 Hz
PPMCM 0.07471 ppm/cm
HZCM 44.72623 Hz/cm

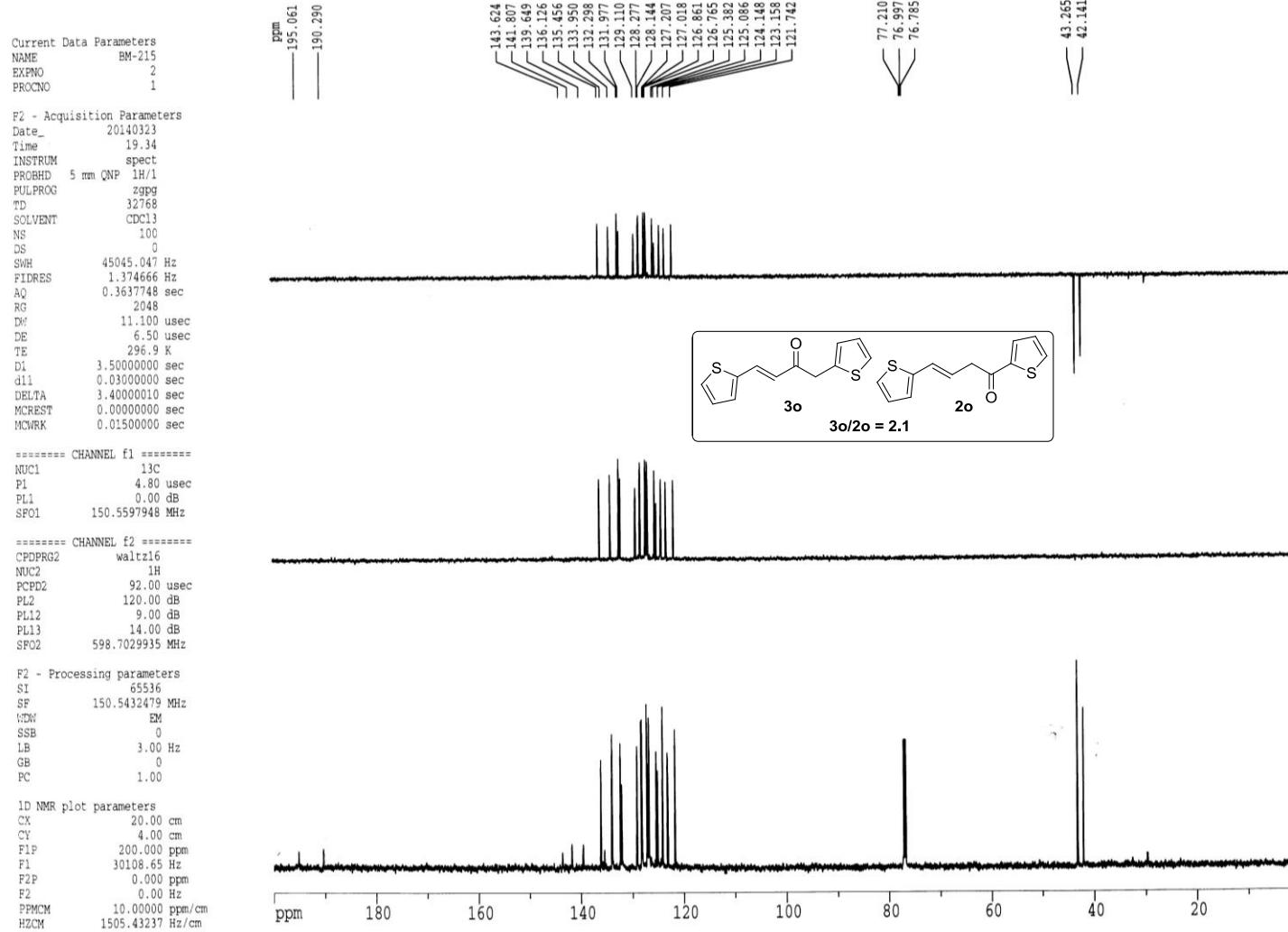












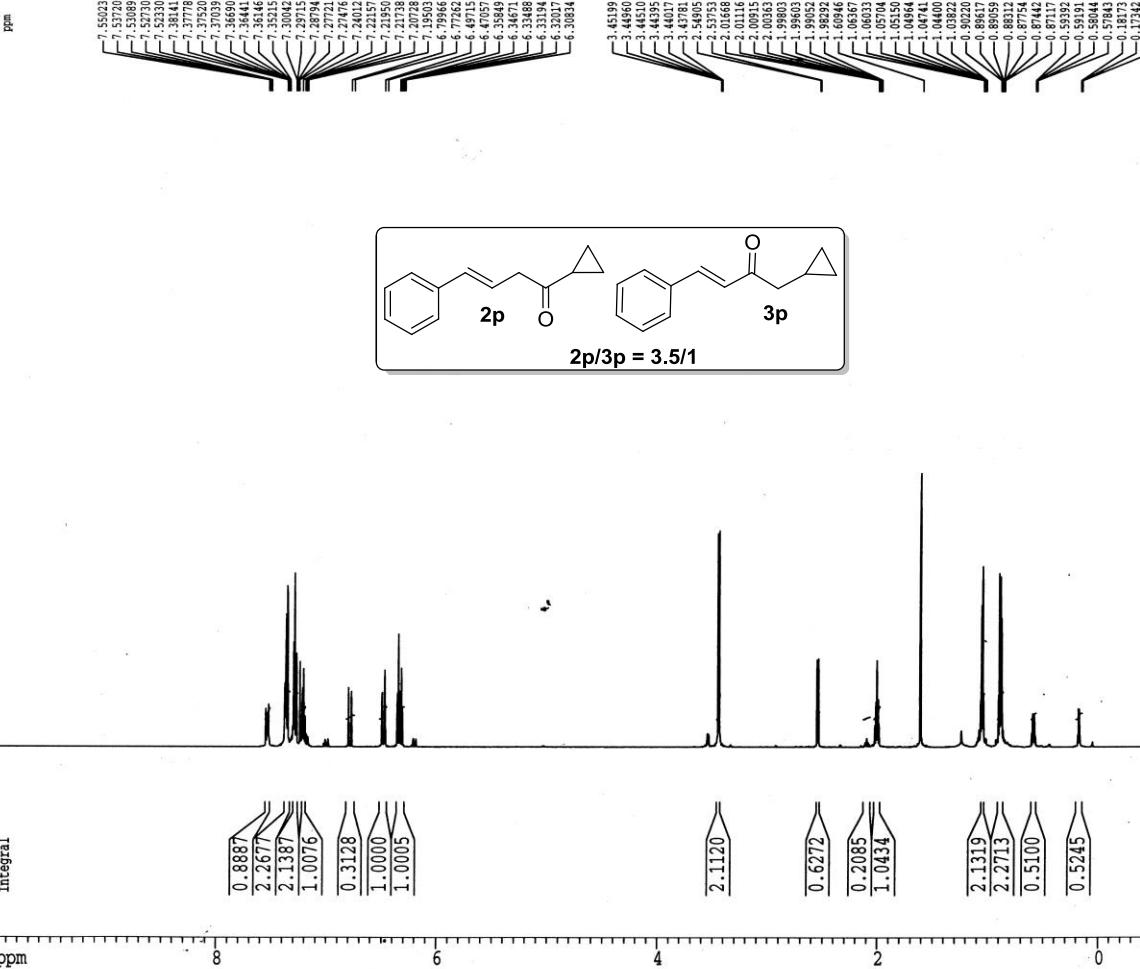
Current Data Parameters
 NAME BM-03-09
 EXPNO 1
 PROCNO 1

F2 - Acquisition Parameters
 Date_ 20140522
 Time 19.17
 INSTRUM spect
 PROBHD 5 mm QNP 1H/1
 PULPROG zg
 TD 32768
 SOLVENT CDCl₃
 NS 16
 DS 0
 SWH 8389.262 Hz
 FIDRES 0.256020 Hz
 AQ 1.9530228 sec
 RG 512
 DW 59.600 usec
 DE 6.50 usec
 TE 298.4 K
 D1 2.0000000 sec
 MCREST 0.0000000 sec
 MWRK 0.0150000 sec

===== CHANNEL f1 ======
 NUC1 1H
 P1 10.00 usec
 PLL 3.00 dB
 SF01 598.7029935 MHz

F2 - Processing parameters
 SI 32768
 SF 598.7000261 MHz
 WDW no
 SSB 0
 LB 0.00 Hz
 GB 0
 PC 1.00

1D NMR plot parameters
 CX 20.00 cm
 CY 5.00 cm
 F1P 10.000 ppm
 F1 5987.00 Hz
 F2P -0.500 ppm
 F2 -299.35 Hz
 PPMCM 0.52500 ppm/cm
 HZCM 314.31750 Hz/cm



Current Data Parameters
NAME BM-03-09
EXPNO 2
PROCNO 1

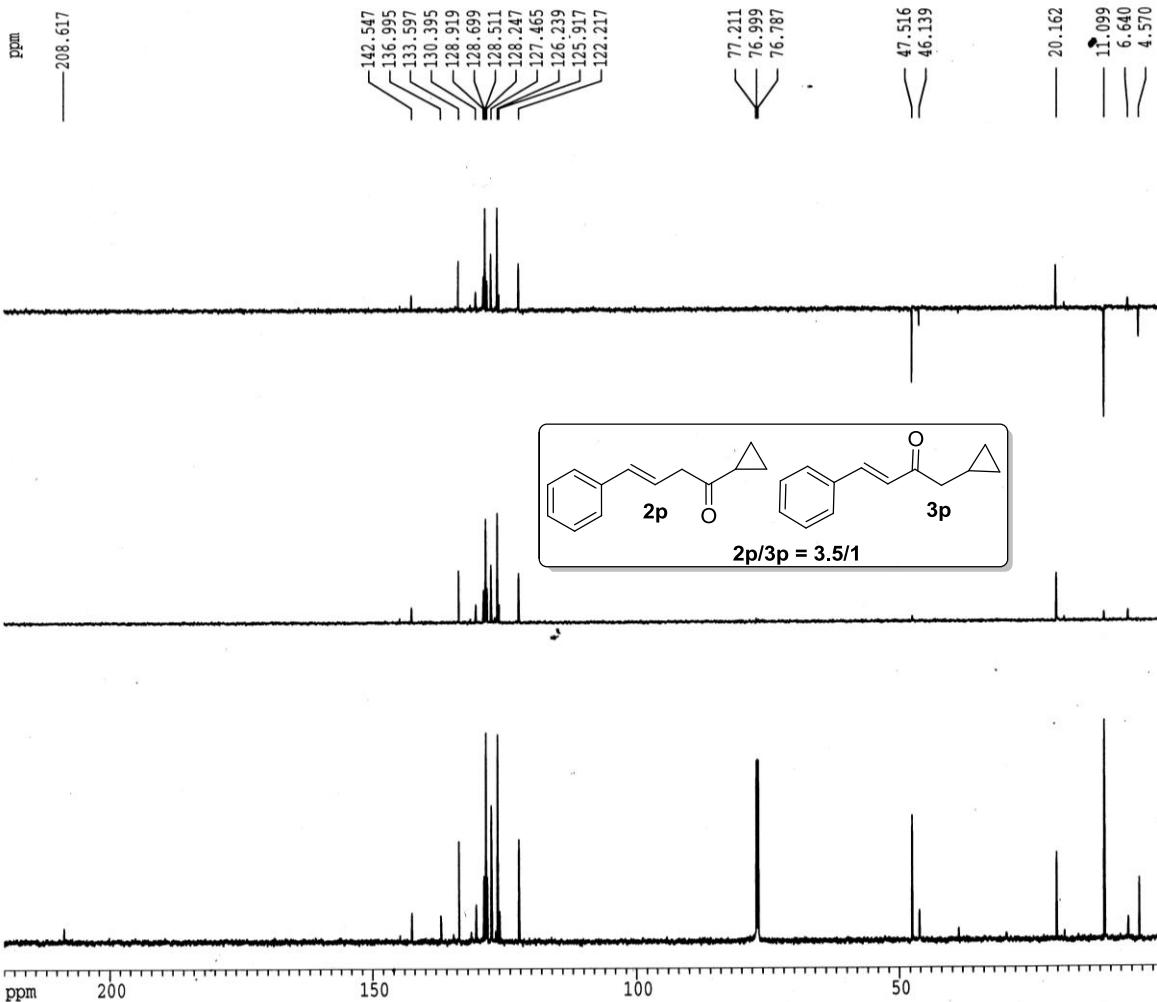
F2 - Acquisition Parameters
Date 20140522
Time 19.19
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT CDCl₃
NS 211
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 299.1 K
D1 3.5000000 sec
d11 0.0300000 sec
DELTA 3.4000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SF01 150.5597948 MHz

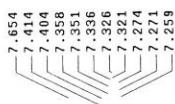
===== CHANNEL f2 =====
CPDPG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SF02 598.7029935 MHz

F2 - Processing parameters
SI 65536
SF 150.5432397 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 0.50

1D NMR plot parameters
CX 20.00 cm
CY 4.00 cm
F1P 220.000 ppm
F1 33119.51 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPMCM 11.00000 ppm/cm
HZCM 1655.97571 Hz/cm

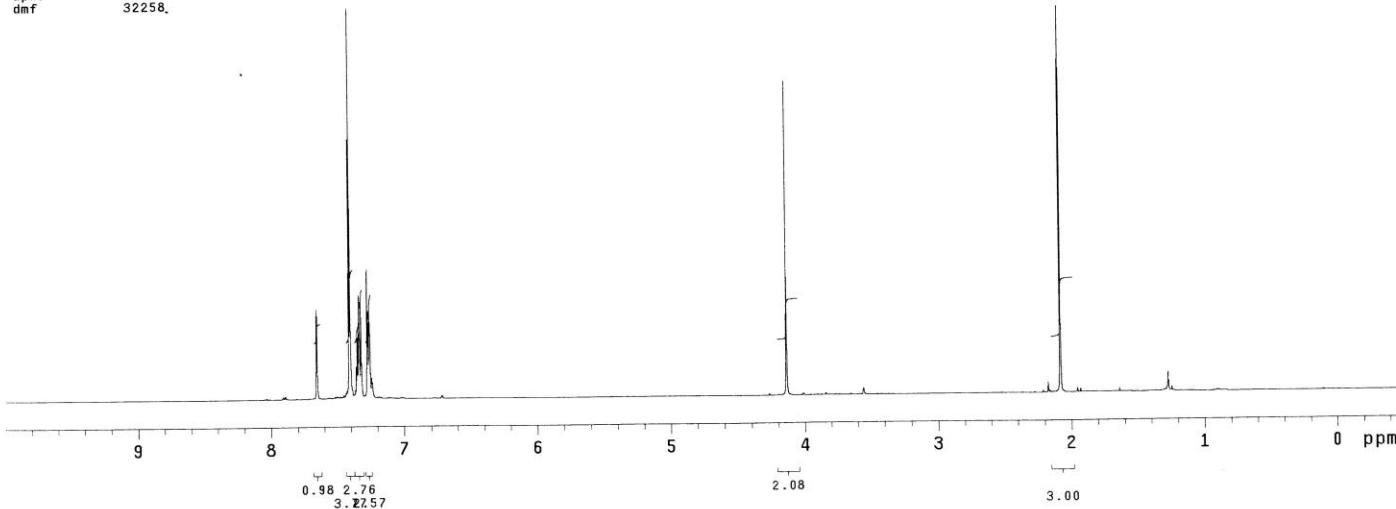
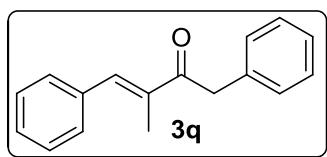


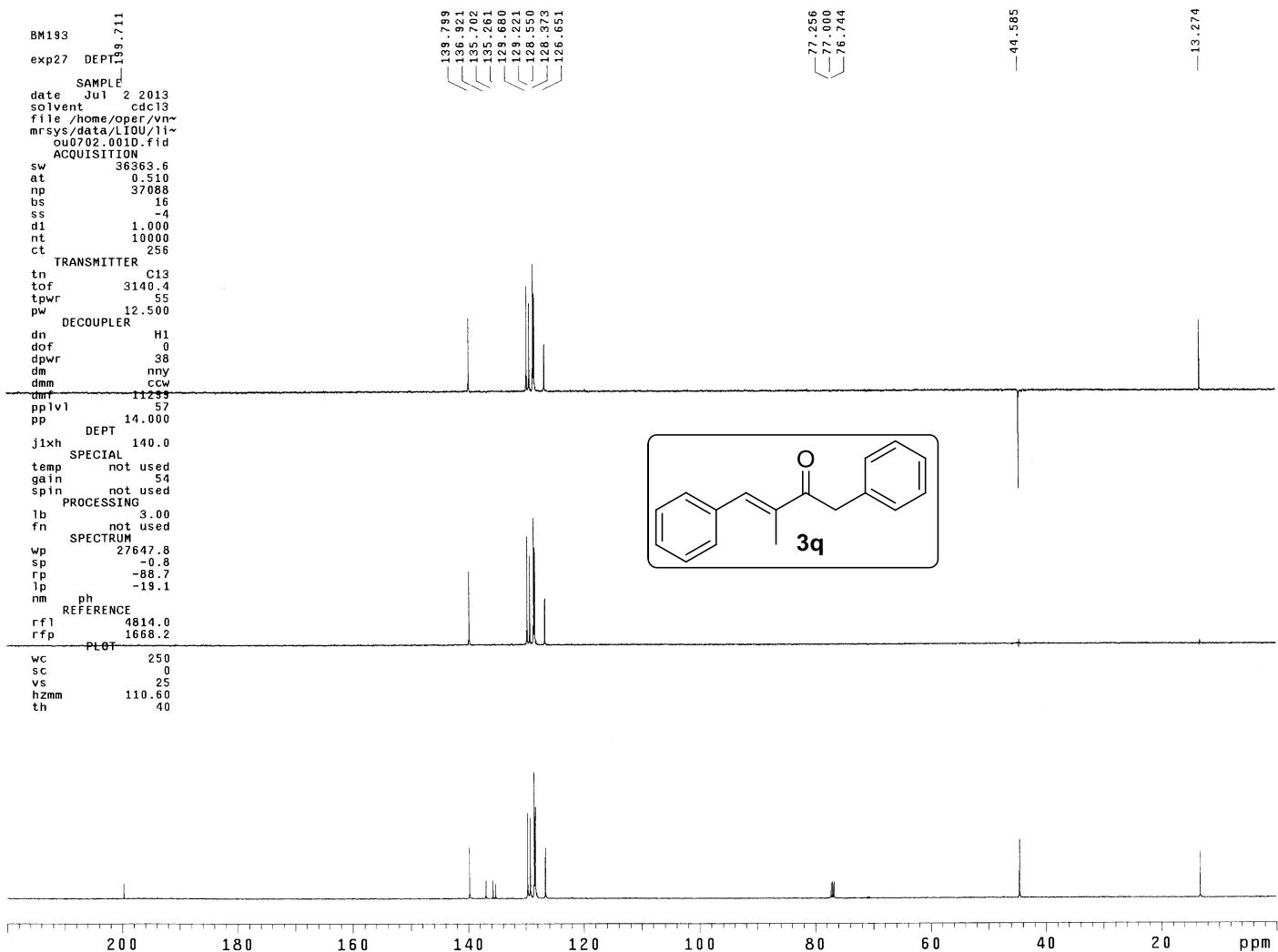
BM193

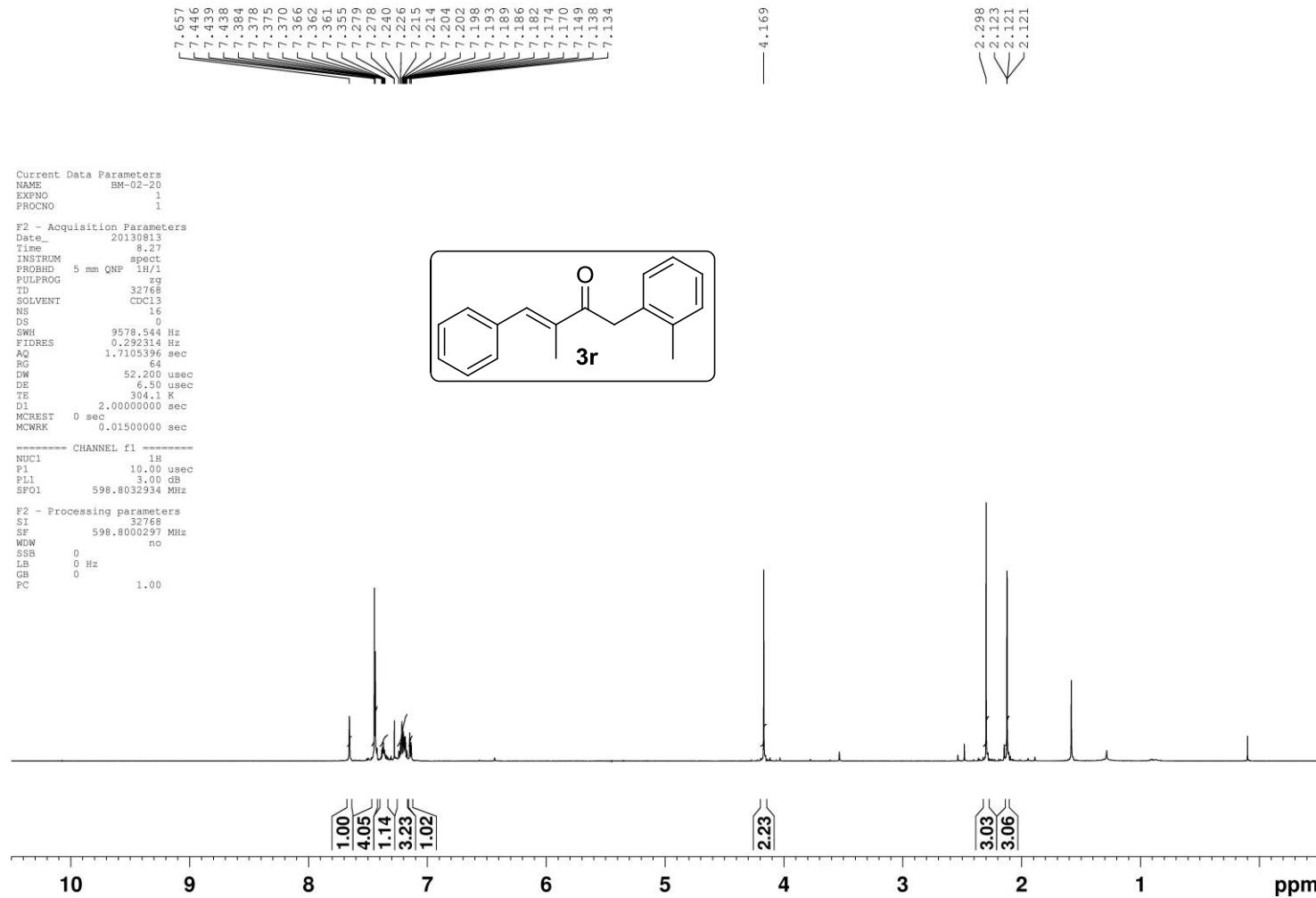


exp24 Proton

SAMPLE SPECIAL
date Jul 2 2013 temp not used
solvent cdc13 gain not used
file /home/oper/vn~ spin not used
mrsys/data/LIOU/l1~ hst 0.008
ou0702.001H.fid pw90 12.600
ACQUISITION alfa 6.600
sw 9000.9 FLAGS
at 2.044 i1 n
np 36788 in n
fb not used dp y
bs not used hs nn
ss 2 PROCESSING
d1 1.000 fn 32768
nt 32 DISPLAY
ct 32 sp -250.1
TRANSMITTER H1 rfp 5247.6
sfrq 499.811 rfp 4632.8
tof 999.5 rp 3618.6
tpwr 55 1p 145.8
pw 6.300 PLOT
DECOUPLER wc 250
dn C13 sc 0
dof 0 vs 70
dm nnn th 4
dmm c nm ph
dpwr 36
dmf 32258.









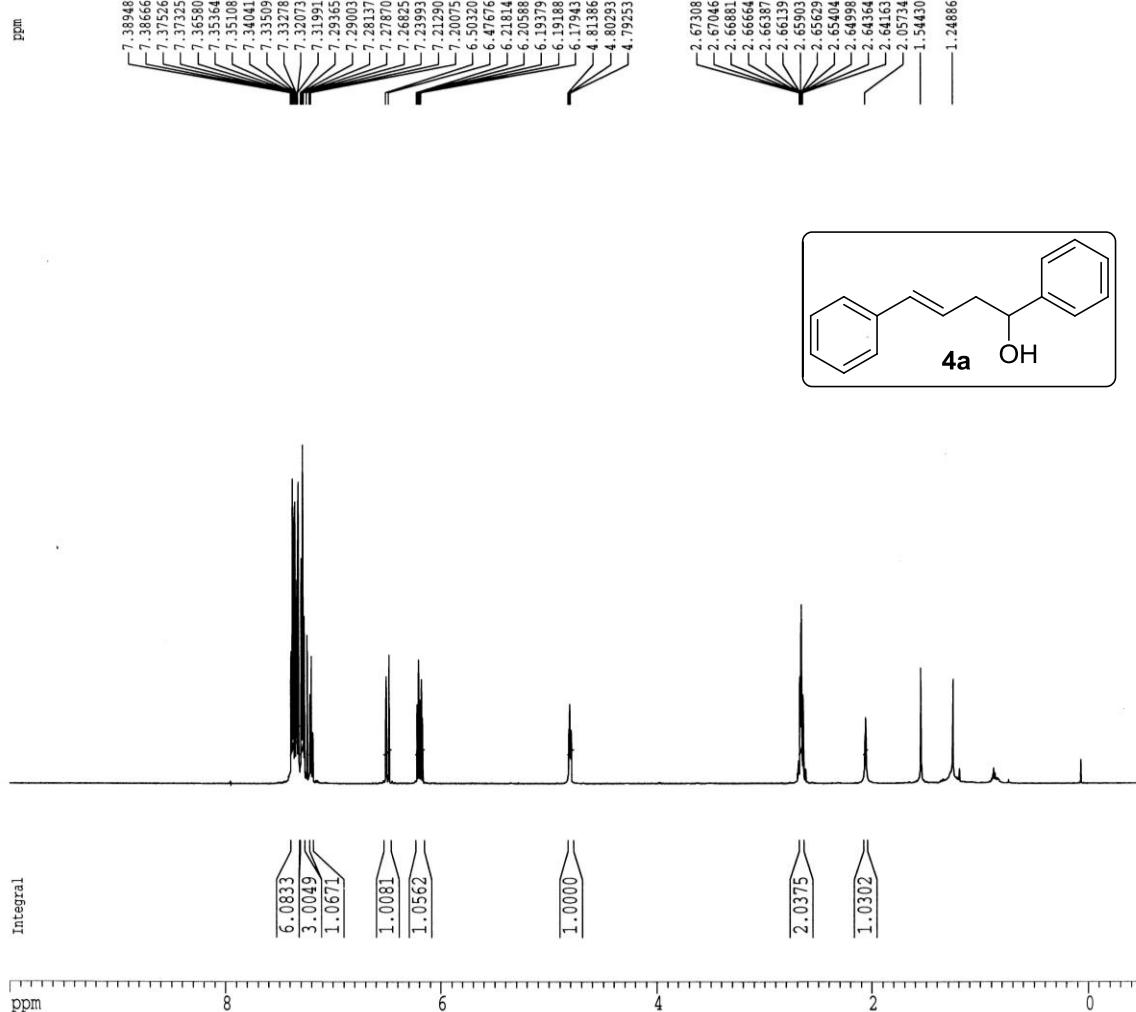
Current Data Parameters
 NAME BM-02-62
 EXPNO 1
 PROCNO 1

F2 - Acquisition Parameters
 Date_ 20130929
 Time 14.17
 INSTRUM spect
 PROBHD 5 mm QNP 1H/1
 PULPROG zg
 TD 32768
 SOLVENT CDCl₃
 NS 16
 DS 0
 SWH 13123.359 Hz
 FIDRES 0.400493 Hz
 AQ 1.2485108 sec
 RG 256
 DW 38.100 usec
 DE 6.50 usec
 TE 304.5 K
 D1 2.0000000 sec
 MCREST 0.0000000 sec
 MCWRK 0.0150000 sec

===== CHANNEL f1 =====
 NUC1 1H
 P1 10.00 usec
 PL1 0.00 dB
 SF01 598.8047904 MHz

F2 - Processing parameters
 SI 32768
 SF 598.8000290 MHz
 WDW no
 SSB 0
 LB 0.00 Hz
 GB 0
 PC 1.00

1D NMR plot parameters
 CX 20.00 cm
 CY 6.00 cm
 F1P 10.000 ppm
 F1 5988.00 Hz
 F2P -0.500 ppm
 F2 -299.40 Hz
 PPMCM 0.52500 ppm/cm
 HZCM 314.37003 Hz/cm



Current Data Parameters
NAME BM-02-62
EXPNO 2
PROCNO 1

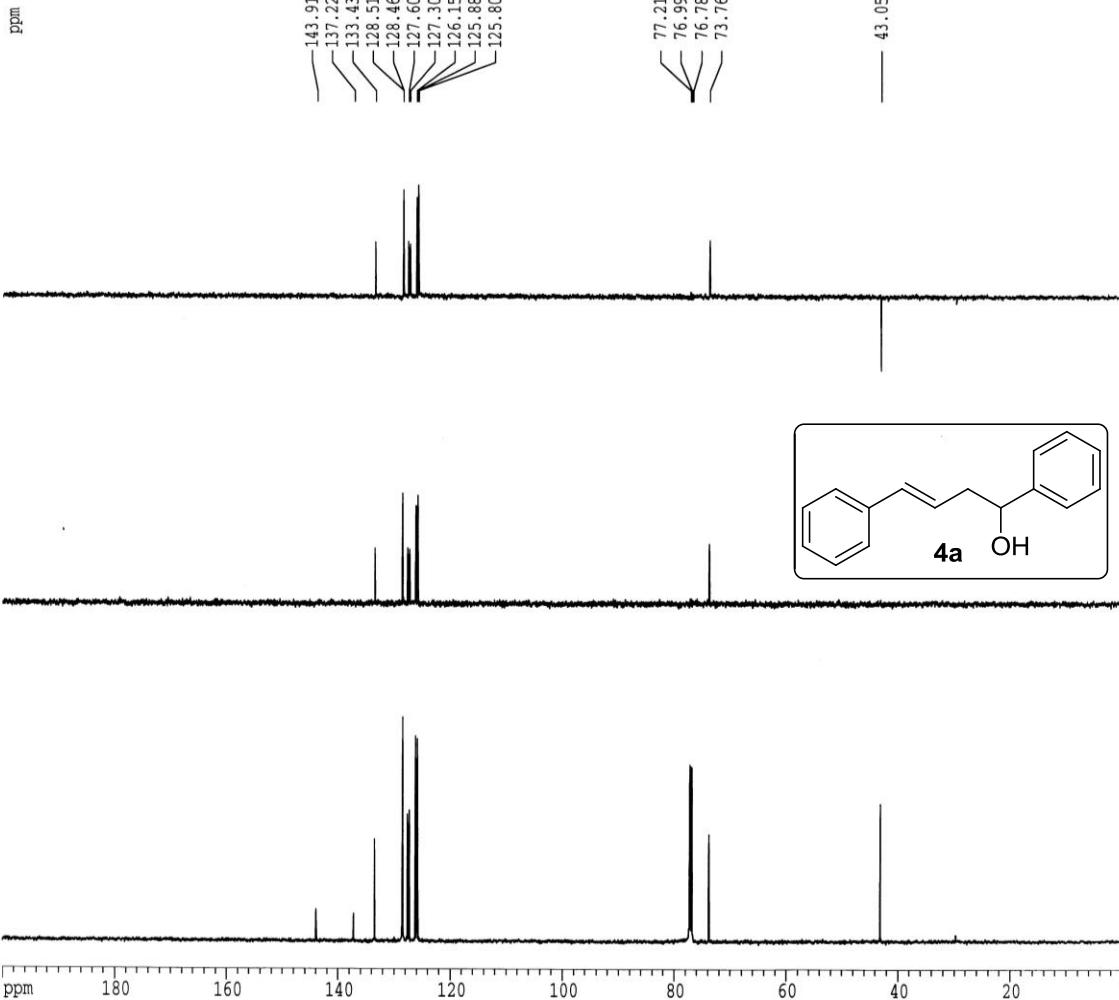
F2 - Acquisition Parameters
Date 20130929
Time 14.50
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpp
TD 32768
SOLVENT CDCl3
NS 500
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 305.2 K
D1 3.5000000 sec
d11 0.0300000 sec
DELTA 3.4000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

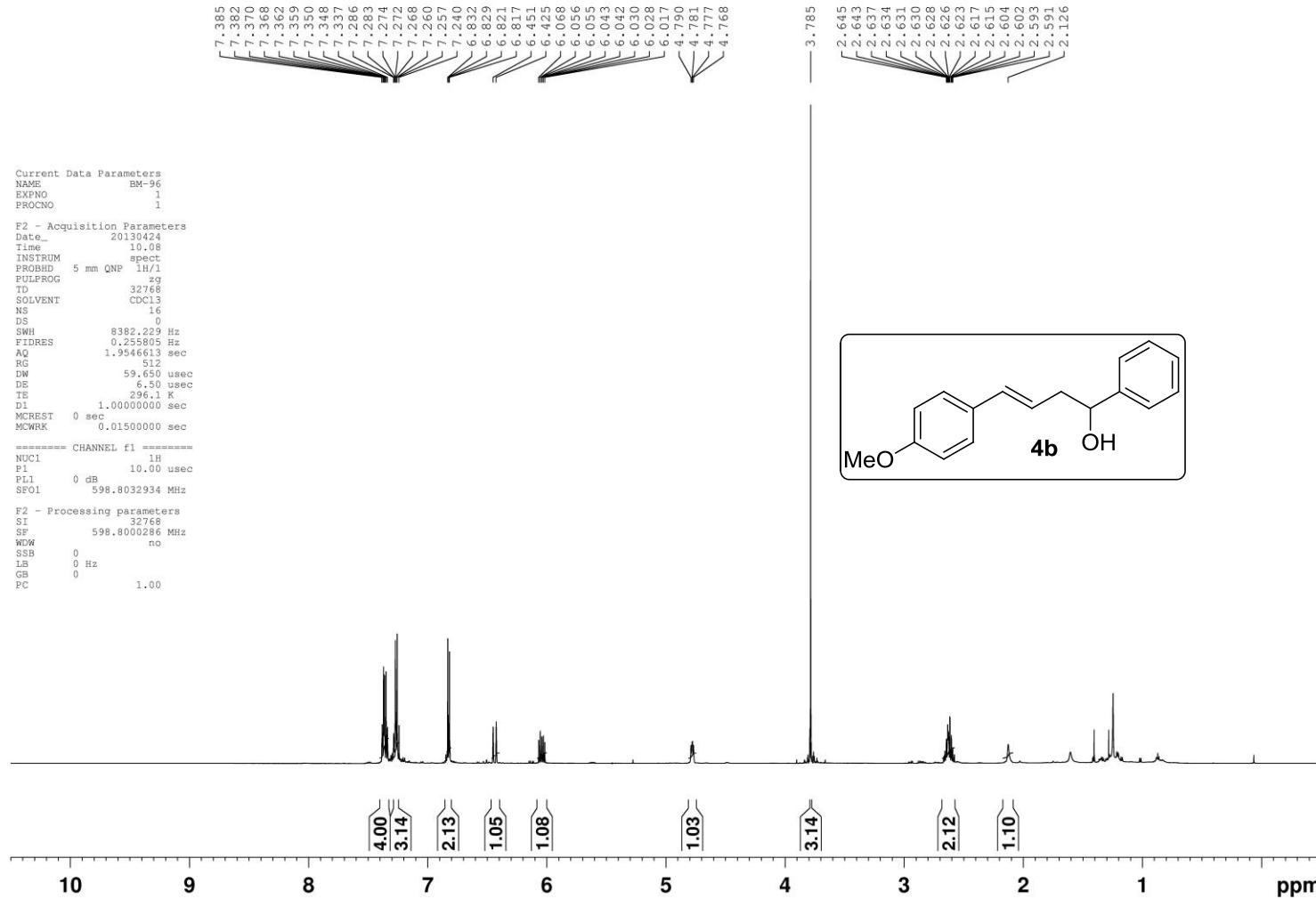
===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SFO1 150.5849425 MHz

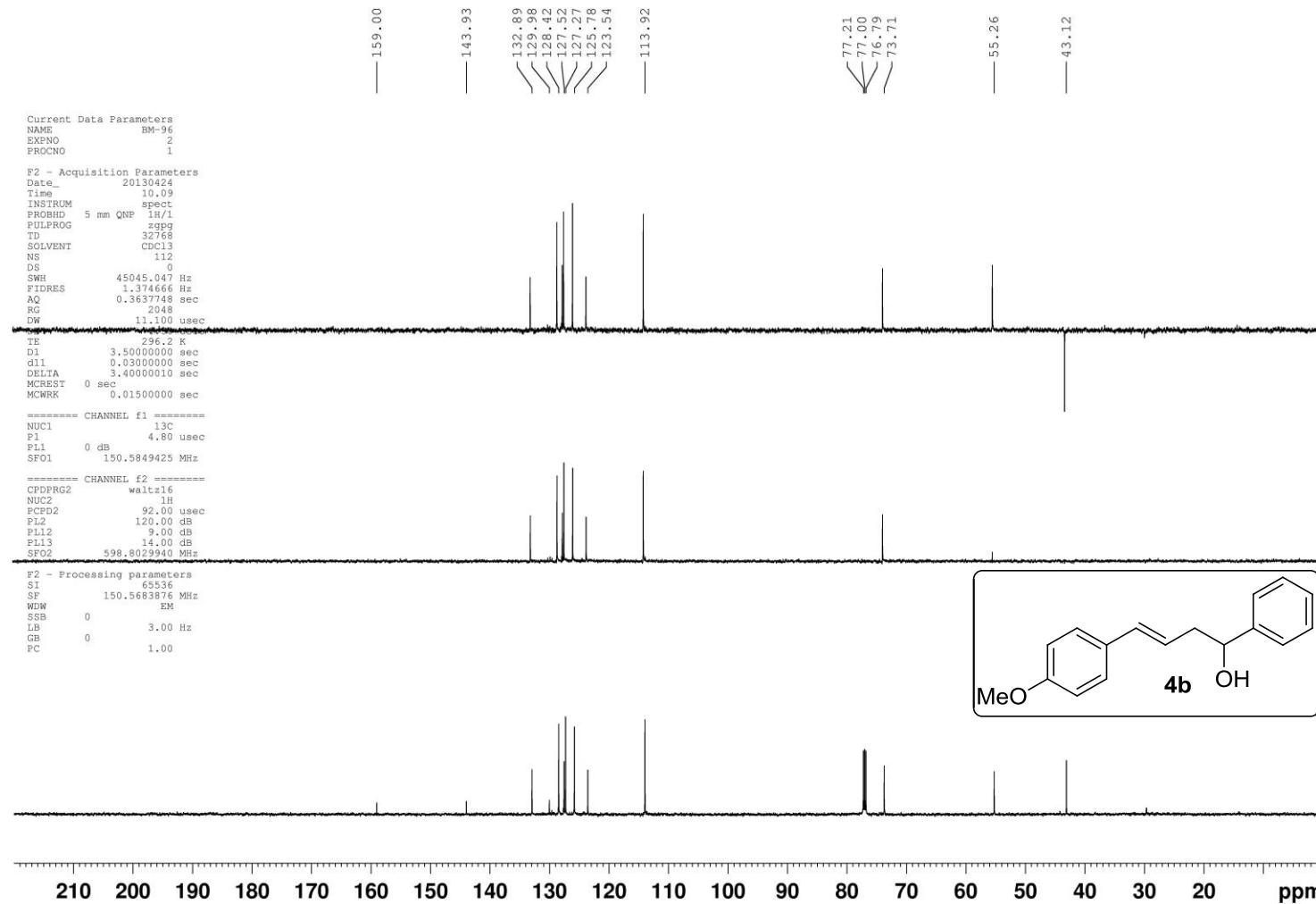
===== CHANNEL f2 =====
CPDPRG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SFO2 598.8029940 MHz

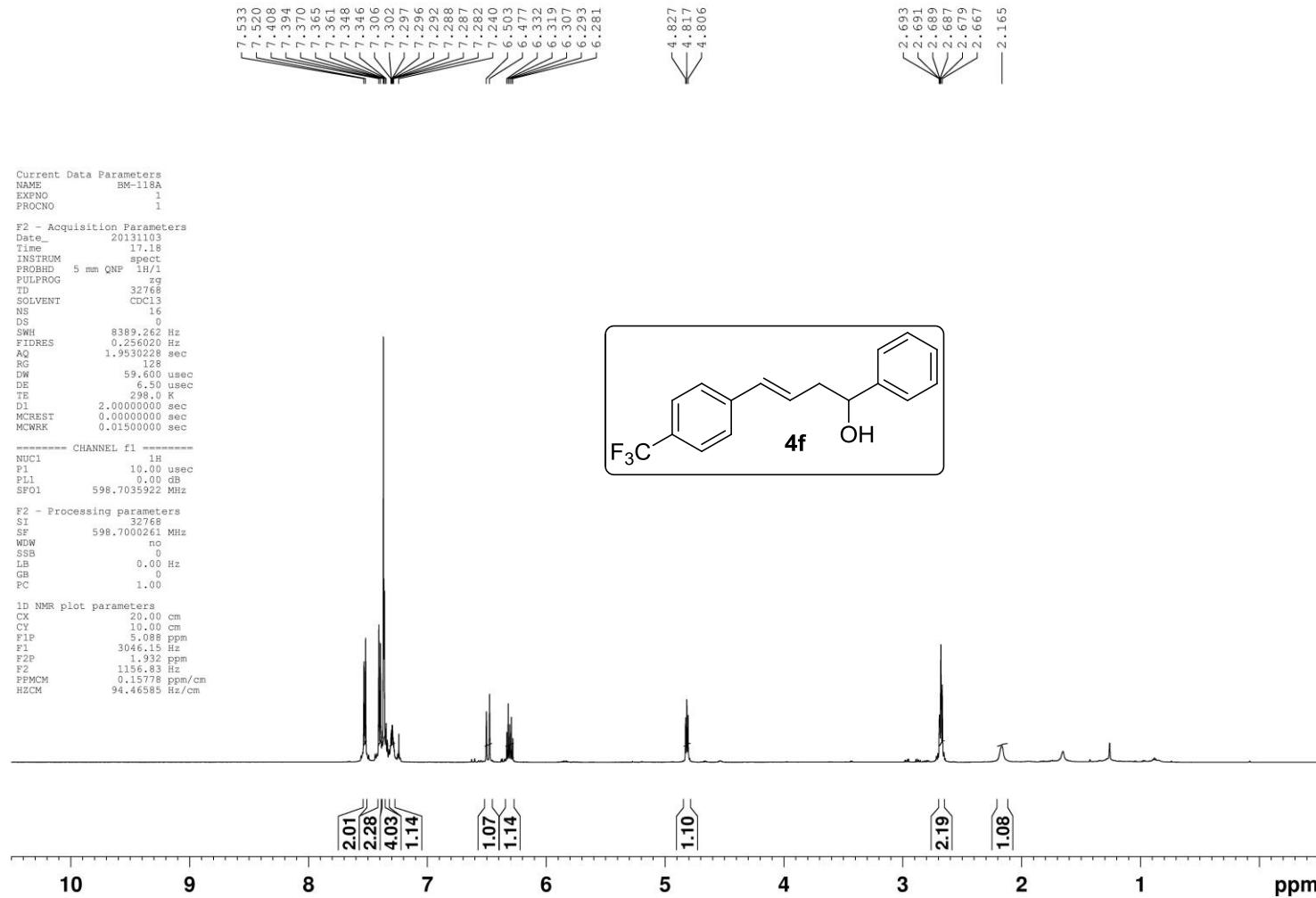
F2 - Processing parameters
SI 65536
SF 150.5683835 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 0.50

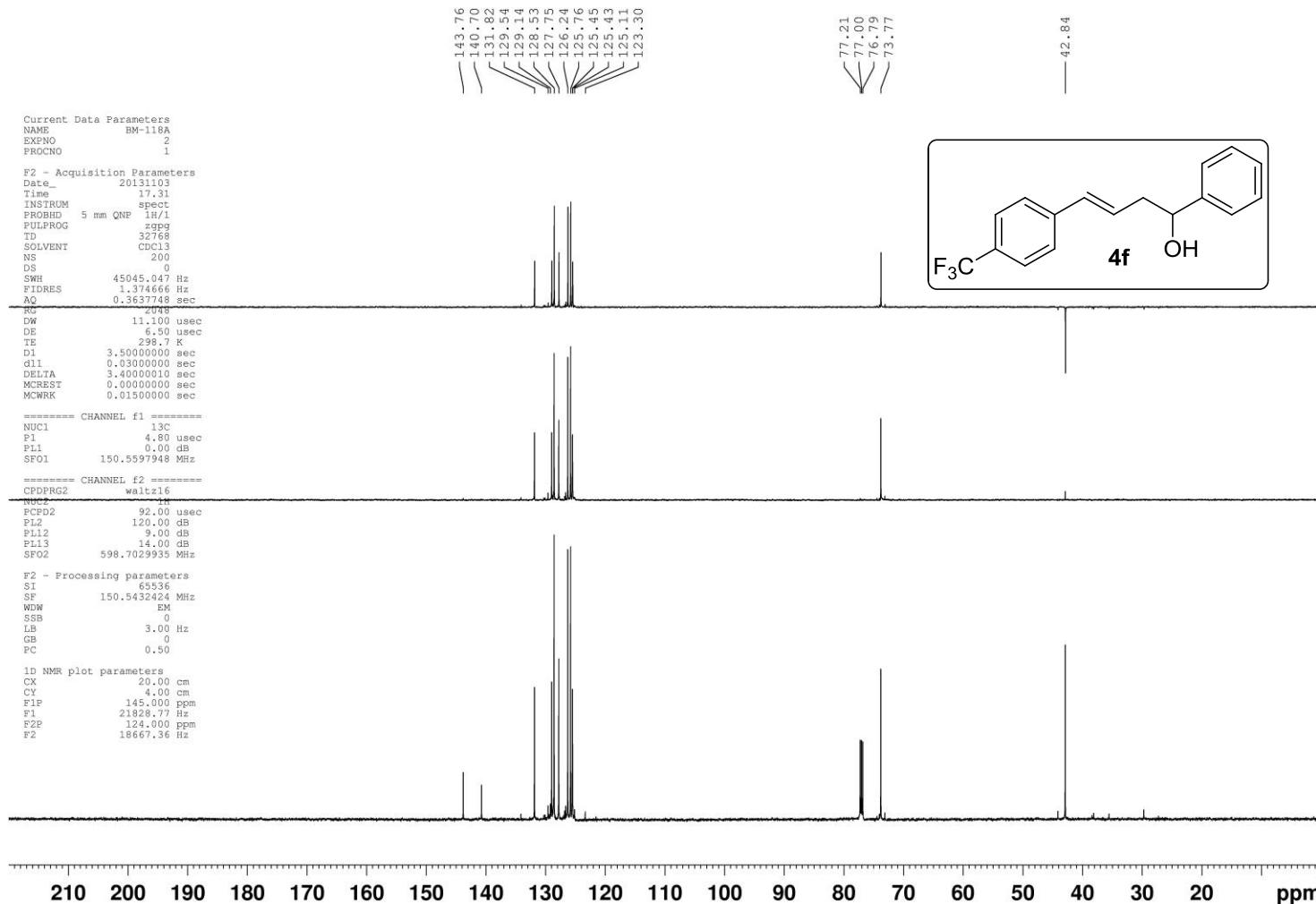
1D NMR plot parameters
CX 20.00 cm
CY 4.00 cm
FLP 200.000 ppm
F1 30113.68 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPMCM 10.00000 ppm/cm
HZCM 1505.68384 Hz/cm











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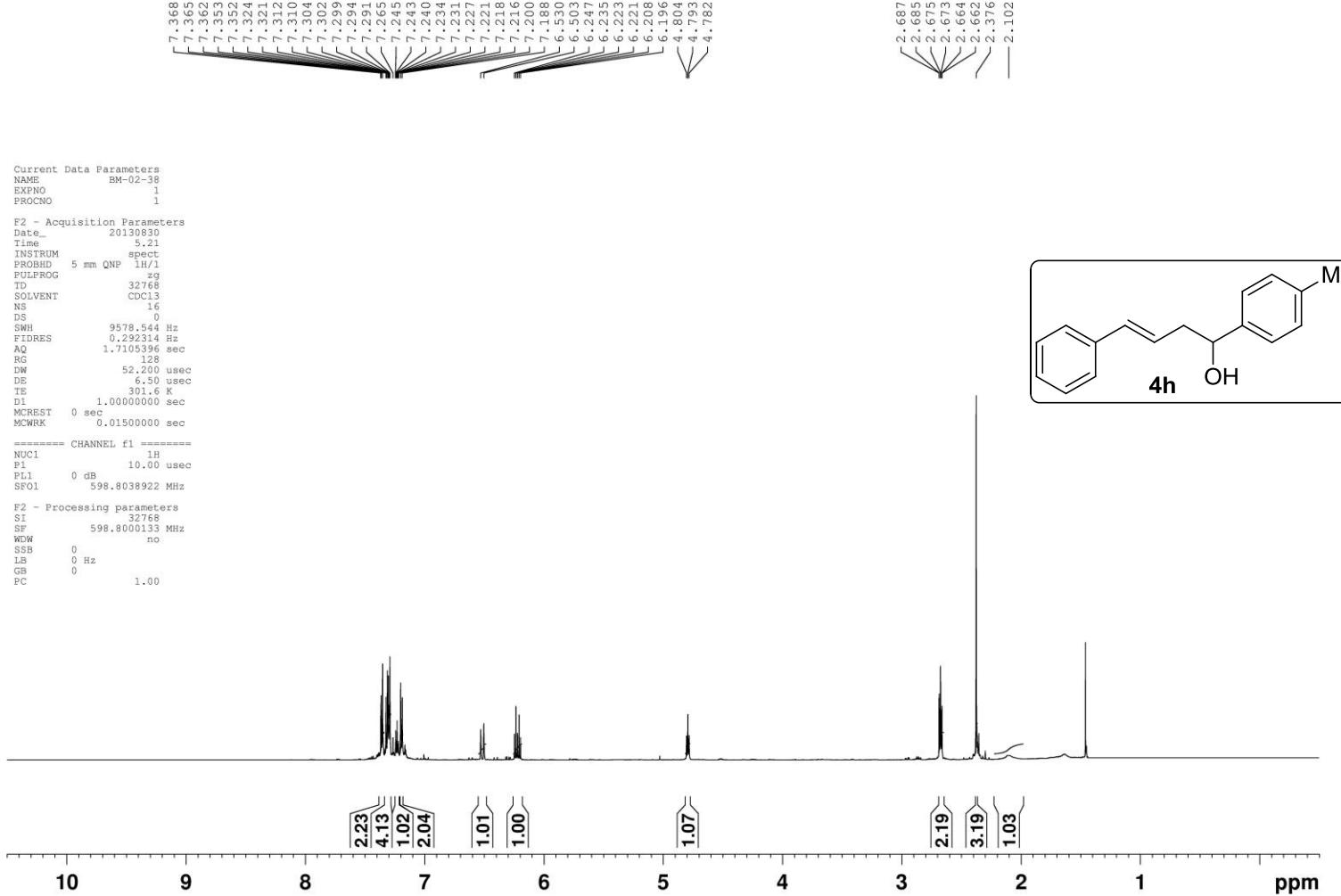
Current Data Parameters
NAME          BM-02-38
EXPNO         1
PROCNO        1

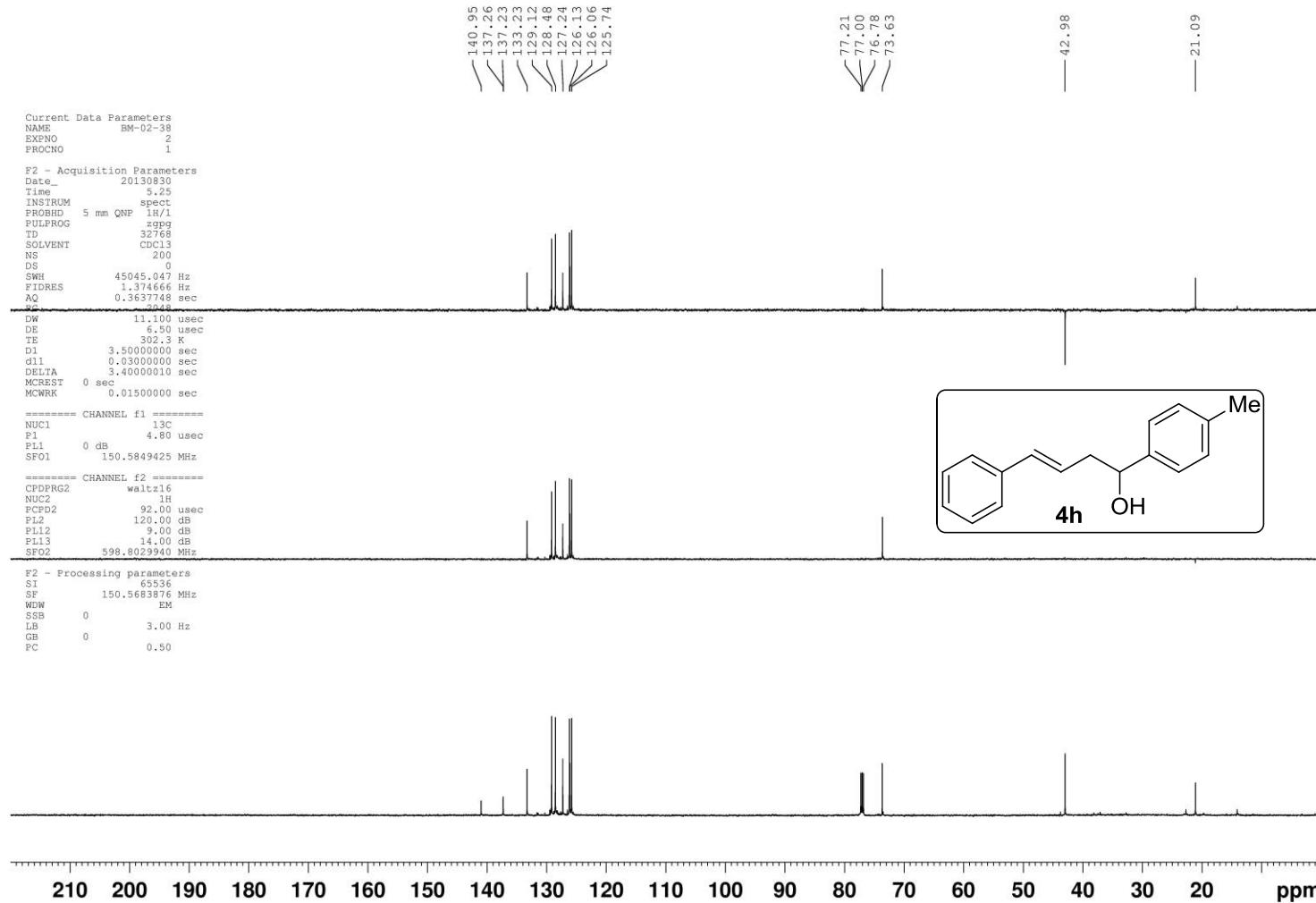
F2 - Acquisition Parameters
Date_       20130830
Time        5.21
INSTRUM     spect
PROBHD      5 mm QNP 1H/1
PULPROG    FID
TD           32768
SOLVENT      CDCl3
NS            16
DS            0
SWH         9578.544 Hz
FIDRES     0.232314 Hz
AQ           1.710536 sec
RG           52,200 usec
DW           52,200 usec
DE           6.50 usec
TE           301.6 K
D1          1.0000000 sec
MCREST      0 sec
MCRWKR     0.01500000 sec

===== CHANNEL f1 =====
NUC1          1H
P1           10.00 usec
PL1          0 dB
SP01        598.8038922 MHz

F2 - Processing parameters
SI            32768
SF          598.8000133 MHz
DW0         no
SSB            0
LB            0 Hz
GB            0
PC           1.00

```





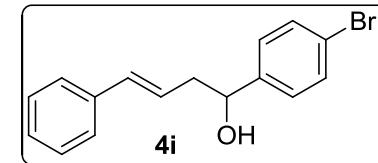
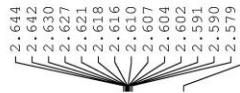
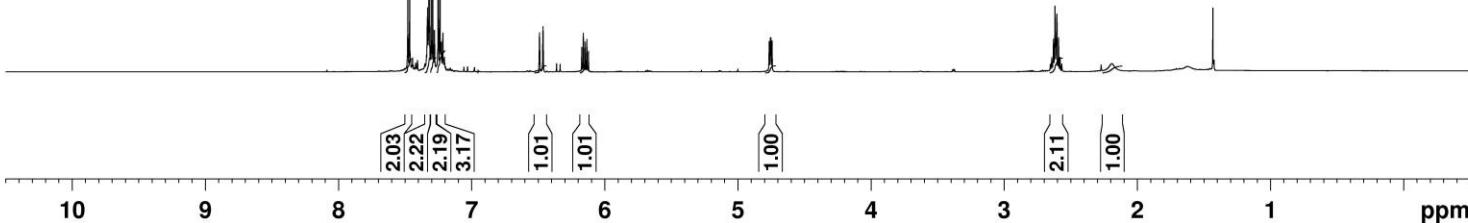
Current Data Parameters
NAME BM-02-40
EXPNO 1
PROCNO 1

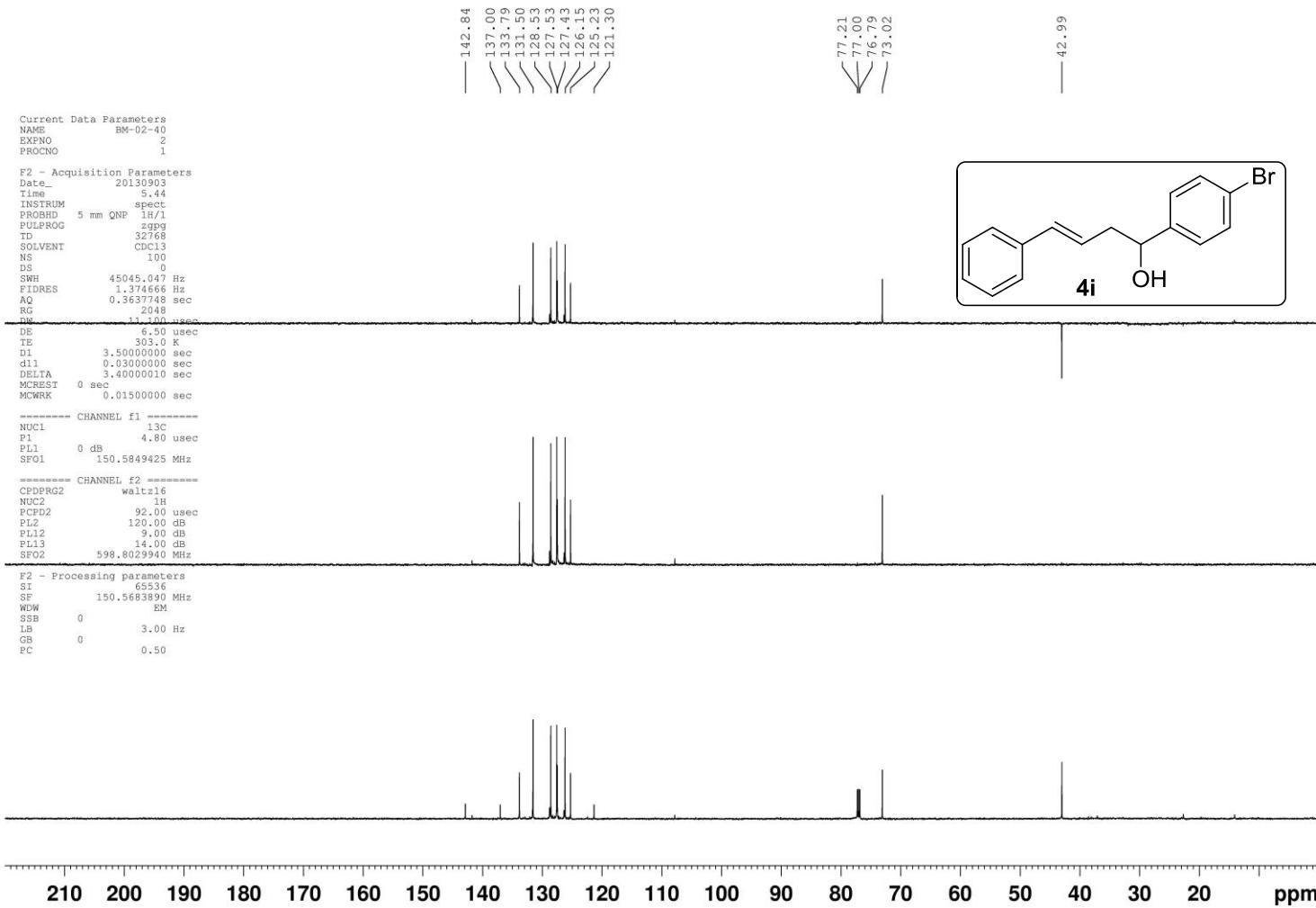
F2 - Acquisition Parameters
Date 20130902
Time 14.37
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 32768
SOLVENT CDCl3
NS 16
DS 0
SWH 9578.544 Hz
FIDRES 0.292314 Hz
AQ 1.7105396 sec
RG 128
DW 52.200 usec
DE 6.5 usec
TE 302.1 K
D1 1.0000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

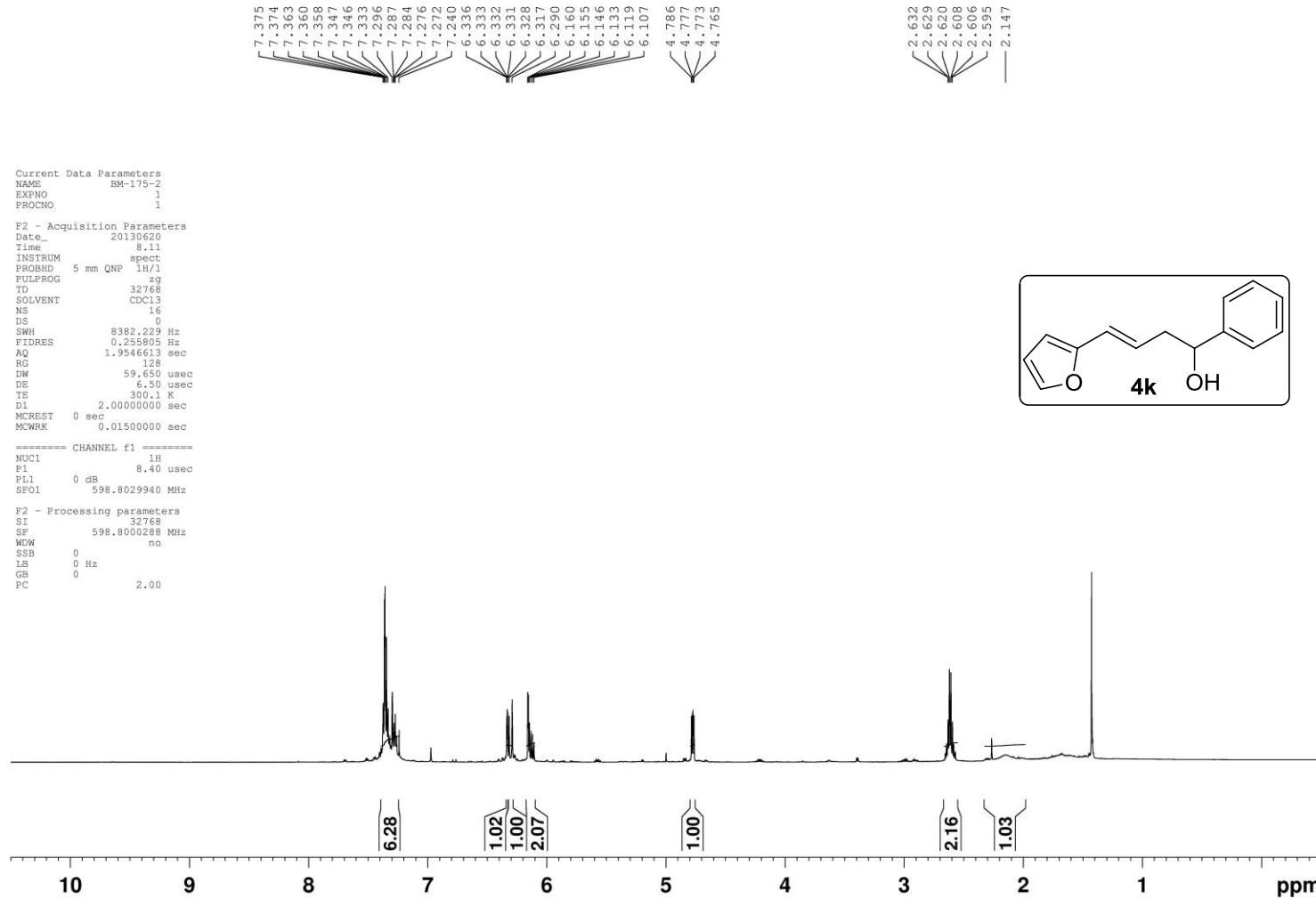
===== CHANNEL f1 =====
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SF01 598.8041916 MHz

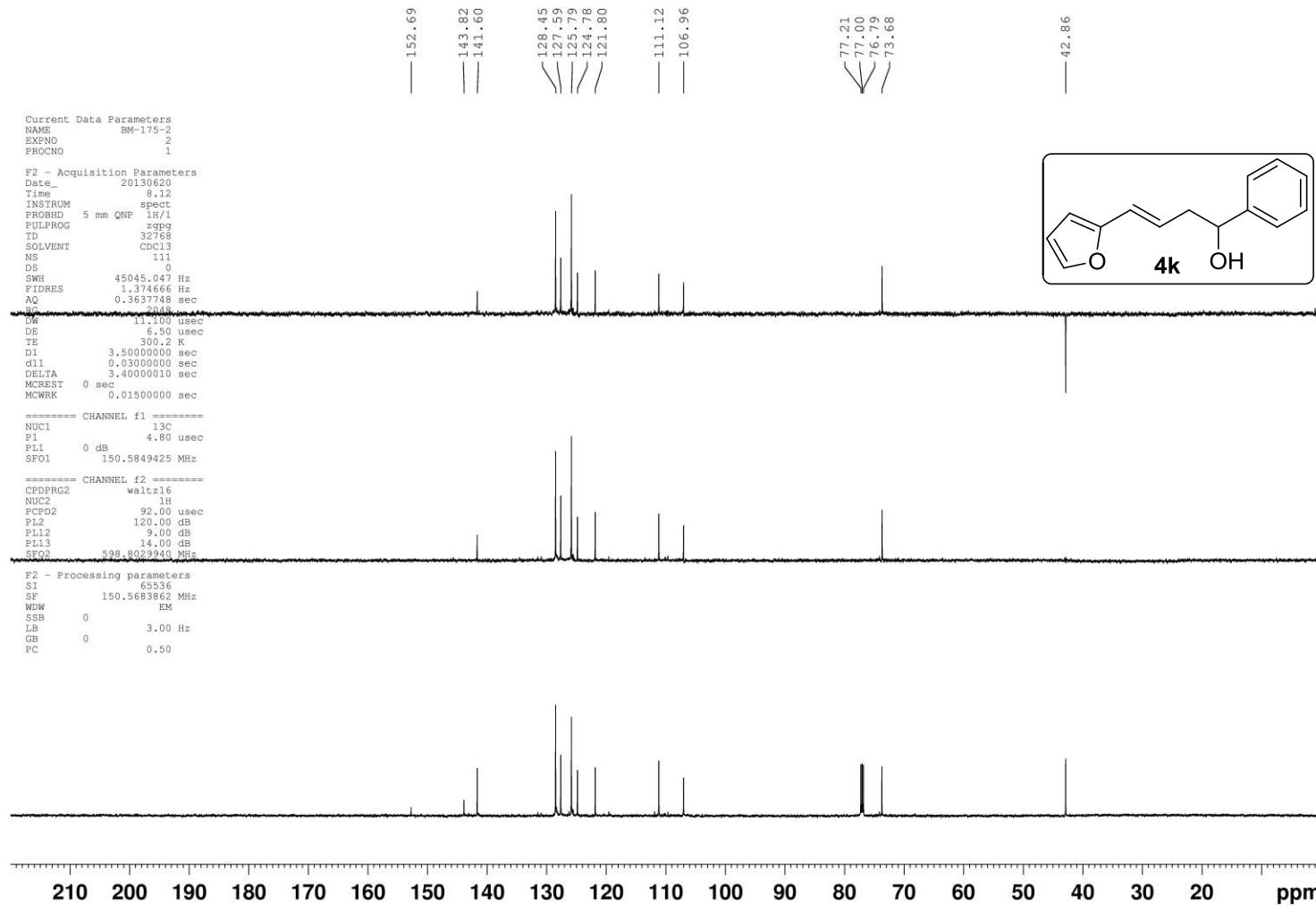
F2 - Processing parameters
SI 32768
SF 598.8000288 MHz
WDW no
SSB 0
LB 0.00 Hz
GB 0
PC 1.00

1D NMR plot parameters
CX 20.00 cm
CY 6.00 cm
F1P 5.088 ppm
F1 3.00 Hz
F2P 2.038 ppm
F2 1220.28 Hz
PPMCM 0.15251 ppm/cm
HZCM 91.32087 Hz/cm









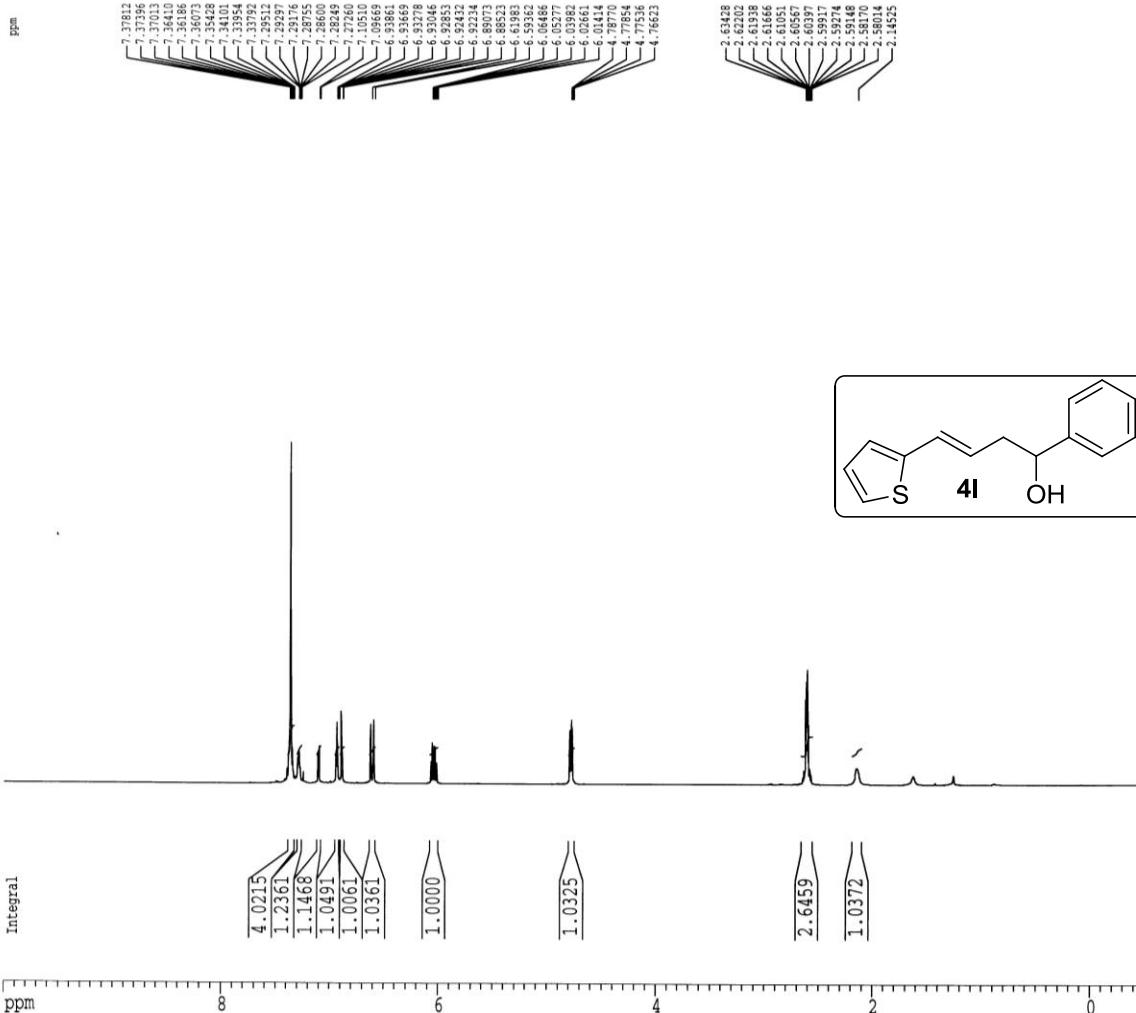
Current Data Parameters
NAME BM-165A
EXPNO 1
PROCNO 1

F2 - Acquisition Parameters
Date_ 20131104
Time 15.49
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zg
TD 32768
SOLVENT CDCl3
NS 16
DS 0
SWH 8389.262 Hz
FIDRES 0.256020 Hz
AQ 1.9530228 sec
RG 128
DW 59.600 usec
DE 6.50 usec
TE 297.7 K
D1 2.0000000 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

===== CHANNEL f1 =====
NUC1 1H
P1 10.00 usec
PL1 0.00 dB
SFO1 598.7029935 MHz

F2 - Processing parameters
SI 32768
SF 598.7000261 MHz
WDW no
SSB 0
LB 0.00 Hz
GB 0
PC 1.00

1D NMR plot parameters
CX 20.00 cm
CY 6.00 cm
F1P 10.000 ppm
F1 5987.00 Hz
F2P -0.500 ppm
F2 -299.35 Hz
PPCM 0.52500 ppm/cm
HZCM 314.31750 Hz/cm



Current Data Parameters
 NAME BM-165A
 EXPNO 2
 PROCNO 1

F2 - Acquisition Parameters
 Date 2011104
 Time 15:56
 INSTRUM spect
 PROBHD 5 mm QNP 1H/1
 PULPROG zgpg
 TD 32768
 SOLVENT CDCl₃
 NS 100
 DS 0
 SWH 45045.047 Hz
 FIDRES 1.374666 Hz
 AQ 0.3637748 sec
 RG 2048
 DW 11.100 usec
 DE 6.50 usec
 TE 298.6 K
 D1 3.5000000 sec
 d11 0.0100000 sec
 DELTA 3.40000010 sec
 MCREST 0.0000000 sec
 MCWRK 0.01500000 sec

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

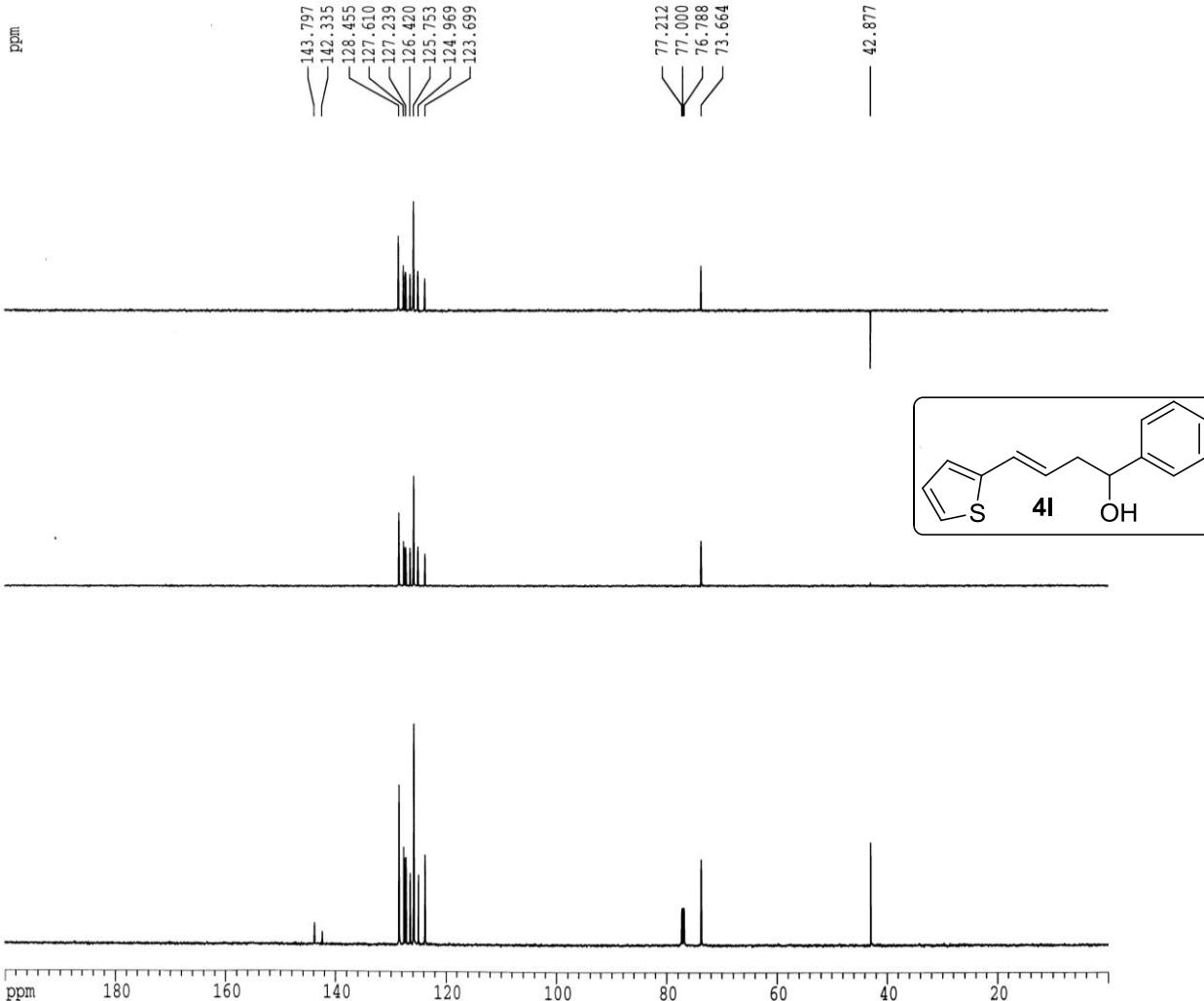
NUC1 13C
 P1 4.80 usec
 PL1 0.00 dB
 SF01 150.5597948 MHz

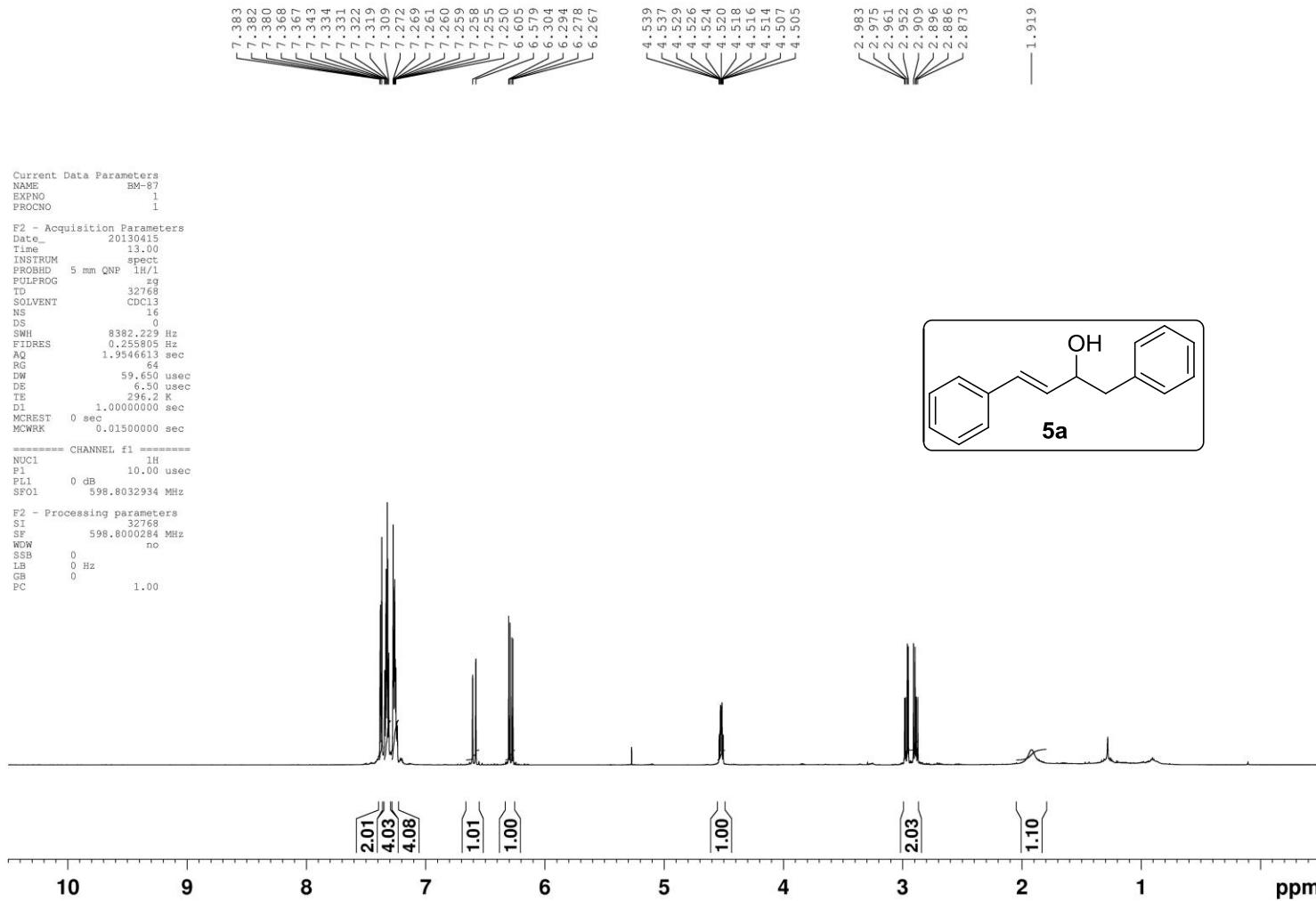
===== CHANNEL f2 ======

CPDPRG2 waltz16
 NUC2 1H
 PCPD2 92.00 usec
 PL2 120.00 dB
 PL12 9.00 dB
 PL13 14.00 dB
 SF02 598.7029935 MHz

F2 - Processing parameters
 SI 65536
 SF 150.5432438 MHz
 WDN EM
 SSB 0
 LB 3.00 Hz
 GB 0
 PC 0.50

1D NMR plot parameters
 CX 20.00 cm
 CY 4.00 cm
 F1P 200.000 ppm
 F1 30108.65 Hz
 F2P 0.000 ppm
 F2 0.00 Hz
 PPMCM 10.00000 ppm/cm
 HZCM 1505.43237 Hz/cm





Current Data Parameters

NAME BM-87
EXPNO 2
PROCNO 1

F2 - Acquisition Parameters

Date 20130414
Time 22.06
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpg
TD 32768
SOLVENT CDCl3
NS 19
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 296.2 K
D1 3.5000000 sec
d11 0.0300000 sec
DELTA 3.40000010 sec
MCREST 0.00000000 sec
MCWRK 0.01500000 sec

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

NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SFO1 150.5849425 MHz

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

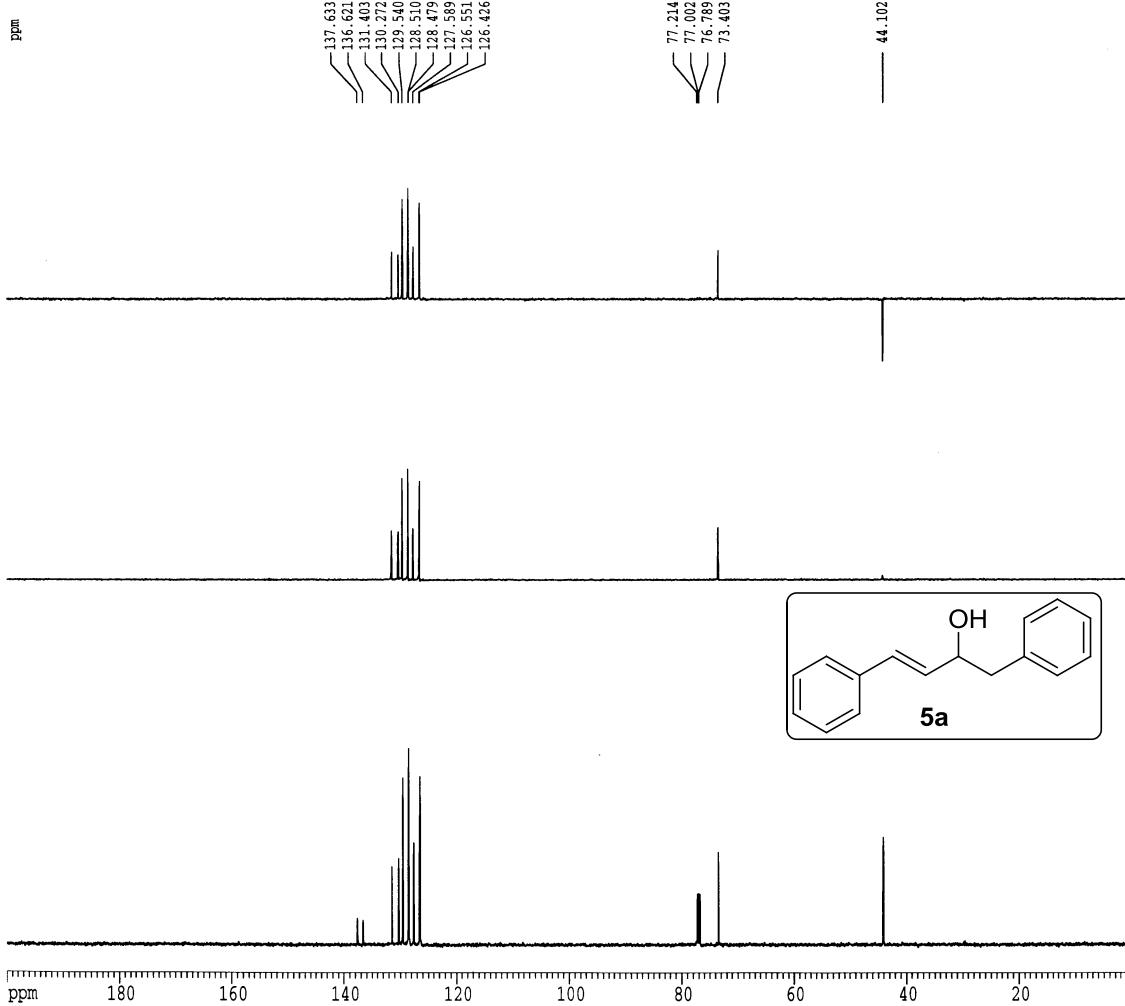
CPPRQ2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SFQ2 598.8029940 MHz

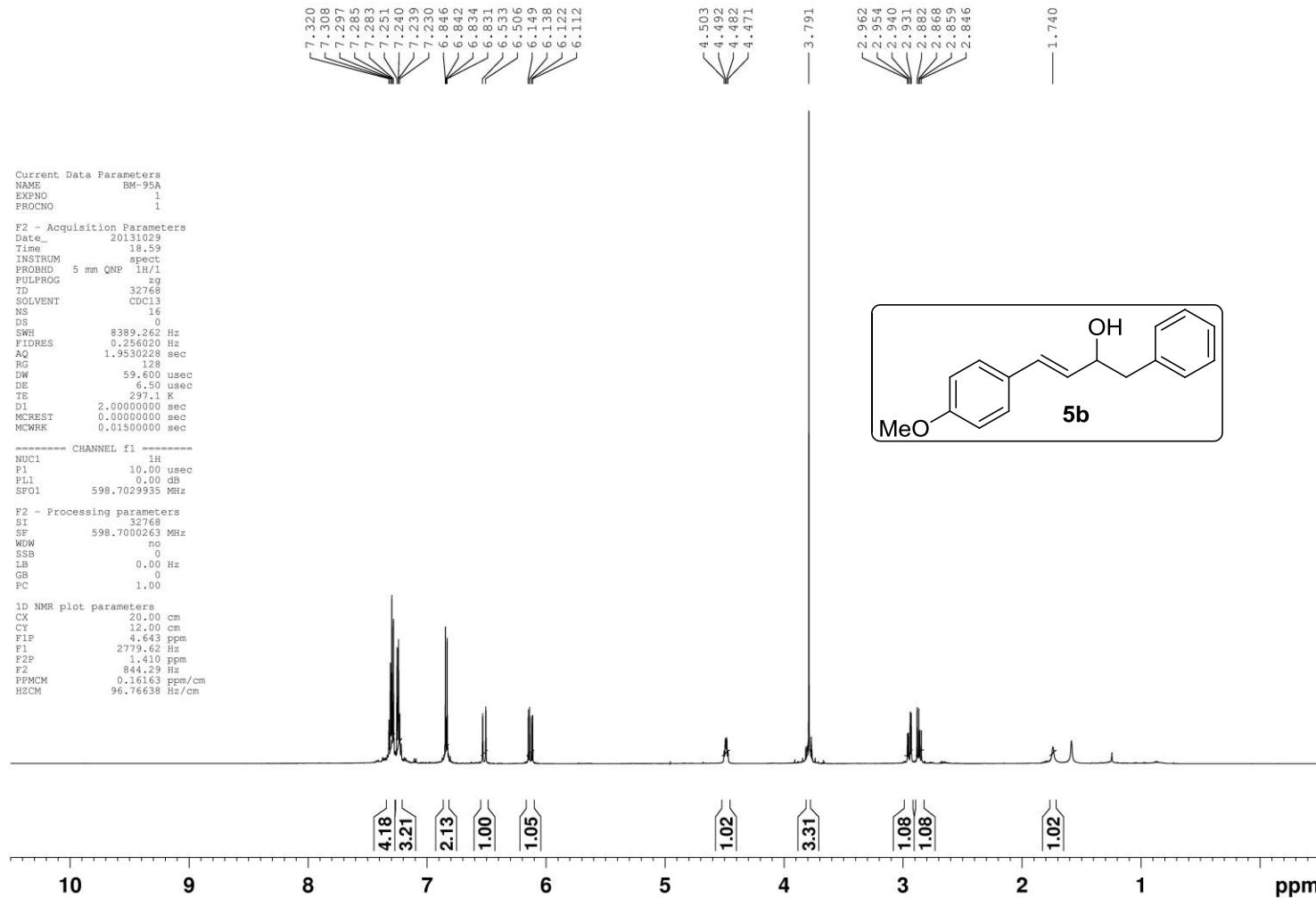
F2 - Processing parameters

SI 65536
SF 150.5683993 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 1.00

1D NMR plot parameters

CX 20.00 cm
CY 3.50 cm
F1P 200.000 ppm
F1 30113.68 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPCM 10.00000 ppm/cm
HZCM 1505.68408 Hz/cm





Current Data Parameters
NAME BM-95A
EXPNO 2
PROCNO 1

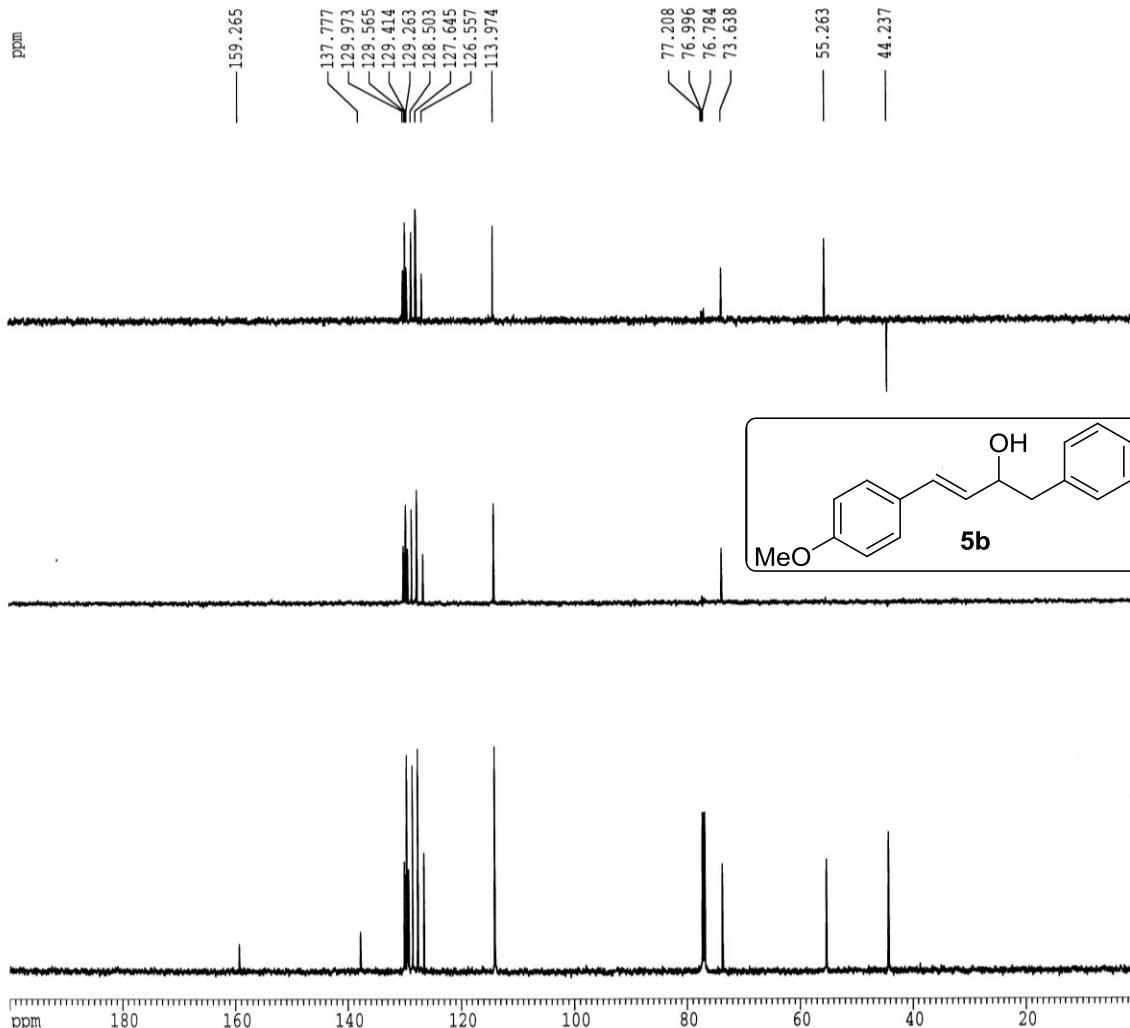
F2 - Acquisition Parameters
Date 20131029
Time 19.04
INSTRUM spect
PROBHD 5 mm QNP 1H/1
PULPROG zgpp
TD 32768
SOLVENT CDCl3
NS 85
DS 0
SWH 45045.047 Hz
FIDRES 1.374666 Hz
AQ 0.3637748 sec
RG 2048
DW 11.100 usec
DE 6.50 usec
TE 298.1 K
D1 3.5000000 sec
d1 0.0300000 sec
DELTA 3.4000010 sec
MCREST 0.0000000 sec
MCWRK 0.0150000 sec

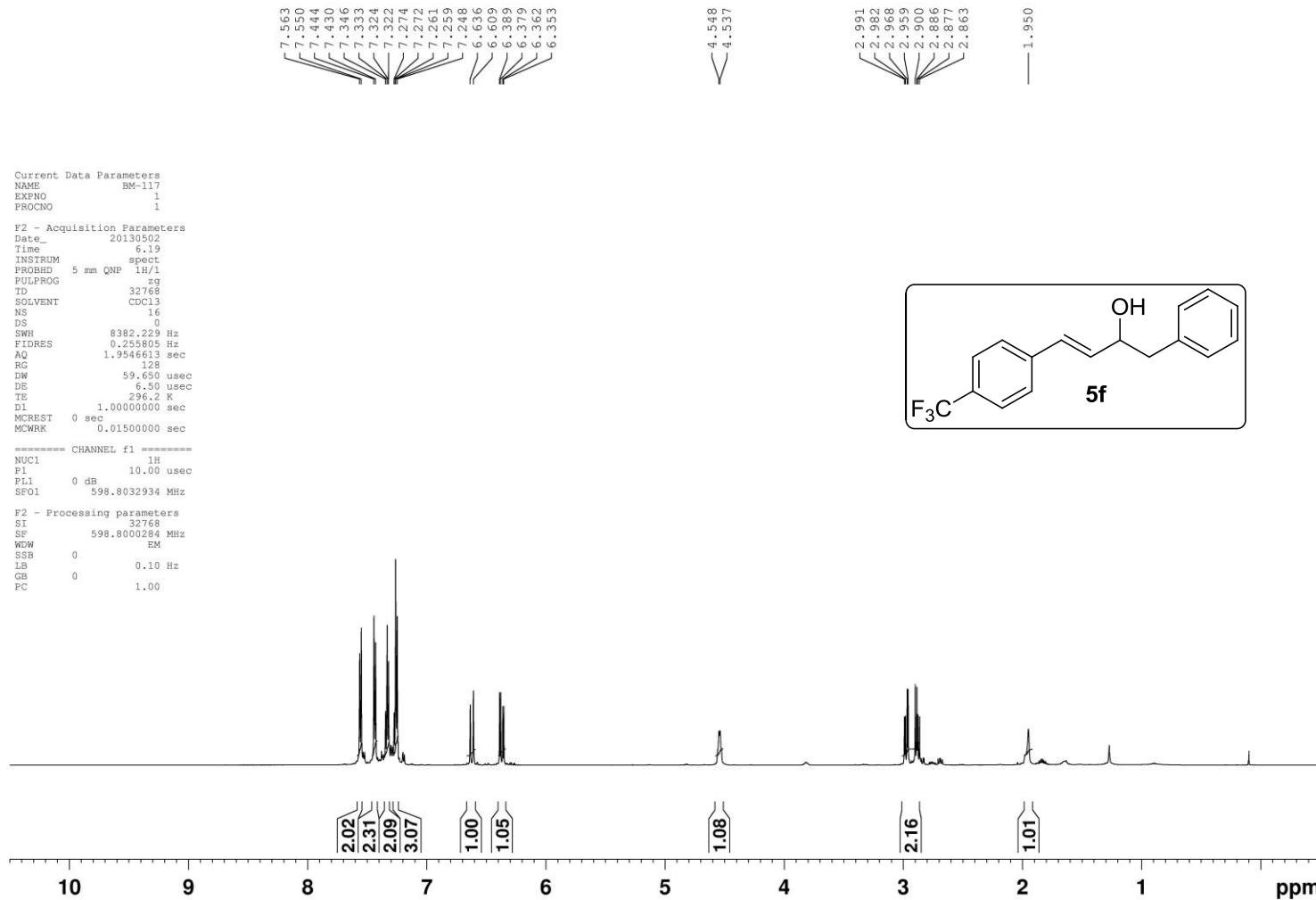
===== CHANNEL f1 =====
NUC1 13C
P1 4.80 usec
PL1 0.00 dB
SF01 150.5597948 MHz

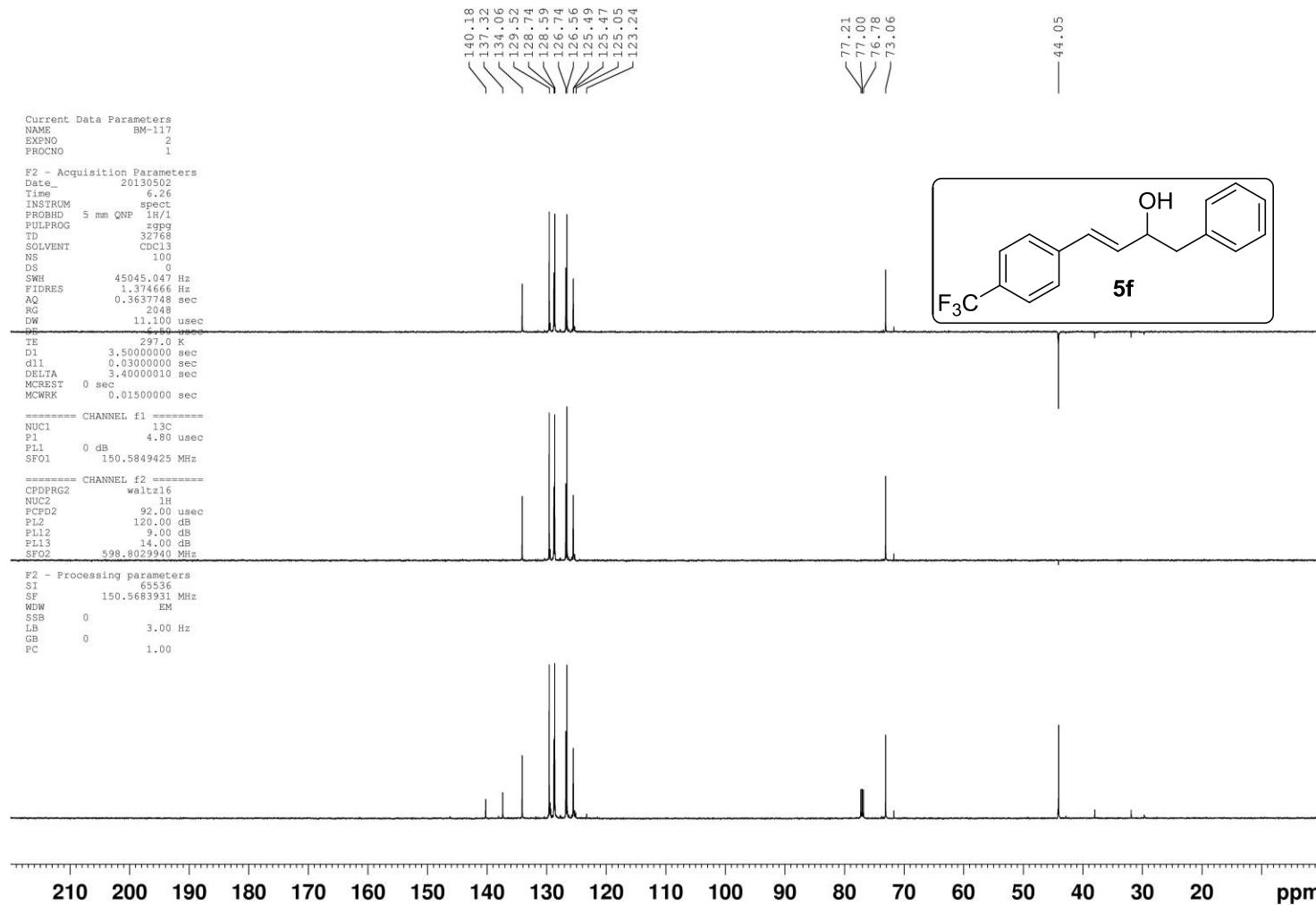
===== CHANNEL f2 =====
CPDPRG2 waltz16
NUC2 1H
PCPD2 92.00 usec
PL2 120.00 dB
PL12 9.00 dB
PL13 14.00 dB
SF02 598.7029935 MHz

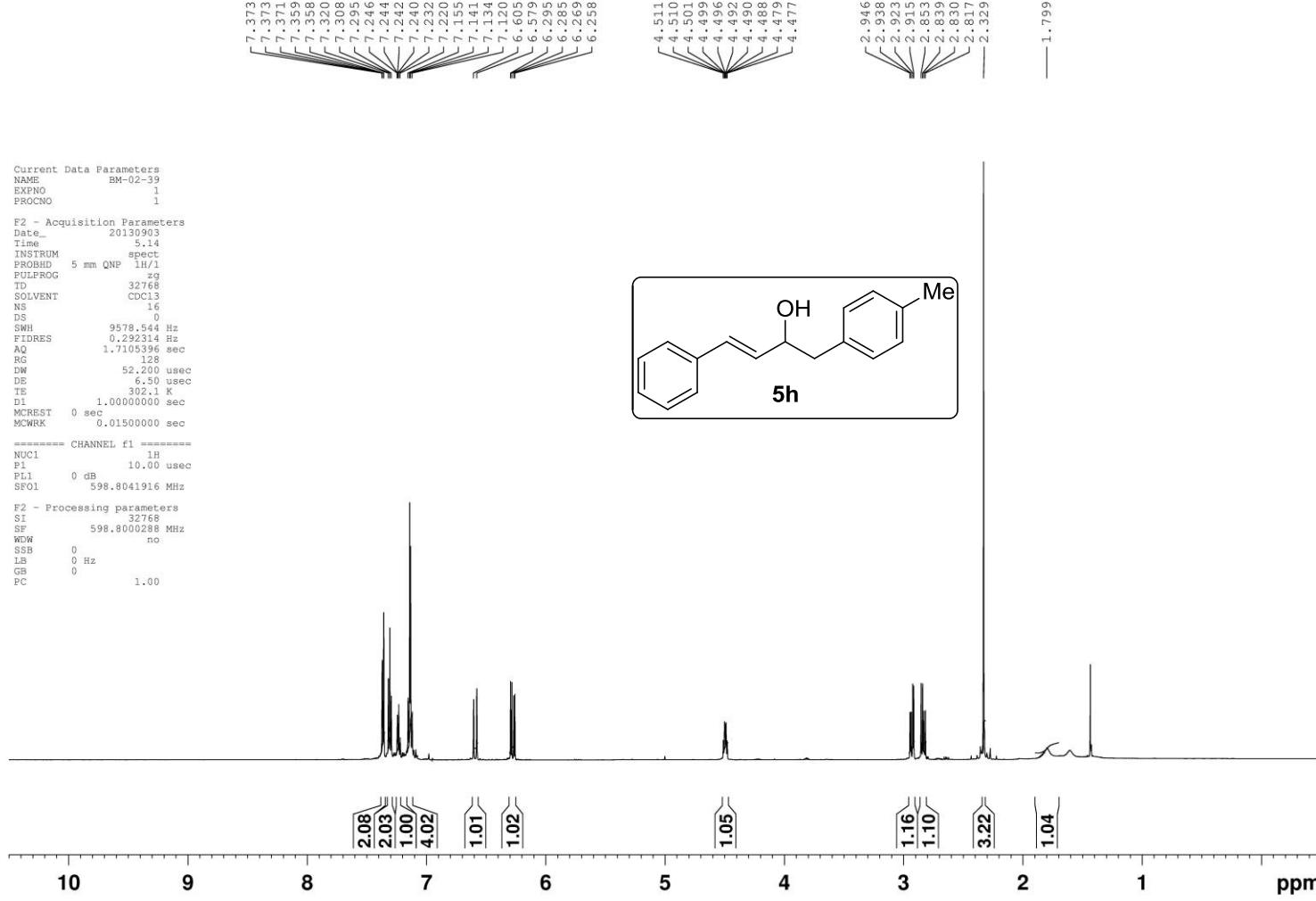
F2 - Processing parameters
SI 65536
SF 150.5432418 MHz
WDW EM
SSB 0
LB 3.00 Hz
GB 0
PC 0.50

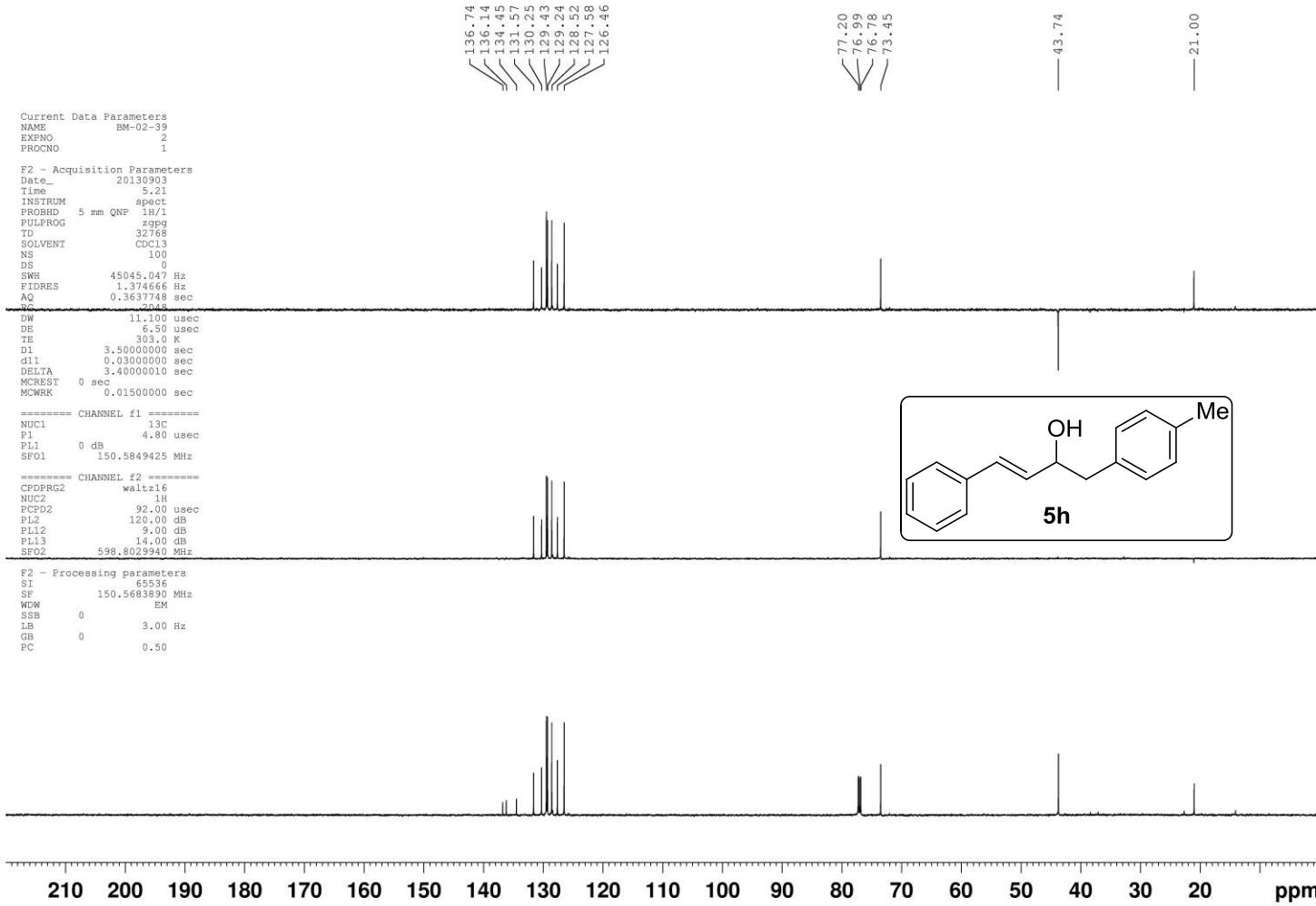
1D NMR plot parameters
CX 20.00 cm
CY 4.00 cm
F1P 200.000 ppm
F1 30108.65 Hz
F2P 0.000 ppm
F2 0.00 Hz
PPCM 10.00000 ppm/cm
HZCM 1505.43237 Hz/cm

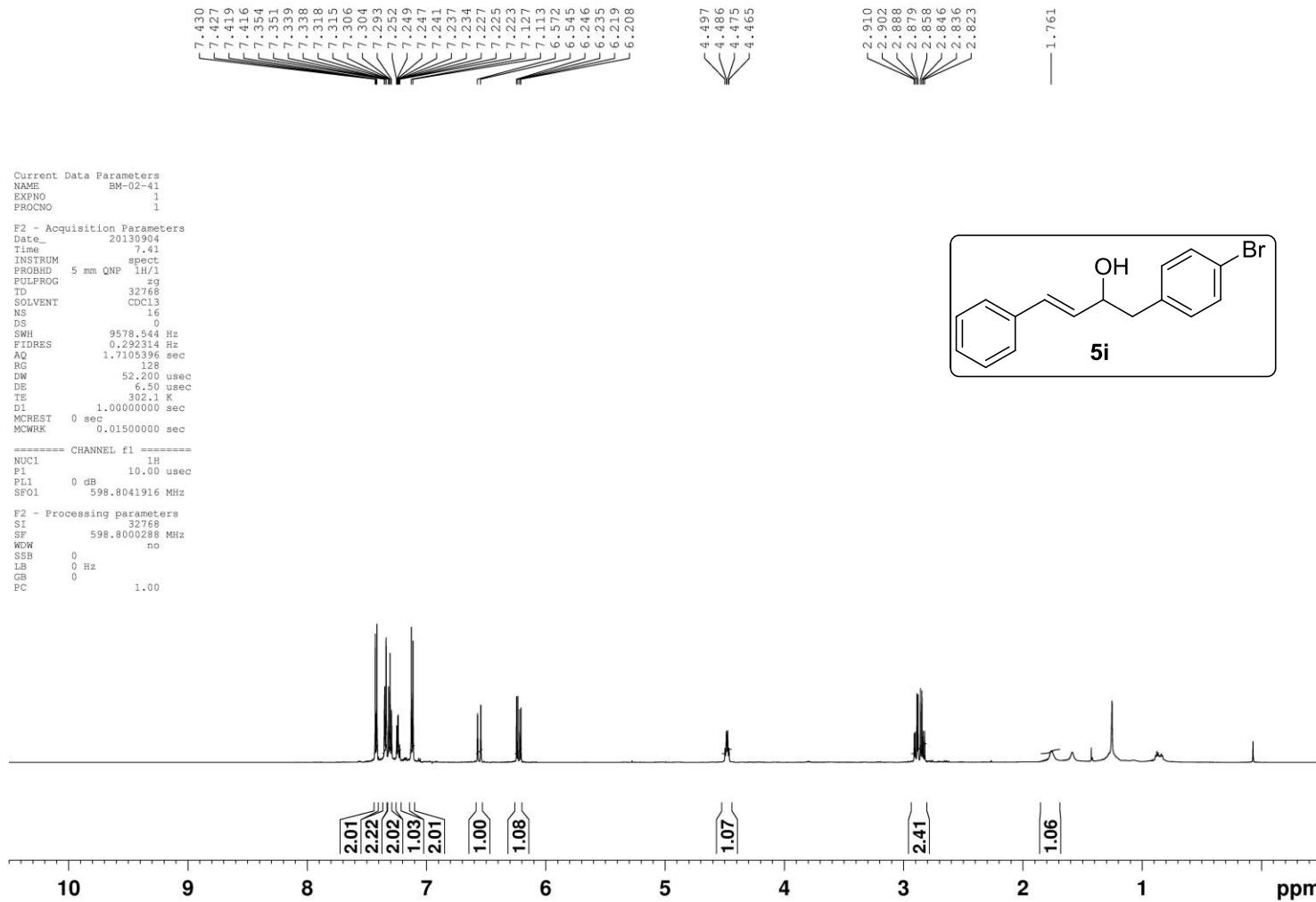


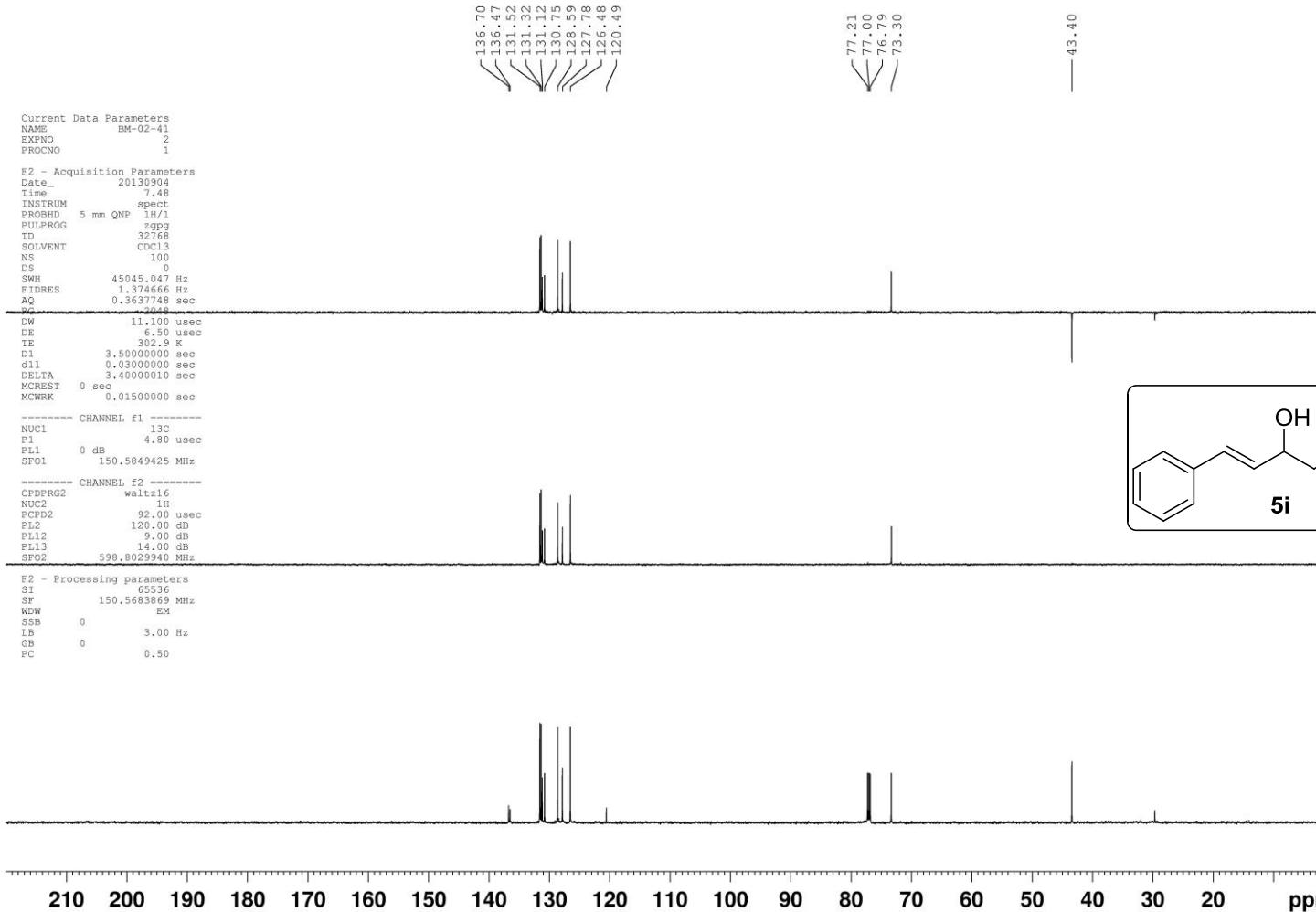


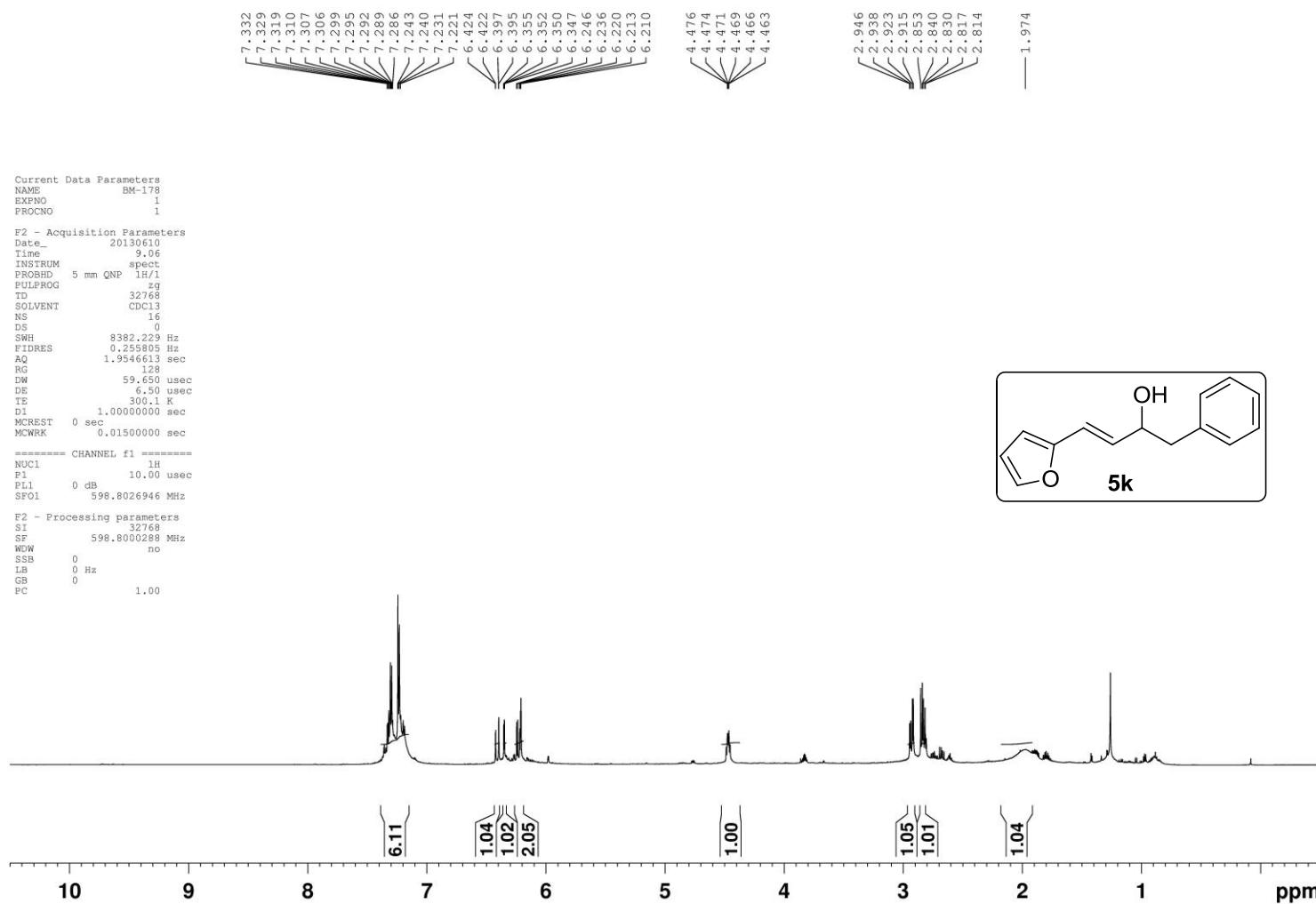


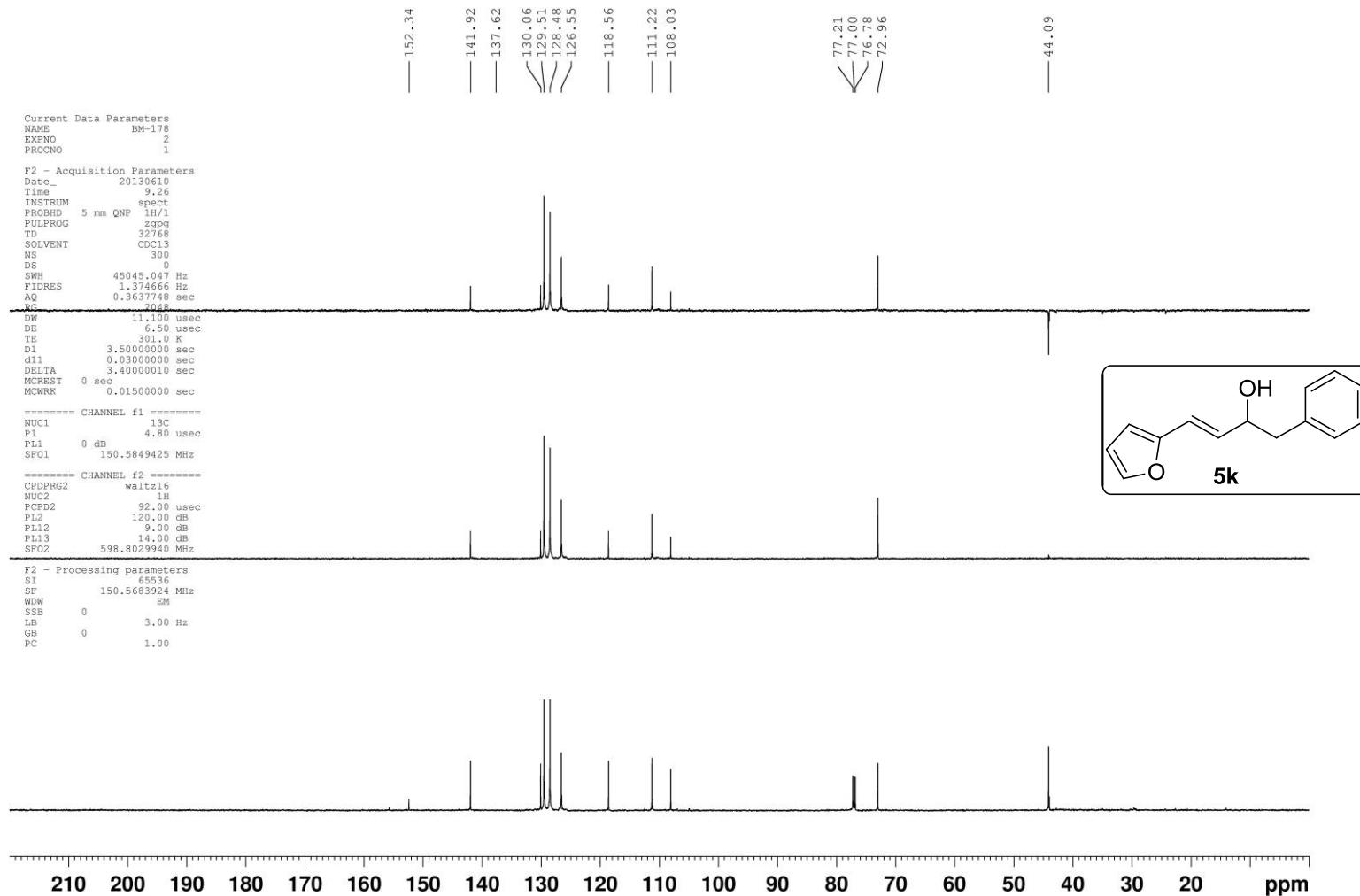












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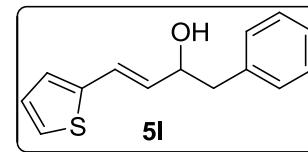
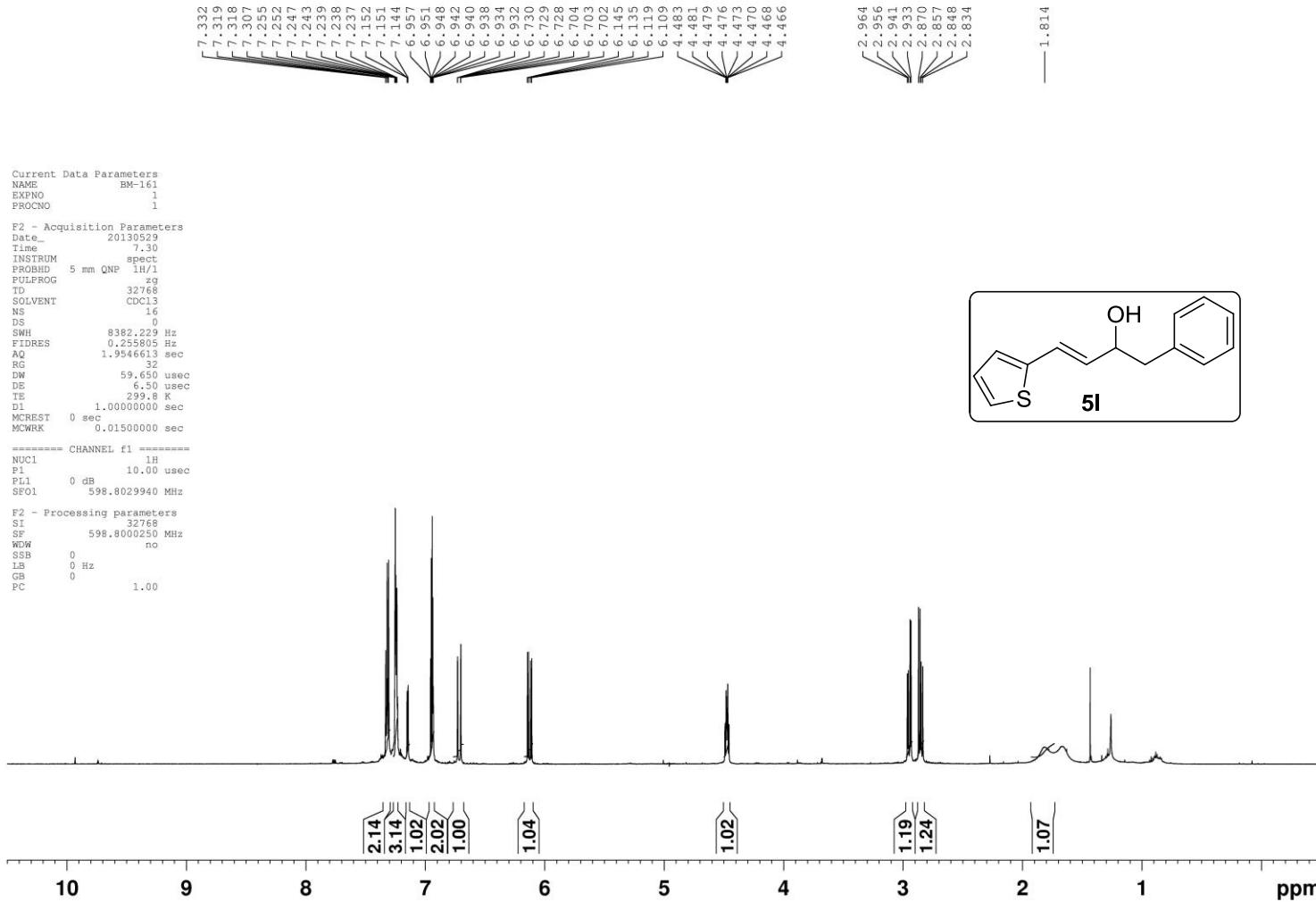
Current Data Parameters
NAME          BM-161
EXPN         1
PROCNO       1

F2 - Acquisition Parameters
Date        20130529
Time        7.30
INSTRUM     spect
PROBHD      5 mm QNP 1H/1
PULPROG    PULFRQG
TD           32768
SOLVENT      CDCl3
NS            16
DS            0
SWH         8382.229 Hz
FIDRES     0.255805 Hz
AQ          1.9546613 sec
RG           1000
DW           59.650 usedc
DE           6.500 usec
TE           299.8 K
D1           1.0000000 sec
MCREST      0 sec
MCRWKR     0.01500000 sec

===== CHANNEL f1 =====
NUC1          1H
P1           10.00 usec
PL1          0 dB
SF01        598.8029990 MHz

F2 - Processing parameters
SI            32768
SF          598.8000250 MHz
WDW        no
SSB           0
LB           0 Hz
GB           0
PC           1.00

```



Current Data Parameters
 NAME BM-161
 EXPNO 2
 PROCN0 1

F2 - Acquisition Parameters
 Date_ 20130528
 Time 16.37
 INSTRUM spect
 PROBHD 5 mm QNP 1H/1
 PULPROG zgpg
 TD 32768
 SOLVENT CDCl3
 NS 100
 DS 0
 SWH 45045.047 Hz
 FIDRES 1.374666 Hz
 AQ 0.3637748 sec
 RG 2048
 DW 11.100 usec
 DE 6.50 usec
 TE 300.4 K
 D1 3.5000000 sec
 d11 0.0300000 sec
 DELTA 3.40000010 sec
 MCREST 0.0900000 sec
 MCWRK 0.0150000 sec

===== CHANNEL f1 =====
 NUC1 13C
 P1 4.80 usec
 PL1 0.00 dB
 SFO1 150.5849425 MHz

===== CHANNEL f2 =====
 CDPRG2 waltz16
 NUC2 1H
 PCPD2 92.00 usec
 PL2 120.00 dB
 PL12 9.00 dB
 PL13 14.00 dB
 SFO2 598.8029940 MHz

F2 - Processing parameters
 SI 65536
 SP 150.5683890 MHz
 NDM EM
 SSB 0
 LB 3.00 Hz
 GB 0
 PC 1.00

1D NMR plot parameters
 CX 20.00 cm
 CY 4.50 cm
 F1P 200.000 ppm
 F1 30113.68 Hz
 F2P 0.000 ppm
 F2 0.00 Hz
 PPMCM 10.00000 ppm/cm
 HZCM 1505.68384 Hz/cm

