

Regioselective hydroarylation and arylation of maleimides with indazoles via a Rh(III)-catalyzed C–H activation

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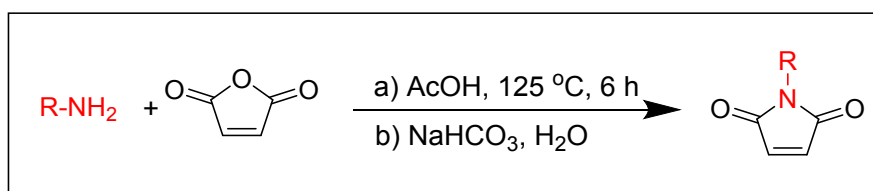
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General Information. All reagents were purchased from commercial sources and used without further purification. ^1H NMR spectra were determined on 400 MHz spectrometer as solutions in CDCl_3 . Chemical shifts are expressed in parts per million (δ) and the signals were reported as s (singlet), d (doublet), t (triplet), m (multiplet) and coupling constants (J) were given in Hz. $^{13}\text{C}\{^1\text{H}\}$ NMR spectra were recorded at 100 MHz in CDCl_3 . Chemical shifts as internal standard are referenced to CDCl_3 ($\delta = 7.26$ for ^1H and $\delta = 77.16$ for $^{13}\text{C}\{^1\text{H}\}$ NMR) as internal standard. TLC was done on silica gel coated glass slide. All solvents were dried and distilled before use. All reactions involving moisture sensitive reactants were executed using oven dried glassware. Melting points (M.p.) were determined after recrystallization of solid compounds from a solution of dichloromethane/petroleum ether (1:3).

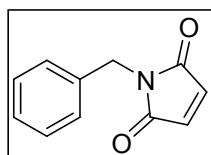
Starting Materials. All 2*H*-indazoles was prepared by the reported method.¹



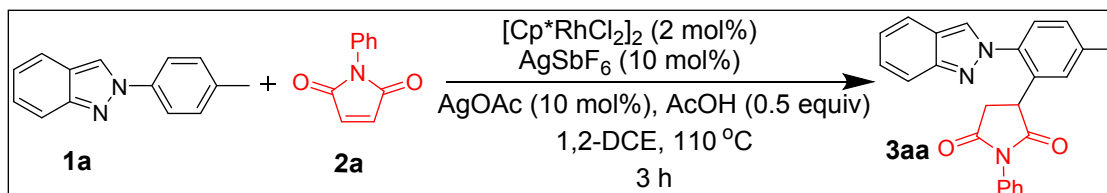
General procedure for the preparation of *N*-aryl & *N*-alkyl maleimides derivatives:² Maleic anhydride (2.0 equiv) were stirred in acetic acid (1.5 mL per mmol of amine) until maleic anhydride was dissolved completely. Then primary amine was added (1.0 equiv) to the reaction mixture and refluxed for 6 h at 125 °C. After completion of the reaction, the reaction mixture was allowed to cool down at room temperature and transfer to a 250 mL beaker. Saturated NaHCO_3 aqueous solution was added to the beaker containing reaction mixture until effervescence of CO_2 stop. Then, the aqueous mixture was extracted with ethyl acetate (3 x 10 mL). The organic phase was dried over anhydrous Na_2SO_4 . The crude residue was obtained after evaporating the solvent in vacuum. Then crude residue was further washed with 1(N) HCl (2 x

20 mL) and brine solution (5 mL) respectively. The aqueous mixture was again extracted with ethyl acetate (3 x 10 mL). Finally, excess solvent was removed under reduced pressure and the residue was purified by column chromatography using ethyl acetate/hexane to get extra pure maleimide with high yields.

2. Characterization data for the synthesized products:

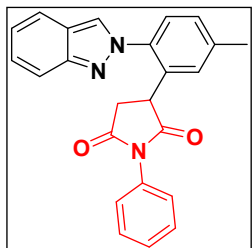


1-Benzyl-1H-pyrrole-2,5-dione (2h): Yellow solid (90%, 33 mg); $R_f = 0.50$ (PE : EA = 80 : 20); ^1H NMR (400 MHz, CDCl_3): δ 7.30-7.20 (m, 5H), 6.64 (s, 2H), 4.62 (s, 2H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 170.5, 136.3, 134.2, 128.8, 128.4, 127.9, 41.5.

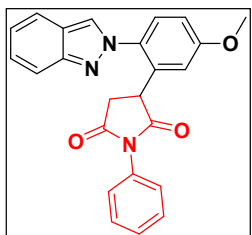


Typical experimental procedure for the synthesized compounds (3aa-3ga): A mixture of 2-(*p*-tolyl)-2*H*-indazole (0.2 mmol, 41.6 mg) (**1a**), $[\text{Cp}^*\text{RhCl}_2]_2$ (2 mol%, 2.4 mg), AgSbF_6 (10 mol%, 6.8 mg), and AgOAc (10 mol%, 3.3 mg) were taken in an oven dried screw-capped reaction tube. Then 1,2-DCE (2 mL) and AcOH (0.5 equiv, 6.0 mg) were added to the mixture and stirred for 5 min at room temperature under open atmosphere. After that, 1-phenyl-1*H*-pyrrole-2,5-dione (**2a**) (0.24 mmol, 41.5 mg) was added, and the resultant mixture was stirred at 110 °C for 3 h. After completion of the reaction (TLC), the reaction was cooled to room temperature and extracted with dichloromethane. The organic phase was dried over anhydrous Na_2SO_4 . The crude residue was obtained after evaporating the solvent in vacuum and was

purified by column chromatography on silica gel using a mixture of petroleum ether and ethyl acetate (72:28) as an eluting solvent to afford the pure product **3aa** (67 mg, 89%) as a light yellow solid.

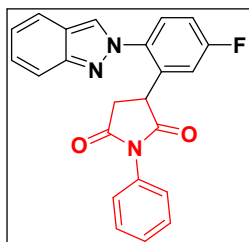


3-(2-(2H-Indazol-2-yl)-5-methylphenyl)-1-phenylpyrrolidine-2,5-dione (3aa): Light yellow solid (89%, 67 mg); $R_f = 0.50$ (PE : EA = 72 : 28); M.p. 143-144 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.24 (s, 1H), 7.71-7.65 (m, 2H), 7.38 (d, $J = 8.0$ Hz, 1H), 7.35-7.28 (m, 5H), 7.23 (s, 1H), 7.15-7.11 (m, 1H), 6.92-6.90 (m, 2H), 4.21 (t, $J = 8.4$ Hz, 1H), 3.26 (d, $J = 8.4$ Hz, 2H), 2.46 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 176.1, 175.1, 149.7, 140.1, 137.7, 132.7, 131.8, 131.2, 129.7, 128.9, 128.4, 127.09, 127.07, 126.3, 125.3, 122.6, 122.4, 120.6, 117.9, 43.9, 38.0, 21.3; Anal. Calcd for $\text{C}_{24}\text{H}_{19}\text{N}_3\text{O}_2$: C, 75.57; H, 5.02; N, 11.02%; Found: C, 75.38; H, 4.98; N, 11.11%.

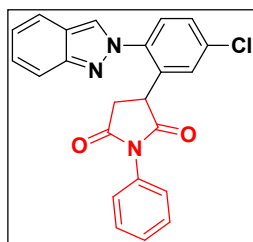


3-(2-(2H-Indazol-2-yl)-5-methoxyphenyl)-1-phenylpyrrolidine-2,5-dione (3ba): Brown solid (73%, 58 mg); $R_f = 0.45$ (PE : EA = 65 : 35); M.p. 148-149 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.21 (s, 1H), 7.71-7.65 (m, 2H), 7.42 (d, $J = 8.8$ Hz, 1H), 7.35-7.28 (m, 4H), 7.13 (t, $J = 7.6$ Hz, 1H), 7.00-6.98 (m, 1H), 6.94 (d, $J = 2.8$ Hz, 1H), 6.89-6.87 (m, 2H), 4.20-4.16 (m, 1H), 3.89 (s,

3H), 3.28-3.25 (m, 2H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 175.9, 175.0, 160.4, 149.7, 134.5, 133.2, 131.8, 129.0, 128.6, 128.5, 127.0, 126.4, 125.6, 122.6, 122.4, 120.6, 117.9, 116.3, 113.4, 55.8, 44.0, 37.9; Anal. Calcd for $\text{C}_{24}\text{H}_{19}\text{N}_3\text{O}_3$: C, 72.53; H, 4.82; N, 10.57; Found: C, 72.66; H, 4.85; N, 10.52%.

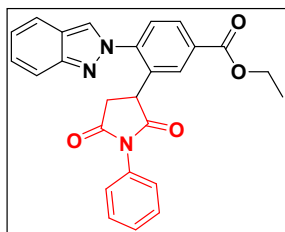


3-(5-Fluoro-2-(2H-indazol-2-yl)phenyl)-1-phenylpyrrolidine-2,5-dione (3ca): White solid (81%, 62 mg); R_f = 0.50 (PE : EA = 67 : 33); M.p. 177-178 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.23 (s, 1H), 7.71-7.65 (m, 2H), 7.50-7.47 (m, 1H), 7.35-7.29 (m, 4H), 7.23-7.13 (m, 3H), 6.89-6.87 (m, 2H), 4.22 (t, J = 8.4 Hz, 1H), 3.27-3.25 (m, 2H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 175.3, 174.6, 162.7 (d, $J_{\text{C-F}}$ = 249.0 Hz), 149.9, 136.4, 136.3, 135.5 (d, $J_{\text{C-F}}$ = 8.0 Hz), 131.6, 129.2 (d, $J_{\text{C-F}}$ = 8.0 Hz), 129.0, 128.6, 127.3, 126.3, 125.5, 122.7 (d, $J_{\text{C-F}}$ = 8.0 Hz), 120.6, 117.9 (d, $J_{\text{C-F}}$ = 10.0 Hz), 117.6, 116.0 (d, $J_{\text{C-F}}$ = 22.0 Hz), 43.9, 37.7; Anal. Calcd for $\text{C}_{23}\text{H}_{16}\text{FN}_3\text{O}_2$: C, 71.68; H, 4.18; N, 10.90%; Found: C, 71.85; H, 4.23; N, 10.79%.

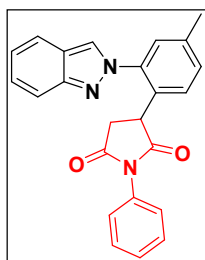


3-(5-Chloro-2-(2H-indazol-2-yl)phenyl)-1-phenylpyrrolidine-2,5-dione (3da): White solid (91%, 73 mg); R_f = 0.50 (PE : EA = 71 : 29); M.p. 158-159 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.23 (s, 1H), 7.71-7.62 (m, 2H), 7.49-7.42 (m, 3H), 7.34-7.27 (m, 4H), 7.16-7.13 (m, 1H), 6.87-

6.83 (m, 2H), 4.25-4.21 (m, 1H), 3.28-3.24 (m, 2H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 175.3, 174.6, 149.9, 138.5, 135.5, 134.5, 131.6, 131.2, 129.2, 128.9, 128.5, 128.3, 127.4, 126.2, 125.2, 122.9, 122.5, 120.6, 117.9, 44.1, 37.7; HRMS (ESI-TOF) m/z : $[\text{M} + \text{Na}]^+$ Calcd for $\text{C}_{23}\text{H}_{16}\text{ClN}_3\text{O}_2\text{Na}$: 424.0823; found: 424.0816.

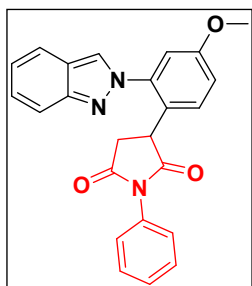


Ethyl-3-(2,5-dioxo-1-phenylpyrrolidin-3-yl)-4-(2H-indazol-2-yl)benzoate (3ea): Light yellow solid (79%, 69 mg); $R_f = 0.45$ (PE : EA = 69 : 31); M.p. 169-170 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.32 (s, 1H), 8.19-8.16 (m, 2H), 7.72 (d, $J = 8.8$ Hz, 1H), 7.63-7.59 (m, 2H), 7.34-7.27 (m, 4H), 7.17-7.13 (m, 1H), 6.88-6.86 (m, 2H), 4.47-4.38 (m, 3H), 3.37-3.34 (m, 2H), 1.46-1.42 (m, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 175.5, 174.9, 165.2, 150.2, 143.3, 133.0, 132.8, 131.7, 131.5, 130.3, 129.0, 128.5, 127.6, 126.9, 126.3, 125.1, 123.1, 122.7, 120.7, 118.0, 61.8, 44.6, 37.9, 14.4; HRMS (ESI-TOF) m/z : $[\text{M} + \text{H}]^+$ Calcd for $\text{C}_{26}\text{H}_{22}\text{N}_3\text{O}_4$: 440.1605; found: 440.1602.

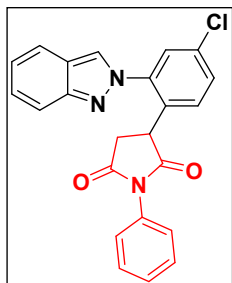


3-(2-(2H-Indazol-2-yl)-4-methylphenyl)-1-phenylpyrrolidine-2,5-dione (3fa): Yellow solid (93%, 71 mg); $R_f = 0.45$ (PE : EA = 74 : 26); M.p. 138-139 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.27 (s, 1H), 7.71-7.66 (m, 2H), 7.34-7.28 (m, 7H), 7.13 (t, $J = 8.0$ Hz, 1H), 6.91-6.89 (m, 2H),

4.21-4.17 (m, 1H), 3.22-3.20 (m, 2H), 2.43 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 176.1, 175.1, 149.7, 139.7, 139.3, 131.8, 130.4, 129.8, 128.8, 128.3, 127.7, 127.0, 126.3, 125.2, 122.5, 122.3, 120.5, 117.9, 43.6, 37.8, 20.9; Anal. Calcd for $\text{C}_{24}\text{H}_{19}\text{N}_3\text{O}_2$: C, 75.57; H, 5.02; N, 11.02%; Found: C, 75.76; H, 5.05; N, 11.10%.

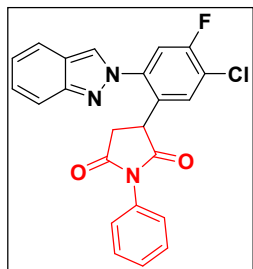


3-(2-(2H-Indazol-2-yl)-4-methoxyphenyl)-1-phenylpyrrolidine-2,5-dione (3ga): Light yellow solid (88%, 69 mg); R_f = 0.45 (PE : EA = 59 : 41); M.p. 147-148 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.25 (s, 1H), 7.68-7.63 (m, 2H), 7.30-7.25 (m, 5H), 7.11 (t, J = 7.6 Hz, 1H), 7.03-7.01 (m, 2H), 6.87 (d, J = 8.0 Hz, 2H), 4.13 (t, J = 8.0 Hz, 1H), 3.81 (s, 3H), 3.16 (d, J = 8.0 Hz, 2H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 176.3, 175.1, 159.7, 149.7, 140.7, 131.8, 131.7, 128.8, 128.3, 127.1, 126.3, 125.2, 124.6, 122.6, 122.3, 120.6, 117.9, 115.4, 112.7, 55.7, 43.4, 37.8; Anal. Calcd for $\text{C}_{24}\text{H}_{19}\text{N}_3\text{O}_3$: C, 72.53; H, 4.82; N, 10.57%; Found: C, 72.75; H, 4.88; N, 10.48%.

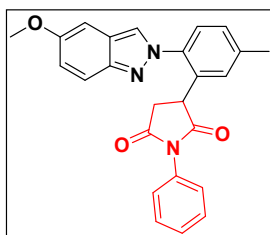


3-(4-Chloro-2-(2H-indazol-2-yl)phenyl)-1-phenylpyrrolidine-2,5-dione (3ha): Brown solid (85%, 68 mg); R_f = 0.45 (PE : EA = 70 : 30); M.p. 153-154 °C; ^1H NMR (400 MHz, CDCl_3): δ

8.25 (s, 1H), 7.70-7.62 (m, 2H), 7.51-7.45 (m, 2H), 7.36-7.27 (m, 5H), 7.14 (t, $J = 8.0$ Hz, 1H), 6.87-6.84 (m, 2H), 4.23 (t, $J = 8.0$ Hz, 1H), 3.21 (d, $J = 7.6$ Hz, 2H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 175.5, 174.7, 149.9, 140.7, 134.4, 132.3, 131.6, 131.2, 129.7, 128.9, 128.4, 127.4, 127.1, 126.2, 125.1, 122.9, 122.4, 120.6, 117.9, 43.8, 37.6; Anal. Calcd for $\text{C}_{23}\text{H}_{16}\text{ClN}_3\text{O}_2$: C, 68.75; H, 4.01; N, 10.46%; Found: C, 68.90; H, 3.97; N, 10.35%.

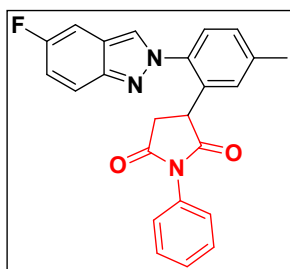


3-(5-Chloro-4-fluoro-2-(2H-indazol-2-yl)phenyl)-1-phenylpyrrolidine-2,5-dione (3ia): White solid (87%, 72 mg); $R_f = 0.50$ (PE : EA = 65 : 35); M.p. 173-174 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.23 (s, 1H), 7.71-7.58 (m, 3H), 7.33-7.25 (m, 5H), 7.15 (t, $J = 8.4$ Hz, 1H), 6.86-6.83 (m, 2H), 4.22 (t, $J = 8.0$ Hz, 1H), 3.25 (d, $J = 8.0$ Hz, 2H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 175.0, 174.4, 158.2 (d, $J_{\text{C-F}} = 252.0$ Hz), 150.0, 136.6 (d, $J_{\text{C-F}} = 4.0$ Hz), 133.5 (d, $J_{\text{C-F}} = 6.0$ Hz), 131.5, 129.5, 129.0, 128.6, 127.6, 126.2, 125.4, 123.2, 122.6, 121.5 (d, $J_{\text{C-F}} = 19.0$ Hz), 120.6, 118.9 (d, $J_{\text{C-F}} = 23.0$ Hz), 117.9, 43.8, 37.6; Anal. Calcd for $\text{C}_{23}\text{H}_{15}\text{ClFN}_3\text{O}_2$: C, 65.80; H, 3.60; N, 10.01%; Found: C, 65.64; H, 3.62; N, 9.92%.

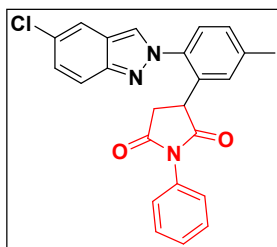


3-(2-(5-Methoxy-2H-indazol-2-yl)-5-methylphenyl)-1-phenylpyrrolidine-2,5-dione (3ja): Brown gummy mass (86%, 70 mg); $R_f = 0.55$ (PE : EA = 64 : 36); ^1H NMR (400 MHz, CDCl_3):

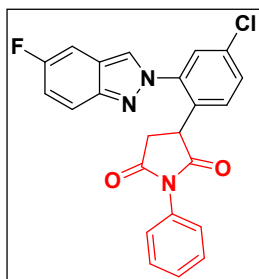
δ 8.07 (s, 1H), 7.54 (d, J = 9.6 Hz, 1H), 7.36-7.25 (m, 5H), 7.21 (s, 1H), 7.01-6.98 (m, 1H), 6.95-6.93 (m, 2H), 6.88 (d, J = 2.4 Hz, 1H), 4.20 (t, J = 8.0 Hz, 1H), 3.82 (s, 3H), 3.20 (d, J = 8.0 Hz, 2H), 2.44 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 176.1, 175.1, 155.6, 146.5, 139.7, 137.5, 132.4, 131.8, 131.4, 129.5, 128.8, 128.3, 126.8, 126.3, 124.1, 122.3, 122.1, 119.2, 96.4, 55.4, 44.0, 37.8, 21.1; Anal. Calcd for $\text{C}_{25}\text{H}_{21}\text{N}_3\text{O}_3$: C, 72.98; H, 5.14; N, 10.21%; Found: C, 72.76; H, 5.19; N, 10.31%.



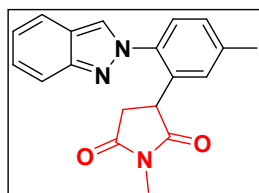
3-(2-(5-Fluoro-2H-indazol-2-yl)-5-methylphenyl)-1-phenylpyrrolidine-2,5-dione (3ka): Yellow solid (76%, 60 mg); R_f = 0.40 (PE : EA = 71 : 29); M.p. 175-176 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.20 (s, 1H), 7.63-7.60 (m, 1H), 7.39-7.24 (m, 7H), 7.12-7.07 (m, 1H), 6.94-6.92 (m, 2H), 4.24-4.20 (m, 1H), 3.27-3.24 (m, 2H), 2.46 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 176.1, 175.1, 160.0, 147.2, 140.3, 137.5, 132.5, 131.6 (d, $J_{\text{C-F}}$ = 36.0 Hz), 129.8, 129.0, 128.5, 127.0, 126.2, 125.4 (d, $J_{\text{C-F}}$ = 8.0 Hz), 121.7 (d, $J_{\text{C-F}}$ = 12.0 Hz), 120.1 (d, $J_{\text{C-F}}$ = 10.0 Hz), 118.9, 118.6, 102.9 (d, $J_{\text{C-F}}$ = 24.0 Hz), 44.0, 37.9, 21.3; Anal. Calcd for $\text{C}_{24}\text{H}_{18}\text{FN}_3\text{O}_2$: C, 72.17; H, 4.54; N, 10.52%; Found: C, 72.31; H, 4.61; N, 10.43%.



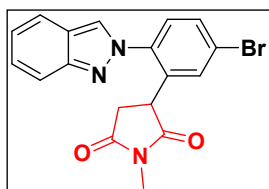
3-(2-(5-Chloro-2H-indazol-2-yl)-5-methylphenyl)-1-phenylpyrrolidine-2,5-dione (3la): White solid (81%, 67 mg); $R_f = 0.50$ (PE : EA = 68 : 32); M.p. 179-180 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.19 (s, 1H), 7.67 (s, 1H), 7.58 (d, $J = 9.2$ Hz, 1H), 7.38-7.29 (m, 5H), 7.24-7.21 (m, 2H), 6.92-6.90 (m, 2H), 4.20 (t, $J = 7.2$ Hz, 1H), 3.25-3.23 (m, 2H), 2.46 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 176.0, 175.0, 148.0, 140.4, 137.3, 132.6, 131.8, 131.4, 129.8, 129.0, 128.55, 128.50, 128.3, 127.0, 126.2, 125.0, 122.7, 119.5, 119.3, 44.0, 37.9, 21.3; Anal. Calcd for $\text{C}_{24}\text{H}_{18}\text{ClN}_3\text{O}_2$: C, 69.31; H, 4.36; N, 10.10%; Found: C, 69.18; H, 4.32; N, 10.04%.



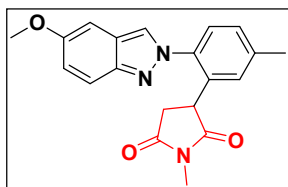
3-(4-Chloro-2-(5-fluoro-2H-indazol-2-yl)phenyl)-1-phenylpyrrolidine-2,5-dione (3ma): Yellow solid (92%, 77 mg); $R_f = 0.50$ (PE : EA = 70 : 30); M.p. 160-161 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.24 (s, 1H), 7.63-7.60 (m, 1H), 7.53-7.49 (m, 2H), 7.40-7.27 (m, 5H), 7.14-7.09 (m, 1H), 6.91 (d, $J = 8.4$ Hz, 2H), 4.29-4.25 (m, 1H), 3.25-3.22 (m, 2H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 175.5, 174.7, 158.9 (d, $J_{\text{C-F}} = 240.0$ Hz), 147.3, 140.5, 134.6, 132.4, 131.6, 131.1, 129.8, 128.9, 128.5, 127.1, 126.1, 125.2 (d, $J_{\text{C-F}} = 9.0$ Hz), 121.8 (d, $J_{\text{C-F}} = 12.0$ Hz), 120.1 (d, $J_{\text{C-F}} = 9.0$ Hz), 119.2 (d, $J_{\text{C-F}} = 29.0$ Hz), 102.9 (d, $J_{\text{C-F}} = 25.0$ Hz), 43.8, 37.6; Anal. Calcd for $\text{C}_{23}\text{H}_{15}\text{ClFN}_3\text{O}_2$: C, 65.80; H, 3.60; N, 10.01%; Found: C, 65.96; H, 3.53; N, 10.10%.



3-(2-(2H-Indazol-2-yl)-5-methylphenyl)-1-methylpyrrolidine-2,5-dione (3ab): Light yellow solid (80%, 51 mg); $R_f = 0.40$ (PE : EA = 69 : 29); M.p. 134-135 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.19 (s, 1H), 7.71-7.64 (m, 2H), 7.36 (d, $J = 8.0$ Hz, 1H), 7.32-7.25 (m, 2H), 7.11 (t, $J = 8.8$ Hz, 2H), 4.03-3.99 (m, 1H), 3.07 (t, $J = 9.6$ Hz, 2H), 2.68 (s, 3H), 2.43 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 177.4, 176.3, 149.5, 140.1, 137.5, 133.1, 130.8, 129.5, 126.95, 126.91, 125.3, 122.5, 122.3, 120.5, 117.7, 43.5, 38.1, 25.0, 21.2; Anal. Calcd for $\text{C}_{19}\text{H}_{17}\text{N}_3\text{O}_2$: C, 71.46; H, 5.37; N, 13.16%; Found: C, 71.26; H, 5.34; N, 13.04%.

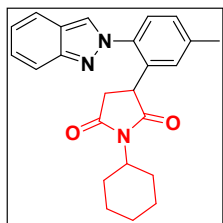


3-(5-Bromo-2-(2H-indazol-2-yl)phenyl)-1-methylpyrrolidine-2,5-dione (3nb): Yellow gummy mass (88%, 67 mg); $R_f = 0.45$ (PE : EA = 69 : 29); ^1H NMR (400 MHz, CDCl_3): δ 8.19 (s, 1H), 7.69 (d, $J = 8.4$ Hz, 1H), 7.65-7.59 (m, 2H), 7.50 (d, $J = 2.4$ Hz, 1H), 7.36-7.30 (m, 2H), 7.12 (t, $J = 8.0$ Hz, 1H), 4.07-4.03 (m, 1H), 3.10-3.07 (m, 2H), 2.66 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 176.5, 175.7, 149.7, 139.0, 135.3, 133.7, 132.1, 128.4, 127.3, 125.2, 123.5, 122.9, 122.5, 120.5, 117.8, 43.5, 37.9, 25.1; Anal. Calcd for $\text{C}_{18}\text{H}_{14}\text{BrN}_3\text{O}_2$: C, 56.27; H, 3.67; N, 10.94%; Found: C, 56.15; H, 3.71; N, 11.01%.

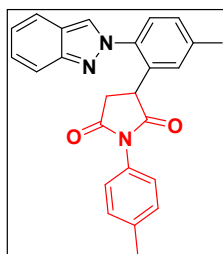


3-(2-(5-Methoxy-2H-indazol-2-yl)-5-methylphenyl)-1-methylpyrrolidine-2,5-dione (3jb): White solid (84%, 58 mg); $R_f = 0.45$ (PE : EA = 68 : 32); M.p. 165-167 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.06 (s, 1H), 7.56 (d, $J = 9.2$ Hz, 1H), 7.36 (d, $J = 8.4$ Hz, 1H), 7.27 (s, 1H), 7.13 (s,

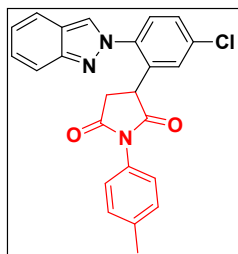
1H), 7.04-7.01 (m, 1H), 6.91 (d, $J=2.4$ Hz, 1H), 4.06-4.02 (m, 1H), 3.85 (s, 3H), 3.10-3.05 (m, 2H), 2.72 (s, 3H), 2.44 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 177.5, 176.3, 155.6, 146.4, 139.8, 137.6, 133.0, 131.0, 129.5, 126.8, 124.2, 122.3, 122.1, 119.1, 96.4, 55.4, 43.6, 38.1, 25.0, 21.2; Anal. Calcd for $\text{C}_{20}\text{H}_{19}\text{N}_3\text{O}_3$: C, 68.75; H, 5.48; N, 12.03%; Found: C, 68.96; H, 5.43; N, 11.92%.



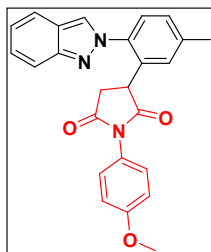
3-(2-(2H-Indazol-2-yl)-5-methylphenyl)-1-cyclohexylpyrrolidine-2,5-dione (3ac): Yellow gummy mass (75%, 58 mg); $R_f = 0.50$ (PE : EA = 75 : 25); ^1H NMR (400 MHz, CDCl_3): δ 8.26 (s, 1H), 7.71 (d, $J = 8.8$ Hz, 2H), 7.35-7.29 (m, 2H), 7.24 (d, $J = 8.0$ Hz, 1H), 7.11 (t, $J = 7.6$ Hz, 1H), 7.02 (s, 1H), 3.94-3.85 (m, 2H), 3.07-3.00 (m, 1H), 2.92-2.86 (m, 1H), 2.42 (s, 3H), 2.14 (s, 1H), 2.06-1.97 (m, 2H), 1.75 (d, $J = 10.8$ Hz, 2H), 1.60 (d, $J = 11.2$ Hz, 1H), 1.40 (d, $J = 9.2$ Hz, 1H), 1.25-1.19 (m, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 177.6, 176.2, 149.6, 140.2, 138.0, 133.7, 129.3, 129.0, 127.1, 126.8, 125.5, 122.4, 122.2, 120.5, 117.8, 52.0, 42.0, 38.0, 28.7, 25.8, 25.0, 21.3; Anal. Calcd for $\text{C}_{24}\text{H}_{25}\text{N}_3\text{O}_2$: C, 74.39; H, 6.50; N, 10.84%; Found: C, 74.62; H, 6.53; N, 10.96%.



3-(2-(2H-Indazol-2-yl)-5-methylphenyl)-1-(p-tolyl)pyrrolidine-2,5-dione (3ad): White solid (83%, 65 mg); $R_f = 0.50$ (PE : EA = 71 : 29); M.p. 155-156 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.24 (s, 1H), 7.71-7.66 (m, 2H), 7.38 (d, $J = 8.0$ Hz, 1H), 7.32-7.27 (m, 2H), 7.22 (s, 1H), 7.16-7.12 (m, 3H), 6.79 (d, $J = 8.4$ Hz, 2H), 4.19 (t, $J = 8.0$ Hz, 1H), 3.25-3.23 (m, 2H), 2.46 (s, 3H), 2.33 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 176.3, 175.2, 149.7, 140.1, 138.4, 137.7, 132.8, 131.1, 129.6, 129.2, 127.08, 127.02, 126.1, 125.3, 122.5, 122.4, 120.6, 117.9, 43.9, 38.0, 21.29, 21.24; Anal. Calcd for $\text{C}_{25}\text{H}_{21}\text{N}_3\text{O}_2$: C, 75.93; H, 5.35; N, 10.63%; Found: C, 76.05; H, 5.32; N, 10.55%.

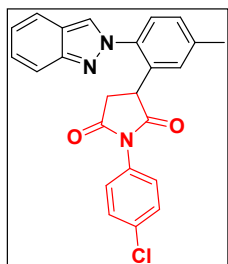


3-(5-Chloro-2-(2H-indazol-2-yl)phenyl)-1-(p-tolyl)pyrrolidine-2,5-dione (3dd): White solid (69%, 57 mg); $R_f = 0.55$ (PE : EA = 70 : 30); M.p. 192-194 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.24 (s, 1H), 7.72-7.64 (m, 2H), 7.50-7.43 (m, 3H), 7.33-7.29 (m, 1H), 7.17-7.11 (m, 3H), 6.74 (d, $J = 8.4$ Hz, 2H), 4.26-4.22 (m, 1H), 3.29-3.26 (m, 2H), 2.32 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 175.4, 174.7, 150.0, 138.6, 135.6, 134.7, 131.0, 129.7, 129.2, 129.0, 128.4, 127.4, 126.1, 125.3, 123.0, 122.6, 120.6, 118.0, 44.0, 37.8, 21.2; Anal. Calcd for $\text{C}_{24}\text{H}_{18}\text{ClN}_3\text{O}_2$: C, 69.31; H, 4.36; N, 10.10%; Found: C, 69.45; H, 4.41; N, 9.98%.



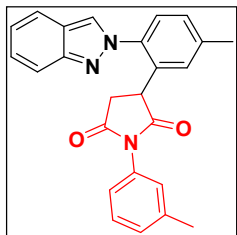
3-(2-(2H-Indazol-2-yl)-5-methylphenyl)-1-(4-methoxyphenyl)pyrrolidine-2,5-dione (3ae):

Yellow solid (82%, 67 mg); $R_f = 0.45$ (PE : EA = 68 : 32); M.p. 188-189 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.23 (s, 1H), 7.71-7.65 (m, 2H), 7.38 (d, $J = 8.0$ Hz, 1H), 7.30 (t, $J = 8.4$ Hz, 2H), 7.22 (s, 1H), 7.13 (t, $J = 7.6$ Hz, 1H), 6.84-6.78 (m, 4H), 4.19 (t, $J = 8.4$ Hz, 1H), 3.77 (s, 3H), 3.26-3.24 (m, 2H), 2.46 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 176.4, 175.4, 159.4, 149.7, 140.1, 137.7, 132.8, 131.2, 129.6, 127.6, 127.1, 127.0, 125.4, 124.5, 122.6, 122.4, 120.6, 117.9, 114.3, 55.5, 43.9, 38.0, 21.3; Anal. Calcd for $\text{C}_{25}\text{H}_{21}\text{N}_3\text{O}_3$: C, 72.98; H, 5.14; N, 10.21%; Found: C, 73.21; H, 5.07; N, 10.11%.

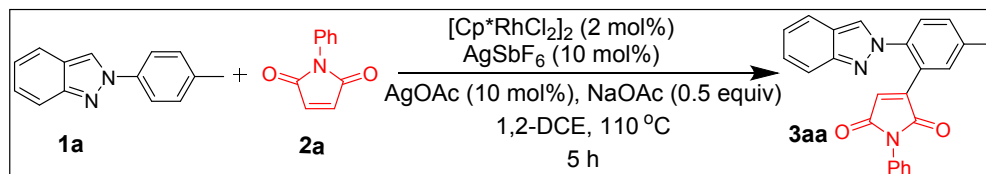


3-(2-(2H-Indazol-2-yl)-5-methylphenyl)-1-(4-chlorophenyl)pyrrolidine-2,5-dione (3af):

White solid (90%, 74 mg); $R_f = 0.50$ (PE : EA = 70 : 30); M.p. 190-191 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.20 (s, 1H), 7.69 (d, $J = 8.4$ Hz, 1H), 7.60 (d, $J = 8.4$ Hz, 1H), 7.38 (d, $J = 8.0$ Hz, 1H), 7.31-7.22 (m, 5H), 7.13 (t, $J = 8.4$ Hz, 1H), 6.76 (d, $J = 8.8$ Hz, 2H), 4.22-4.18 (m, 1H), 3.37-3.19 (m, 2H), 2.46 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 175.7, 174.8, 149.7, 140.0, 137.4, 134.0, 132.3, 132.0, 130.3, 129.8, 129.0, 127.5, 127.1, 127.0, 125.2, 122.7, 122.4, 120.6, 117.9, 44.5, 37.9, 21.2; Anal. Calcd for $\text{C}_{24}\text{H}_{18}\text{ClN}_3\text{O}_2$: C, 69.31; H, 4.36; N, 10.10%; Found: C, 69.43; H, 4.38; N, 10.04%.

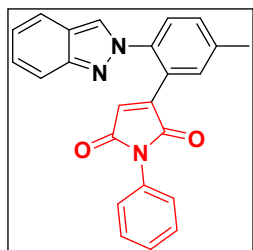


3-(2-(2H-Indazol-2-yl)-5-methylphenyl)-1-(m-tolyl)pyrrolidine-2,5-dione (3ag): Yellow gummy mass (87%, 68 mg); $R_f = 0.45$ (PE : EA = 75 : 25); ^1H NMR (400 MHz, CDCl_3): δ 8.24 (s, 1H), 7.72-7.66 (m, 2H), 7.38 (d, $J = 8.0$ Hz, 1H), 7.30 (t, $J = 8.4$ Hz, 2H), 7.23-7.19 (m, 2H), 7.16-7.09 (m, 2H), 6.72 (d, $J = 8.0$ Hz, 1H), 6.55 (s, 1H), 4.22-4.18 (m, 1H), 3.38-3.21 (m, 2H), 2.46 (s, 3H), 2.26 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 176.1, 175.2, 149.7, 140.0, 139.0, 137.6, 132.7, 131.6, 131.5, 129.6, 129.3, 128.7, 127.1, 127.0, 125.3, 123.4, 122.6, 122.4, 120.6, 118.0, 44.1, 38.1, 21.3, 21.2; Anal. Calcd for $\text{C}_{25}\text{H}_{21}\text{N}_3\text{O}_2$: C, 75.93; H, 5.35; N, 10.63%; Found: C, 76.13; H, 5.31; N, 10.55%.

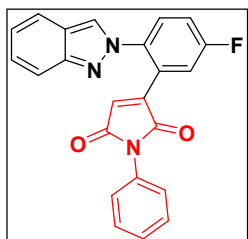


Typical experimental procedure for the synthesized compounds (4aa-4ga): A mixture of 2-(p-tolyl)-2H-indazole (0.2 mmol, 41.6 mg) (**1a**), $[\text{Cp}^*\text{RhCl}_2]_2$ (2 mol%, 2.4 mg), AgSbF_6 (10 mol%, 6.8 mg), AgOAc (10 mol%, 3.3 mg), and NaOAc (0.5 equiv, 8.2 mg) were taken in an oven dried screw-capped reaction tube. Then 1,2-DCE (2 mL) was added to the mixture and stirred for 5 min at room temperature under open atmosphere. After that, 1-phenyl-1H-pyrrole-2,5-dione (**2a**) (0.24 mmol, 41.5 mg) was added, and the resultant mixture was stirred at 110 °C for 5 h. After completion of the reaction (TLC), the reaction was cooled to room temperature and extracted with dichloromethane. The organic phase was dried over anhydrous Na_2SO_4 . The crude

residue was obtained after evaporating the solvent in vacuum and was purified by column chromatography on silica gel using a mixture of petroleum ether and ethyl acetate (79:21) as an eluting solvent to afford the pure product **4aa** (50 mg, 67%) as a yellow solid.

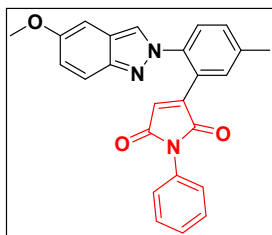


3-(2-(2H-Indazol-2-yl)-5-methylphenyl)-1-phenyl-1H-pyrrole-2,5-dione (4aa): Yellow solid (67%, 50 mg); $R_f = 0.50$ (PE : EA = 79 : 21); M.p. 125-126 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.29 (s, 1H), 7.74-7.68 (m, 2H), 7.55 (d, $J = 8.4$ Hz, 2H), 7.45 (d, $J = 8.0$ Hz, 1H), 7.34-7.29 (m, 3H), 7.26-7.25 (m, 1H), 7.14 (t, $J = 8.0$ Hz, 1H), 7.00-6.98 (m, 2H), 6.48 (s, 1H), 2.51 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 169.2, 167.8, 149.7, 145.4, 139.4, 137.3, 132.1, 131.9, 131.4, 129.0, 127.7, 127.1, 126.7, 126.0, 125.4, 124.7, 124.5, 122.9, 122.8, 120.7, 117.6, 21.2; Anal. Calcd for $\text{C}_{24}\text{H}_{17}\text{N}_3\text{O}_2$: C, 75.98; H, 4.52; N, 11.08%; Found: C, 75.86; H, 4.56; N, 11.03%.



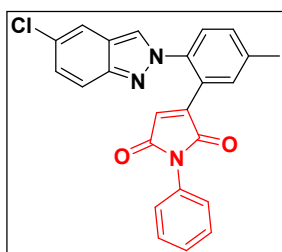
3-(5-Fluoro-2-(2H-indazol-2-yl)phenyl)-1-phenyl-1H-pyrrole-2,5-dione (4ca): Light yellow solid (52%, 39 mg); $R_f = 0.50$ (PE : EA = 78 : 22); M.p. 129-130 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.27 (s, 1H), 7.74-7.69 (m, 2H), 7.66-7.62 (m, 1H), 7.57-7.55 (m, 1H), 7.38-7.27 (m, 5H), 7.18-7.14 (m, 1H), 7.01-6.99 (m, 2H), 6.42 (s, 1H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ

167.5, 163.4, 160.9, 149.9, 143.3, 131.2, 129.3, 129.1, 127.9, 127.8, 127.7, 127.5, 127.1, 126.0, 124.8, 123.0 (d, $J_{\text{C-F}} = 19.0$ Hz), 120.7, 118.6, 118.4, 118.3 (d, $J_{\text{C-F}} = 14.0$ Hz), 117.7; Anal. Calcd for $\text{C}_{23}\text{H}_{14}\text{FN}_3\text{O}_2$: C, 72.06; H, 3.68; N, 10.96%; Found: C, 72.24; H, 3.66; N, 11.03%.



3-(2-(5-Methoxy-2H-indazol-2-yl)-5-methylphenyl)-1-phenyl-1H-pyrrole-2,5-dione (4ja):

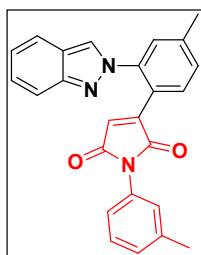
Brown solid (71%, 58 mg); $R_f = 0.45$ (PE : EA = 72 : 28); M.p. 162-163 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.15 (s, 1H), 7.60-7.52 (m, 3H), 7.45-7.43 (m, 1H), 7.34-7.30 (m, 2H), 7.27-7.24 (m, 1H), 7.04-7.00 (m, 3H), 6.91 (d, $J = 2.0$ Hz, 1H), 6.49 (s, 1H), 3.85 (s, 3H), 2.50 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 169.3, 167.9, 155.7, 146.7, 145.7, 139.1, 137.4, 132.1, 131.9, 131.5, 129.0, 127.7, 126.5, 126.1, 125.2, 124.5, 123.4, 123.0, 122.3, 119.0, 96.5, 55.4, 21.2; Anal. Calcd for $\text{C}_{25}\text{H}_{19}\text{N}_3\text{O}_3$: C, 73.34; H, 4.68; N, 10.26%; Found: C, 73.17; H, 4.71; N, 10.14%.



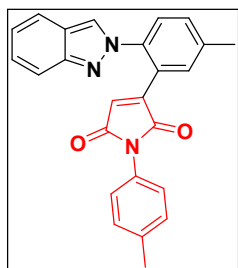
3-(2-(5-Chloro-2H-indazol-2-yl)-5-methylphenyl)-1-phenyl-1H-pyrrole-2,5-dione (4la):

Yellow solid (69%, 56 mg); $R_f = 0.45$ (PE : EA = 72 : 28); M.p. 141-142 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.25 (s, 1H), 7.70 (s, 1H), 7.64 (d, $J = 9.2$ Hz, 1H), 7.54 (d, $J = 8.0$ Hz, 2H), 7.47-7.45 (m, 1H), 7.35-7.31 (m, 2H), 7.28-7.24 (m, 2H), 6.99 (d, $J = 7.6$ Hz, 2H), 6.52 (s, 1H),

2.51 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 169.1, 167.8, 148.0, 145.5, 139.7, 137.0, 132.2, 132.0, 131.4, 129.1, 128.6, 128.4, 127.8, 126.8, 126.0, 125.3, 124.7, 124.1, 123.2, 119.4, 119.2, 21.2; Anal. Calcd for $\text{C}_{24}\text{H}_{16}\text{ClN}_3\text{O}_2$: C, 69.65; H, 3.90; N, 10.15%; Found: C, 69.80; H, 3.92; N, 10.09%.

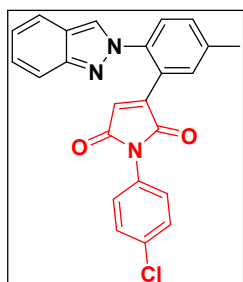


3-(2-(2H-Indazol-2-yl)-4-methylphenyl)-1-(m-tolyl)-1H-pyrrole-2,5-dione (4ag): Light yellow liquid (64%, 50 mg); R_f = 0.45 (PE : EA = 79 : 21); ^1H NMR (400 MHz, CDCl_3): δ 8.30 (s, 1H), 7.74-7.70 (m, 2H), 7.55 (t, J = 4.8 Hz, 2H), 7.45 (d, J = 7.2 Hz, 1H), 7.35-7.31 (m, 1H), 7.21-7.13 (m, 2H), 7.05 (d, J = 7.6 Hz, 1H), 6.82 (d, J = 7.6 Hz, 1H), 6.66 (s, 1H), 6.51 (s, 1H), 2.51 (s, 3H), 2.22 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 169.3, 167.9, 149.7, 145.5, 139.4, 139.0, 137.4, 132.1, 131.9, 131.3, 128.8, 128.6, 127.1, 126.8, 126.6, 125.4, 124.9, 124.7, 123.2, 122.9, 122.8, 120.8, 117.7, 21.3, 21.2; Anal. Calcd for $\text{C}_{25}\text{H}_{19}\text{N}_3\text{O}_2$: C, 76.32; H, 4.87; N, 10.68%; Found: C, 76.44; H, 4.84; N, 10.71%.



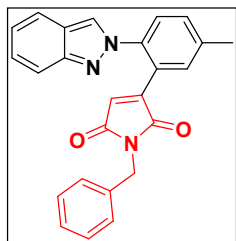
3-(2-(2H-Indazol-2-yl)-5-methylphenyl)-1-(p-tolyl)-1H-pyrrole-2,5-dione (4ad): Yellow solid (76%, 59 mg); R_f = 0.50 (PE : EA = 77 : 23); M.p. 135-136 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.28 (s, 1H), 7.71 (t, J = 8.4 Hz, 2H), 7.57-7.53 (m, 2H), 7.45-7.43 (m, 1H), 7.34-7.30 (m, 1H),

7.15-7.10 (m, 3H), 6.87 (d, $J = 8.4$ Hz, 2H), 6.45 (s, 1H), 2.50 (s, 3H), 2.30 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 169.4, 168.0, 149.7, 145.2, 139.4, 137.8, 137.4, 132.1, 131.9, 129.7, 128.8, 127.1, 126.7, 126.0, 125.5, 124.9, 124.5, 122.9, 122.8, 120.7, 117.7, 21.26, 21.20; HRMS (ESI-TOF) m/z : $[\text{M} + \text{H}]^+$ Calcd for $\text{C}_{25}\text{H}_{20}\text{N}_3\text{O}_2$: 394.1550; found: 394.1548.



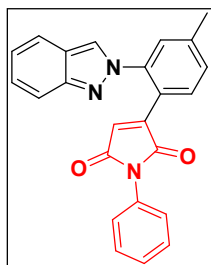
3-(2-(2H-Indazol-2-yl)-5-methylphenyl)-1-(4-chlorophenyl)-1H-pyrrole-2,5-dione (4af):

Yellow solid (77%, 63 mg); $R_f = 0.55$ (PE : EA = 70 : 30); M.p. 149-150 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.31 (s, 1H), 7.74-7.65 (m, 2H), 7.55 (t, $J = 8.0$ Hz, 2H), 7.46 (d, $J = 9.6$ Hz, 1H), 7.34-7.32 (m, 1H), 7.30-7.24 (m, 2H), 7.16-7.14 (m, 1H), 6.93-6.90 (m, 2H), 6.52 (s, 1H), 2.50 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 168.9, 167.4, 149.7, 145.9, 139.4, 137.3, 133.3, 132.3, 131.9, 130.0, 129.2, 127.2, 127.1, 126.8, 126.5, 125.2, 124.6, 122.9, 122.8, 120.7, 117.5, 21.2; Anal. Calcd for $\text{C}_{24}\text{H}_{16}\text{ClN}_3\text{O}_2$: C, 69.65; H, 3.90; N, 10.15%; Found: C, 69.47; H, 3.92; N, 10.12%.

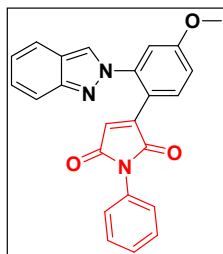


3-(2-(2H-Indazol-2-yl)-5-methylphenyl)-1-benzyl-1H-pyrrole-2,5-dione (4ah): Yellow gummy mass (81%, 63 mg); $R_f = 0.45$ (PE : EA = 79 : 21); ^1H NMR (400 MHz, CDCl_3): δ 8.22 (s, 1H), 7.68 (d, $J = 8.4$ Hz, 1H), 7.56-7.48 (m, 3H), 7.42-7.40 (m, 1H), 7.30-7.26 (m, 1H), 7.21-7.20 (m,

3H), 7.13-7.10 (m, 3H), 6.30 (s, 1H), 4.49 (s, 2H), 2.48 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 170.0, 168.9, 149.6, 145.0, 139.3, 137.4, 136.2, 131.9, 128.7, 128.2, 127.7, 127.0, 126.9, 125.9, 124.9, 124.2, 122.7, 122.6, 120.5, 117.9, 41.6, 21.2; Anal. Calcd for $\text{C}_{25}\text{H}_{19}\text{N}_3\text{O}_2$: C, 76.32; H, 4.87; N, 10.68%; Found: C, 76.54; H, 4.92; N, 10.73%.

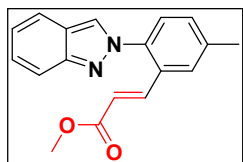


3-(2-(2H-Indazol-2-yl)-4-methylphenyl)-1-phenyl-1H-pyrrole-2,5-dione (4fa): Yellow solid (79%, 59 mg); R_f = 0.50 (PE : EA = 79 : 21); M.p. 116-117 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.20 (s, 1H), 7.64-7.59 (m, 3H), 7.38 (s, 1H), 7.30 (d, J = 7.6 Hz, 1H), 7.24-7.19 (m, 3H), 7.16 (t, J = 4.4 Hz, 1H), 7.07-7.03 (m, 1H), 6.92-6.90 (m, 2H), 6.28 (s, 1H) 2.42 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3): δ 169.3, 168.0, 149.8, 144.9, 142.5, 139.6, 131.5, 131.4, 129.9, 129.0, 127.7, 127.2, 126.5, 126.2, 126.1, 124.5, 122.9, 122.2, 120.7, 117.7, 21.5; HRMS (ESI-TOF) m/z : $[\text{M} + \text{Na}]^+$ Calcd for $\text{C}_{24}\text{H}_{17}\text{N}_3\text{O}_2\text{Na}$: 402.1213; found: 402.1215.

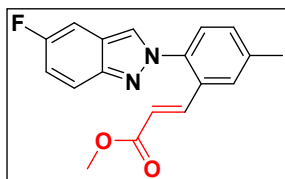


3-(2-(2H-Indazol-2-yl)-4-methoxyphenyl)-1-phenyl-1H-pyrrole-2,5-dione (4ga): Light yellow solid (55%, 43 mg); R_f = 0.50 (PE : EA = 70 : 30); M.p. 136-137 °C; ^1H NMR (400 MHz, CDCl_3): δ 8.28 (s, 1H), 7.81 (d, J = 8.4 Hz, 1H), 7.74-7.71 (m, 2H), 7.36-7.27 (m, 4H), 7.18-7.10 (m, 3H), 7.03 (d, J = 7.2 Hz, 2H), 6.18 (s, 1H), 3.93 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3):

δ 169.4, 168.3, 162.0, 149.8, 144.1, 141.2, 132.9, 131.5, 129.0, 127.7, 127.4, 126.1, 125.2, 124.5, 123.0, 122.9, 120.7, 117.8, 117.4, 114.7, 112.3, 56.0; Anal. Calcd for $C_{24}H_{17}N_3O_3$: C, 72.90; H, 4.33; N, 10.63%; Found: C, 72.68; H, 4.29; N, 10.71%.



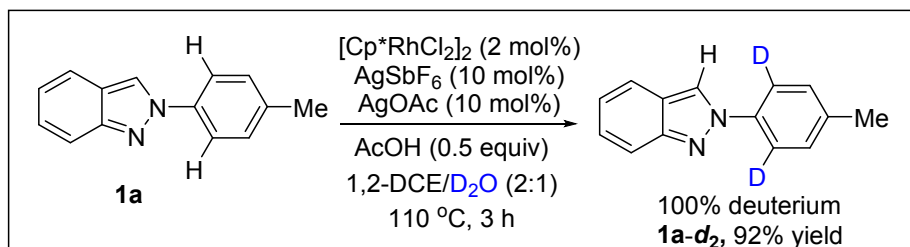
Methyl (E)-3-(2-(2H-indazol-2-yl)-5-methylphenyl)acrylate (4ai): Yellow liquid (81%, 47 mg); R_f = 0.50 (PE : EA = 78 : 22); 1H NMR (400 MHz, $CDCl_3$): δ 8.09 (s, 1H), 7.80 (d, J = 9.2 Hz, 1H), 7.72 (d, J = 8.4 Hz, 1H), 7.56-7.44 (m, 3H), 7.37-7.32 (m, 2H), 7.16-7.13 (m, 1H), 6.41 (d, J = 16.0 Hz, 1H), 3.71 (s, 3H), 2.46 (s, 3H); $^{13}C\{^1H\}$ NMR (100 MHz, $CDCl_3$): δ 166.8, 149.8, 139.9, 139.4, 138.0, 131.4, 130.0, 128.0, 127.0, 126.8, 125.3, 122.5, 122.4, 120.8, 120.4, 118.1, 51.8, 21.3; Anal. Calcd for $C_{18}H_{16}N_2O_2$: C, 73.95; H, 5.52; N, 9.58%; Found: C, 74.16; H, 5.55; N, 9.49%.



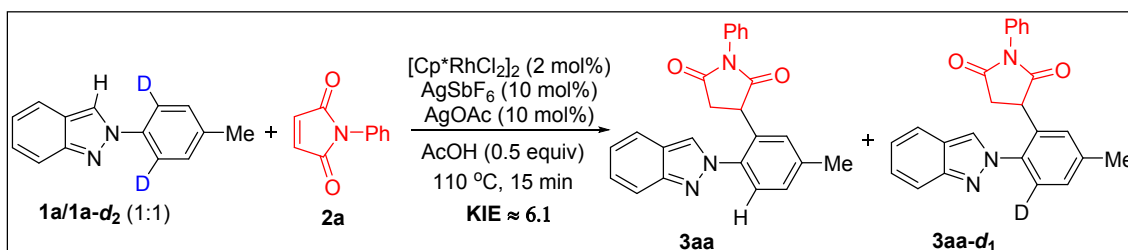
Methyl (E)-3-(2-(5-fluoro-2H-indazol-2-yl)-5-methylphenyl)acrylate (4ki): Yellow liquid (53%, 32 mg); R_f = 0.45 (PE : EA = 71 : 29); 1H NMR (400 MHz, $CDCl_3$): δ 8.05 (s, 1H), 7.78-7.75 (m, 1H), 7.56 (s, 1H), 7.50-7.42 (m, 2H), 7.34-7.27 (m, 2H), 7.18-7.12 (m, 1H), 6.40 (d, J = 16.0 Hz, 1H), 3.72 (s, 3H), 2.47 (s, 3H); $^{13}C\{^1H\}$ NMR (100 MHz, $CDCl_3$): δ 166.8, 158.8 (d, J_{C-F} = 240.0 Hz), 147.3, 139.7 (d, J_{C-F} = 7.0 Hz), 137.8, 131.5, 130.1, 128.1, 126.9, 125.4 (d, J_{C-F} = 9.0 Hz), 121.7, 121.6, 120.5 (d, J_{C-F} = 70.0 Hz), 120.2, 118.6 (d, J_{C-F} = 29.0 Hz), 102.7 (d, J_{C-F}

= 23.0 Hz), 51.9, 21.3; Anal. Calcd for C₁₈H₁₅FN₂O₂: C, 69.67; H, 4.87; N, 9.03%; Found: C, 69.43; H, 4.91; N, 8.91%.

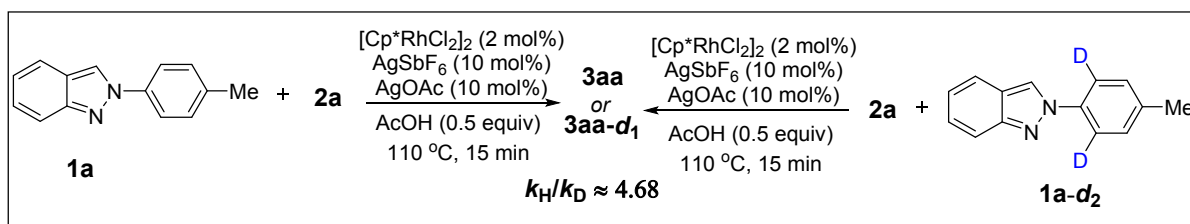
3. Mechanistic investigations:



Preparation of 2-(4-methylphenyl)-2,6-d₂-2H-indazole (1a-d₂**).**³ A mixture of 2-(*p*-tolyl)-2H-indazole (0.2 mmol, 41.6 mg) (**1a**), [Cp*RhCl₂]₂ (2 mol%, 2.4 mg), AgSbF₆ (10 mol%, 6.8 mg), and AgOAc (10 mol%, 3.3 mg) were taken in an oven dried screw-capped reaction tube. Then 1,2-DCE (2 mL) and AcOH (0.5 equiv, 6.0 mg) was added to the mixture and stirred for 5 min at room temperature under open atmosphere. After that, D₂O (1 mL) was added, and the resultant mixture was stirred at 110 °C for 3 h. Then, the reaction was cooled to room temperature and extracted with dichloromethane. The organic phase was dried over anhydrous Na₂SO₄. The crude residue was obtained after evaporating the solvent in vacuum and was purified by column chromatography on silica gel using a mixture of petroleum ether and ethyl acetate (85:15) as an eluting solvent to afford the pure product **1a-d₂** as a white solid. The deuterium incorporation was determined using 400 MHz ¹H NMR as 100%.

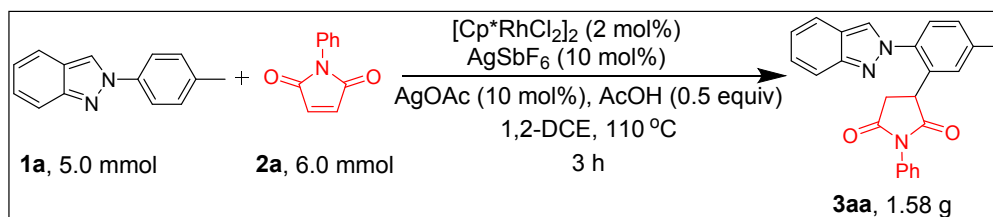
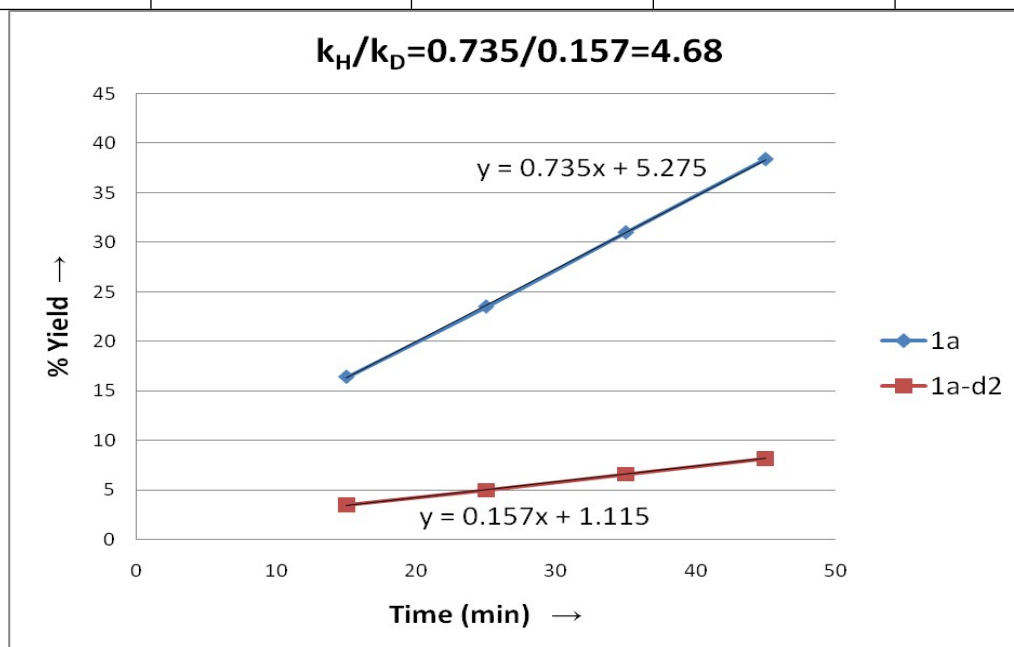


Intermolecular experiment.^{4a} 1-Phenyl-1*H*-pyrrole-2,5-dione (**2a**) (0.24 mmol, 41.5 mg) was reacted with 2-(*p*-tolyl)-2*H*-indazole (**1a**) (0.1 mmol, 20.8 mg) and 2-(4-methylphenyl-2,6-*d*₂)-2*H*-indazole (**1a-d**₂) (0.1 mmol, 21.0 mg) for 15 min under standard reaction condition. The resulting solution was then diluted with dichloromethane (3 x 10 mL) and washed with brine (2 x 5 mL) and water (5 mL). The organic phase was dried over anhydrous Na₂SO₄. The crude residue was obtained after evaporating the solvent in vacuum and was purified by column chromatography on silica gel using a mixture of petroleum ether and ethyl acetate (72:28) as an eluting solvent to afford **3aa** and a mixture of unreacted **1a** and **1a-d**₂ as a light yellow solid. The KIE was found to be 6.1 after 15 min at ~20% conversion, based on 400 MHz ¹H NMR of the product **3aa** and **3aa-d**₁.



Parallel Experiment study.⁴ In a set of two experiments: in first set, 1-phenyl-1*H*-pyrrole-2,5-dione (**2a**) (0.12 mmol, 20.7 mg) was reacted with 2-(*p*-tolyl)-2*H*-indazole (**1a**) (0.1 mmol, 20.8 mg) under standard reaction conditions. Whereas in another set, 2-(4-methylphenyl-2,6-*d*₂)-2*H*-indazole (**1a-d**₂) (0.1 mmol, 21.0 mg) was used instead of **1a** in the reaction with 1-phenyl-1*H*-pyrrole-2,5-dione (**2a**) under the standard reaction conditions. The two reactions were allowed to stir at 110 °C with stirring for 15, 25, 35 and 45 minutes. After cooling down, the mixture was concentrated in vacuo and purified by column chromatography, affording the product **3aa** or **3aa-d**₁. The KIE value was determined to be 4.68.

t/min.	15	25	35	45
yield				
3aa	16.4%	23.5%	31%	38.4%
3aa-<i>d</i>₁	3.5%	5.0%	6.6%	8.2%



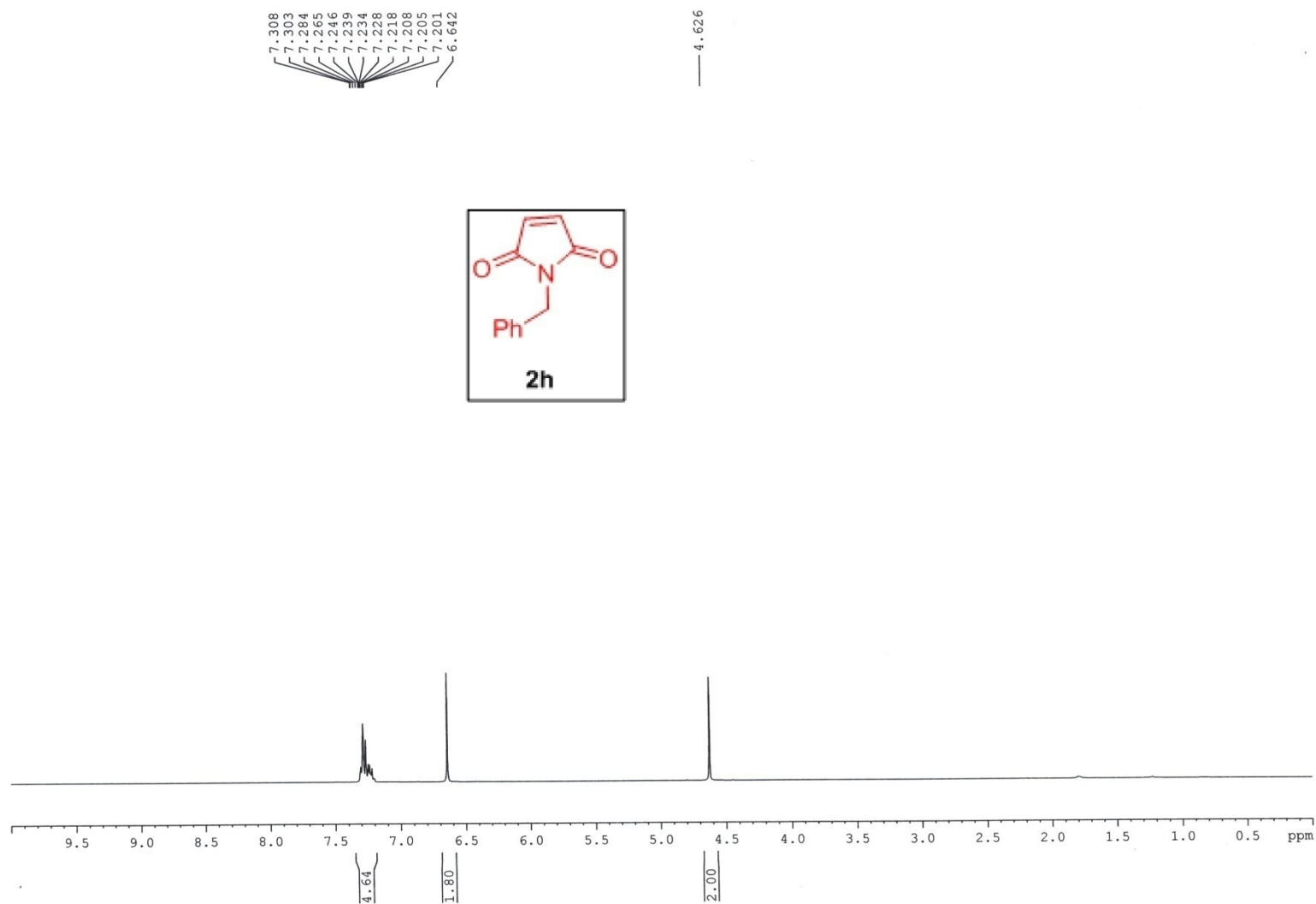
Experimental procedure of 3aa on 5 mmol scale: A mixture of 2-(*p*-tolyl)-2*H*-indazole (5.0 mmol, 1.04 g) (**1a**), [Cp*RhCl₂]₂ (2 mol%, 61.8 mg), AgSbF₆ (10 mol%, 171.8 mg), AgOAc (10 mol%, 83.5 mg), and AcOH (0.5 equiv, 150.0 mg) were taken in an oven dried round-bottom pressure flask (100 mL). Then 1,2-DCE (50 mL) was added to the mixture and stirred for 5 min at room temperature under open atmosphere. After that, 1-phenyl-1*H*-pyrrole-2,5-dione (**2a**) (6.0 mmol, 1.03 g) was added, and the resultant mixture was stirred at 110 °C for 3 h. After completion of the reaction (TLC), the reaction was cooled to room temperature and extracted with dichloromethane. The organic phase was dried over anhydrous Na₂SO₄. The crude residue

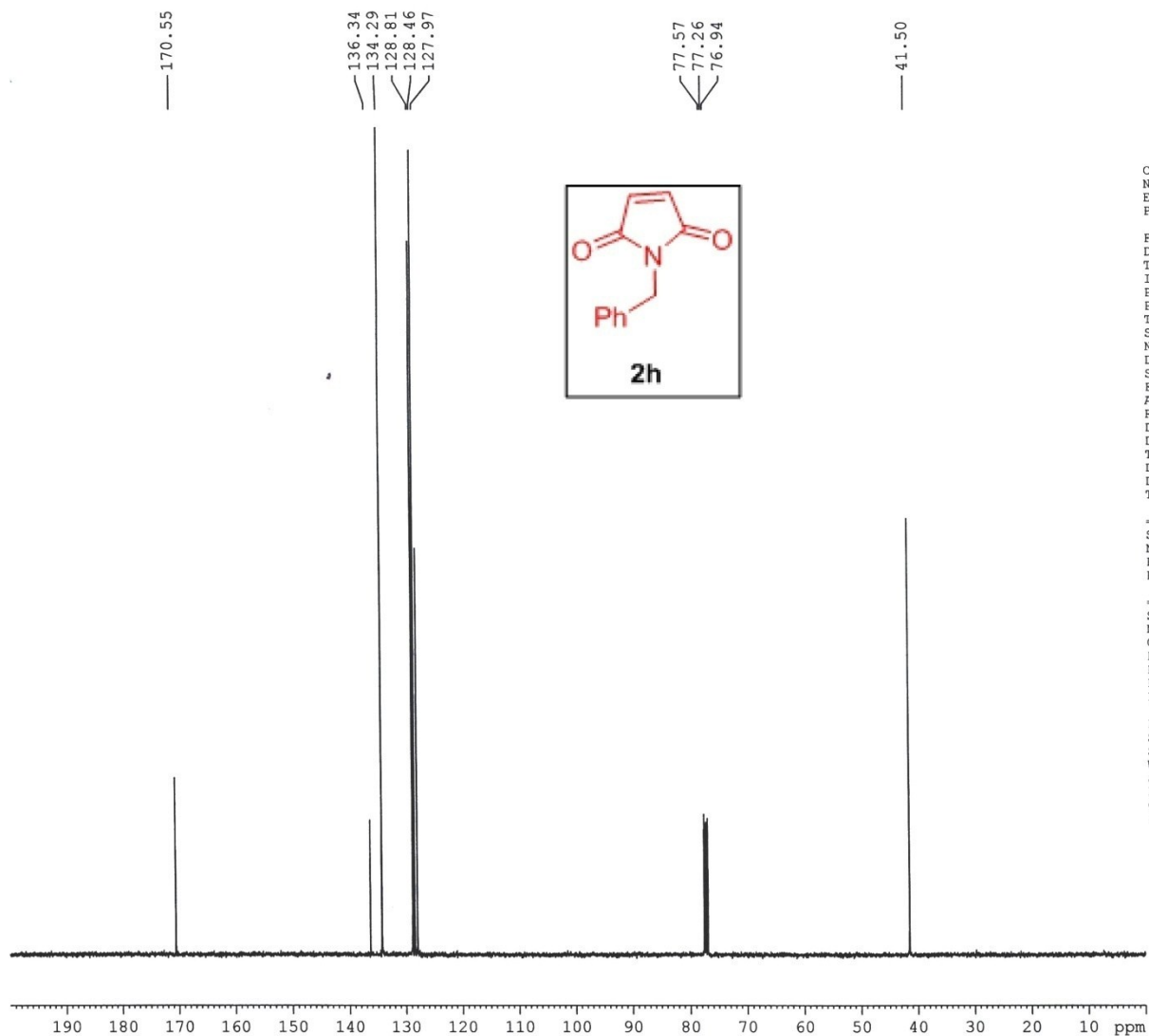
was obtained after evaporating the solvent in vacuum and was purified by column chromatography on silica gel using a mixture of petroleum ether and ethyl acetate (72:28) as an eluting solvent to afford the pure product **3aa** (1.58 g, 83%) as a light yellow solid.

4. References:

1. (a) M. R. Kumar, A. Park, N. Park and S. Lee, *Org. Lett.*, 2011, **13**, 3542-3545; (b) G. Bogonda, H. Y. Kim and K. Oh, *Org. Lett.*, 2018, **20**, 2711-2715.
2. G. B. Deshmukha, N. S. Patila, V. B. Gaikwada, A. D. Bholeb and S. V. Patil, *J. Chem. Pharm. Res.*, 2014, **6**, 393-399.
3. S. Samanta, S. Mondal, D. Ghosh and A. Hajra, *Org. Lett.*, 2019, **21**, 4905-4909.
4. (a) A. Modi, P. Sau, N. Chakraborty and B. K. Patel, *Adv. Synth. Catal.*, 2019, **361**, 1368–1375. (b) A. Maji, S. Guin, S. Feng, A. Dahiya, V. K. Singh, P. Liu and D. Maiti, *Angew. Chem., Int. Ed.*, 2017, **56**, 14903-14907.

5. NMR spectra for the synthesized products





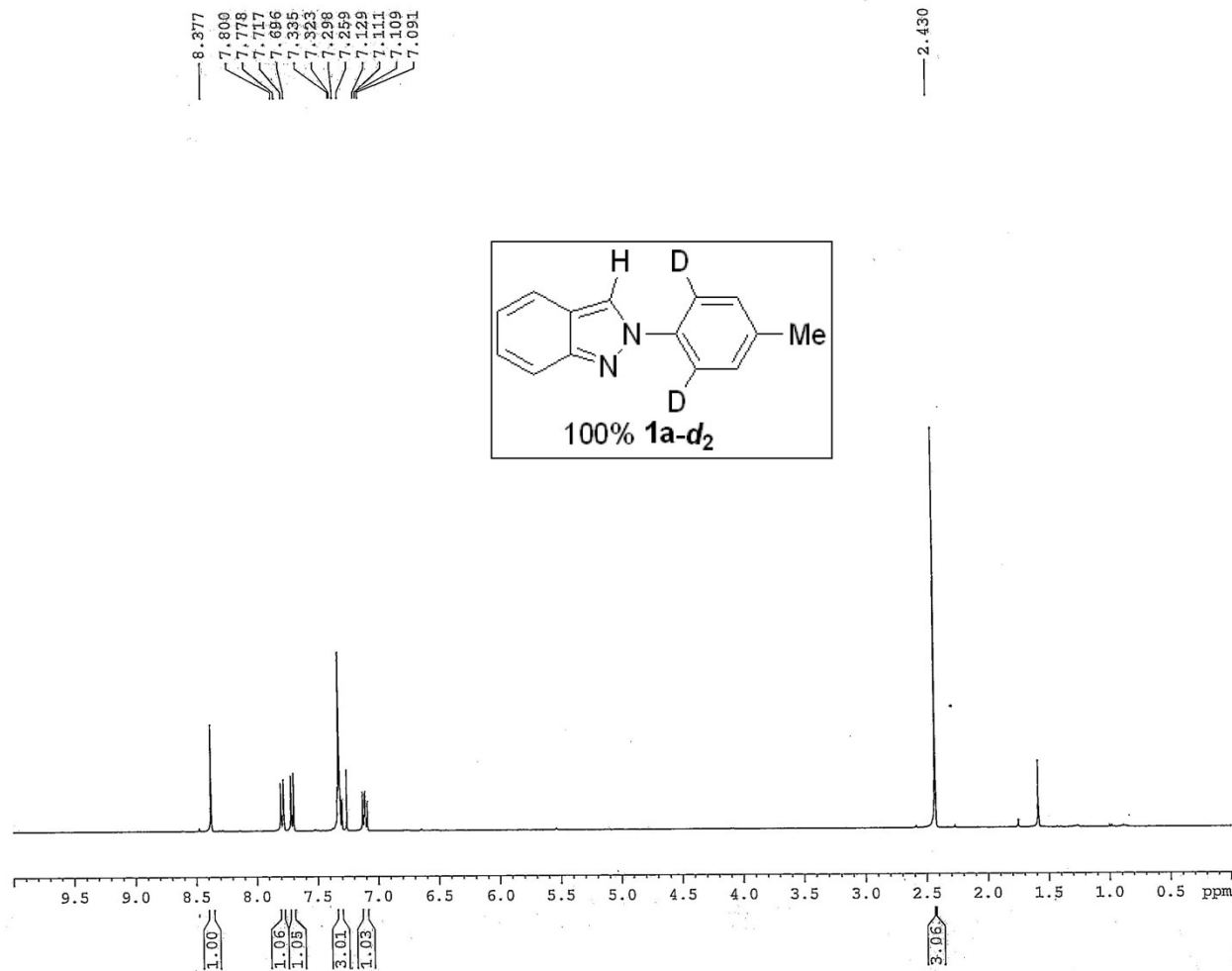
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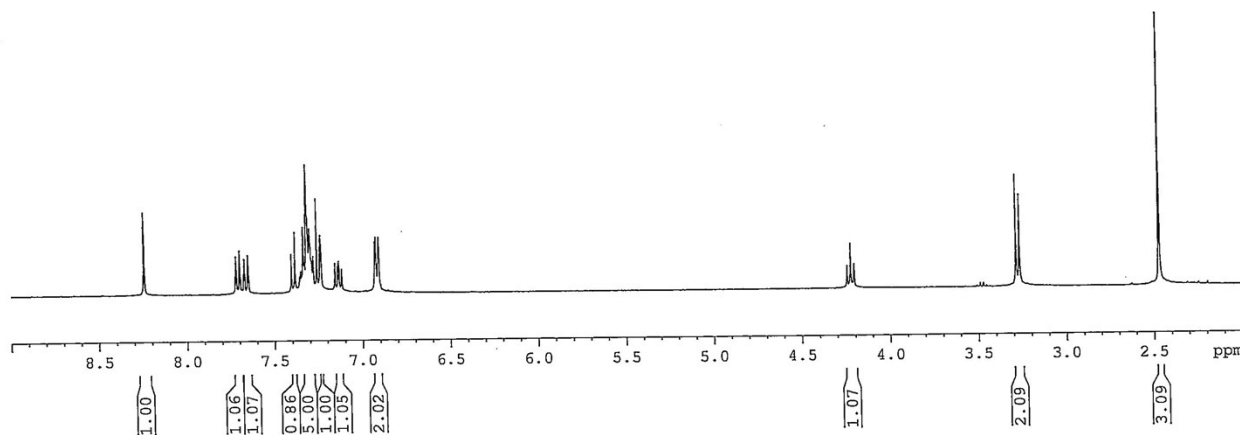
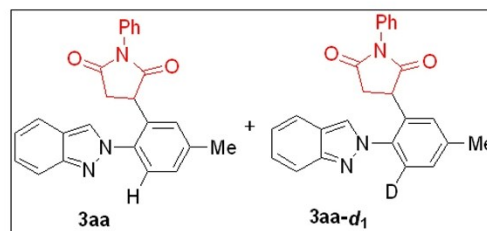


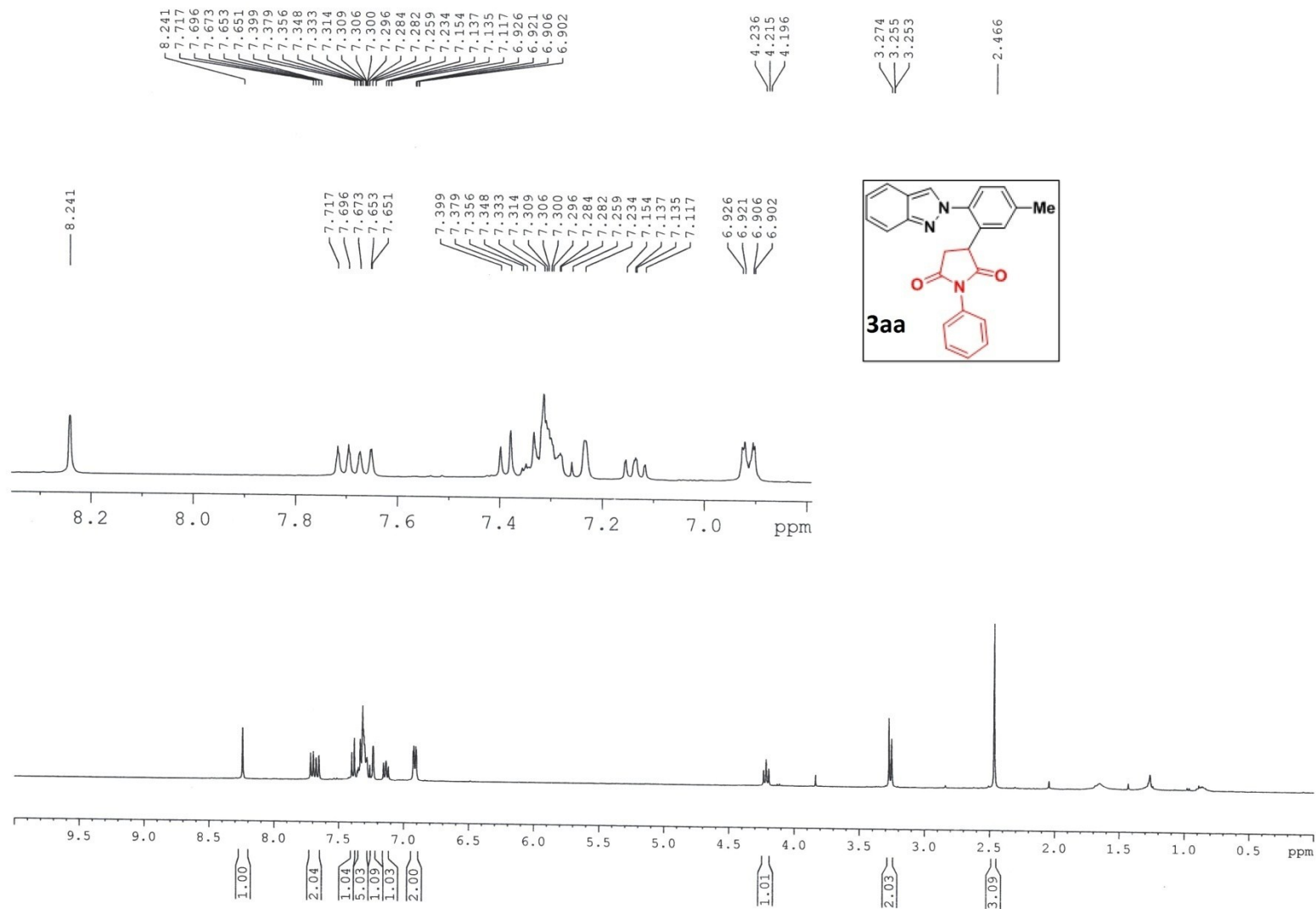
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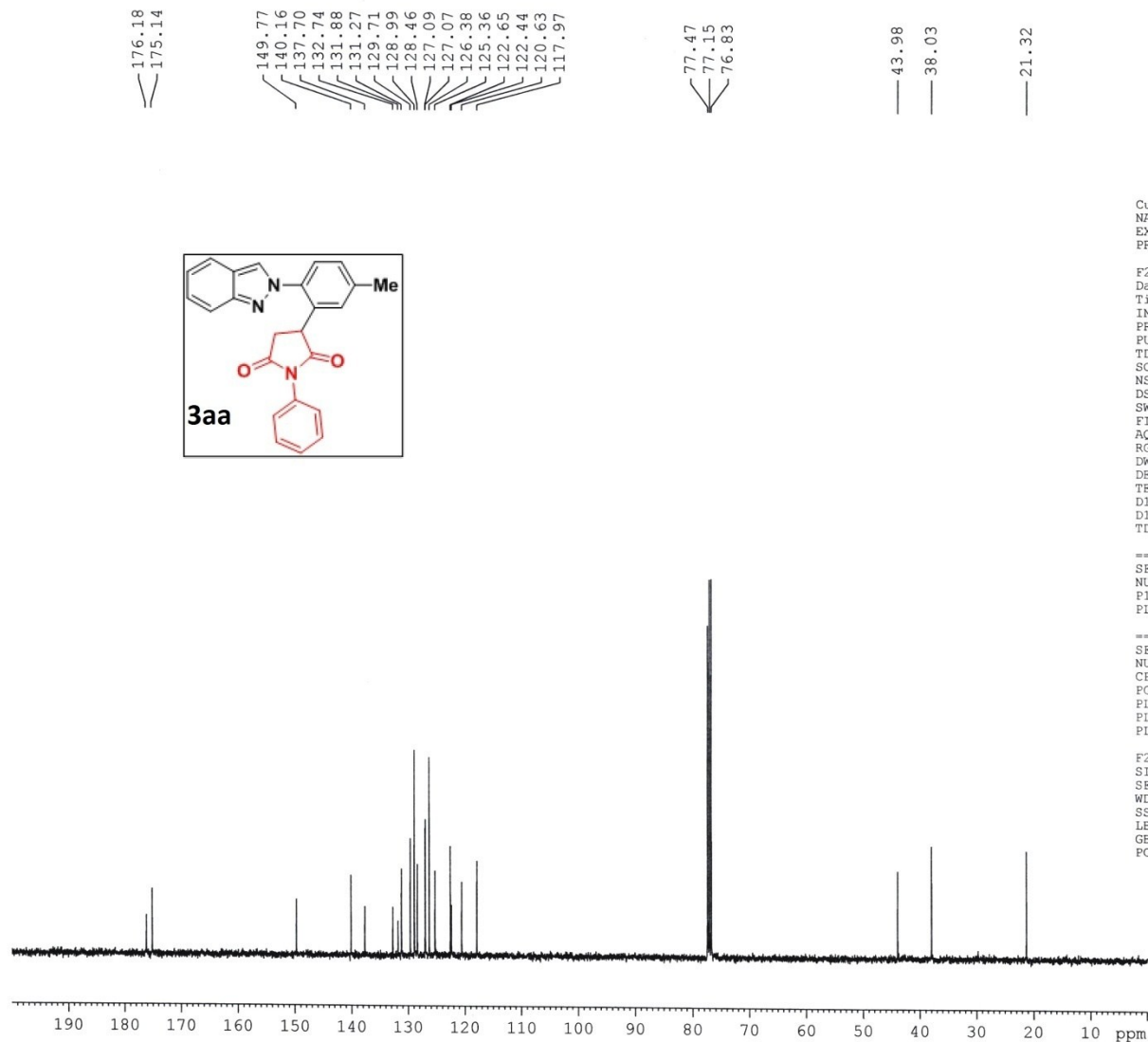
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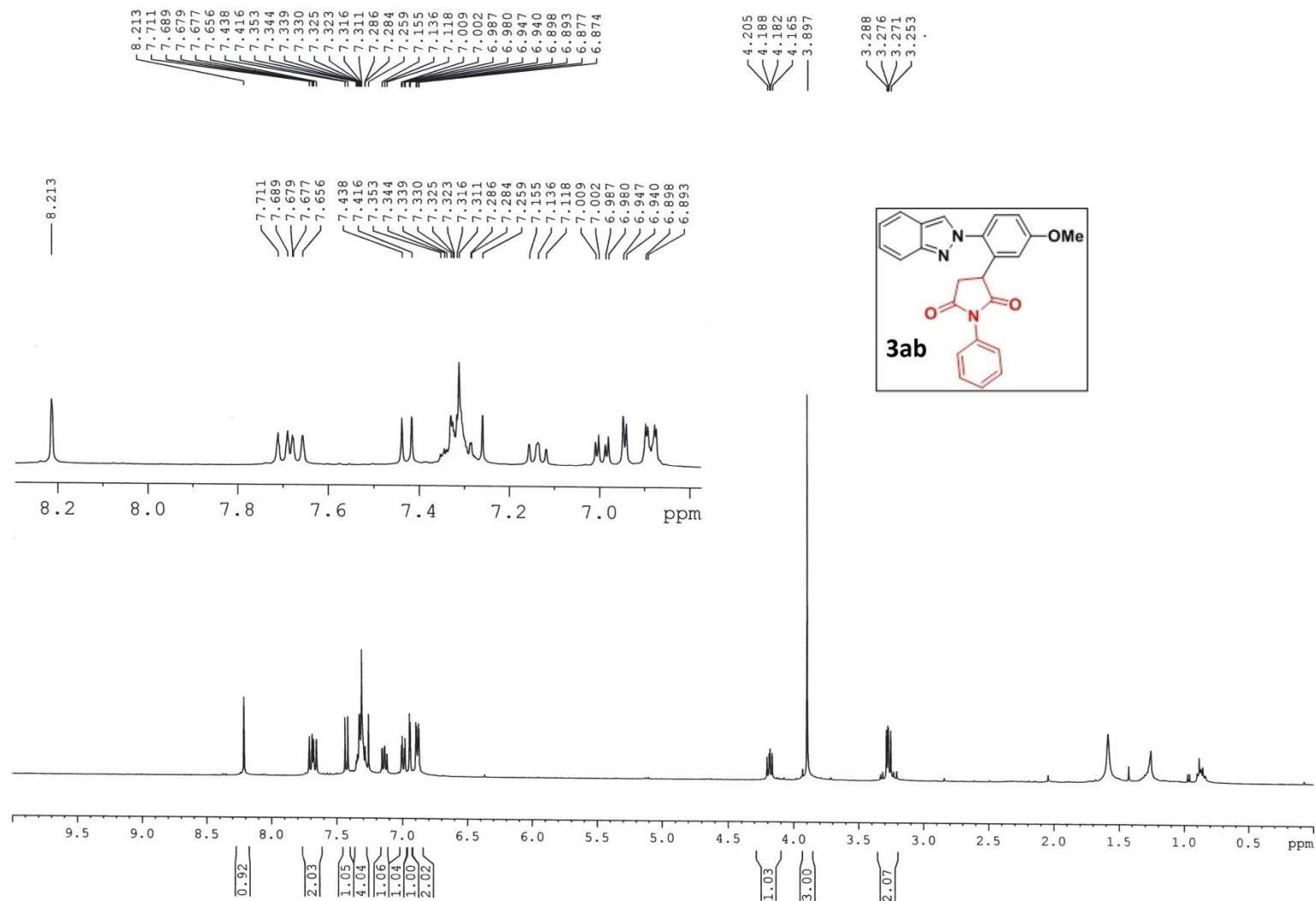
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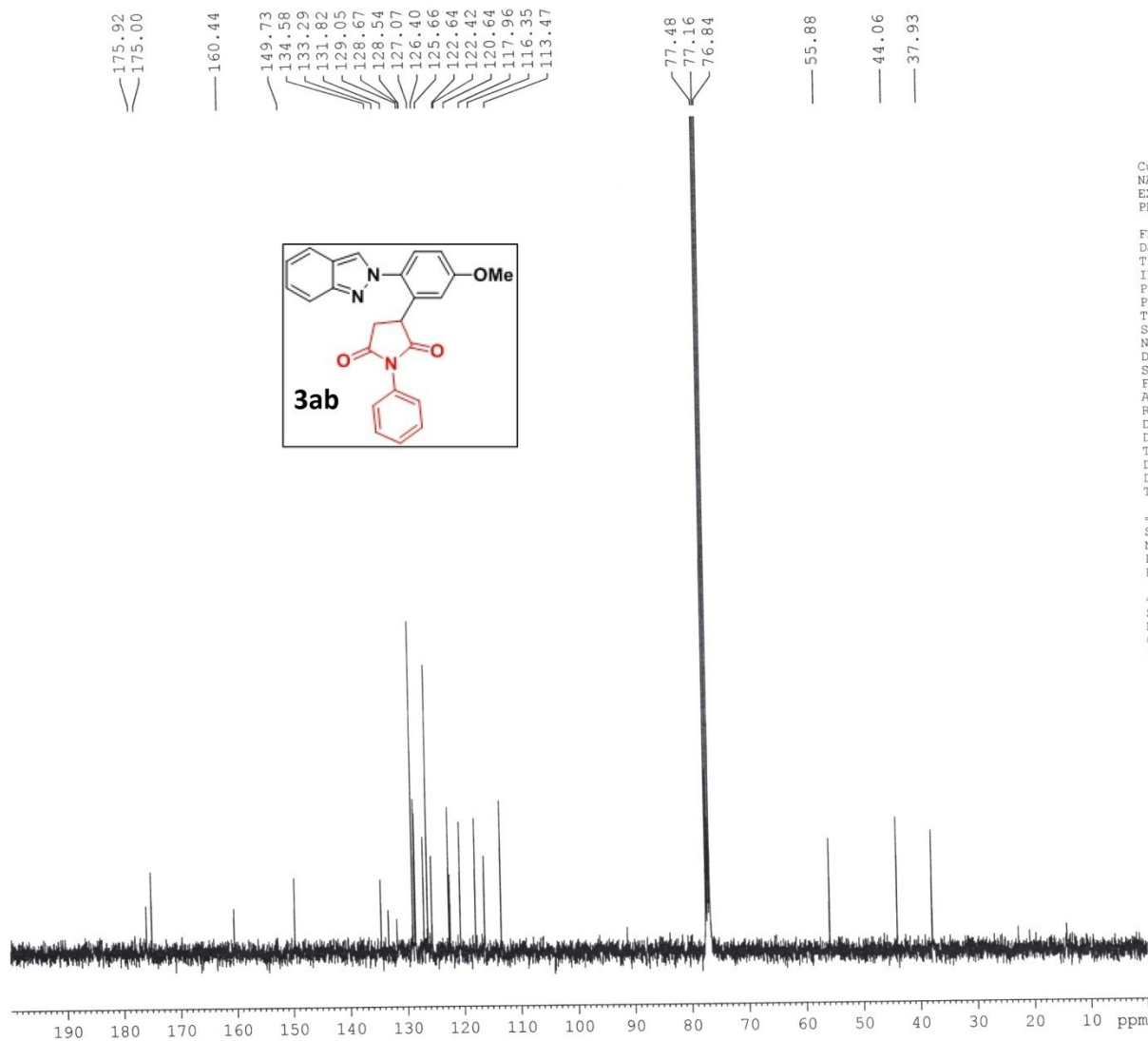
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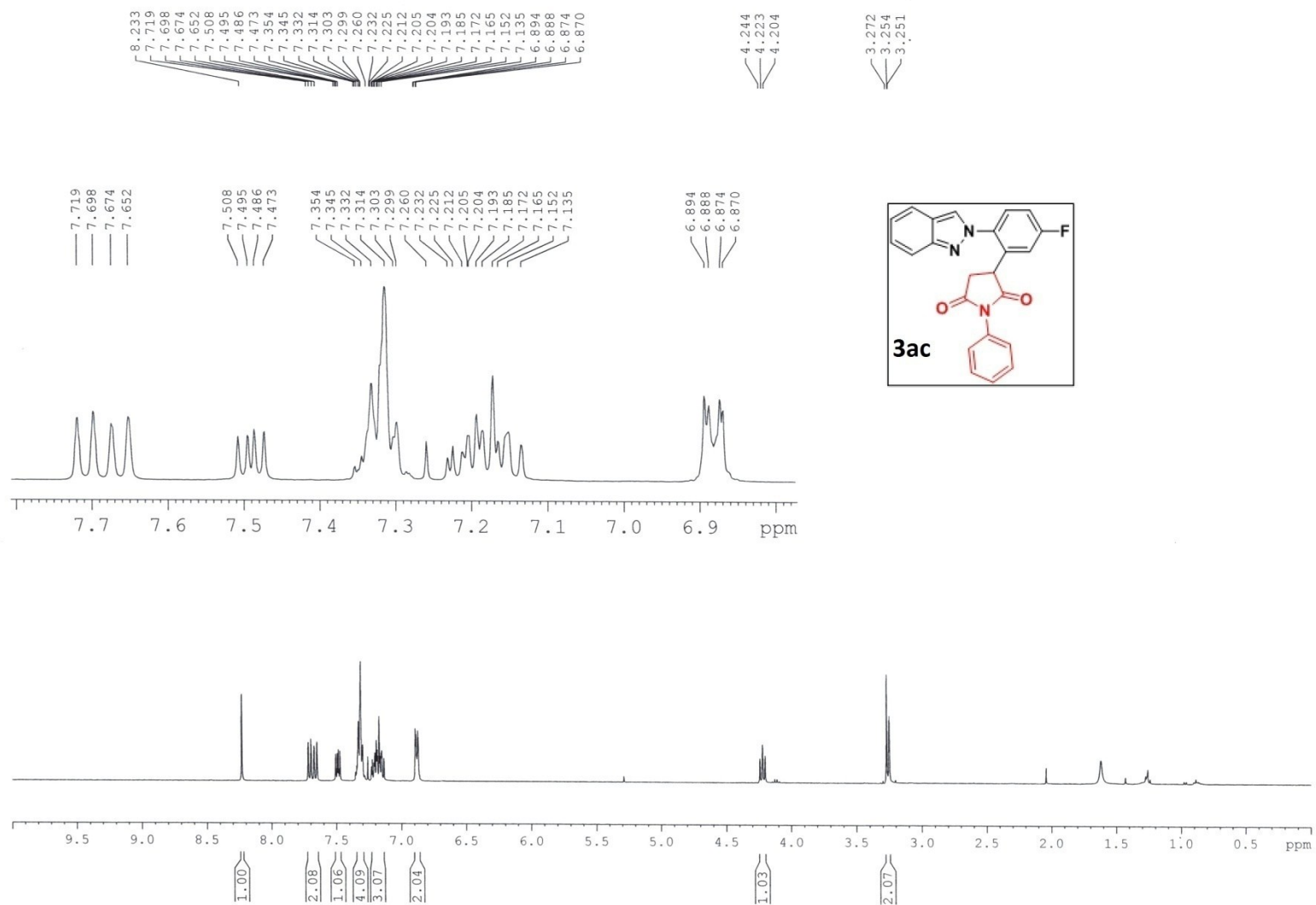
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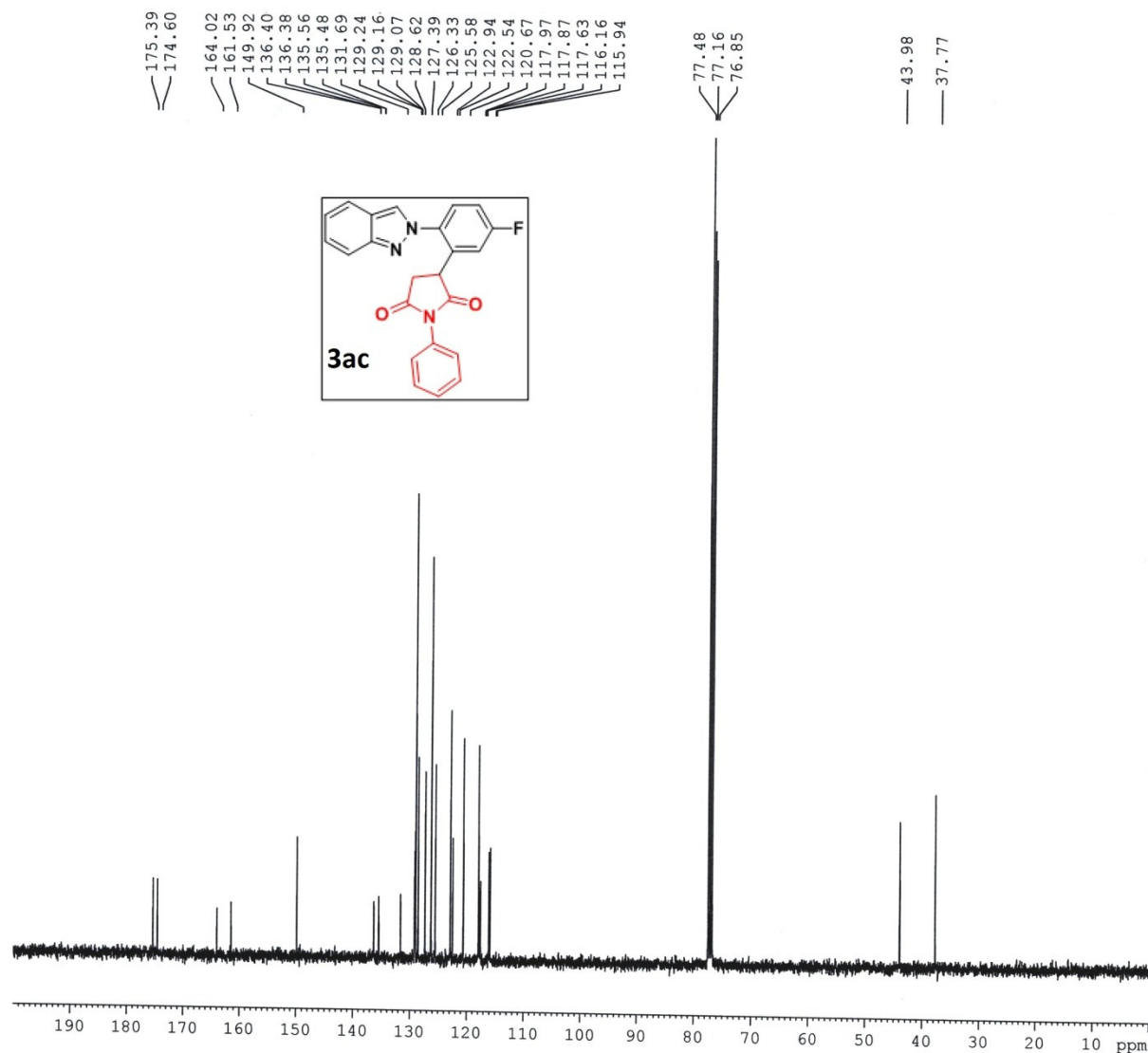
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 Time 9.03
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDCl3
 NS 624
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 297.9 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TDO 1

===== CHANNEL f1 =====
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

===== CHANNEL f2 =====
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG[2] waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177839 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40





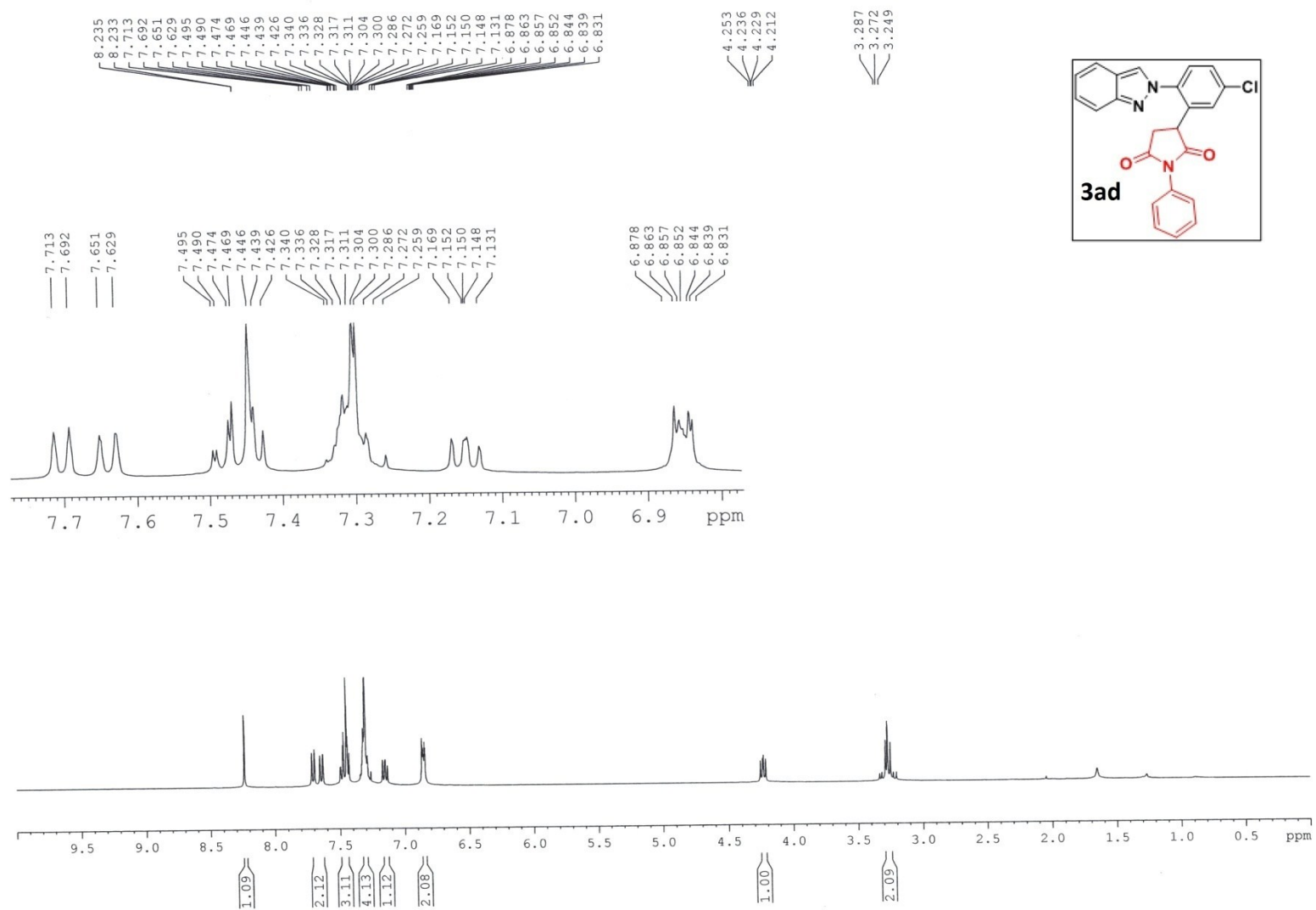
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 402
PROCNO 1

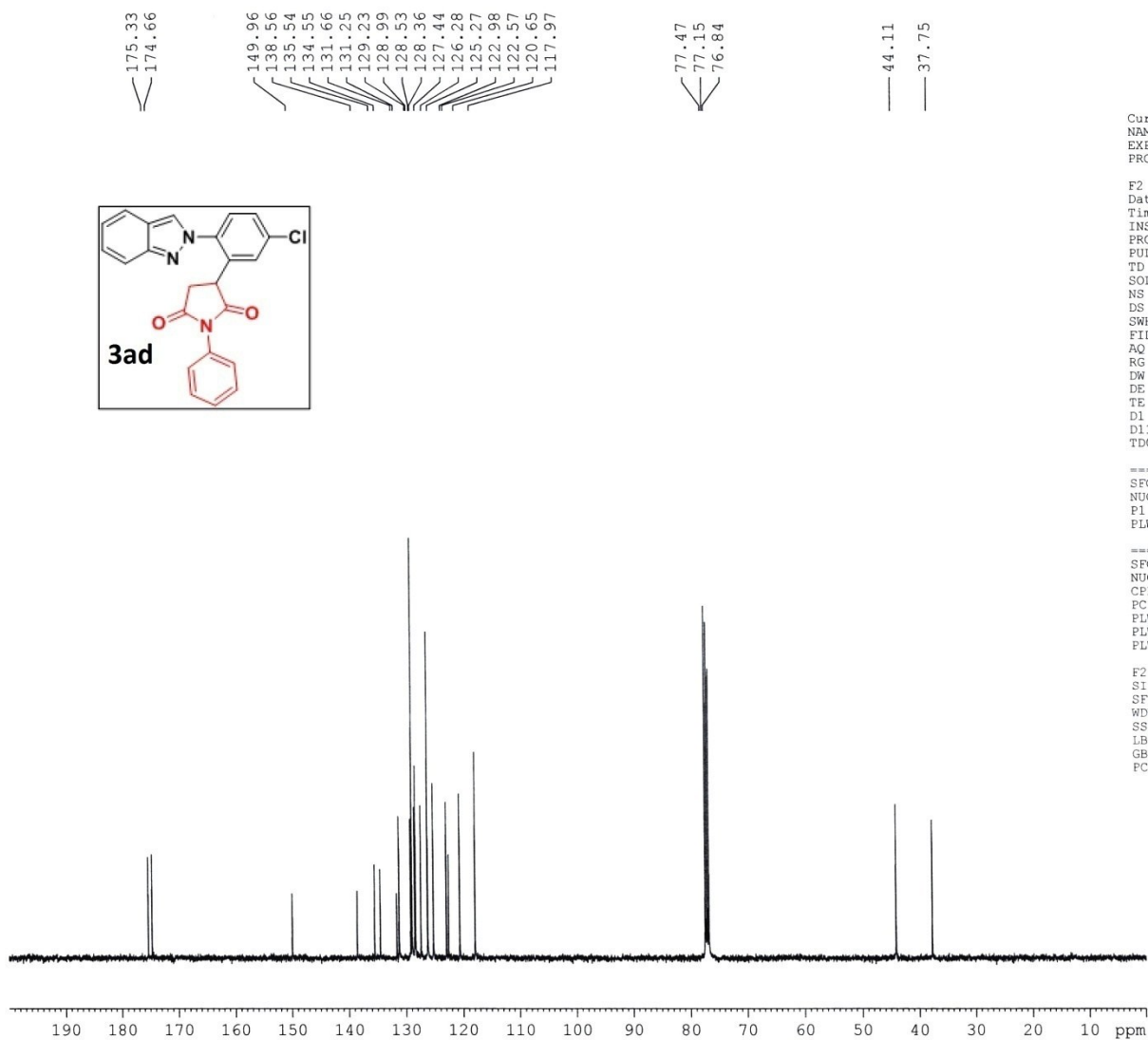
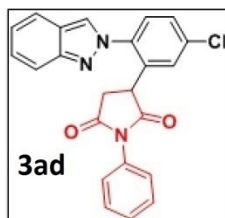
F2 - Acquisition Parameters
Date_ 20191003
Time_ 13.03
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDC13
NS 420
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 186.42
DW 20.800 usec
DE 6.50 usec
TE 299.2 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177854 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
FC 1.40





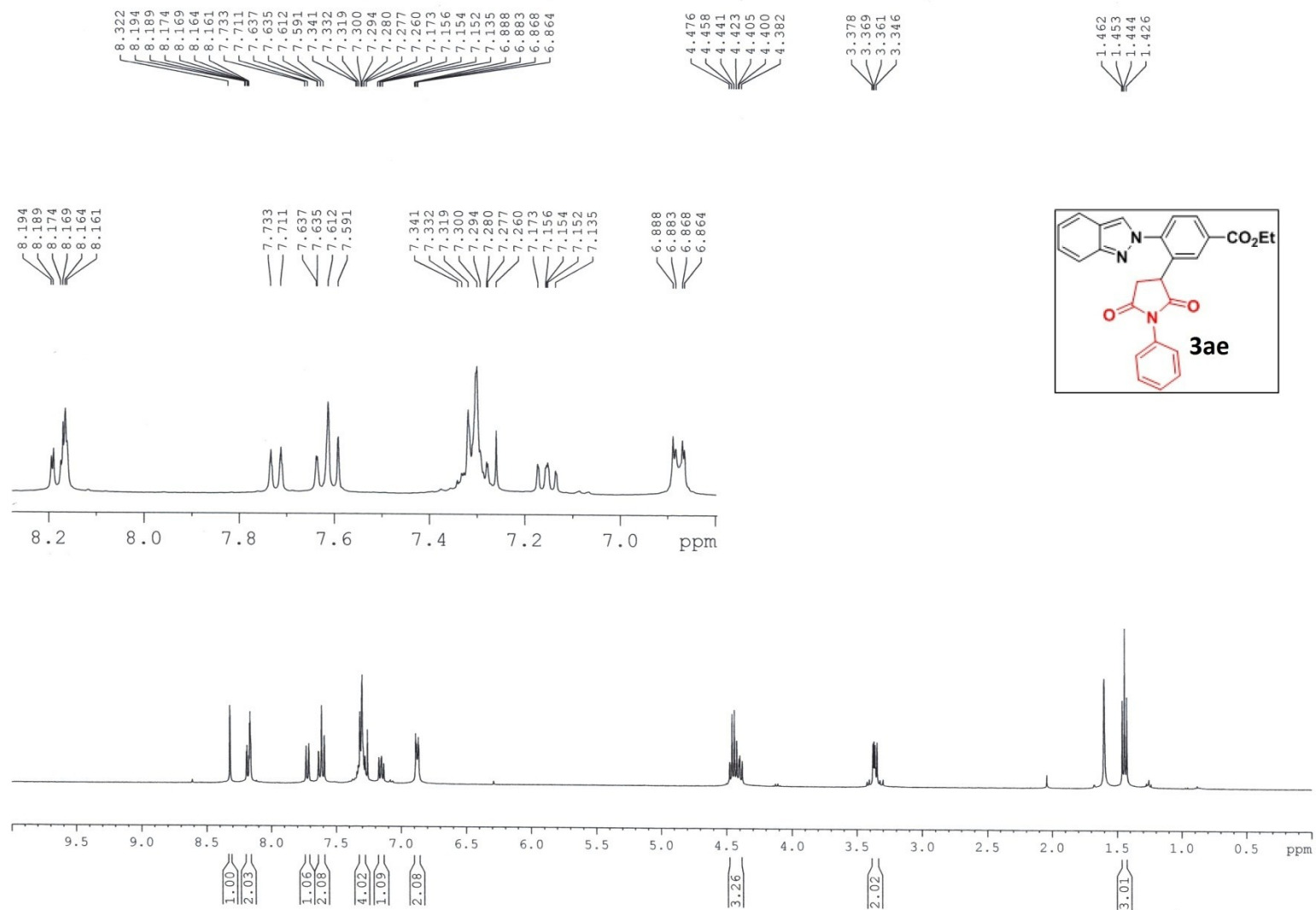
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 412
PROCNO 1

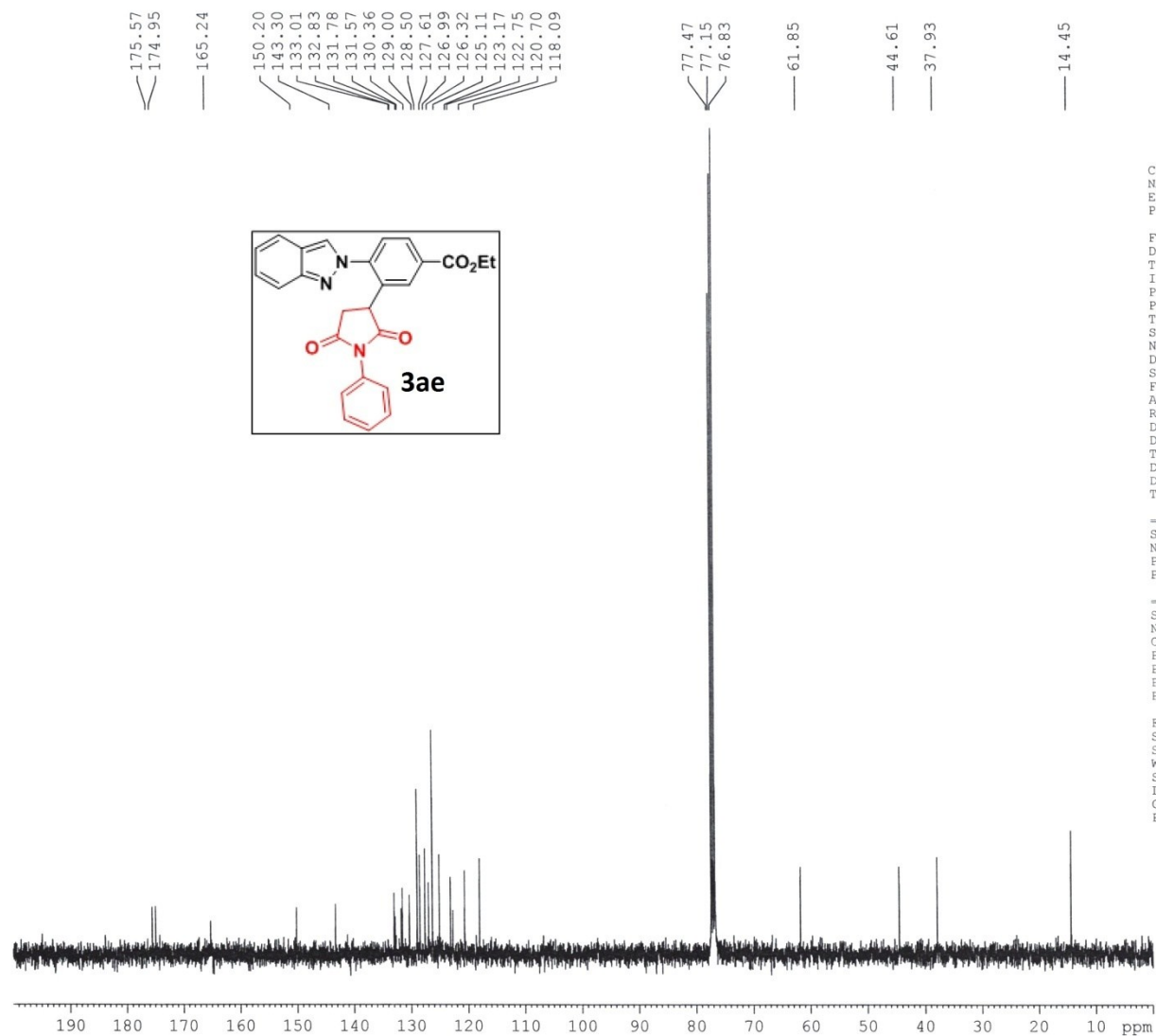
F2 - Acquisition Parameters
Date_ 20191017
Time_ 9.36
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDC13
NS 512
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 186.42
DW 20.800 usec
DE 6.50 usec
TE 296.6 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
PCPDG12 waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177903 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40





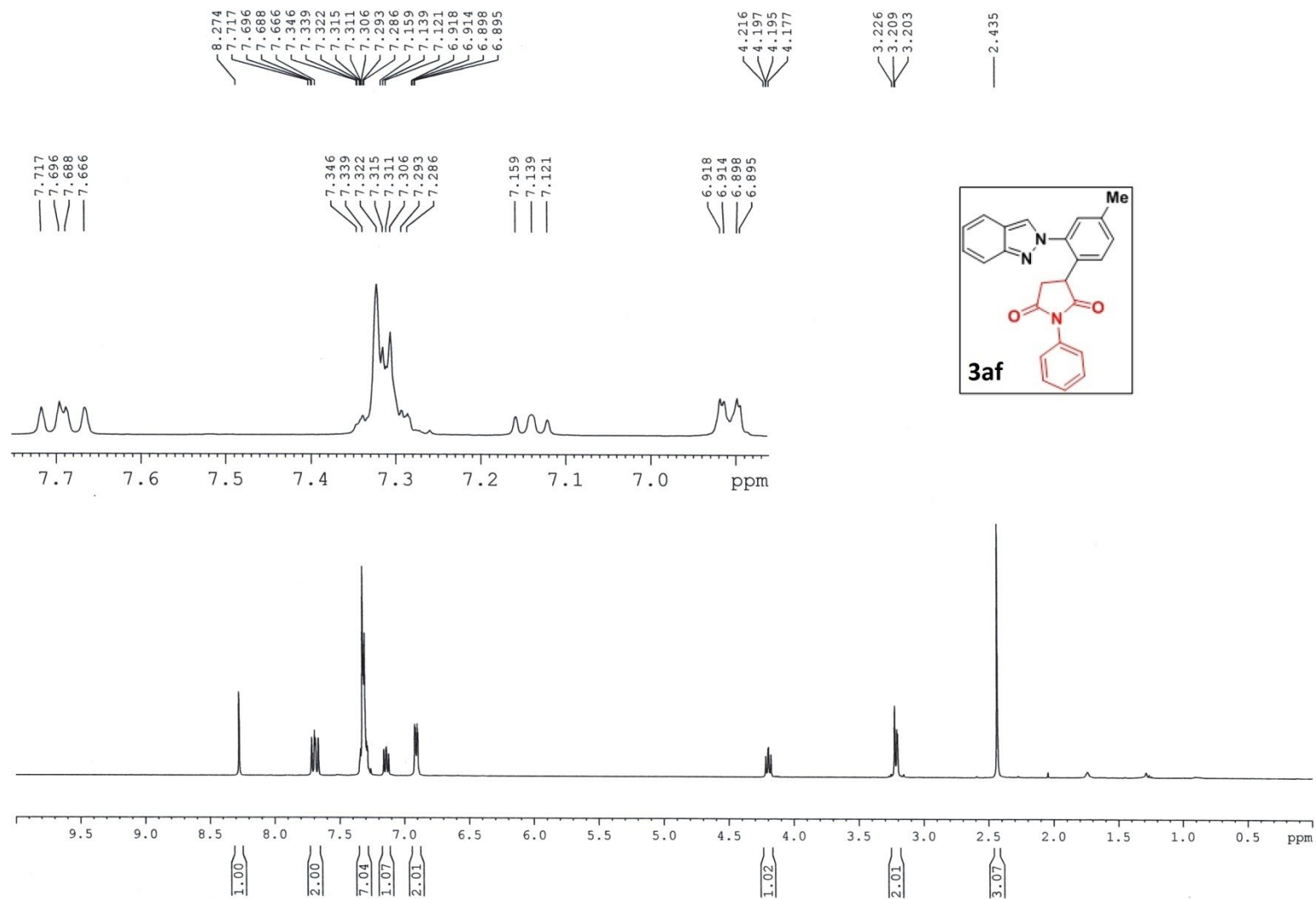
Current Data Parameters
 NAME Dr. A HAJRA-2019-13C
 EXPNO 478
 PROCNO 1

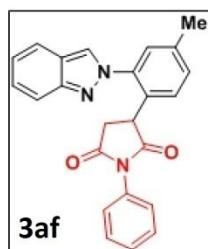
F2 - Acquisition Parameters
 Date_ 20191030
 Time_ 16.43
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDC13
 NS 220
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 297.3 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TD0 1

===== CHANNEL f1 =====
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

===== CHANNEL f2 =====
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG[2] waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177855 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40





176.19
175.10

149.70
139.78
139.34
131.84
130.49
129.83
128.89
128.33
127.77
127.02
126.31
125.24
122.58
122.36
120.59
117.91

77.47
77.15
76.83

43.60
37.88

20.95



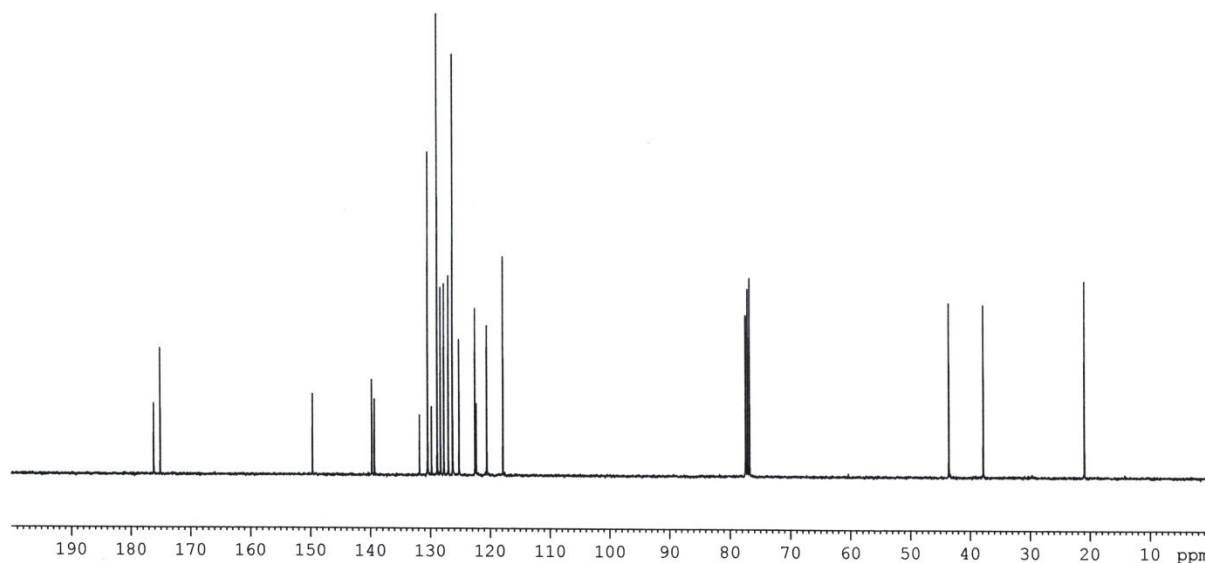
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 448
PROCNO 1

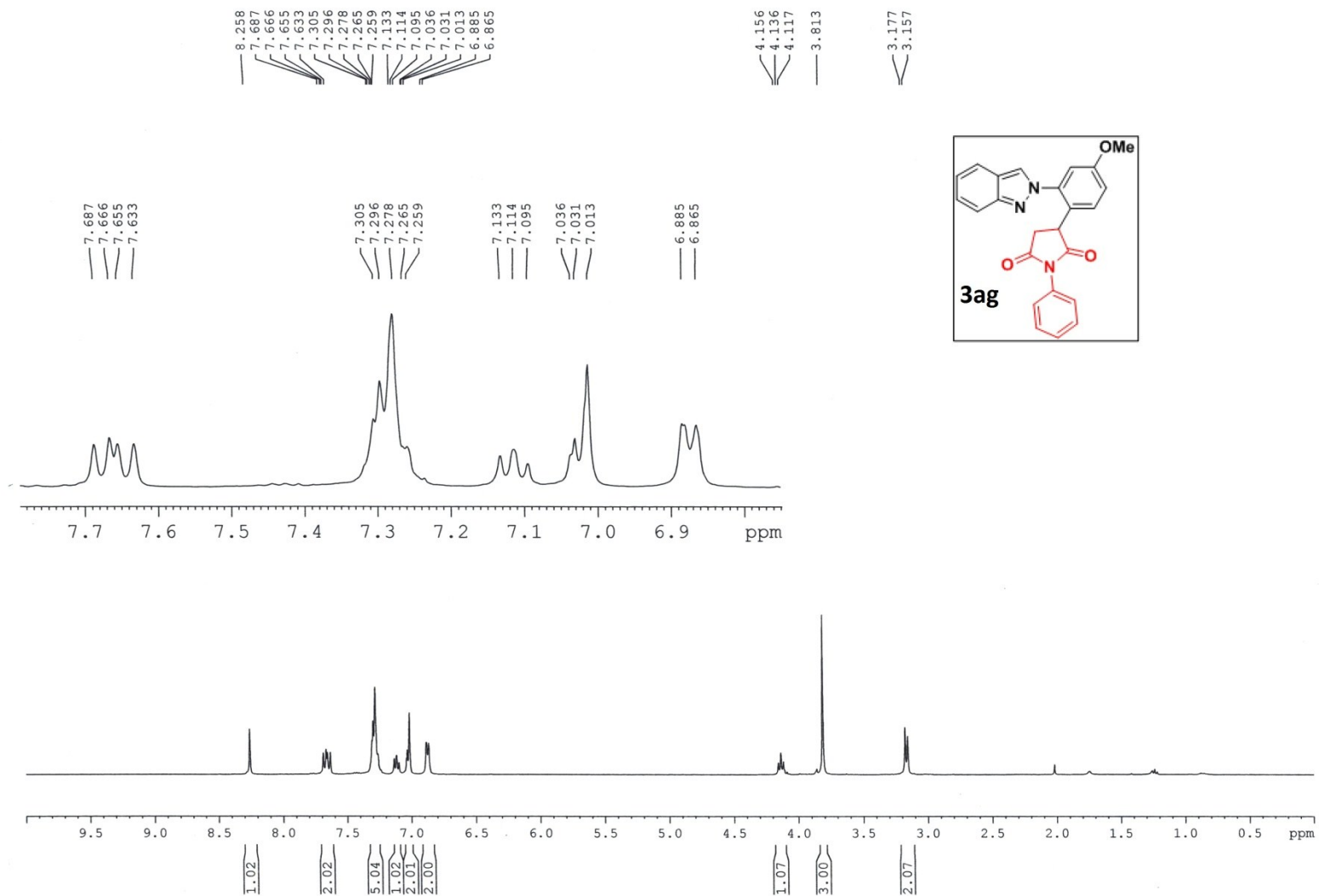
F2 - Acquisition Parameters
Date 20191023
Time 19.31
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 550
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 186.42
DW 20.800 usec
DE 6.50 usec
TE 299.4 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

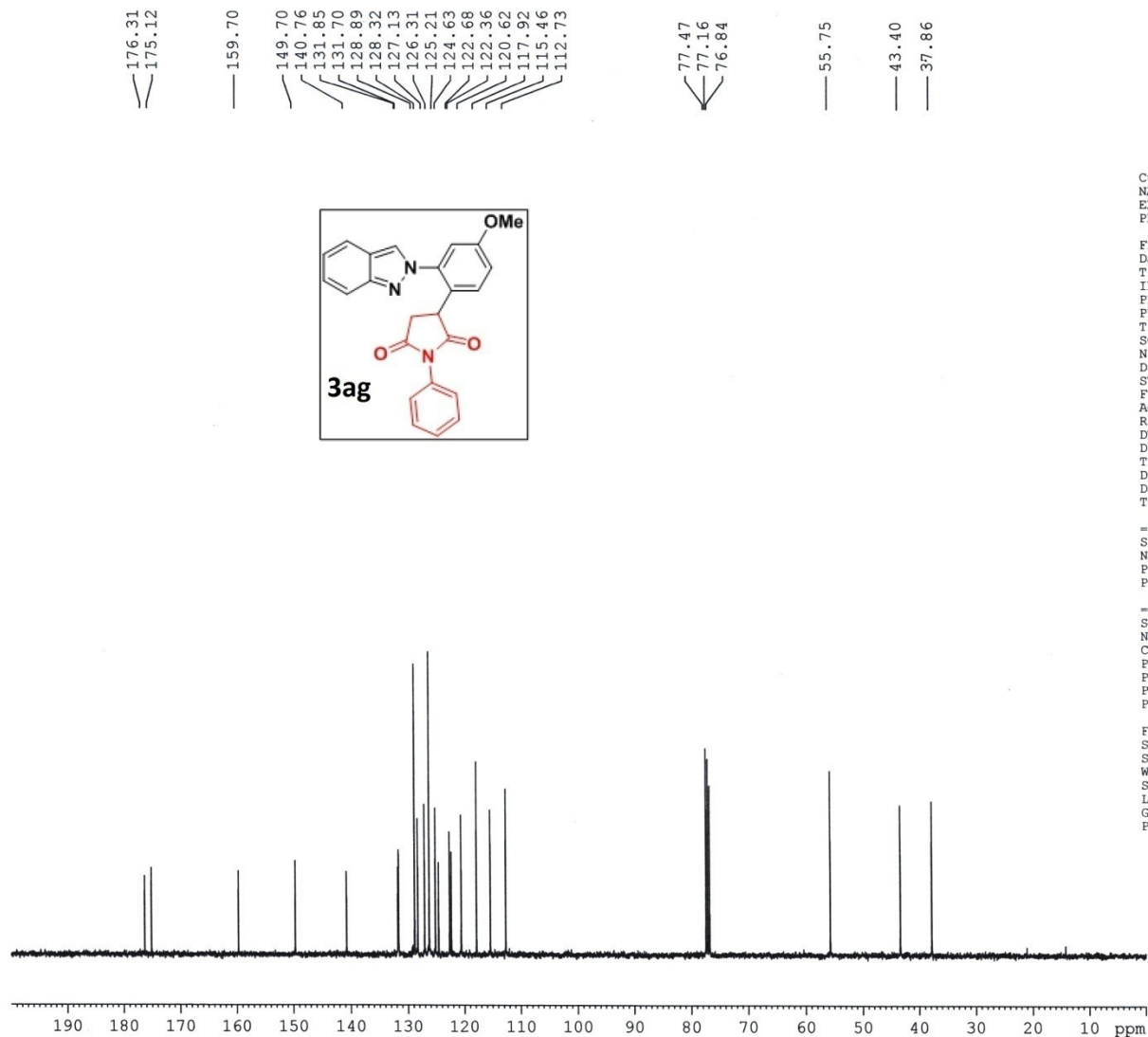
===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177976 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40







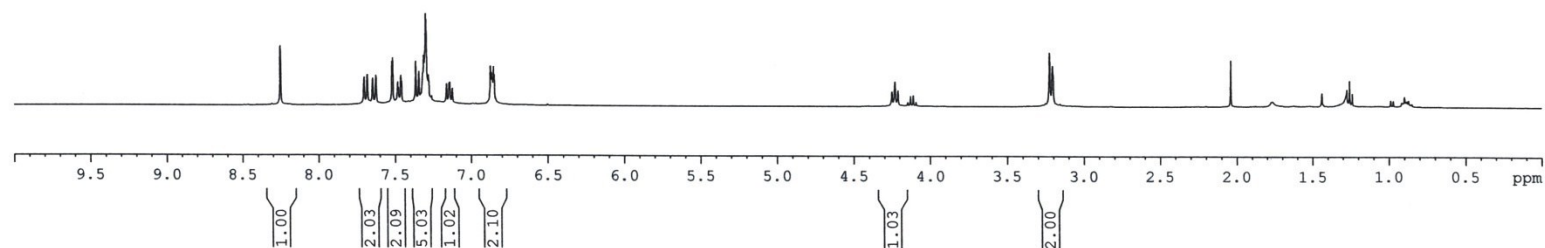
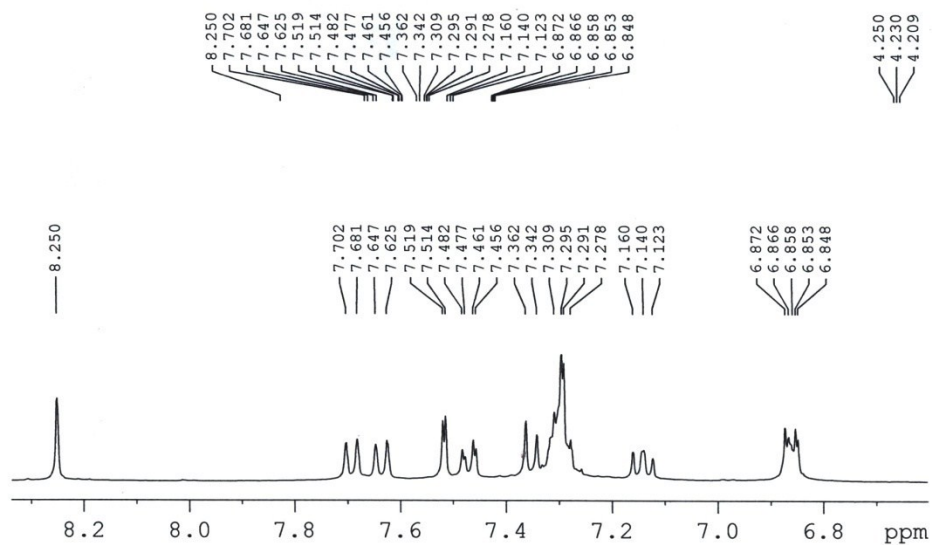
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 485
PROCNO 1

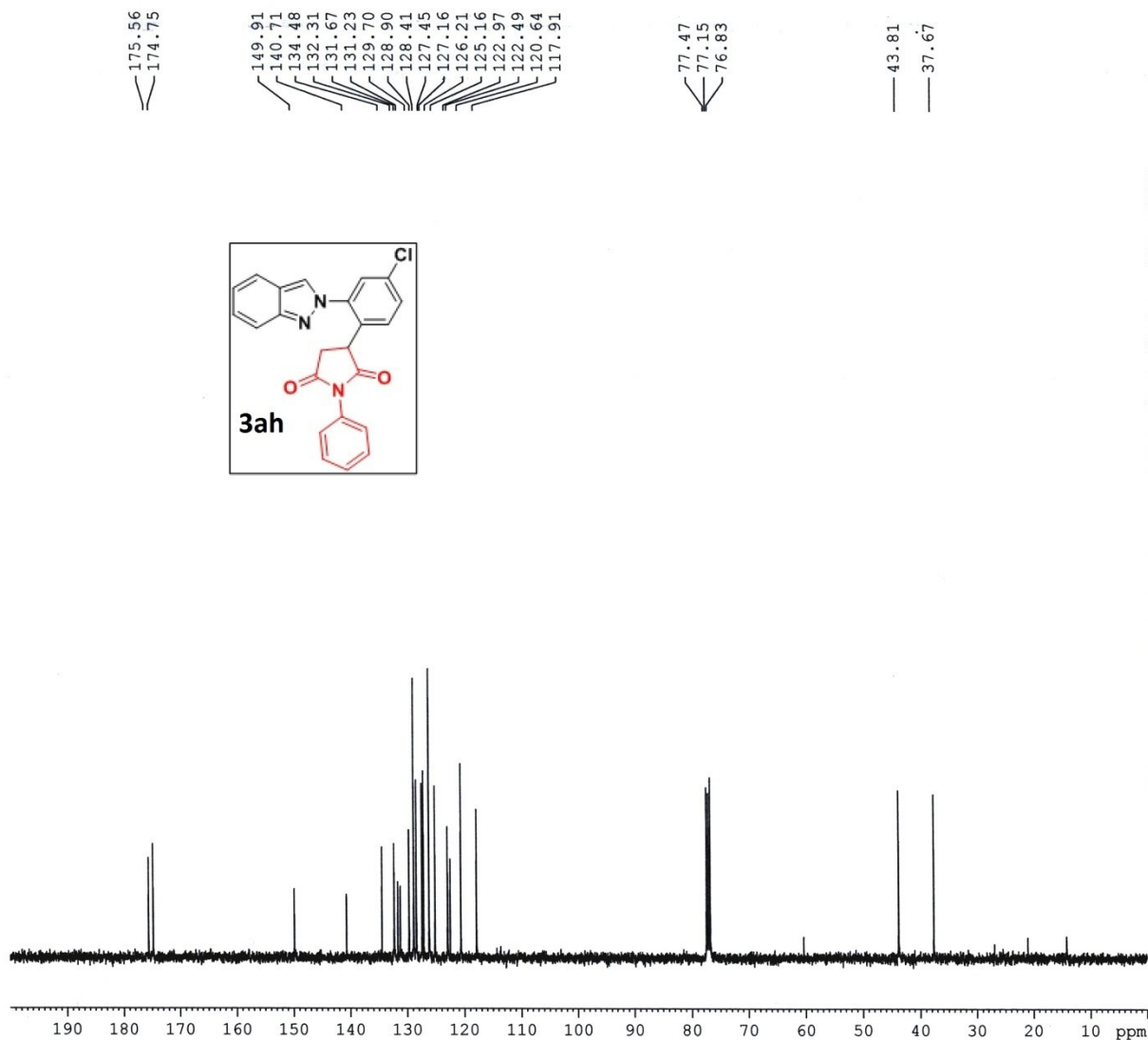
F2 - Acquisition Parameters
Date_ 20191101
Time 10.10
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 120
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 186.42
DW 20.800 usec
DE 6.50 usec
TE 298.4 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177976 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40





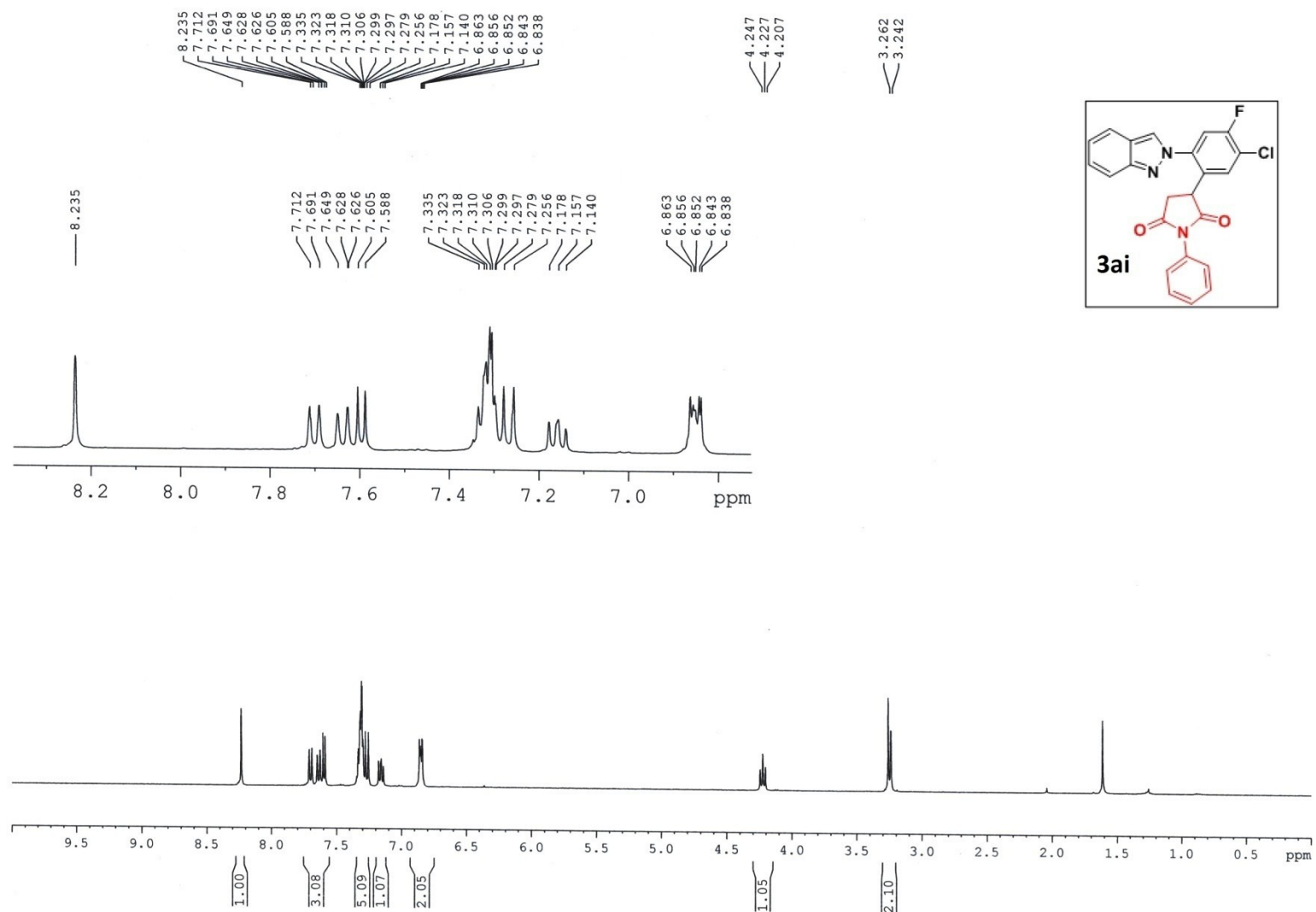
Current Data Parameters
 NAME Dr. A HAJRA-2019-13C
 EXPNO 474
 PROCNO 1

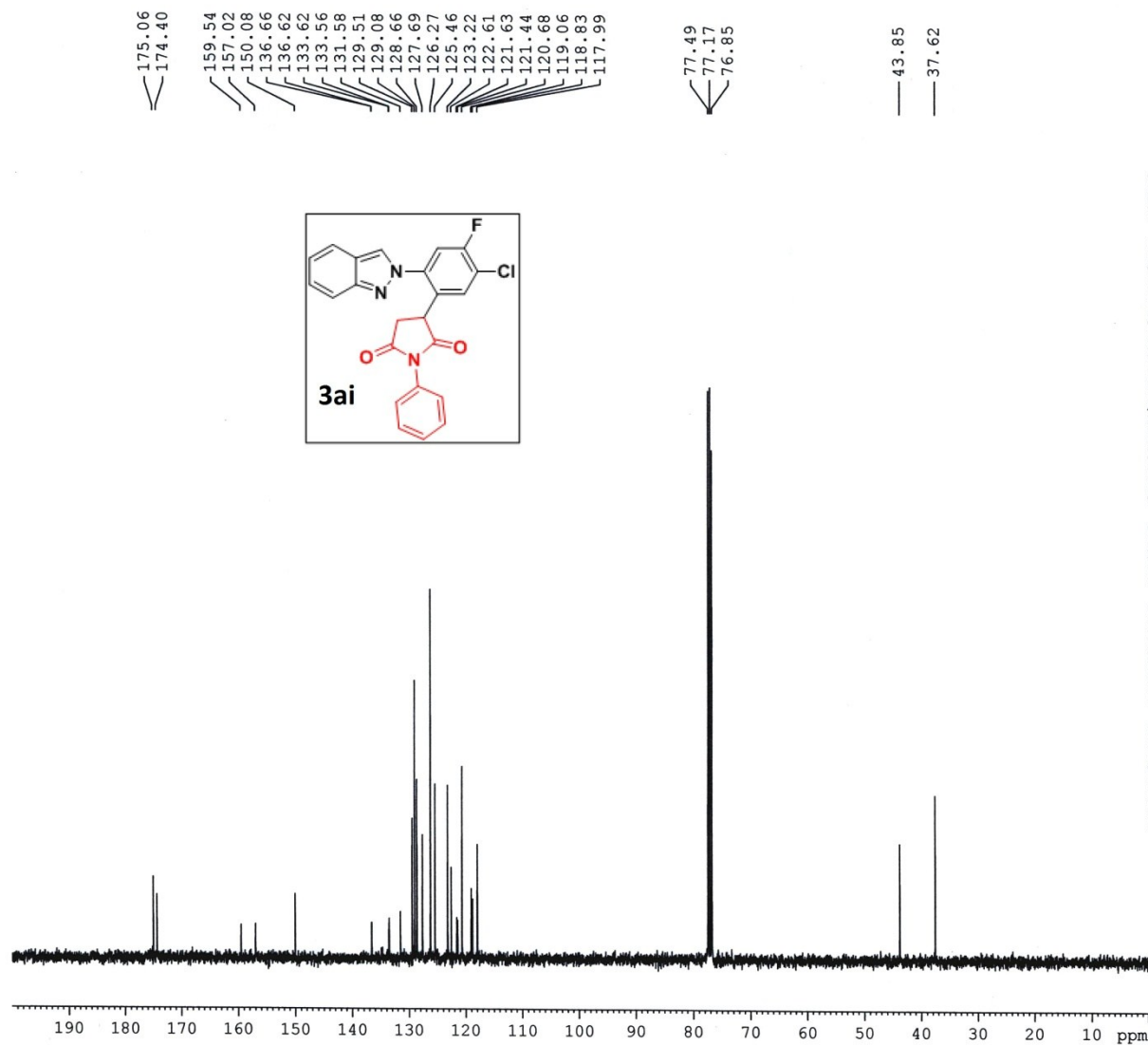
F2 - Acquisition Parameters
 Date_ 20191028
 Time 17.09
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDCl3
 NS 50
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 296.9 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TDO 1

----- CHANNEL f1 -----
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

----- CHANNEL f2 -----
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG[2] waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177985 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40





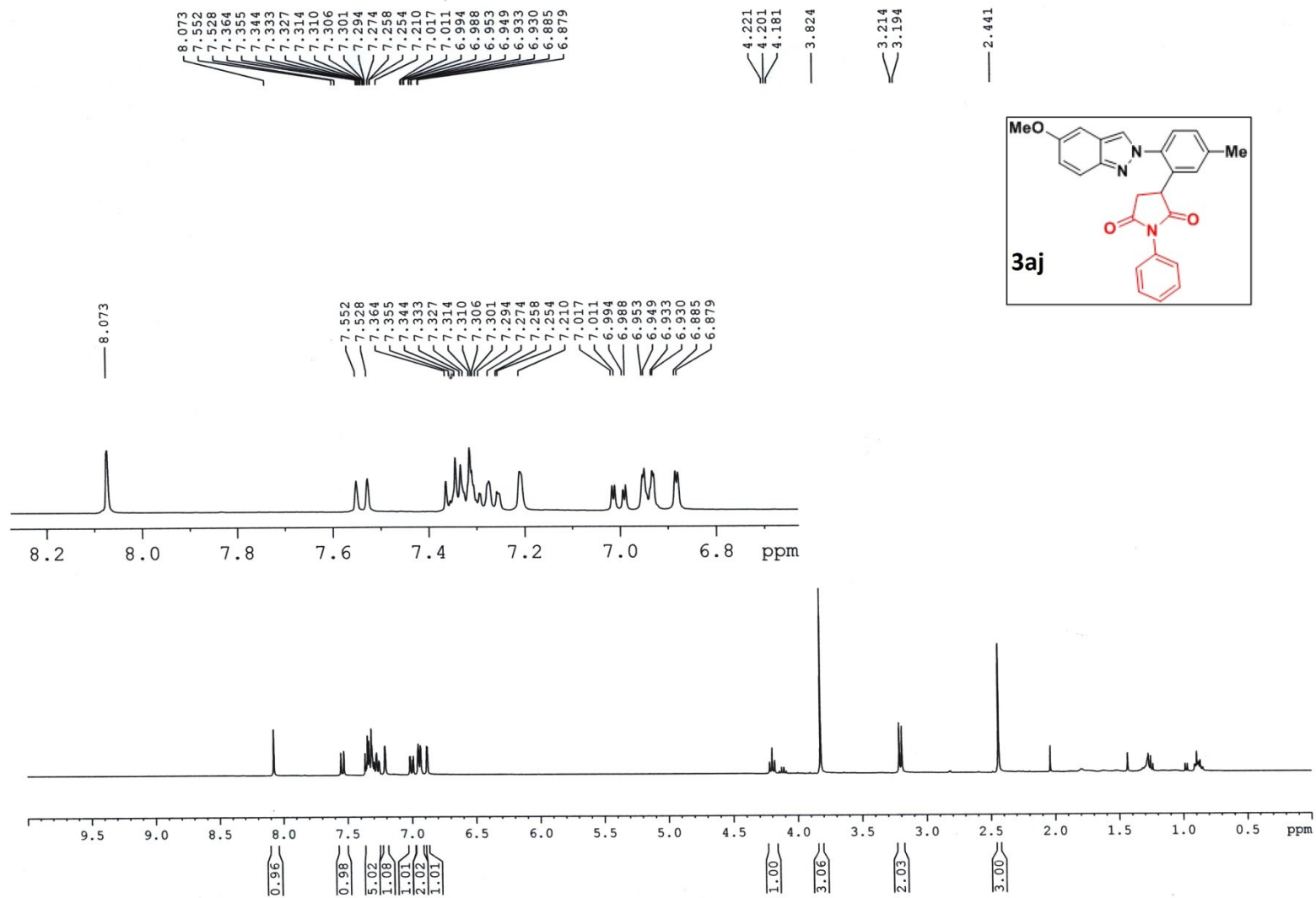
Current Data Parameters
 NAME Dr. A HAJRA-2019-13C
 EXPNO 479
 PROCNO 1

F2 - Acquisition Parameters
 Date_ 20191030
 Time_ 18.33
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDCl3
 NS 220
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 297.1 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TDO 1

===== CHANNEL f1 =====
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.0000000 W

===== CHANNEL f2 =====
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG12 waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177855 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40



176.16
175.11

155.61
146.58
139.75
137.59
132.47
131.86
131.47
129.59
128.87
128.34
126.89
126.30
124.19
122.39
122.10
119.24

96.45

77.47
77.15
76.83

55.40

44.09

37.83

21.18



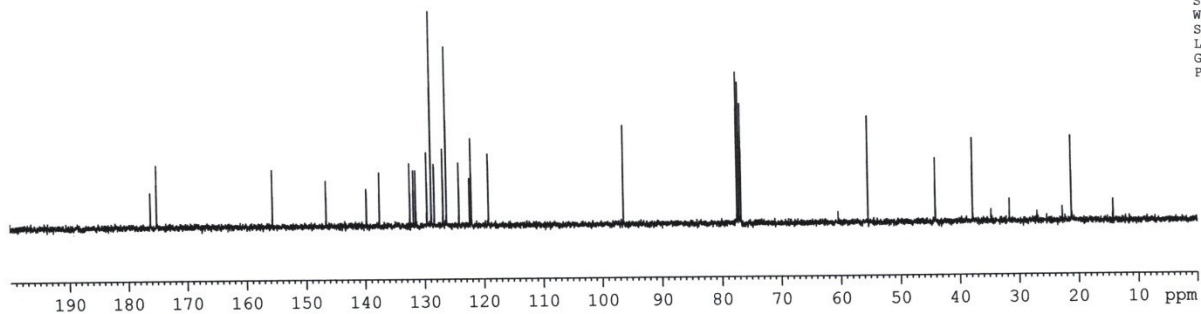
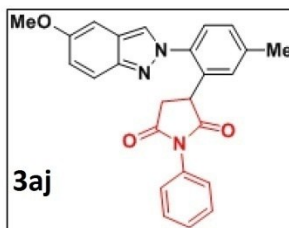
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 452
PROCNO 1

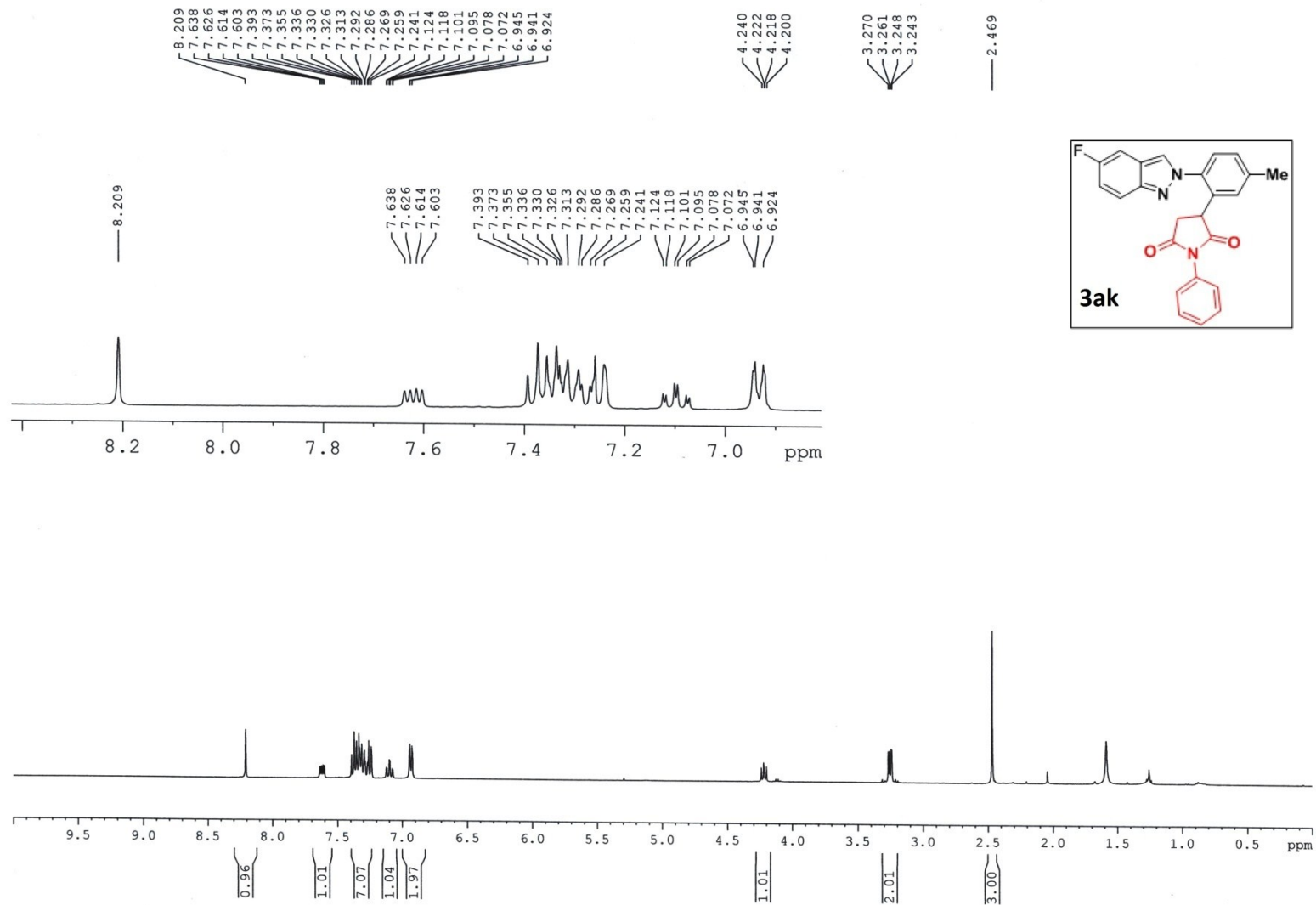
F2 - Acquisition Parameters
Date_ 20191024
Time 12.04
INSTRUM spect
PROBHD 5 mm F4BBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDC13
NS 30
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 186.42
DW 20.800 usec
DE 6.50 usec
TE 297.7 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

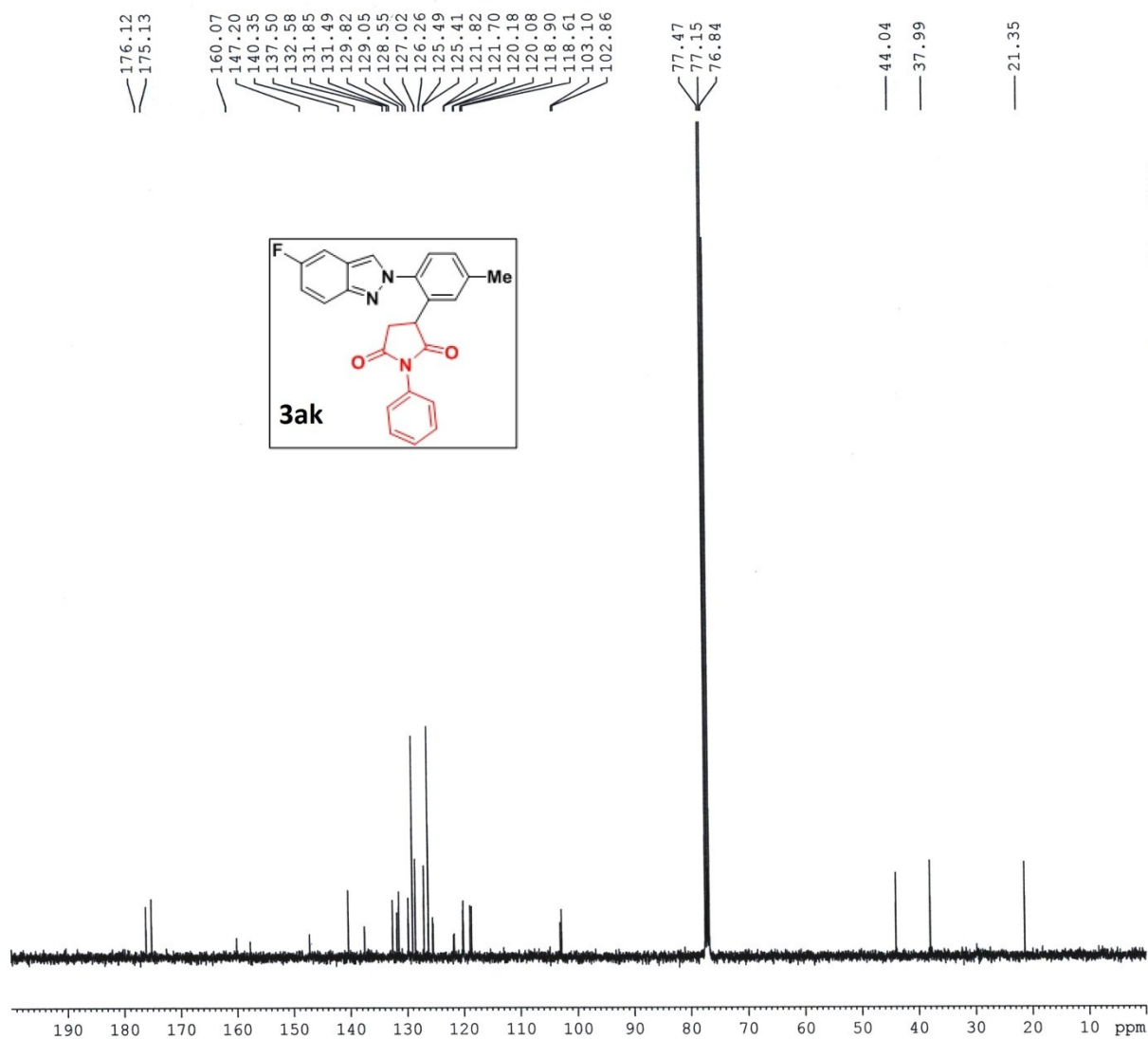
===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177968 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40







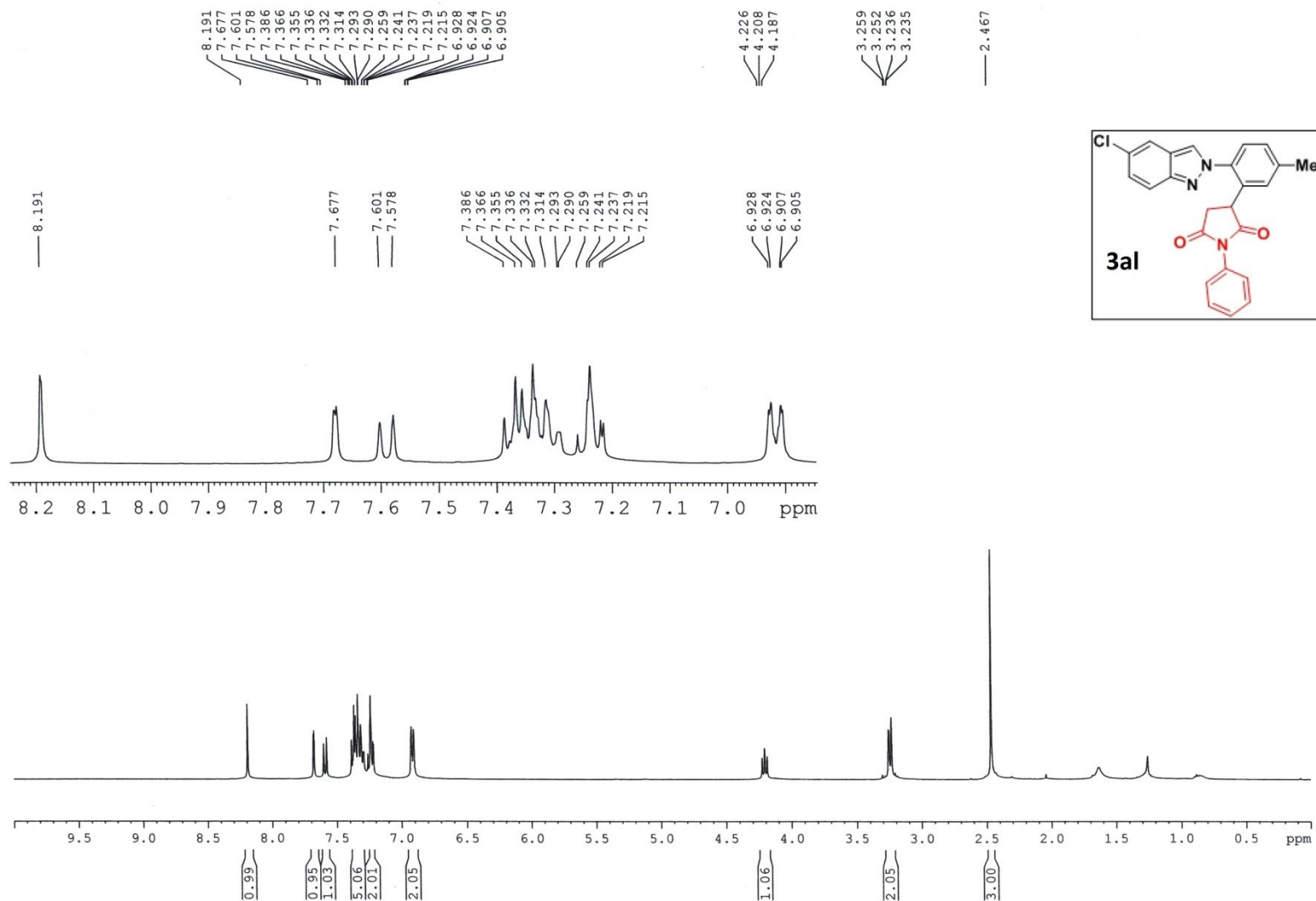
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 467
PROCNO 1

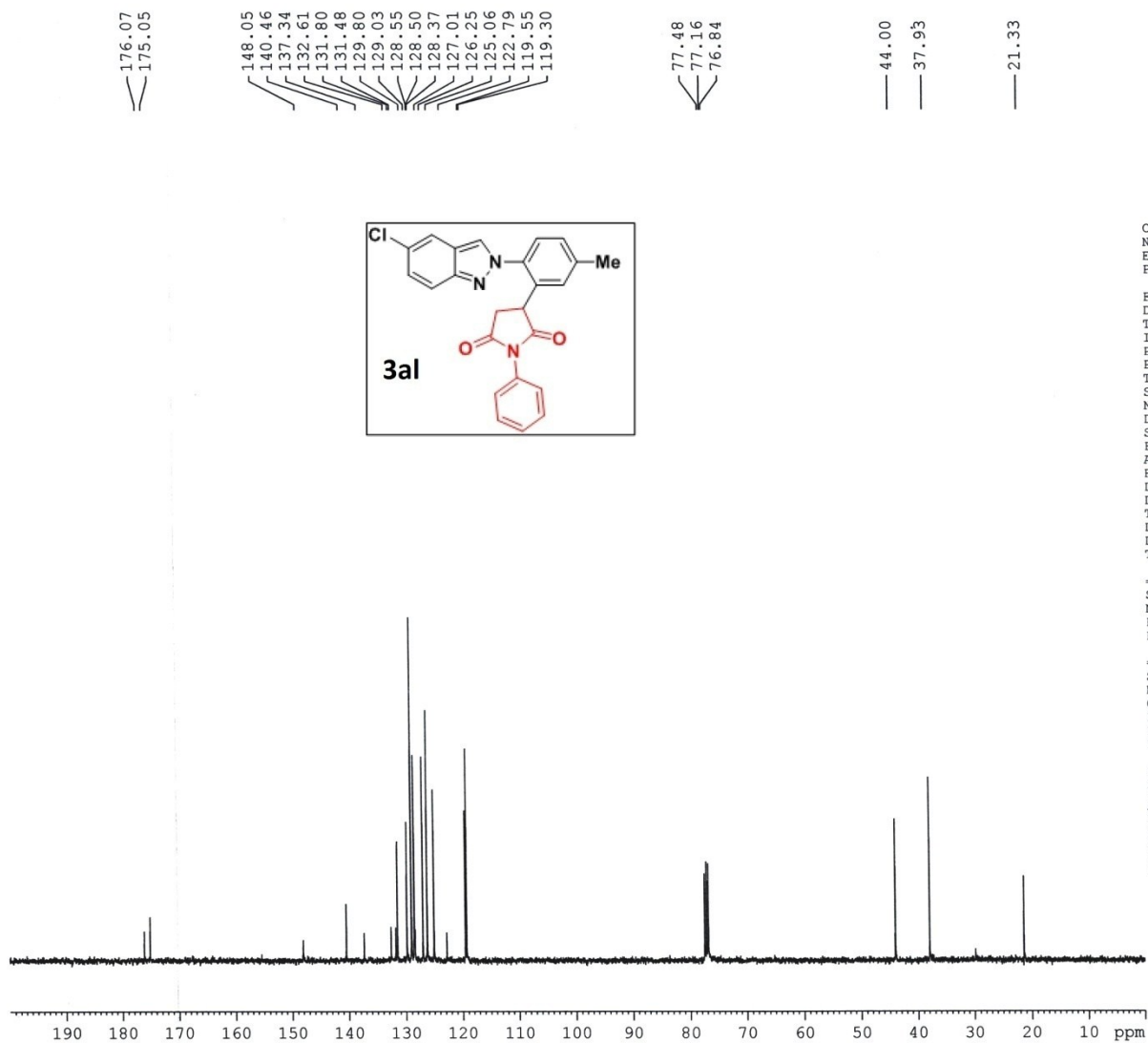
F2 - Acquisition Parameters
Date_ 20191027
Time 12.07
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 650
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 186.42
DW 20.800 usec
DE 6.50 usec
TE 297.6 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177844 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40





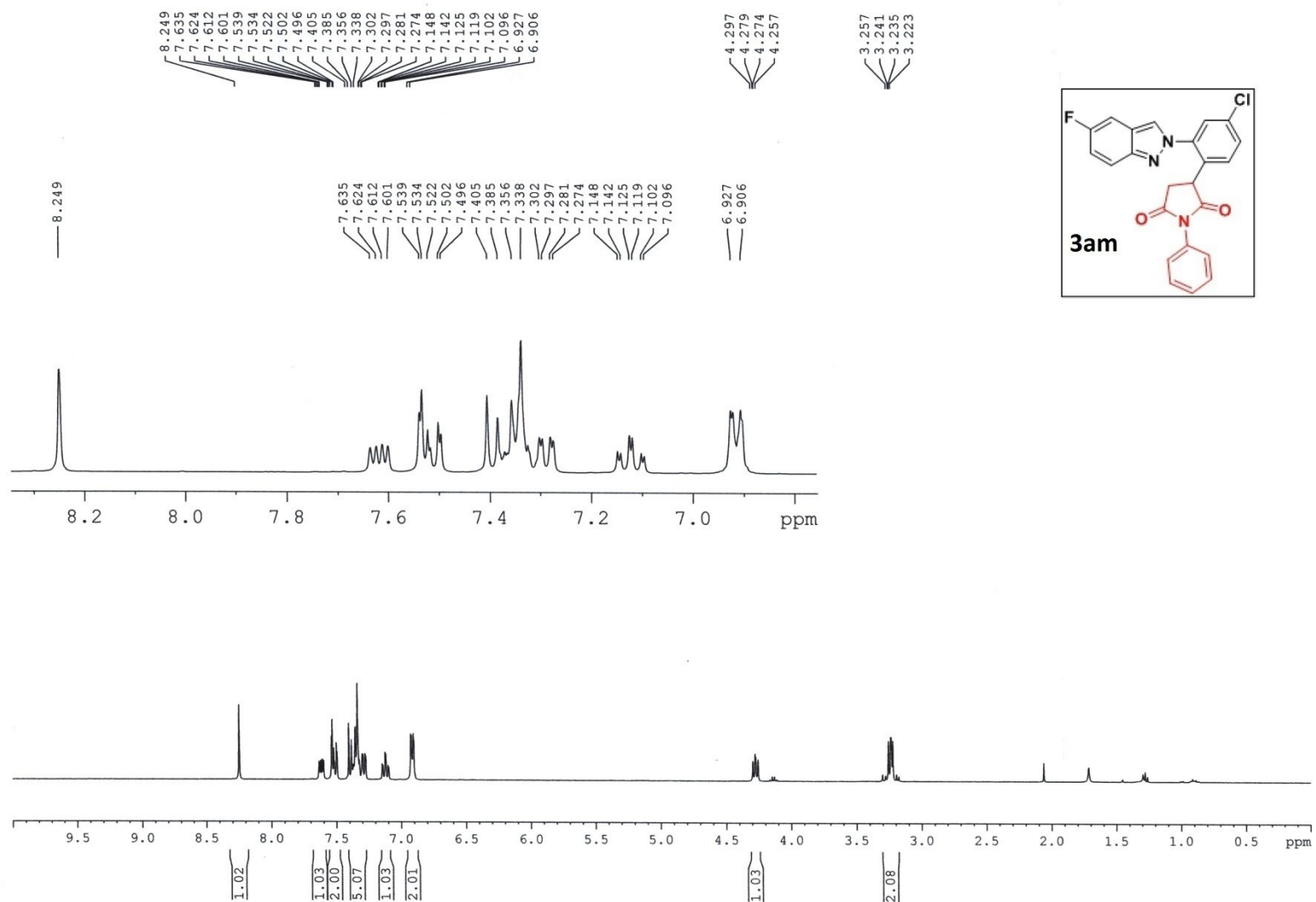
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 419
PROCNO 1

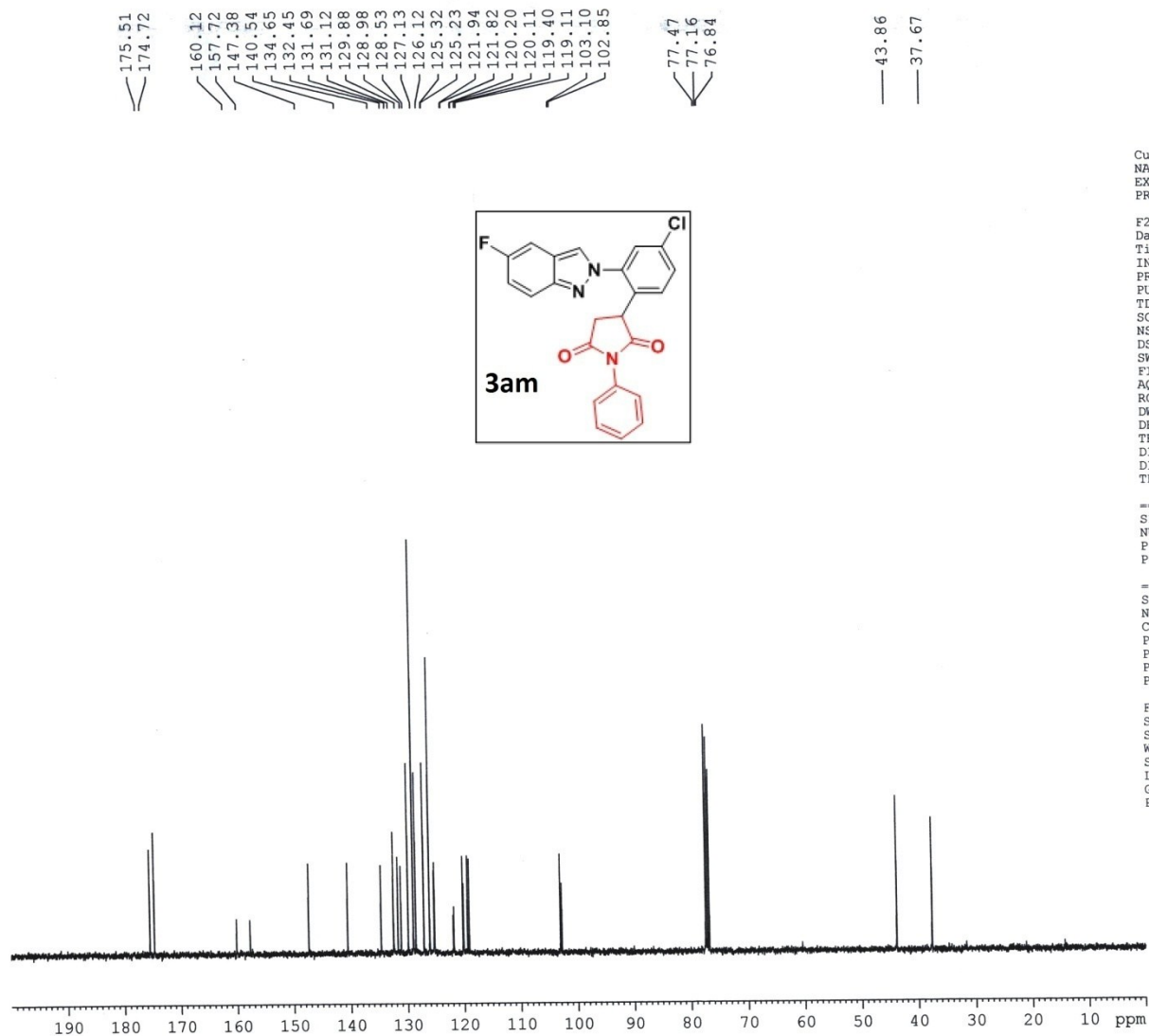
F2 - Acquisition Parameters
Date_ 20191017
Time 20.45
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgdc
TD 32768
SOLVENT CDCl₃
NS 420
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 186.42
DW 20.800 usec
DE 6.50 usec
TE 298.2 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W

F2 - Processing parameters
SI 16384
SF 100.6177867 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00





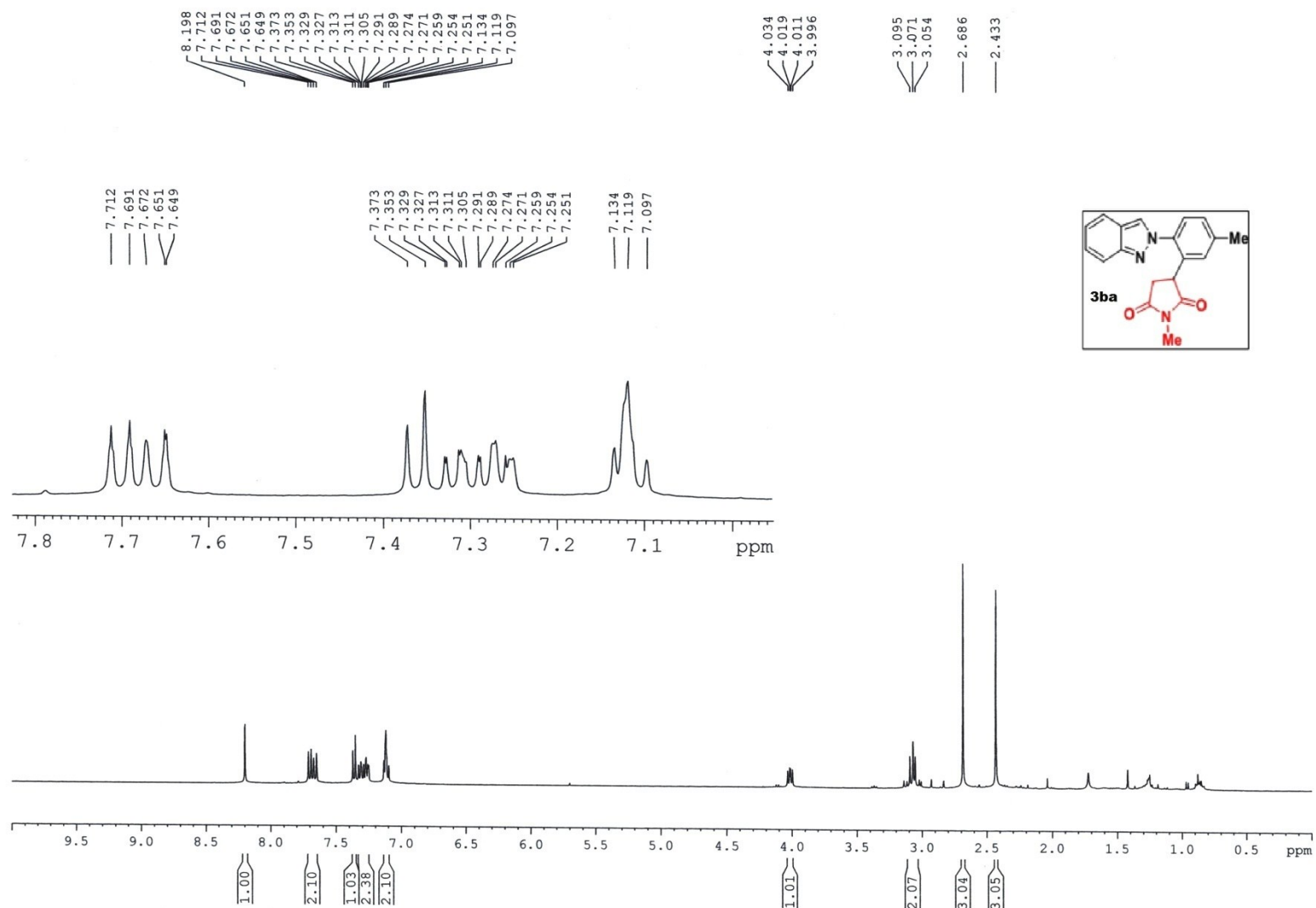
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 481
PROCNO 1

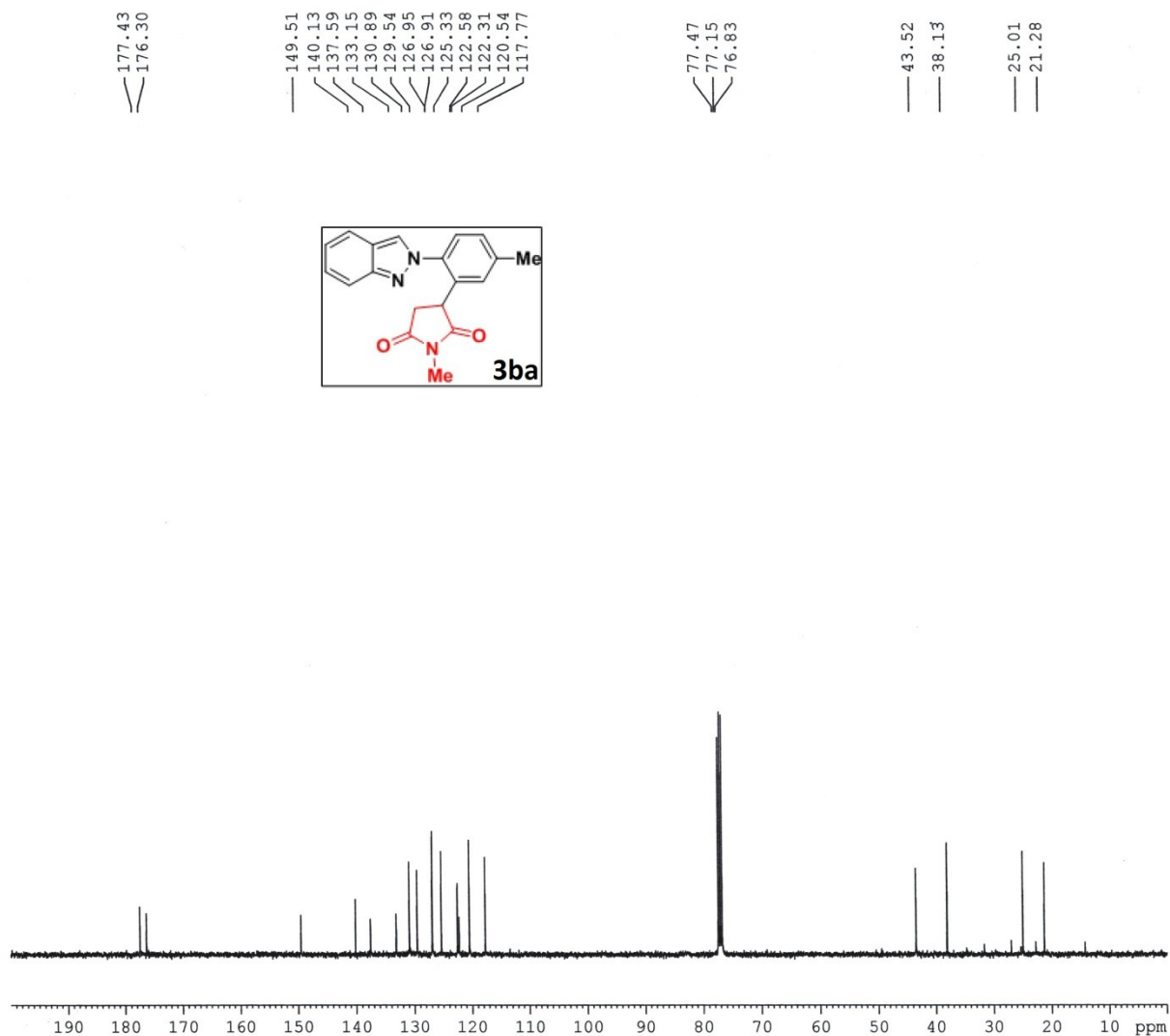
F2 - Acquisition Parameters
Date_ 20191031
Time 10.06
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 120
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 186.42
DW 20.800 usec
DE 6.50 usec
TE 298.3 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG(2) waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177931 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40





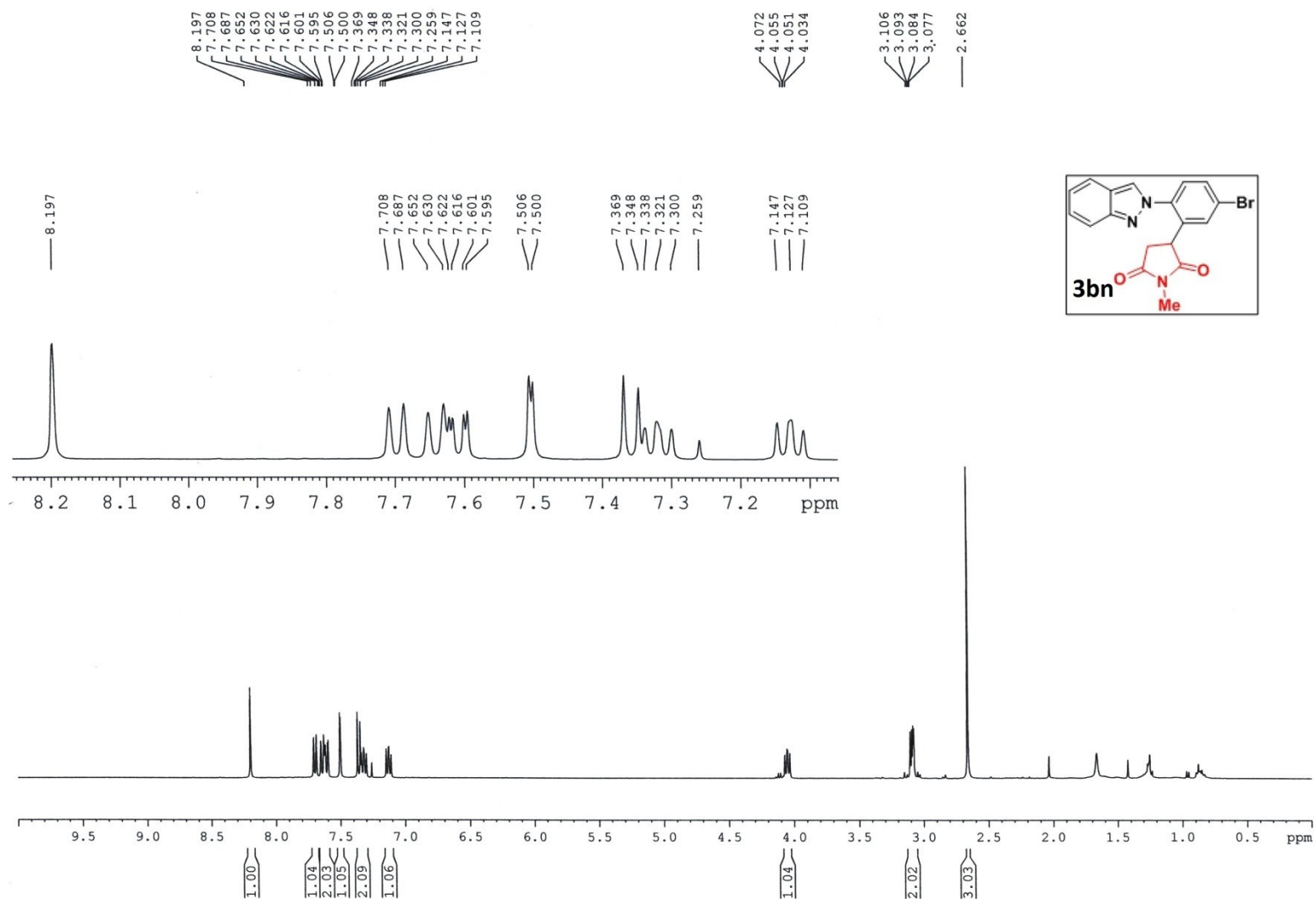
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 520
PROCNO 1

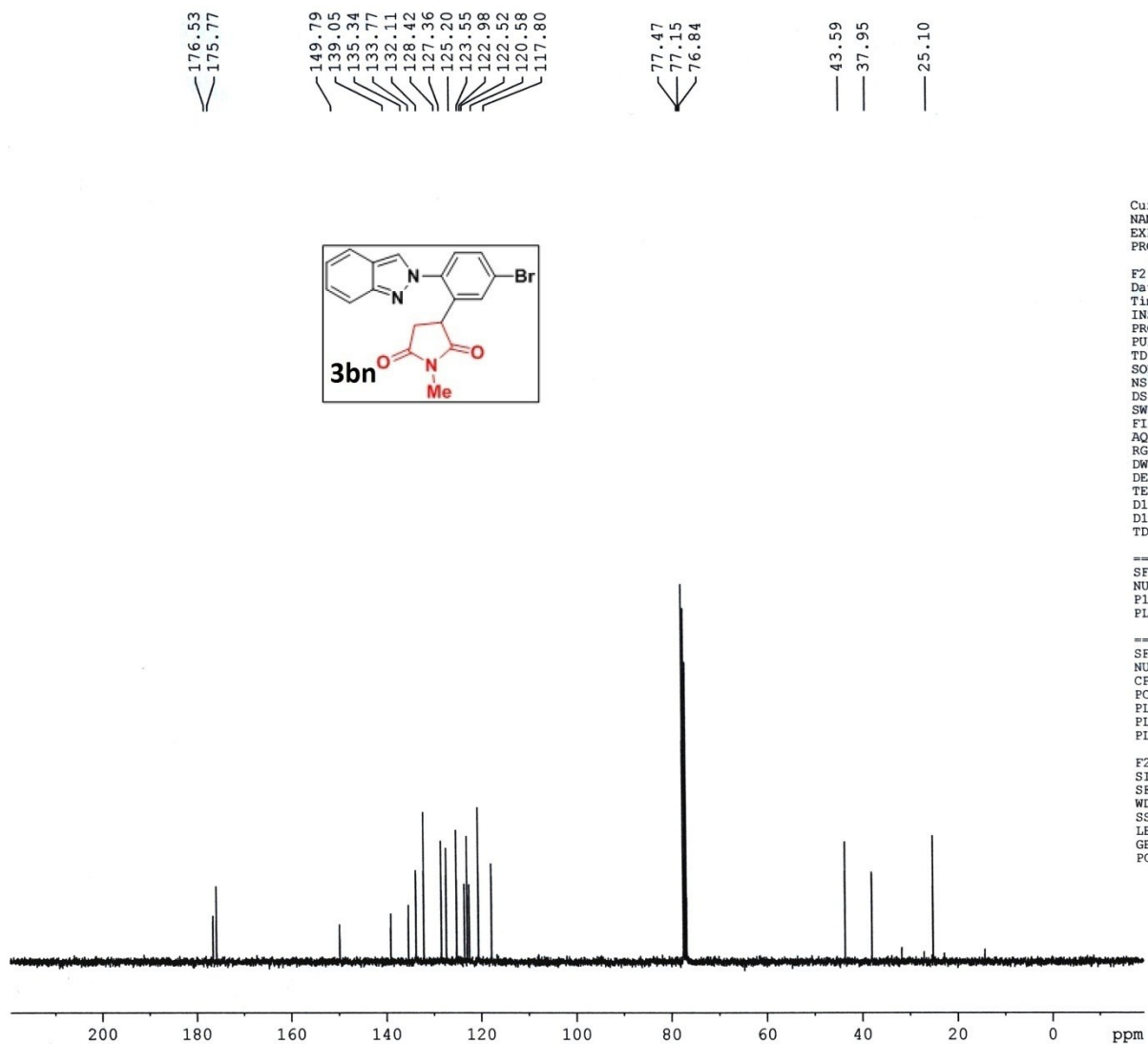
F2 - Acquisition Parameters
Date_ 20191113
Time 10.16
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl₃
NS 300
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 67.81
DW 20.800 usec
DE 6.50 usec
TE 296.7 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG2 waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177886 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40





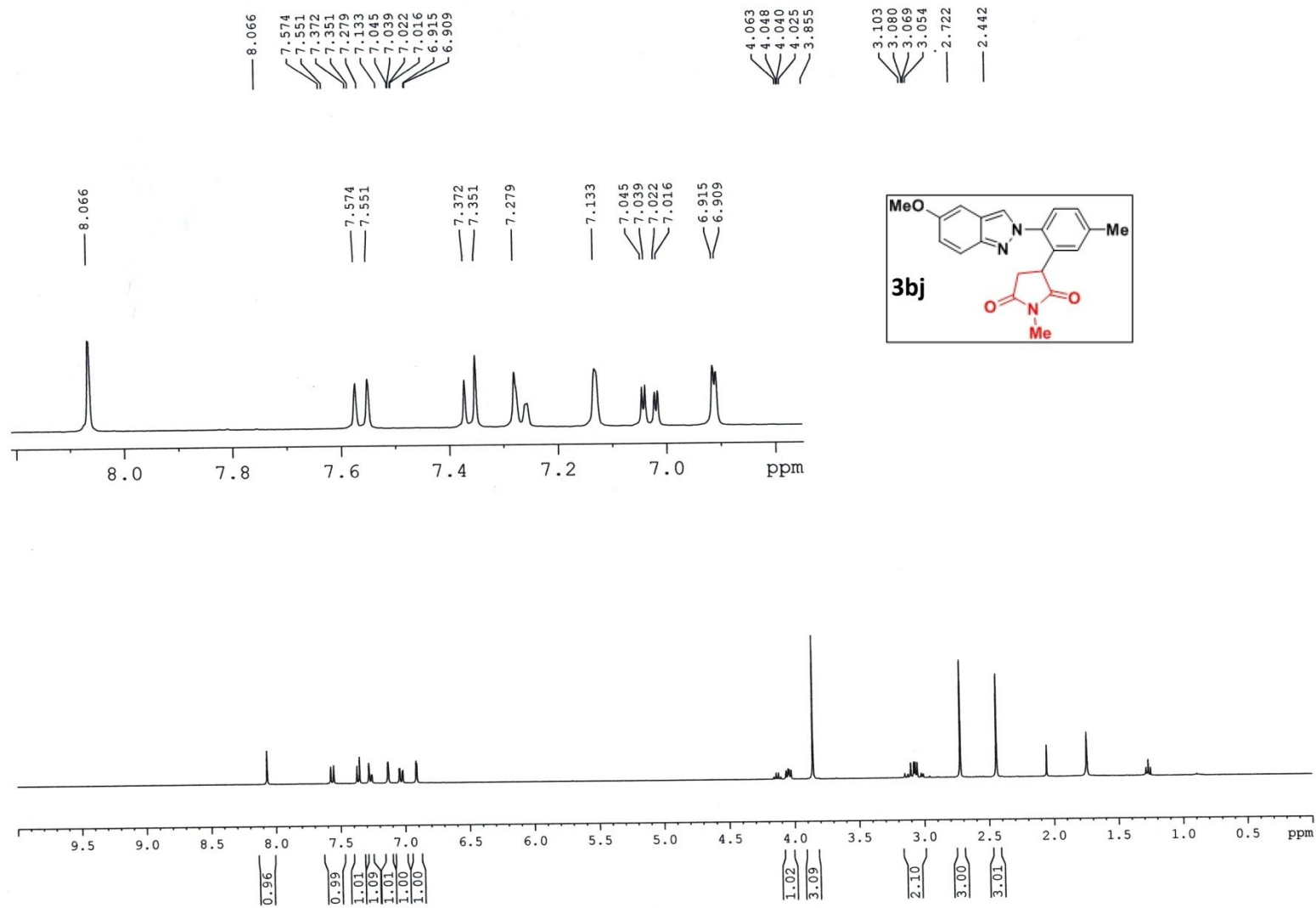
Current Data Parameters
 NAME Dr. A HAJRA-2019-13C
 EXPNO 422
 PROCNO 1

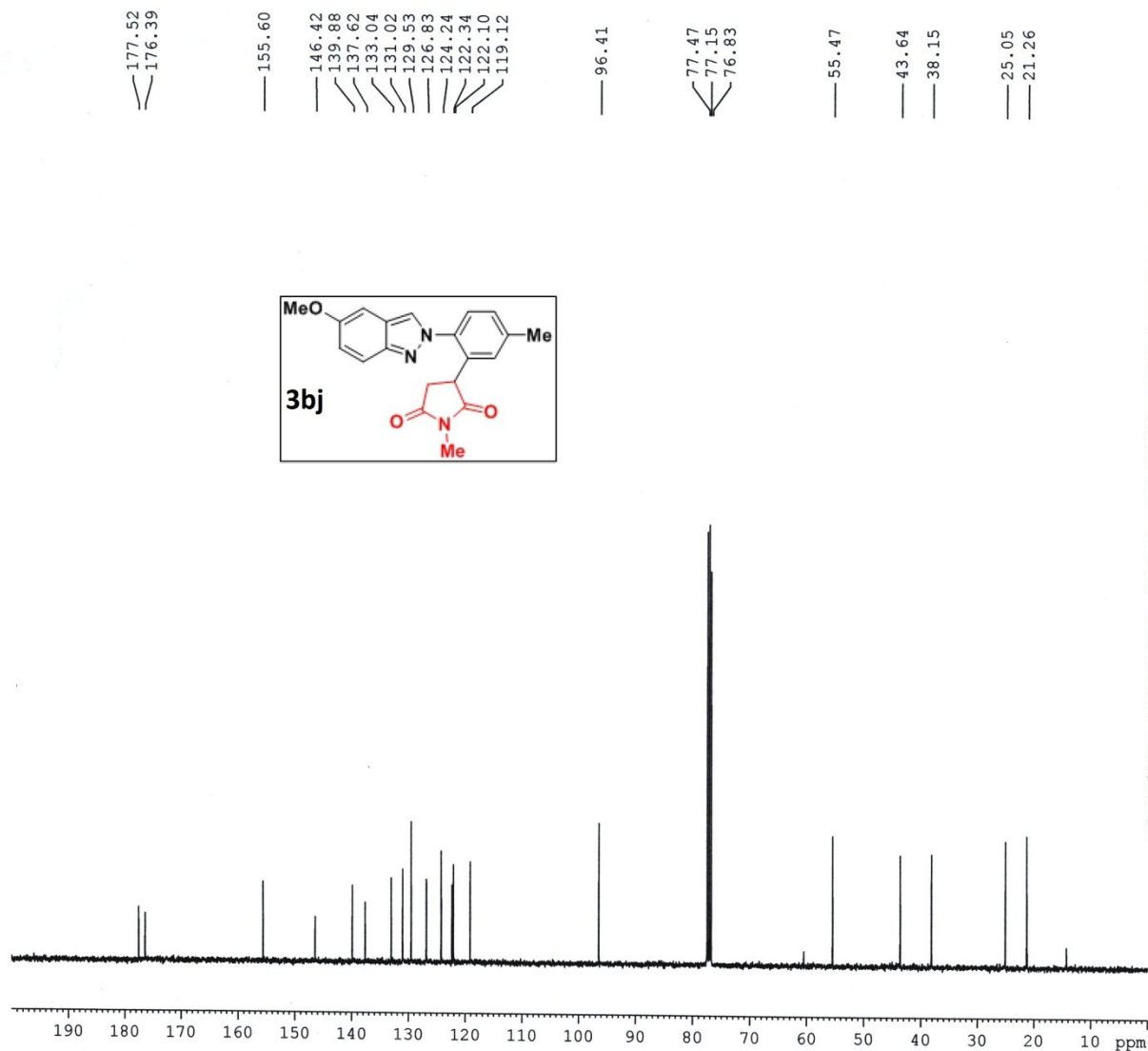
F2 - Acquisition Parameters
 Date_ 20191018
 Time 11.06
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDC13
 NS 240
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 300.9 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TD0 1

===== CHANNEL f1 =====
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

===== CHANNEL f2 =====
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG2 waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177858 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40





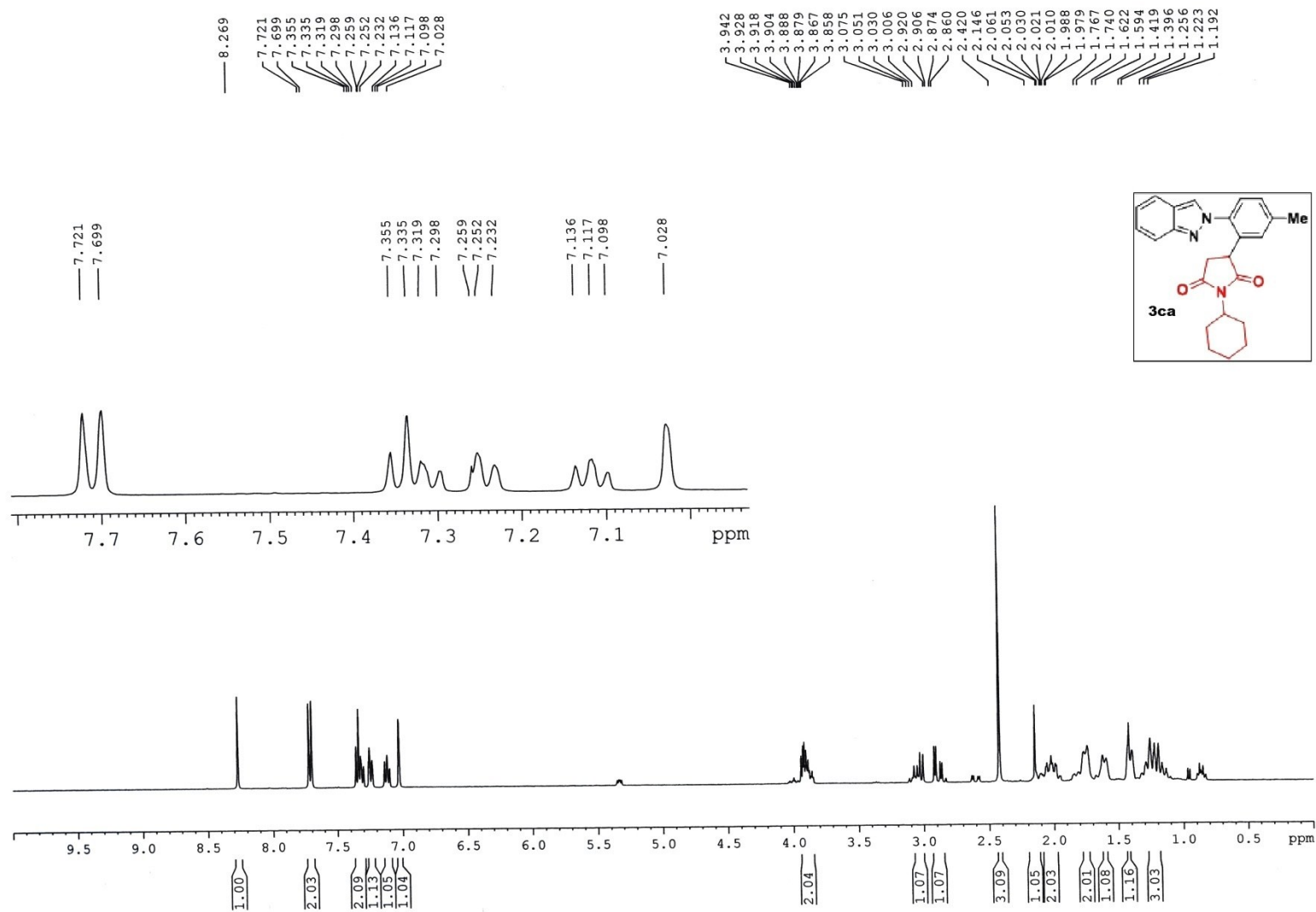
Current Data Parameters
 NAME Dr. A HAJRA-2019-13C
 EXPNO 473
 PROCNO 1

F2 - Acquisition Parameters
 Date_ 20191028
 Time 16.07
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDC13
 NS 512
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 296.8 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TD0 1

===== CHANNEL f1 =====
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

===== CHANNEL f2 =====
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG[2] waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177873 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40



177.65
176.29

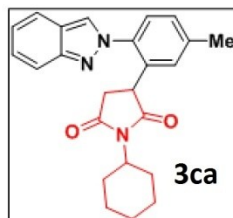
149.66
140.28
138.02
133.79
129.34
129.08
127.18
126.83
125.55
122.46
122.26
120.55
117.88

77.46
77.15
76.83

52.07

42.02
38.04

28.71
25.87
25.04
21.38



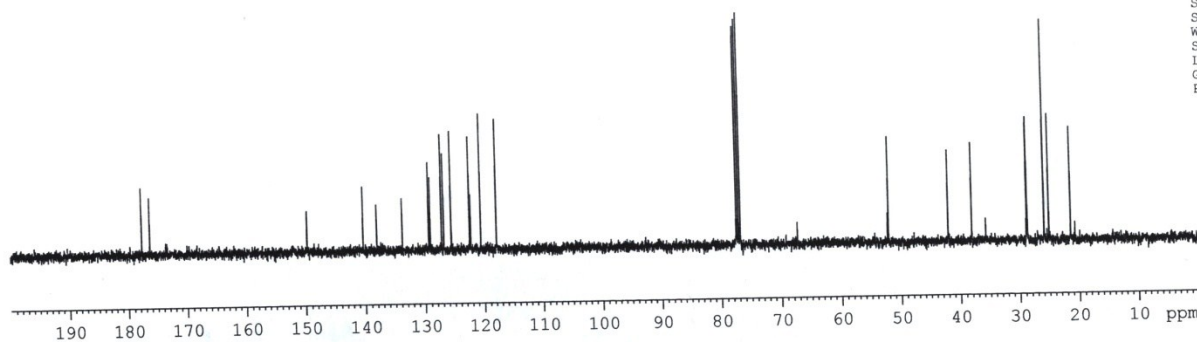
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 527
PROCNO 1

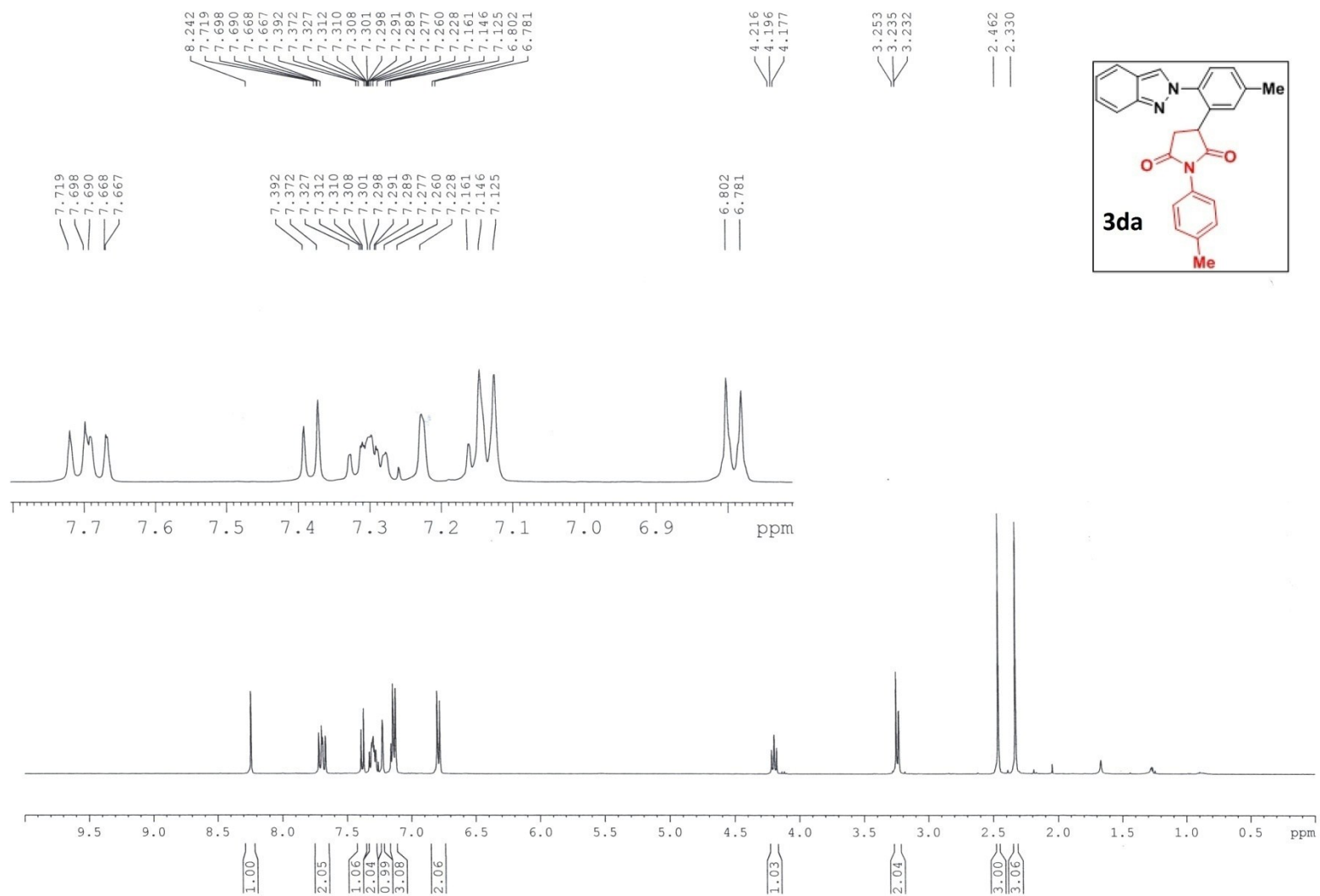
F2 - Acquisition Parameters
Date_ 20191115
Time 12.18
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 50
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 62.69
DW 20.800 usec
DE 6.50 usec
TE 297.1 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

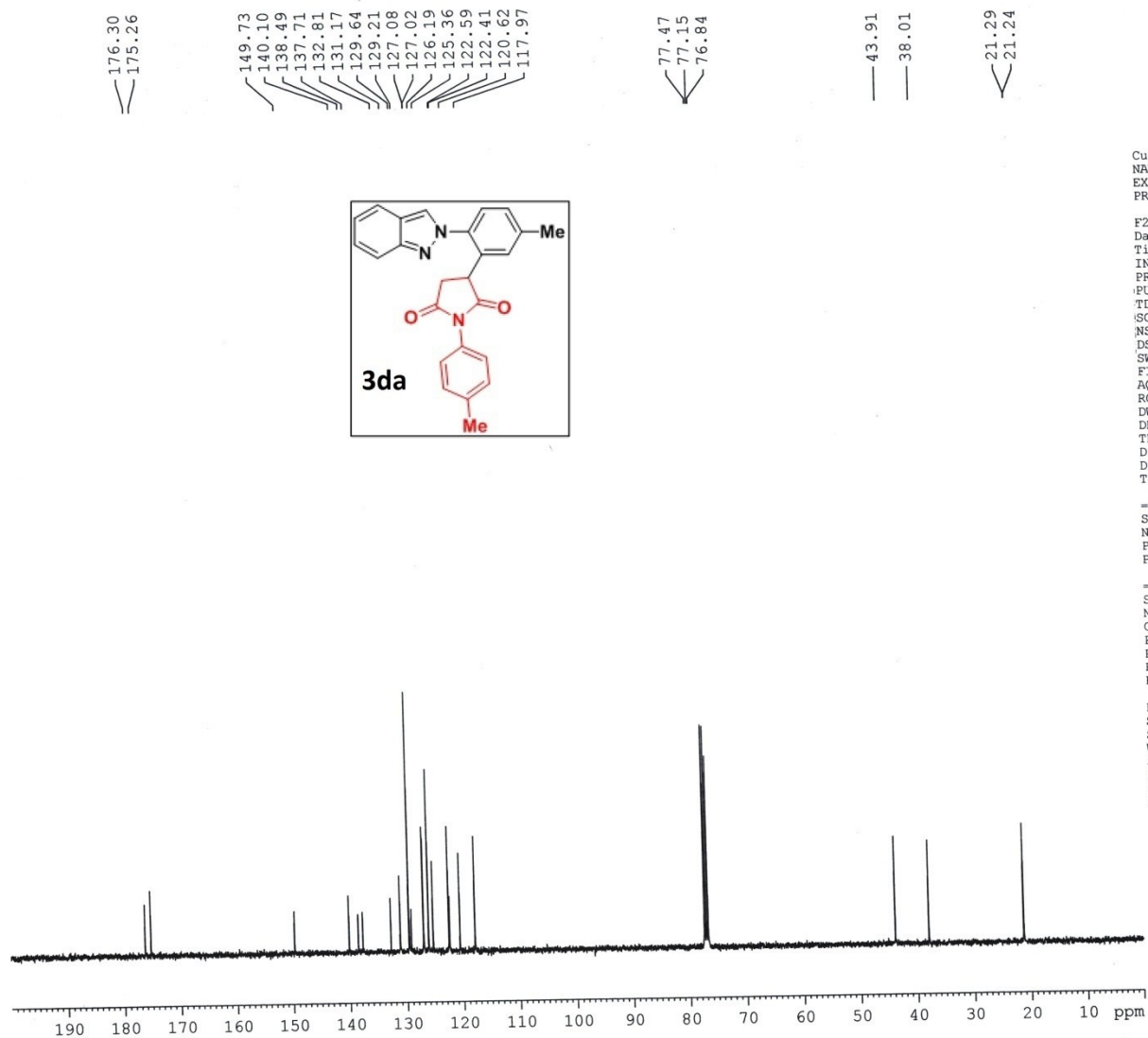
===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
PCPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177901 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40







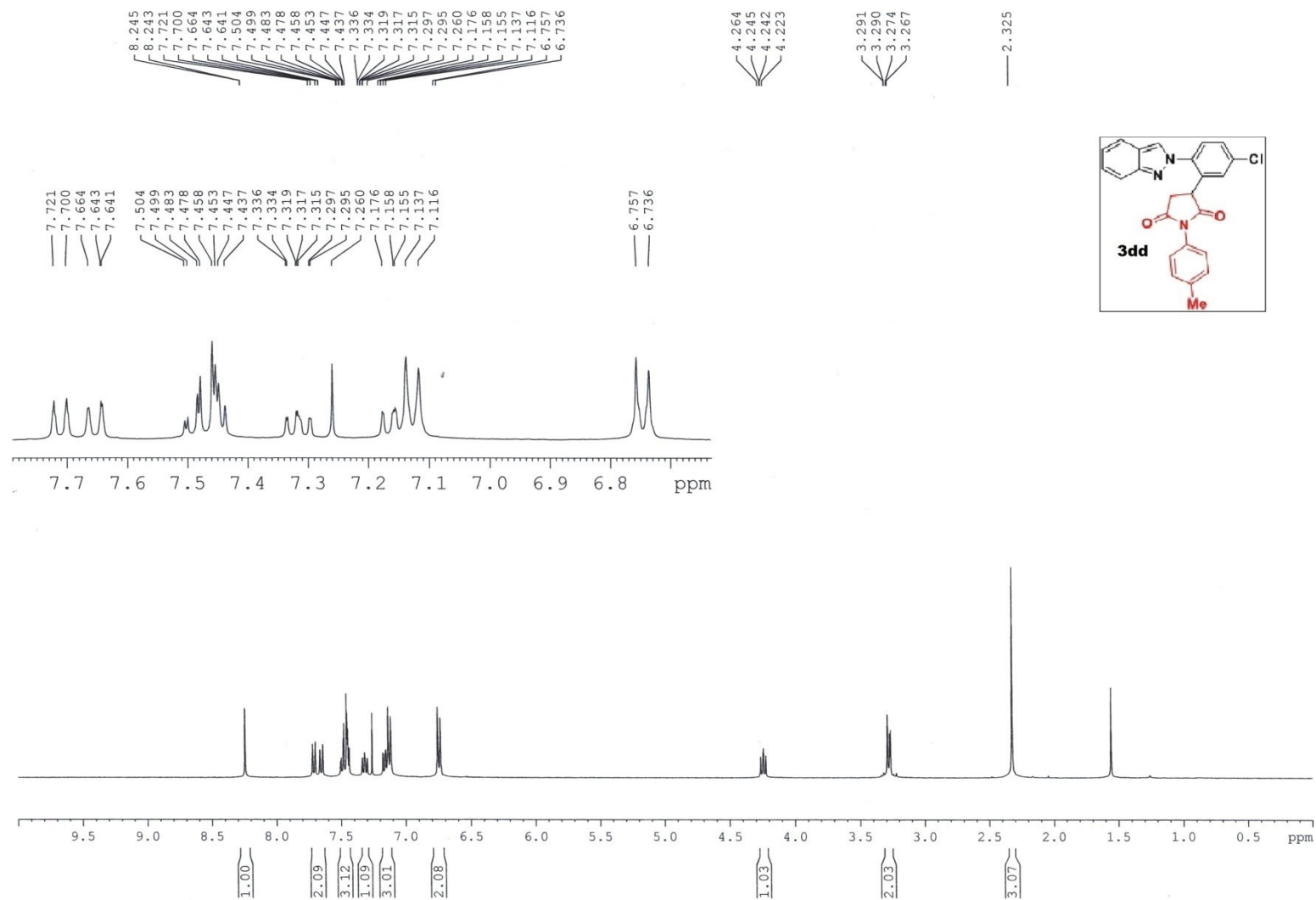
Current Parameters
 NAME 2019-13C
 EXPNO 444
 PROCNO 1

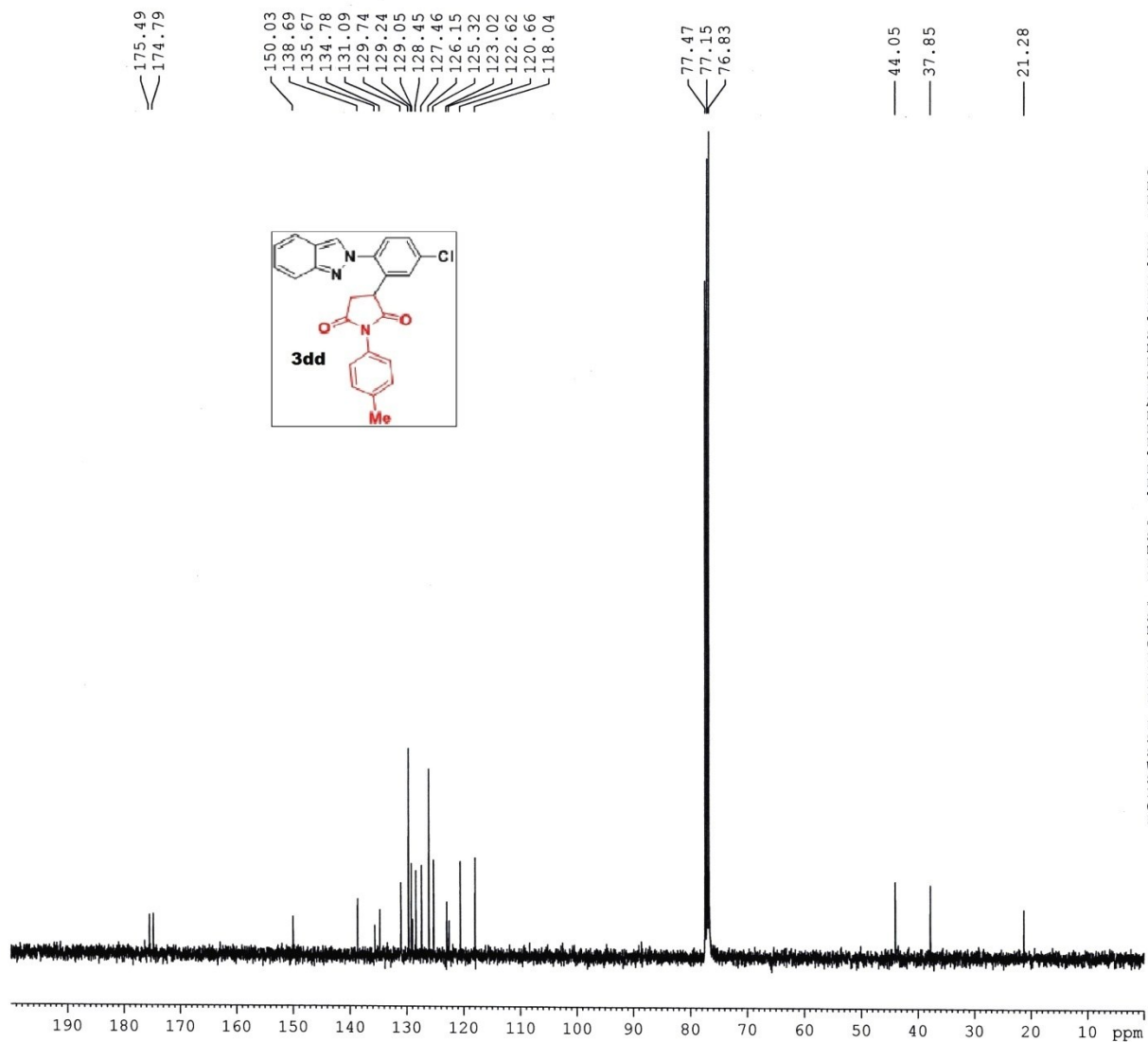
F2 - Acquisition Parameters
 Date 20191022
 Time 15.27
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDCl3
 NS 220
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 299.8 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TD0 1

===== CHANNEL f1 =====
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

===== CHANNEL f2 =====
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG2 waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177903 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40





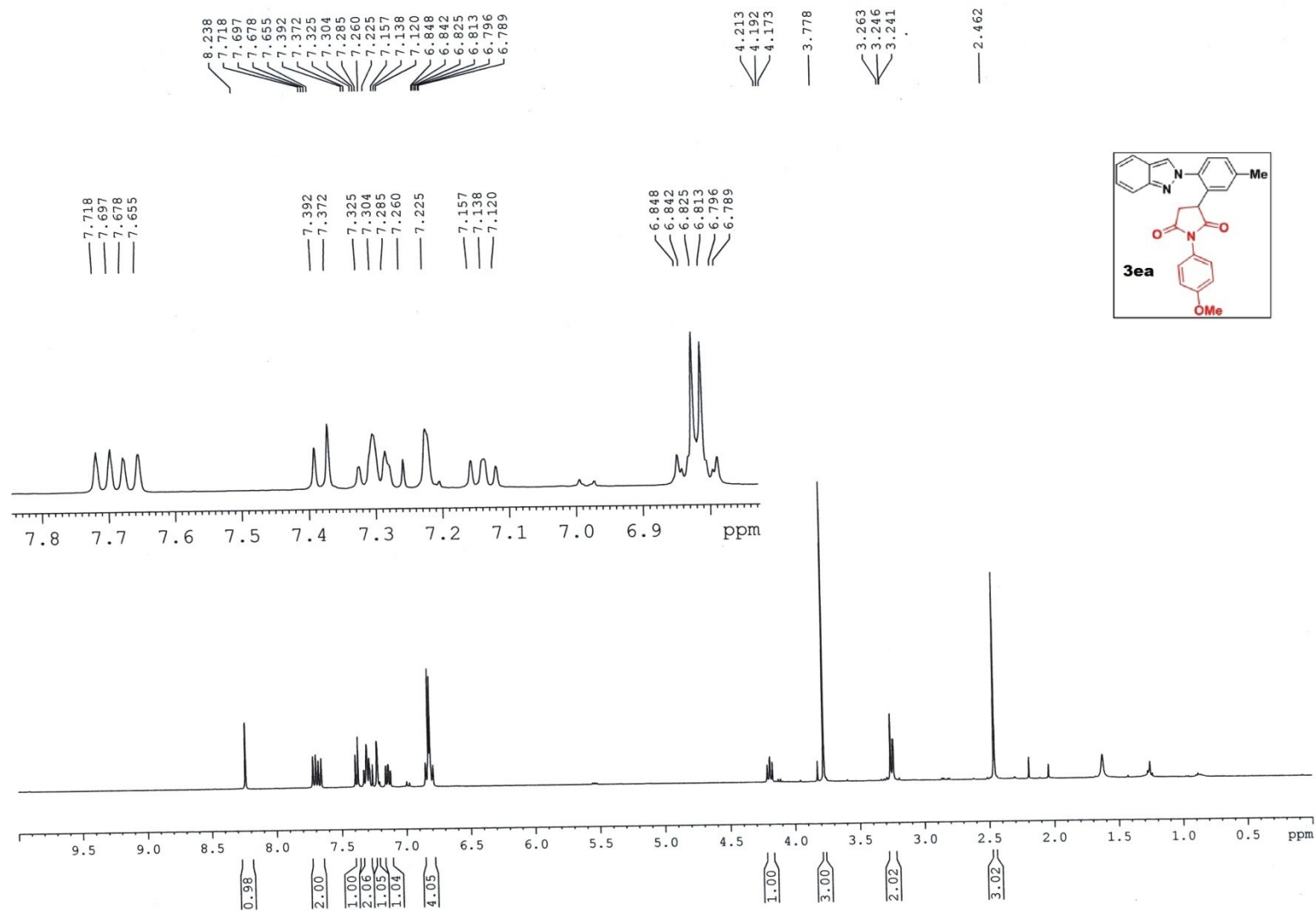
Current Data Parameters
 NAME Dr. A HAJRA-2019-13C
 EXPNO 434
 PROCNO 1

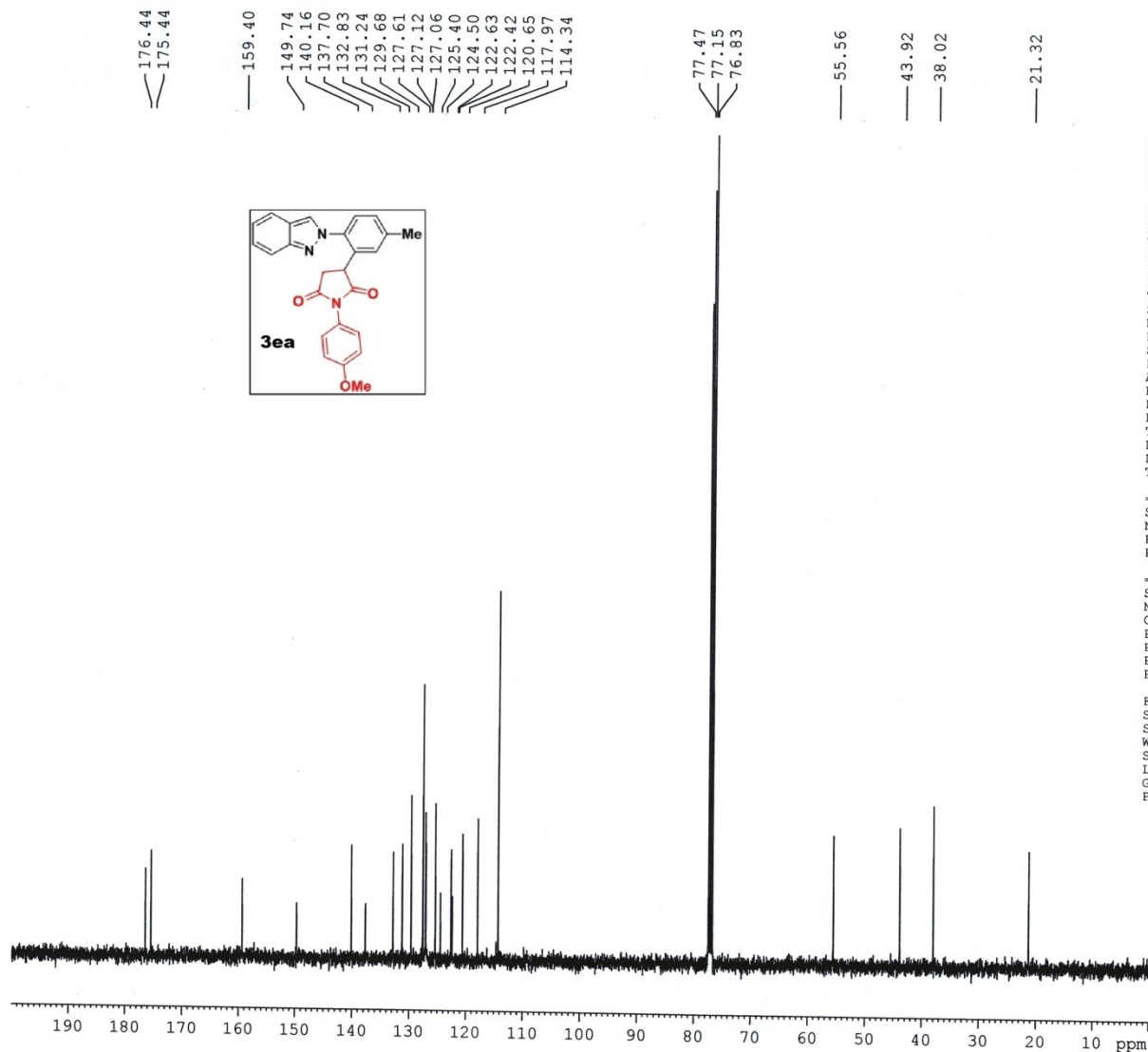
F2 - Acquisition Parameters
 Date_ 20191021
 Time_ 15.49
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDCl3
 NS 512
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 299.0 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TD0 1

===== CHANNEL f1 =====
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

===== CHANNEL f2 =====
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG[2] waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177844 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40





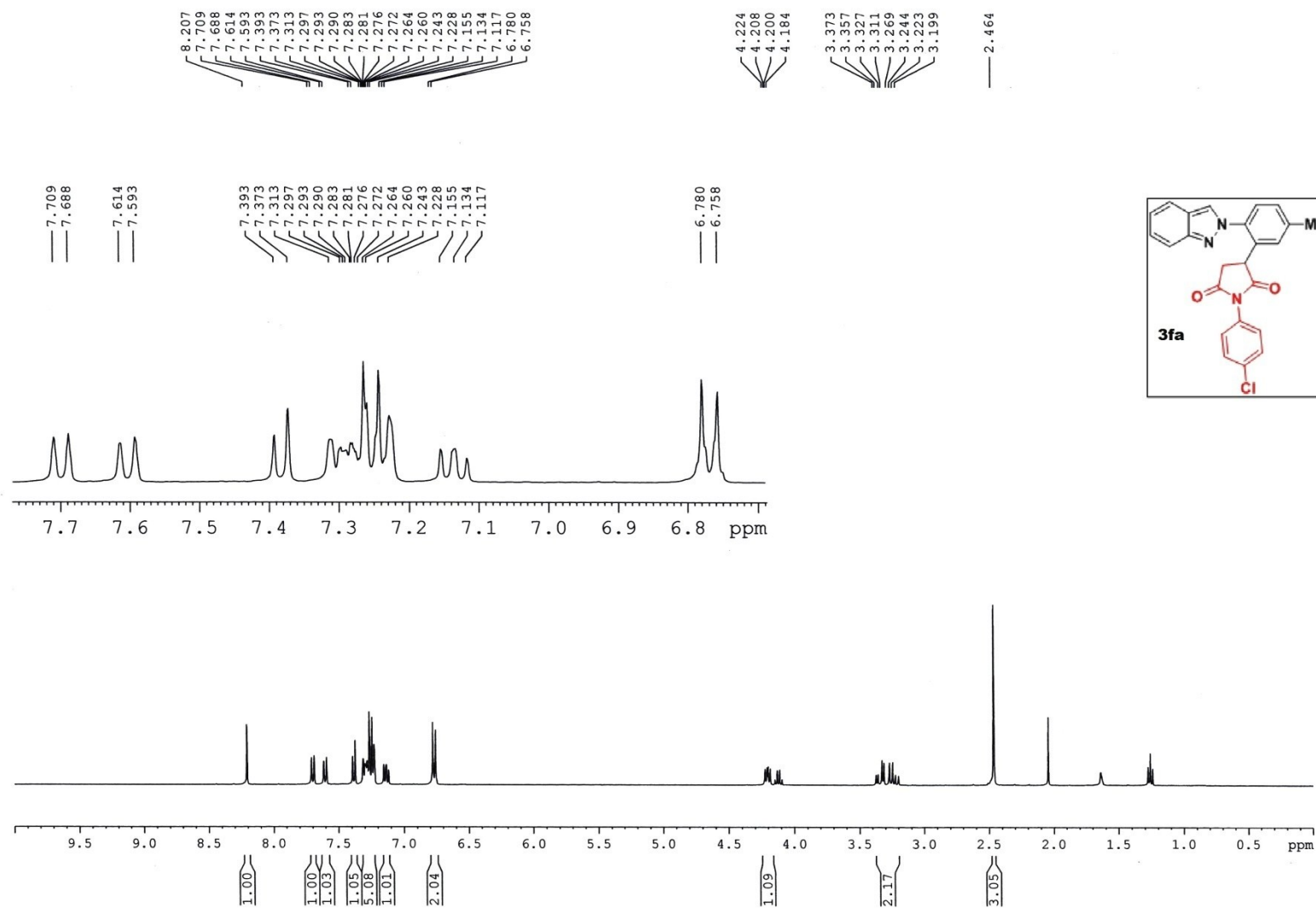
Current Data Parameters
 NAME Dr. A HAJRA-2019-13C
 EXPNO 458
 PROCNO 1

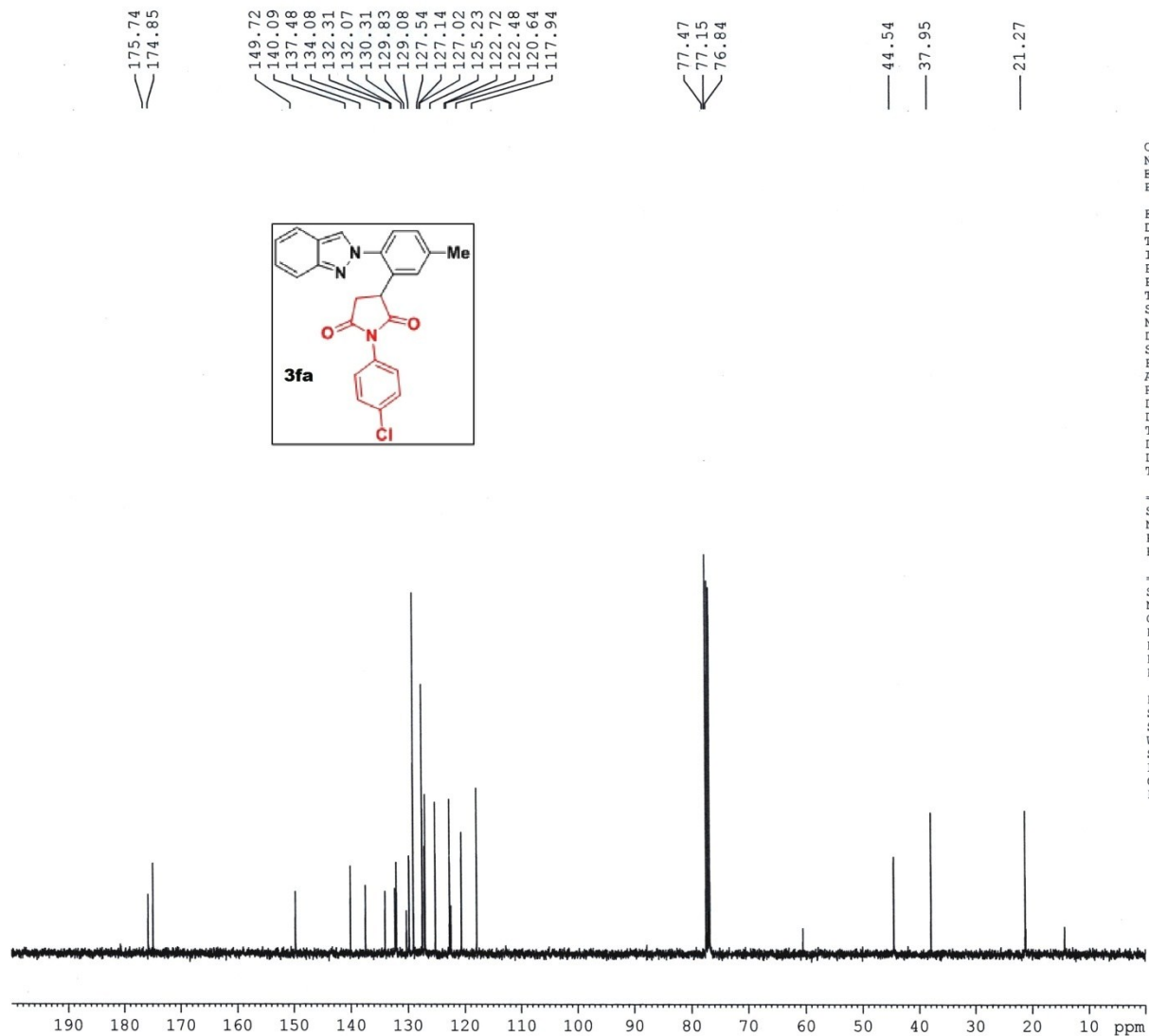
F2 - Acquisition Parameters
 Date_ 20191025
 Time 17.48
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDC13
 NS 250
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 297.8 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TDO 1

----- CHANNEL f1 -----
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

----- CHANNEL f2 -----
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG2 waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177871 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40





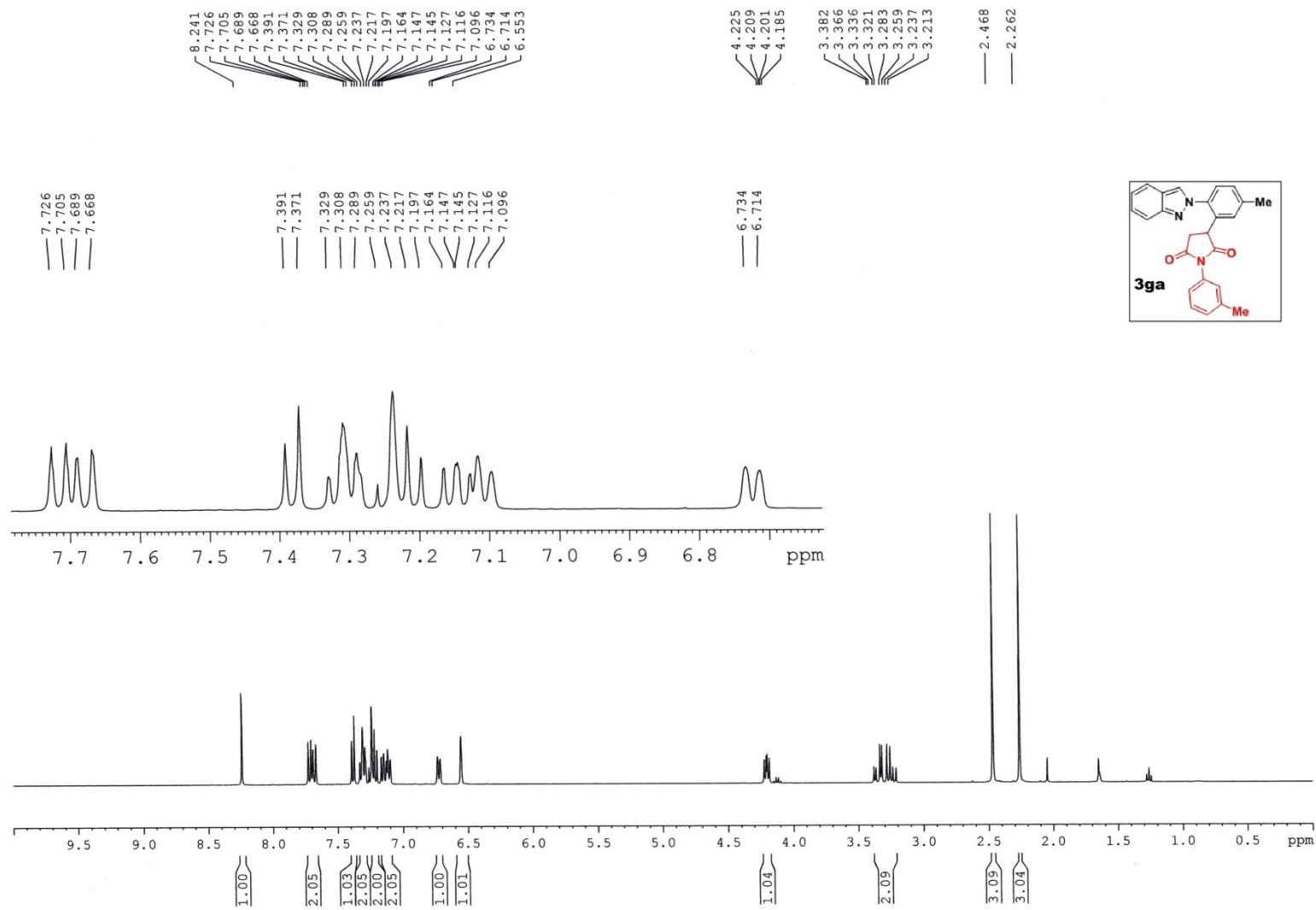
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 509
PROCNO 1

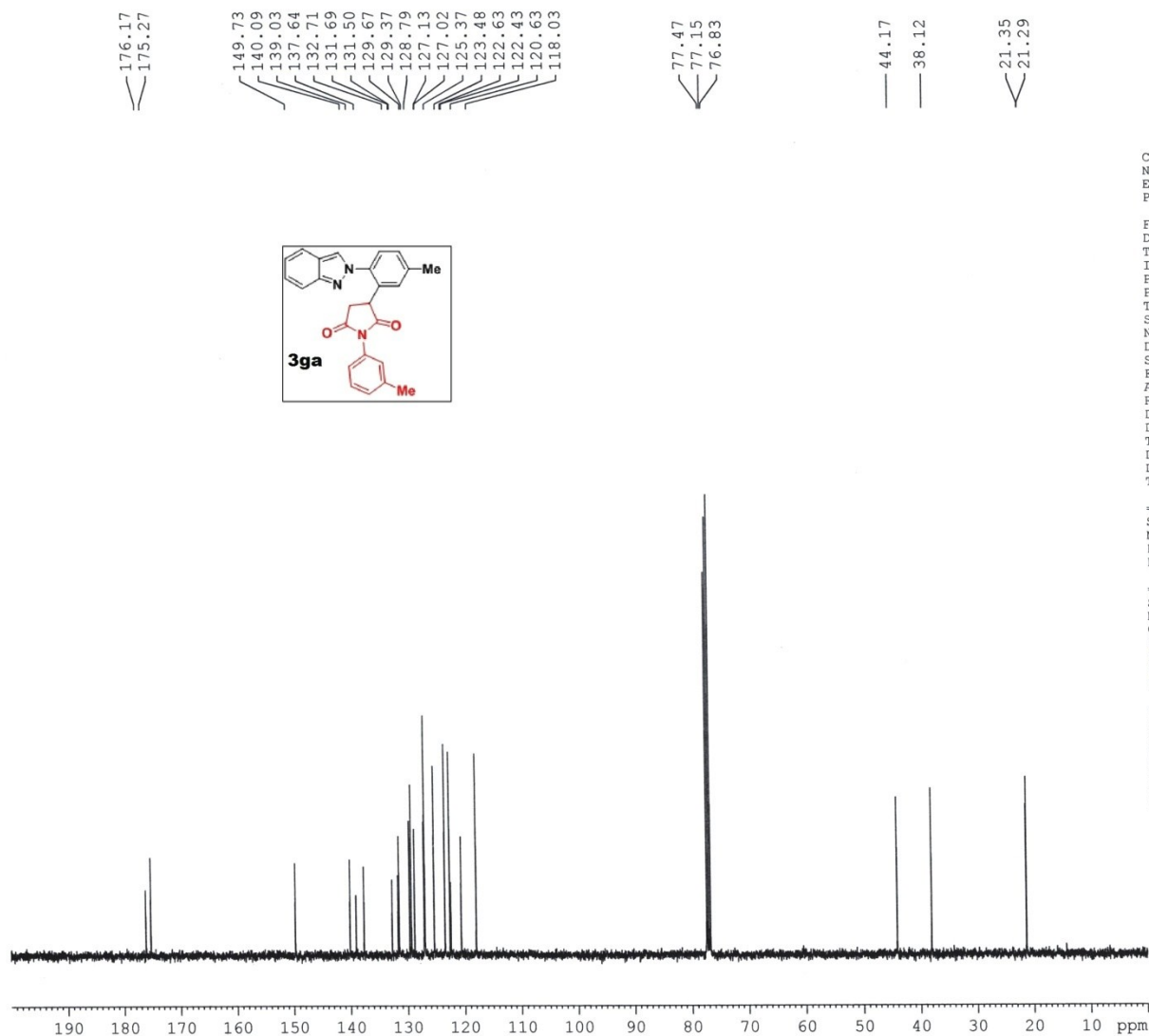
F2 - Acquisition Parameters
Date_ 20191107
Time 10.59
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 205
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 77.59
DW 20.800 usec
DE 6.50 usec
TE 298.7 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177875 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40





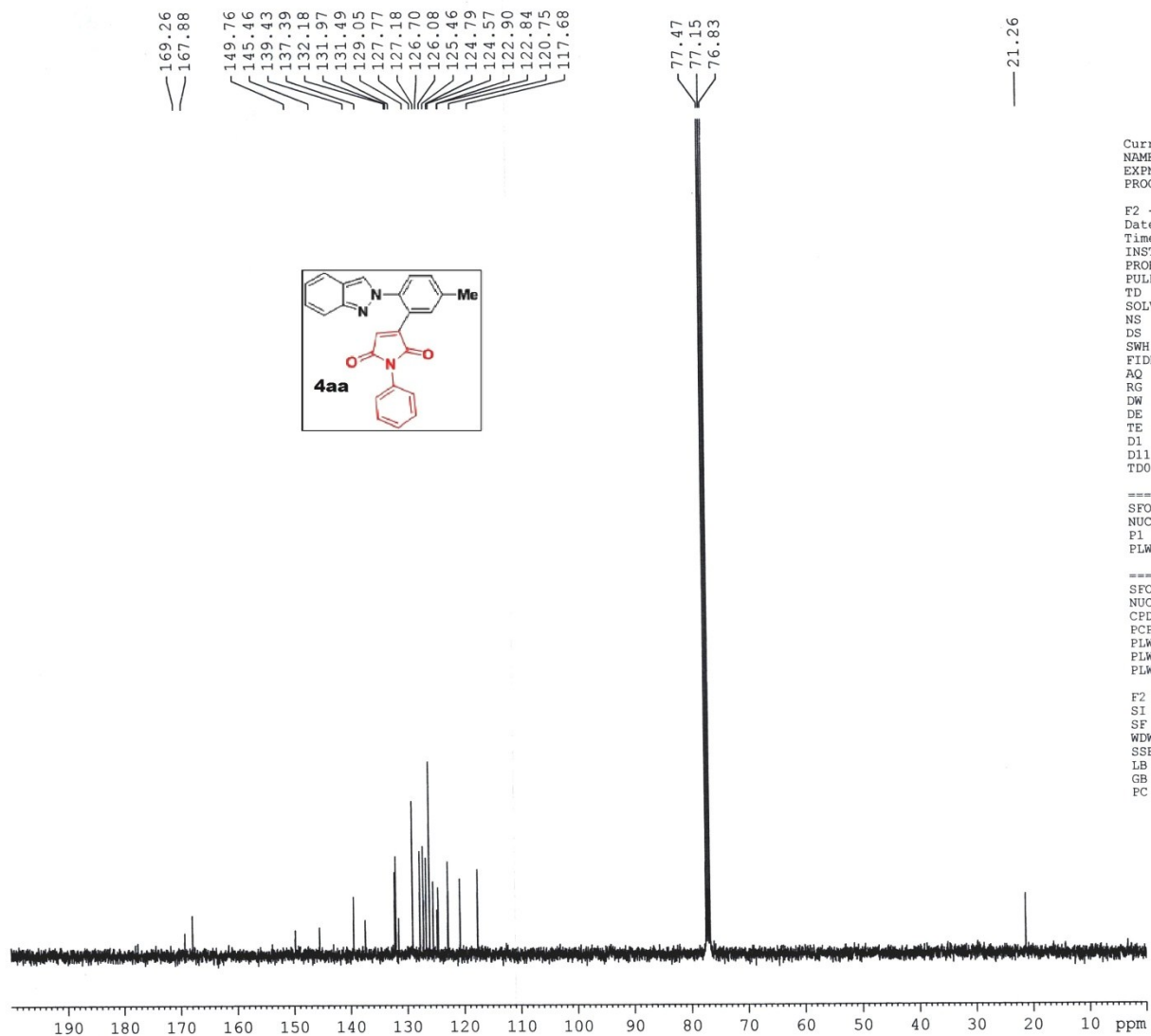
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 499
PROCNO 1

F2 - Acquisition Parameters
Date_ 20191104
Time 20.00
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDC13
NS 220
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 67.81
DW 20.800 usec
DE 6.50 usec
TE 296.8 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177902 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40



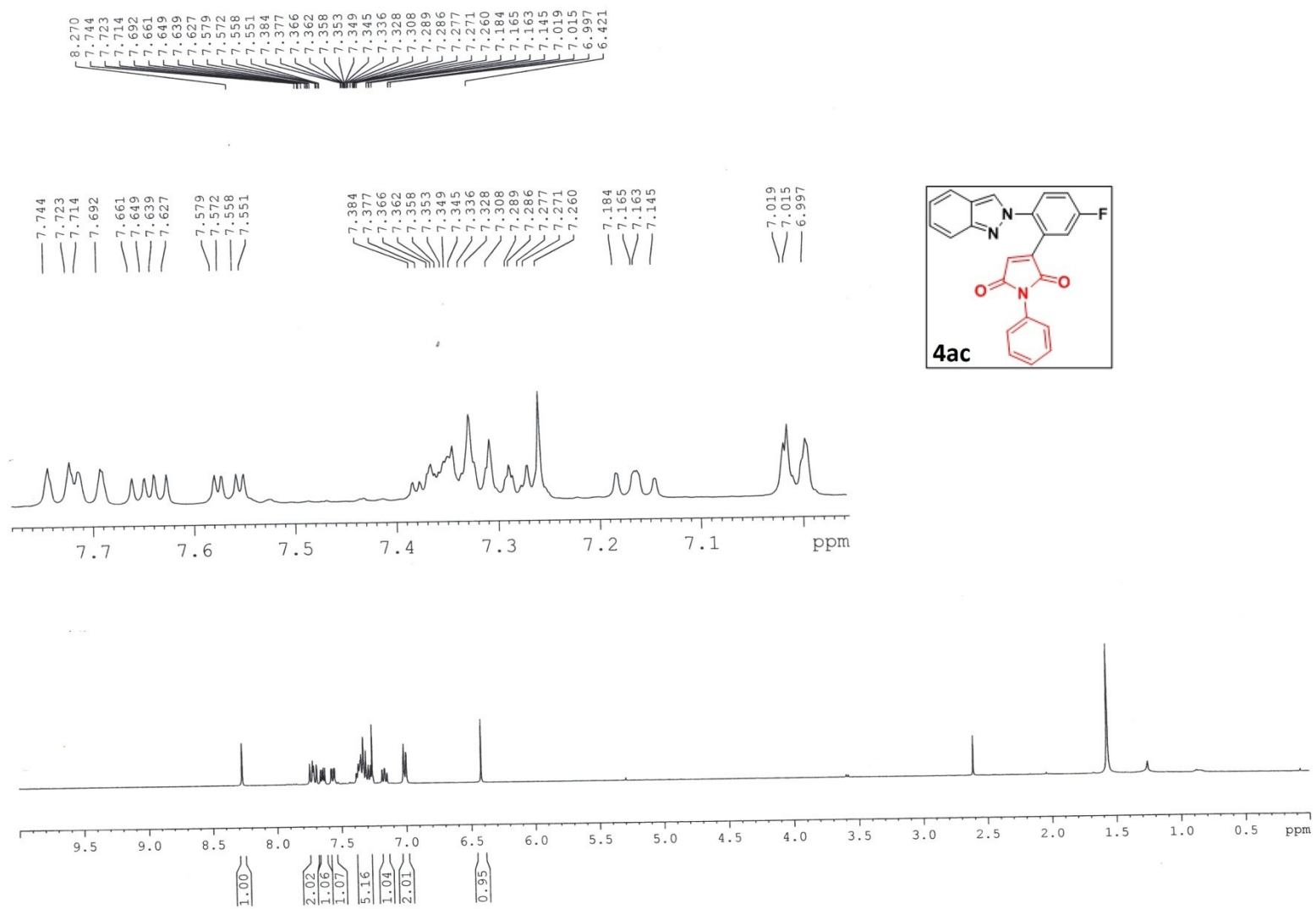
Current Data Parameters
 NAME Dr. A HAJRA-2019-13C
 EXPNO 543
 PROCNO 1

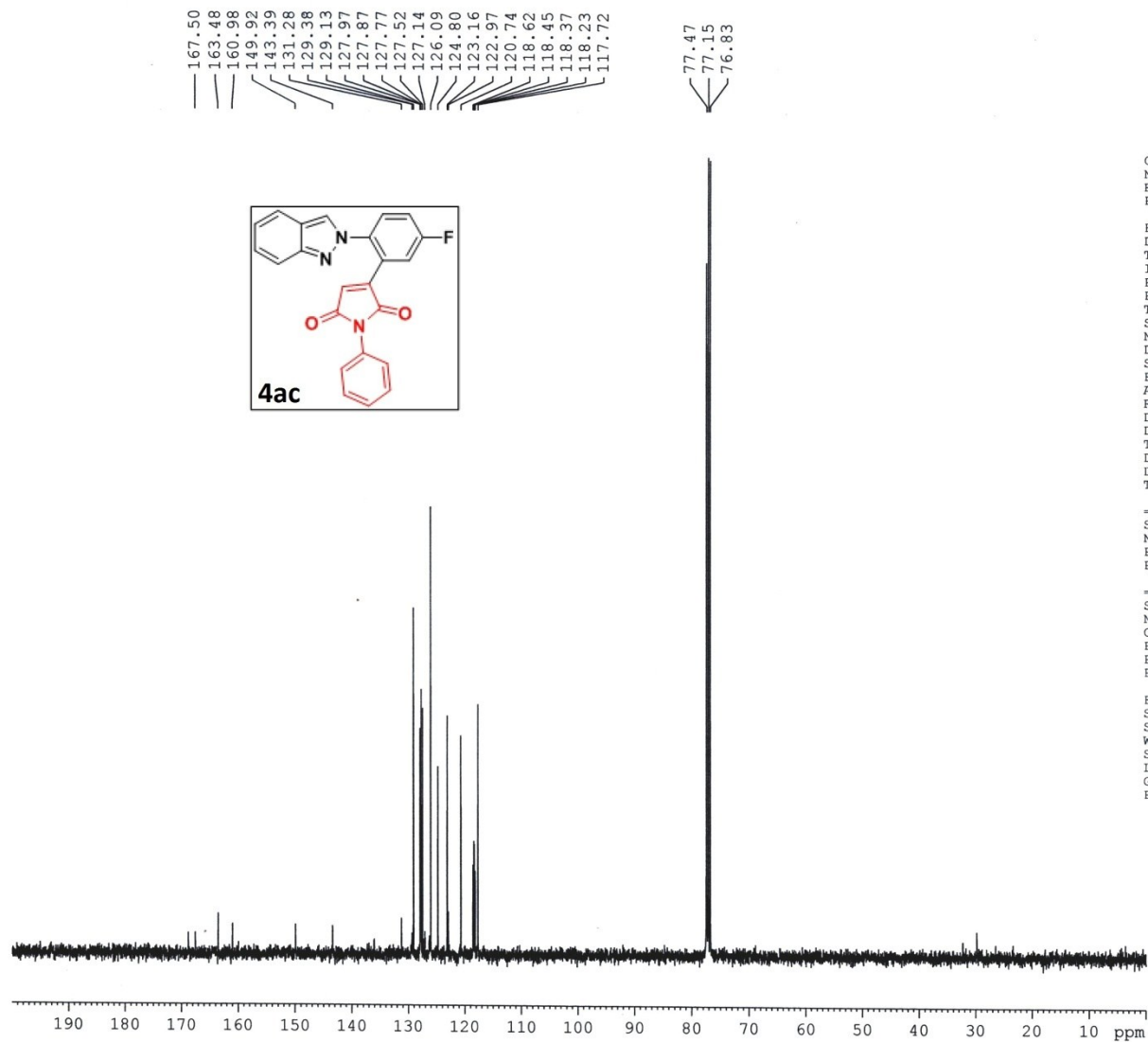
F2 - Acquisition Parameters
 Date_ 20191124
 Time 18.30
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDCl3
 NS 512
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 295.7 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TDO 1

===== CHANNEL f1 =====
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

===== CHANNEL f2 =====
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG[2] waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177858 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40





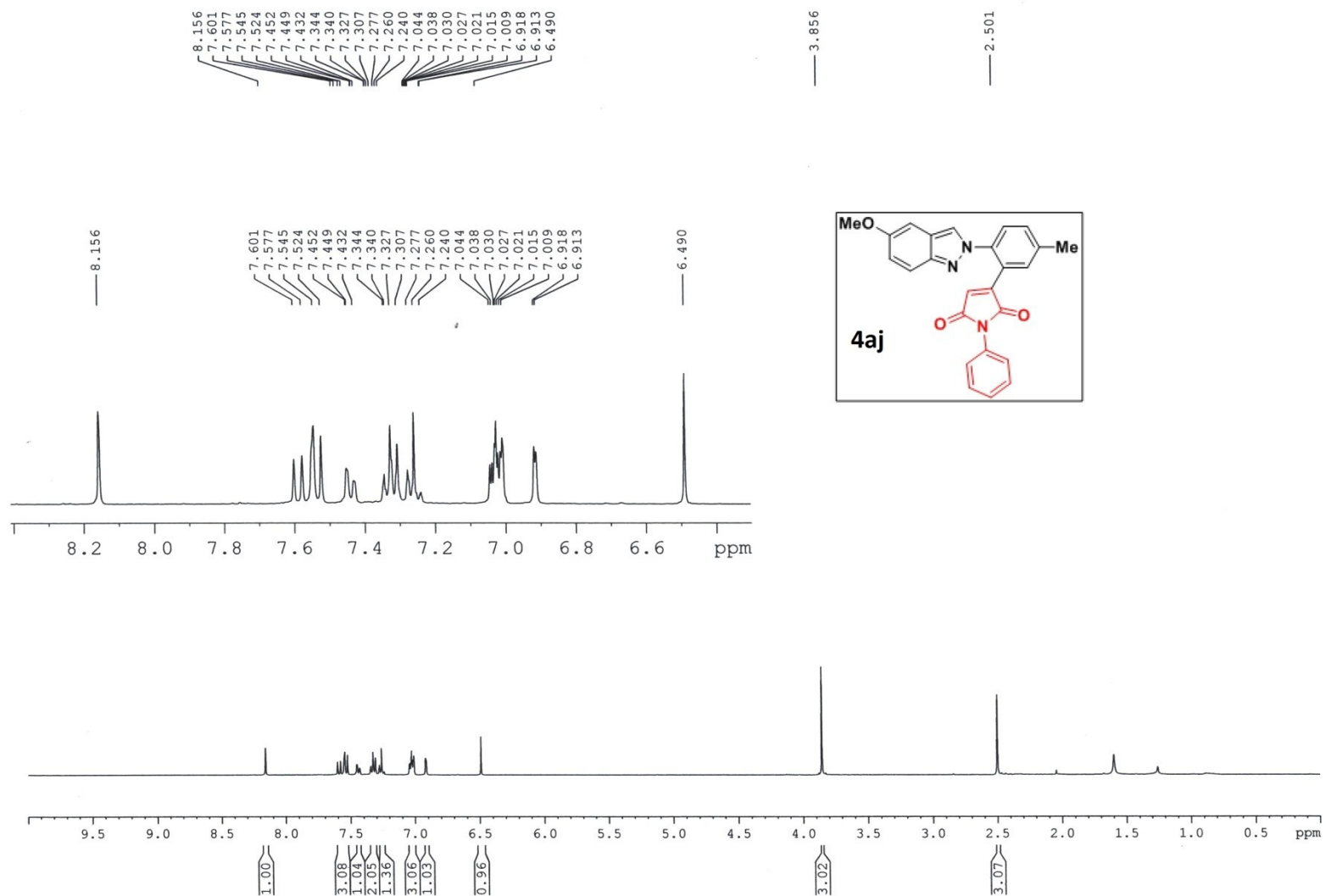
Current Data Parameters
 NAME Dr. A HAJRA-2019-13C
 EXPNO 470
 PROCNO 1

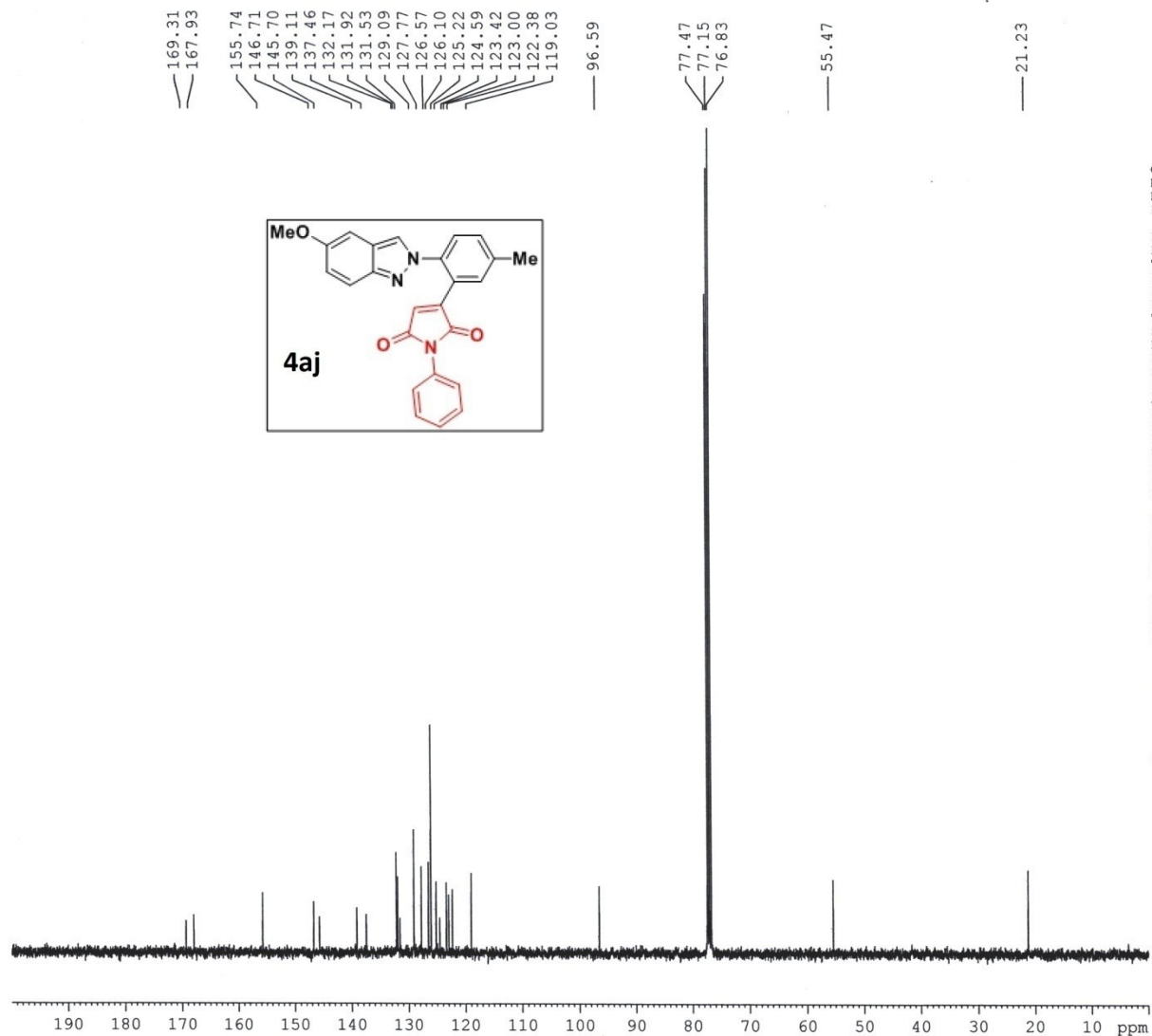
F2 - Acquisition Parameters
 Date_ 20191027
 Time 18.37
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgdc
 TD 32768
 SOLVENT CDCl3
 NS 2048
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 297.1 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TD0 1

===== CHANNEL f1 =====
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

===== CHANNEL f2 =====
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG2 waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177844 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.00





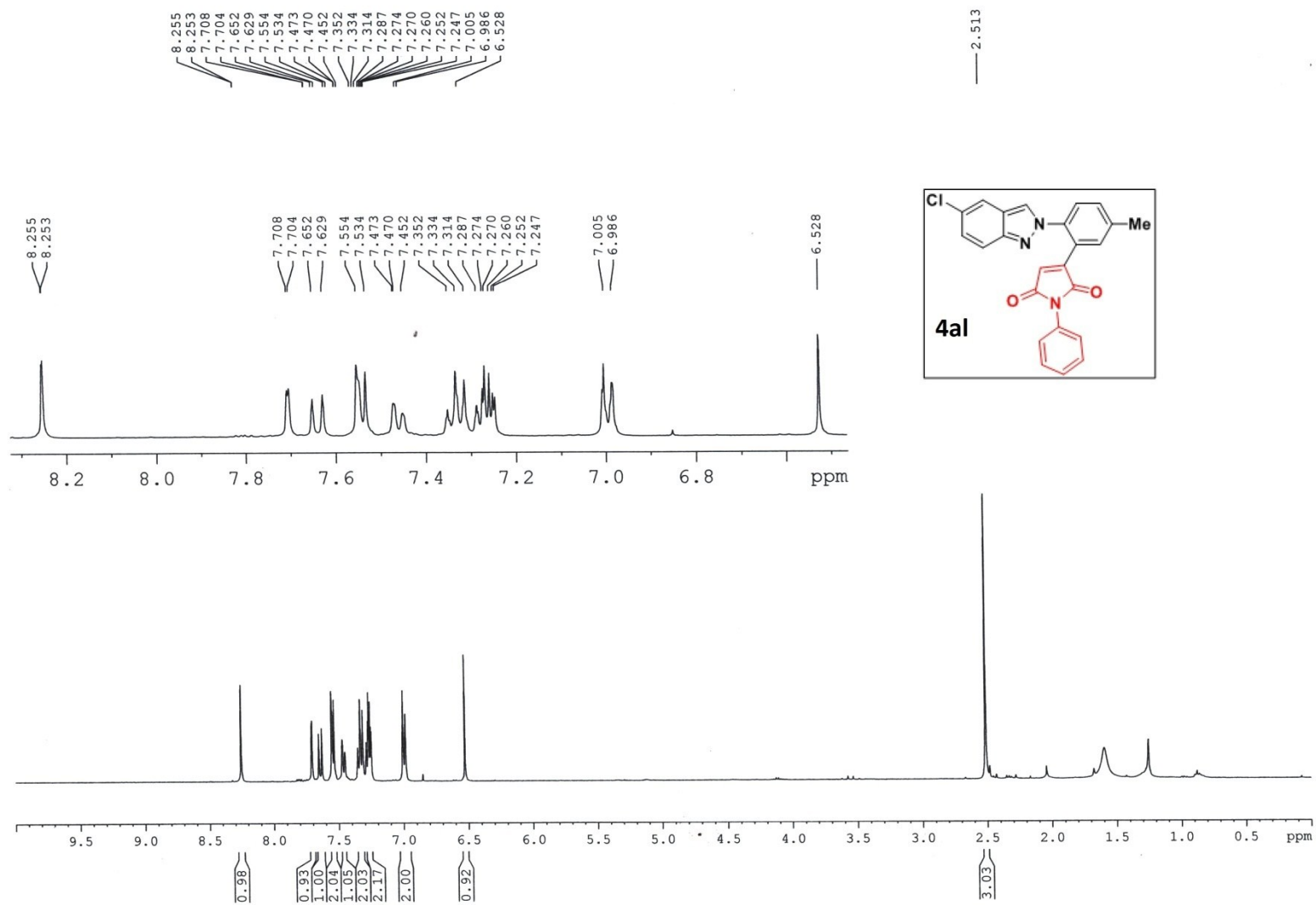
Current Data Parameters
 NAME Dr. A HAJRA-2019-13C
 EXPNO 538
 PROCNO 1

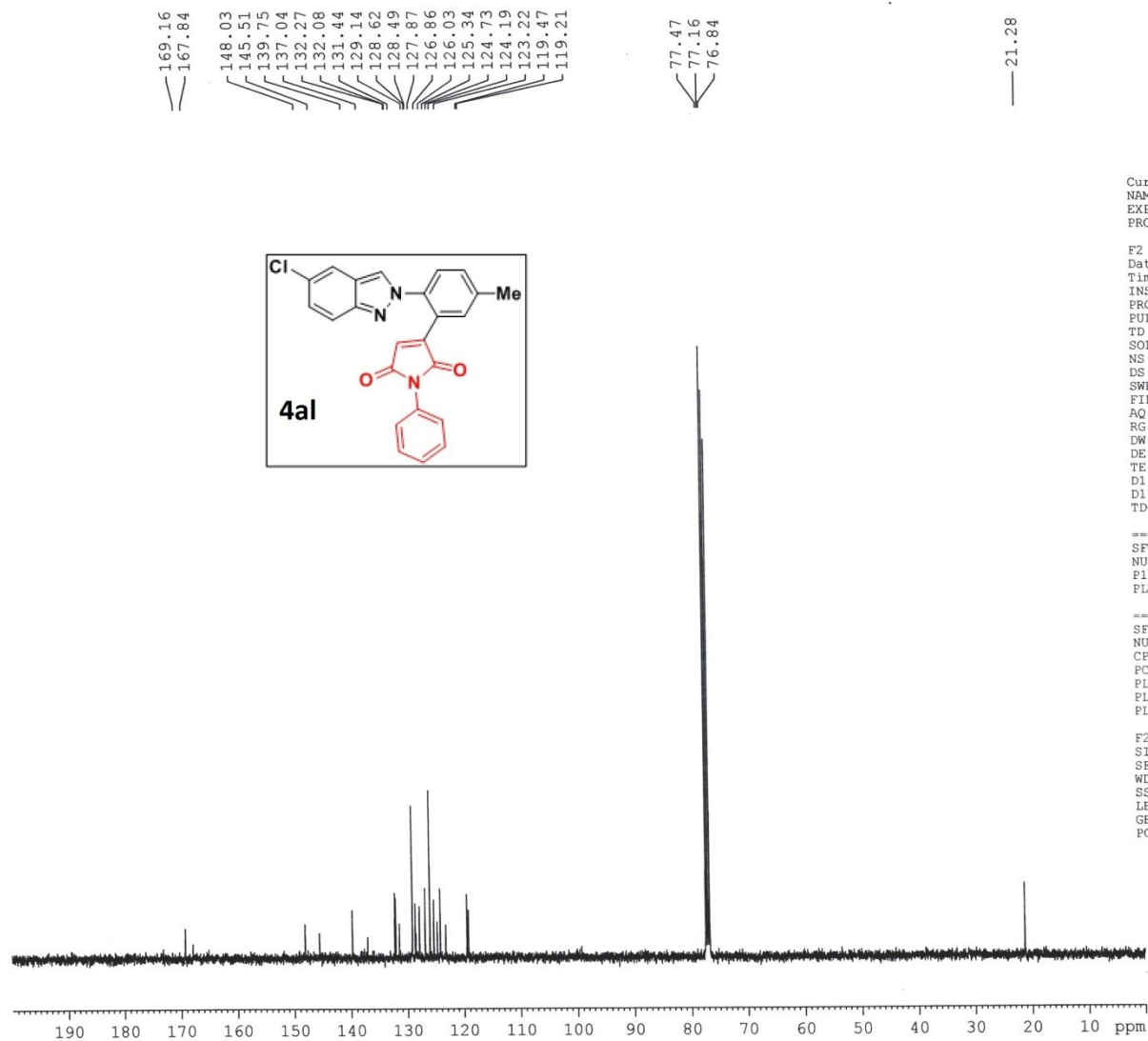
F2 - Acquisition Parameters
 Date_ 20191121
 Time 19.31
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDCl3
 NS 512
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 296.8 K
 D1 2.0000000 sec
 D11 0.03000000 sec
 TDO 1

===== CHANNEL f1 =====
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

===== CHANNEL f2 =====
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG[2] waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177857 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40





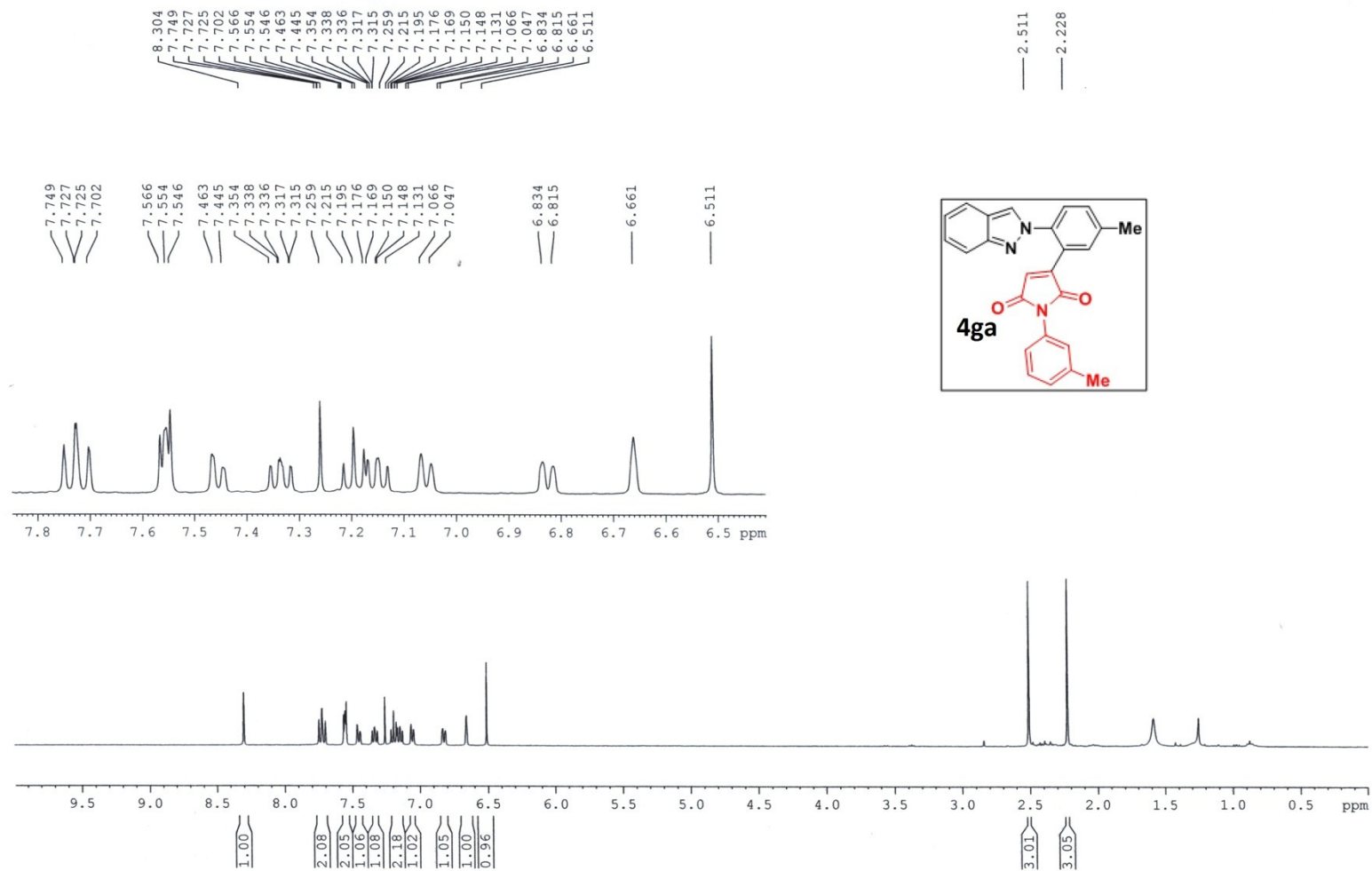
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 420
PROCNO 1

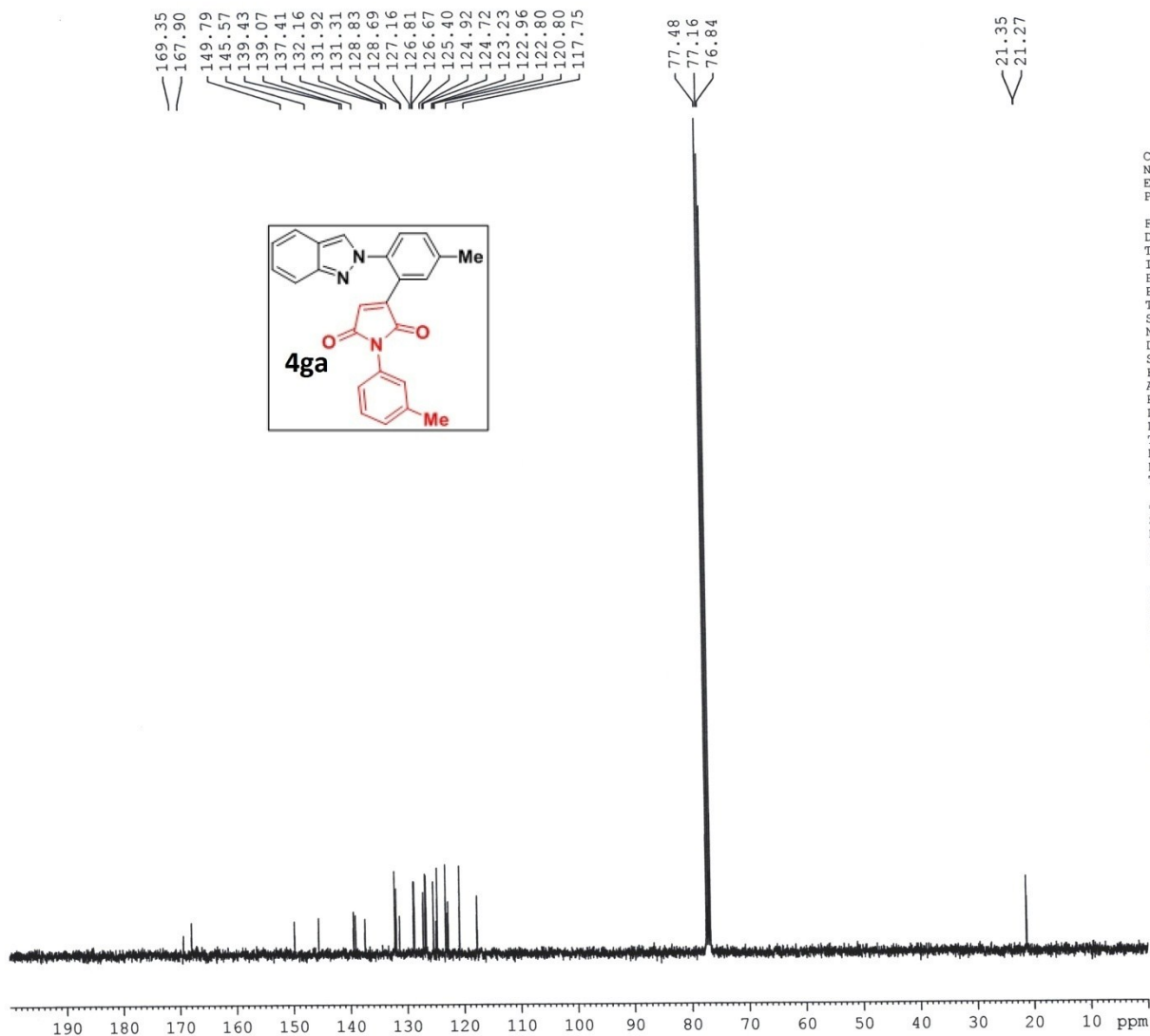
F2 - Acquisition Parameters
Date_ 20191017
Time_ 21.29
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 512
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 186.42
DW 20.800 usec
DE 6.50 usec
TE 297.9 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177843 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40





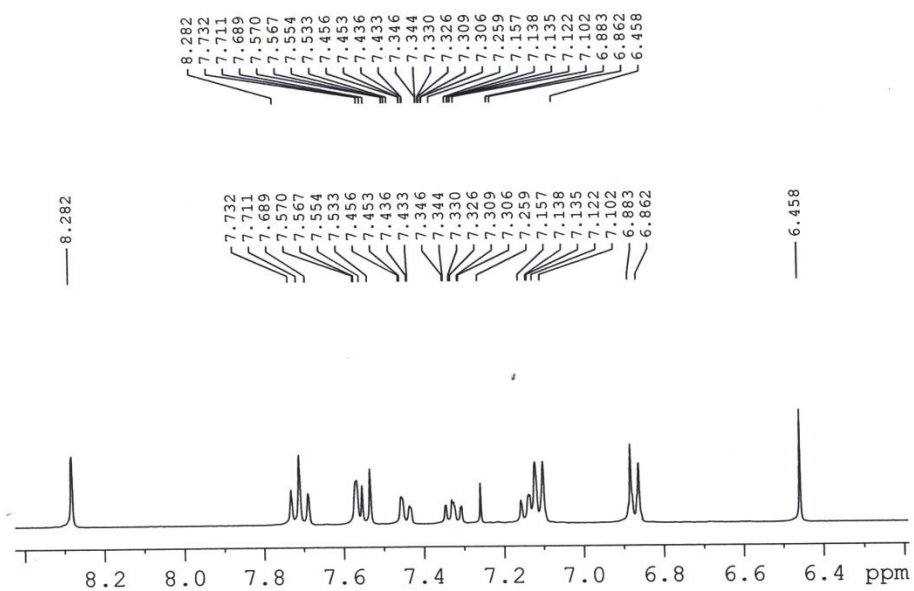
Current Data Parameters
 NAME Dr. A HAJRA-2019-13C
 EXPNO 501
 PROCNO 1

F2 - Acquisition Parameters
 Date_ 20191105
 Time 13.26
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDC13
 NS 640
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 297.1 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TD0 1

===== CHANNEL f1 =====
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

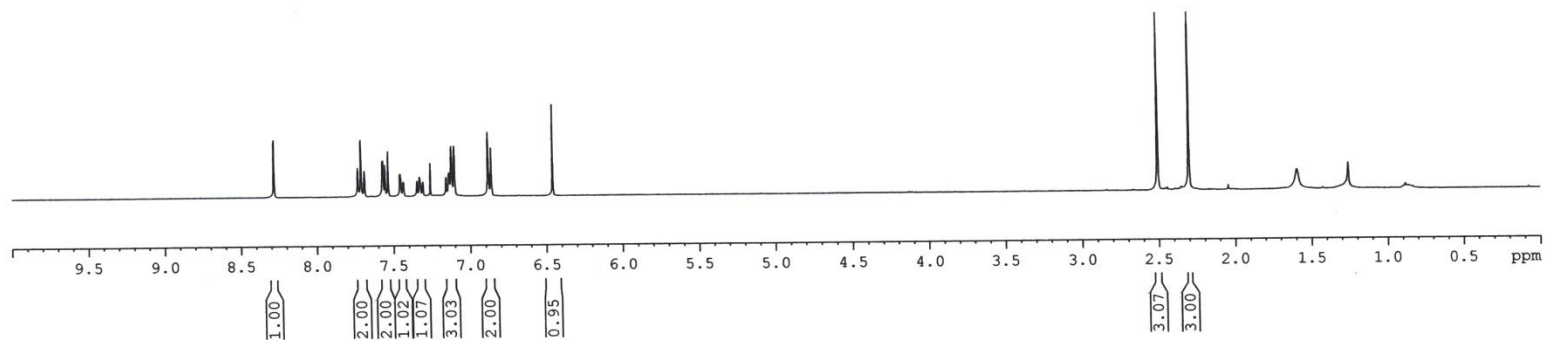
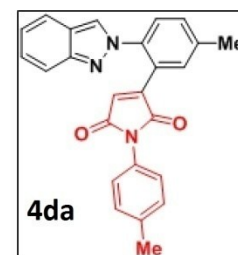
===== CHANNEL f2 =====
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG[2] waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

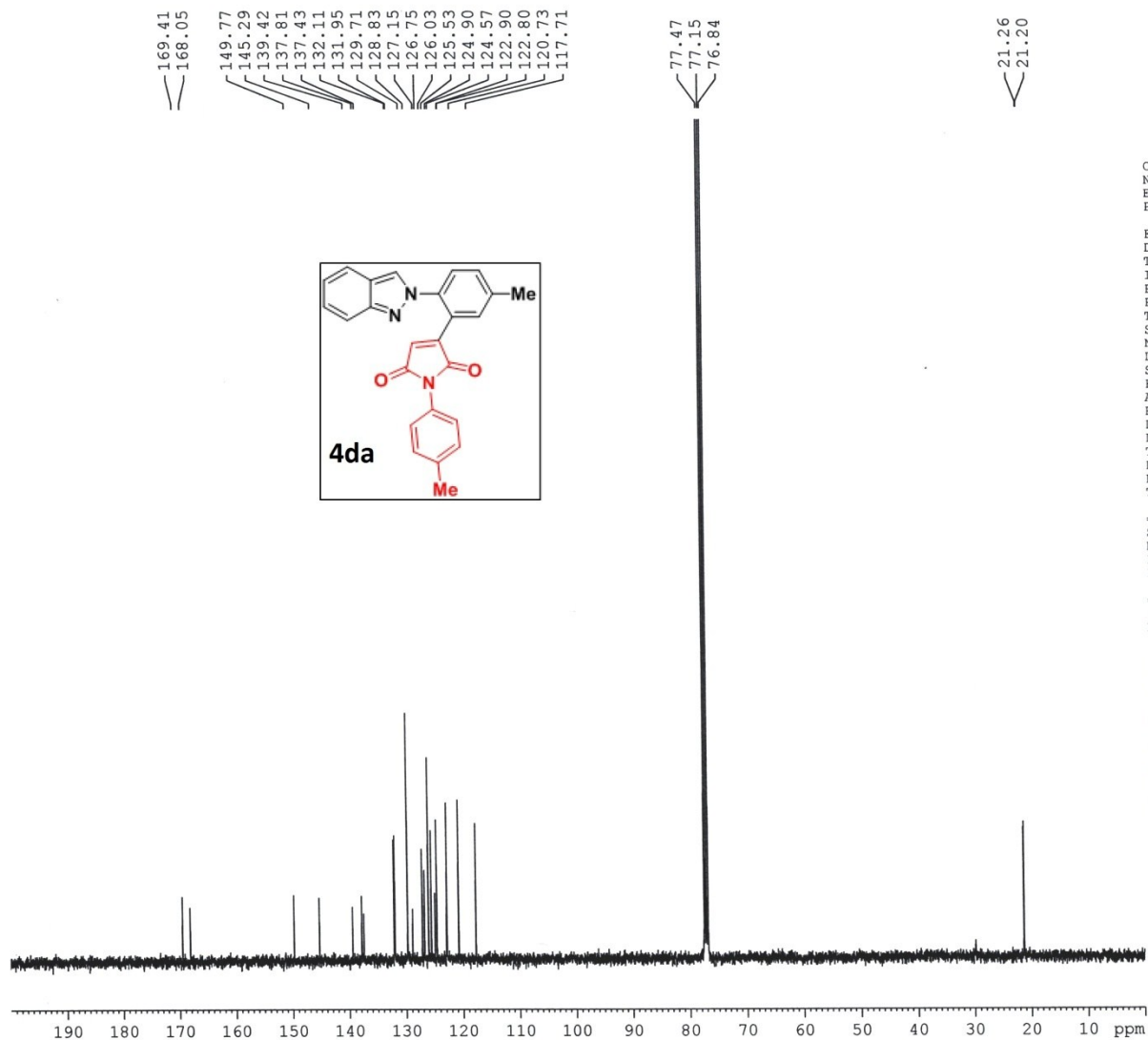
F2 - Processing parameters
 SI 16384
 SF 100.6177842 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40



2.506
2.301

0.074





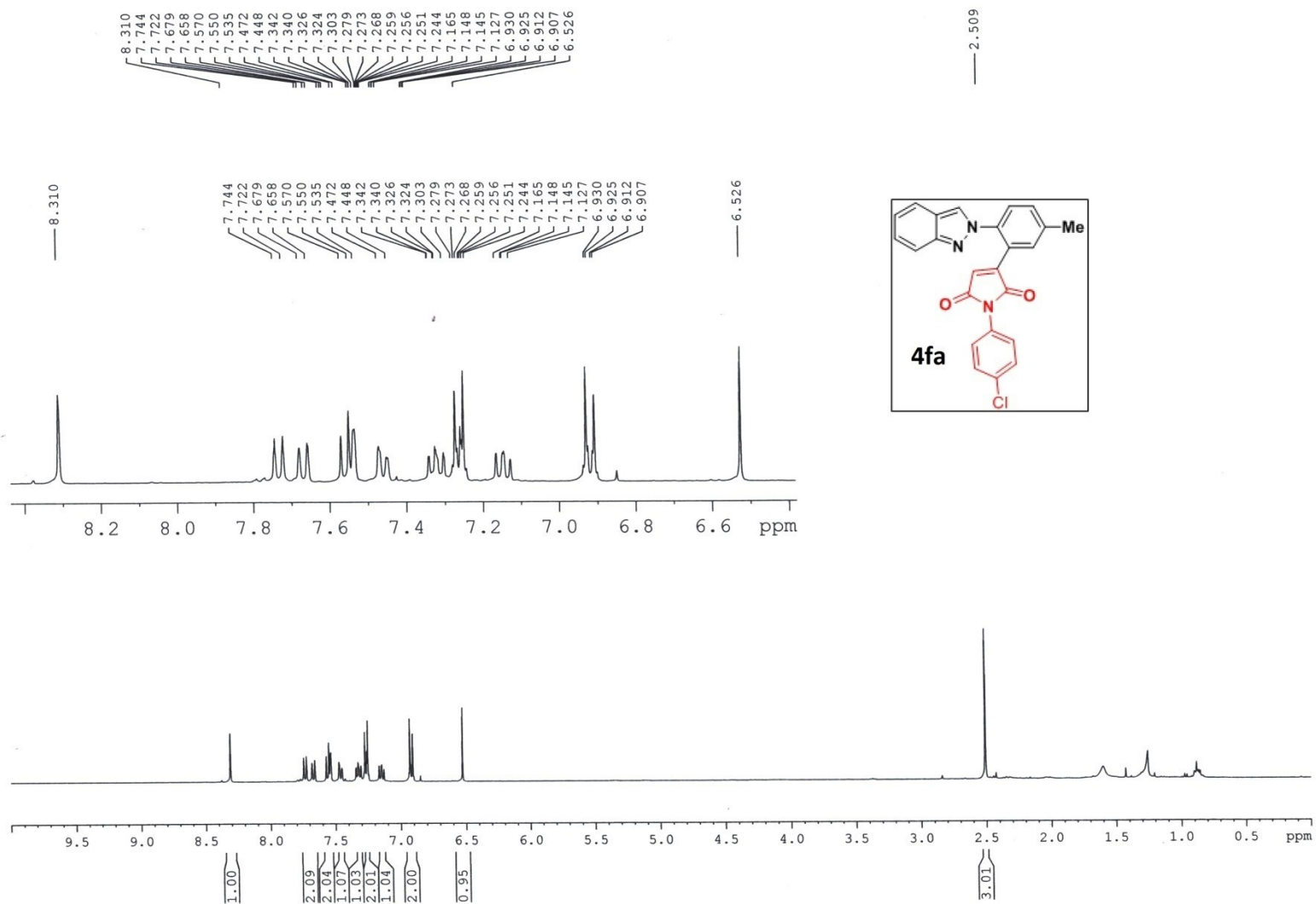
Current Data Parameters
 NAME Dr. A HAJRA-2019-13C
 EXPNO 442
 PROCNO 1

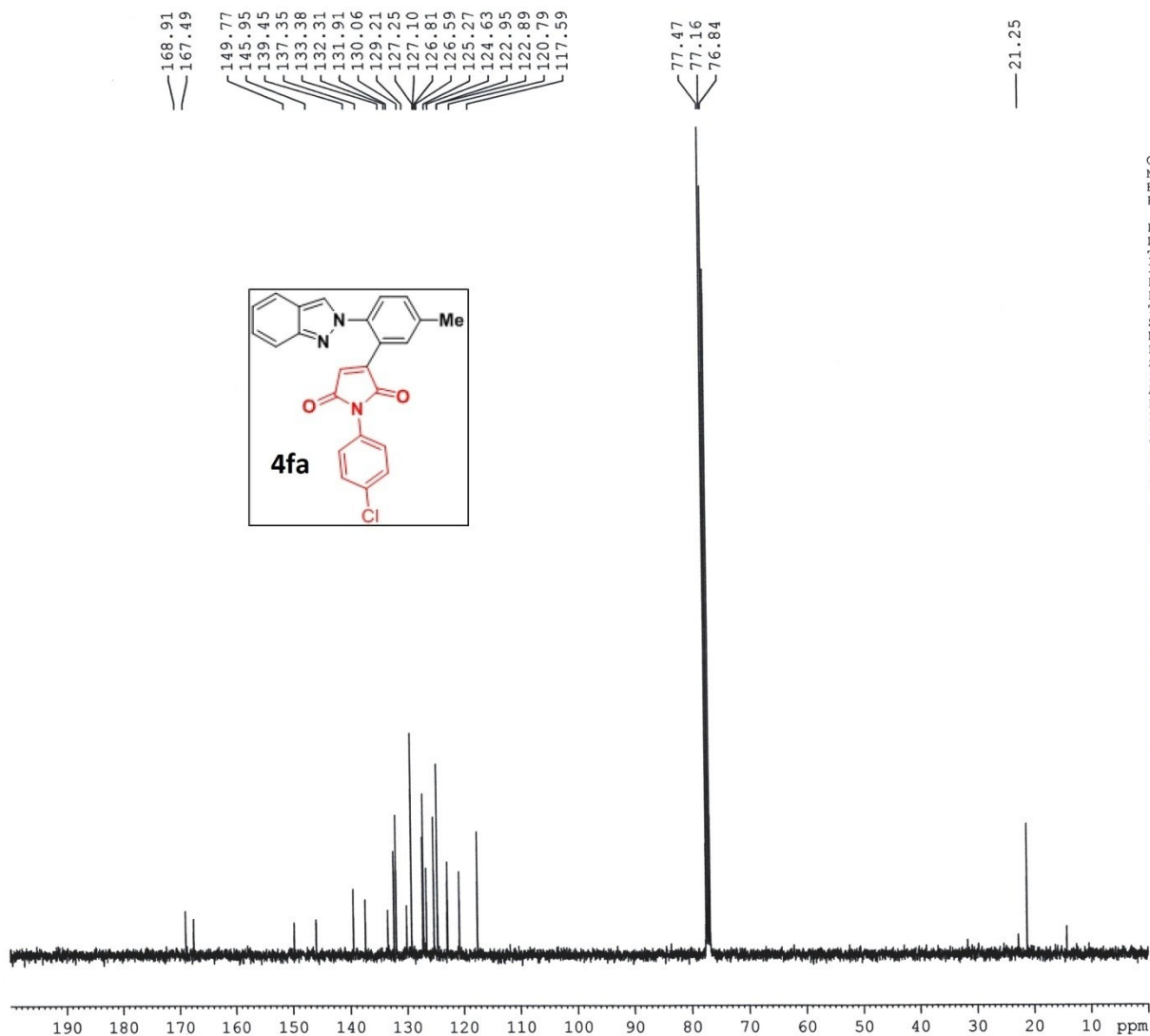
F2 - Acquisition Parameters
 Date_ 20191022
 Time_ 11.00
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDCl₃
 NS 824
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 299.3 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TD0 1

===== CHANNEL f1 =====
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

===== CHANNEL f2 =====
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG[2] waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177847 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40





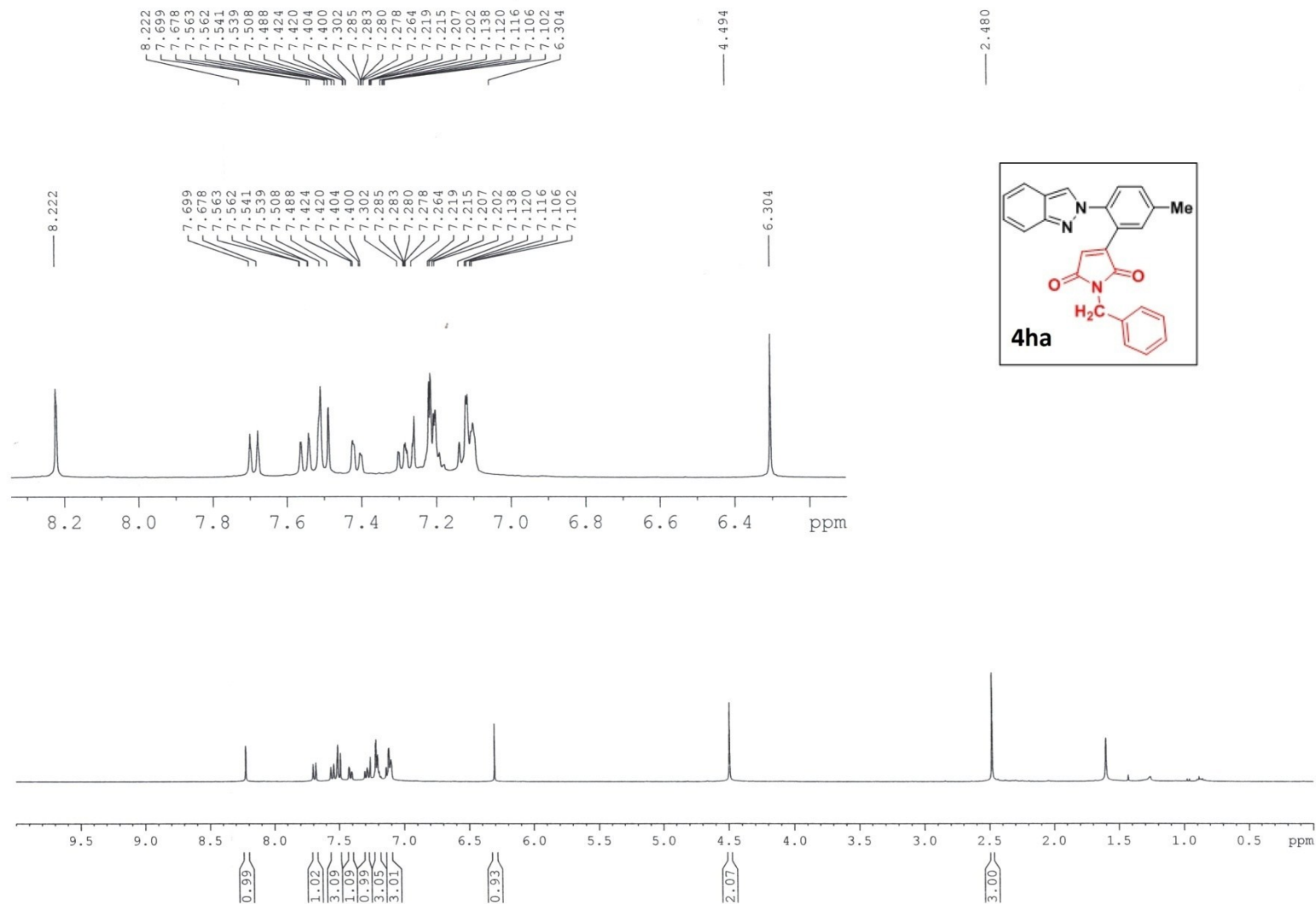
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 508
PROCNO 1

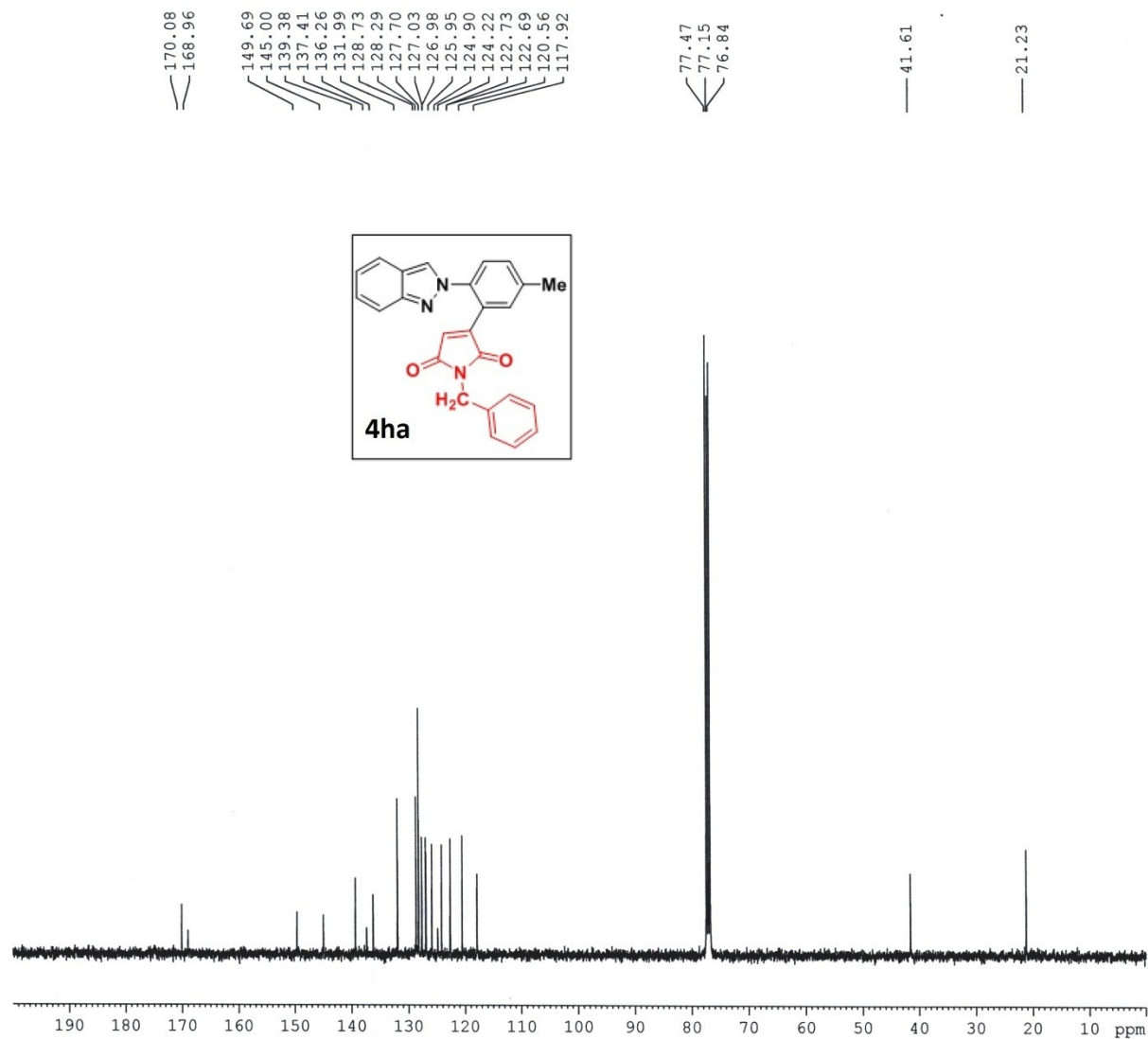
F2 - Acquisition Parameters
Date 20191107
Time 10.17
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 512
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 120.16
DW 20.800 usec
DE 6.50 usec
TE 299.4 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177843 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40





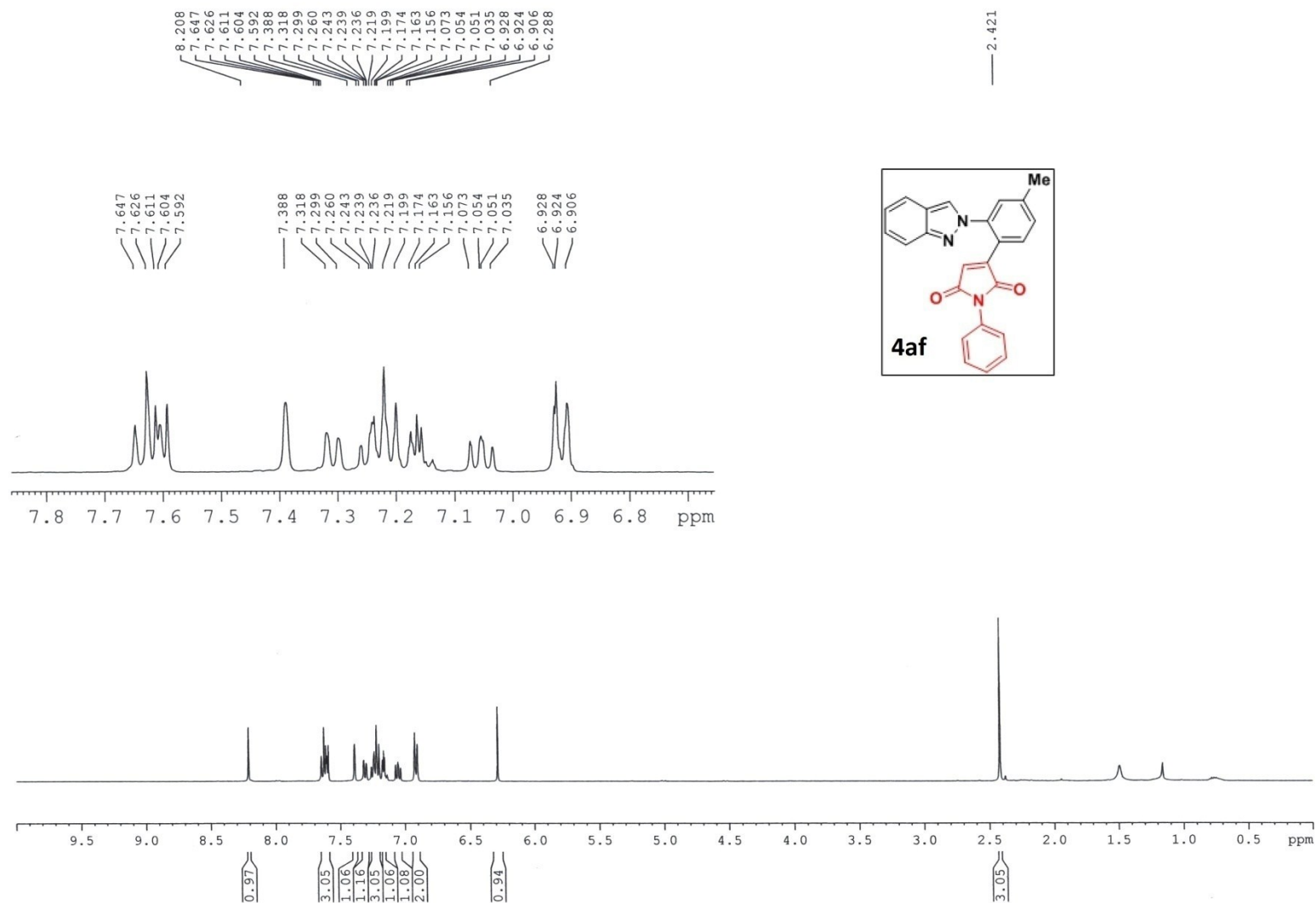
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 462
PROCNO 1

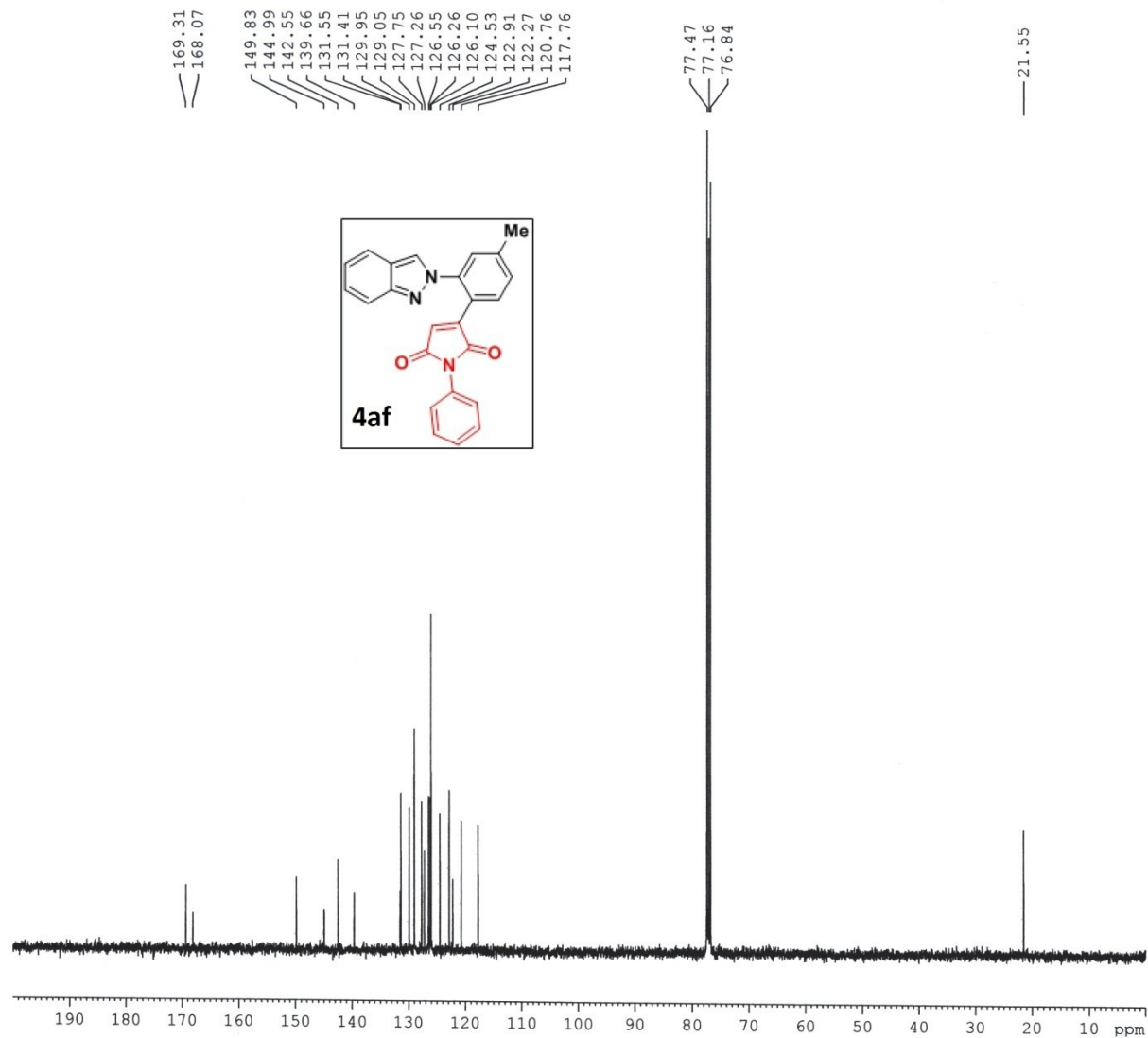
F2 - Acquisition Parameters
Date_ 20191026
Time 9.39
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 640
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 186.42
DW 20.800 usec
DE 6.50 usec
TE 299.1 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG2 waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177848 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40





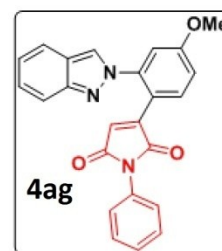
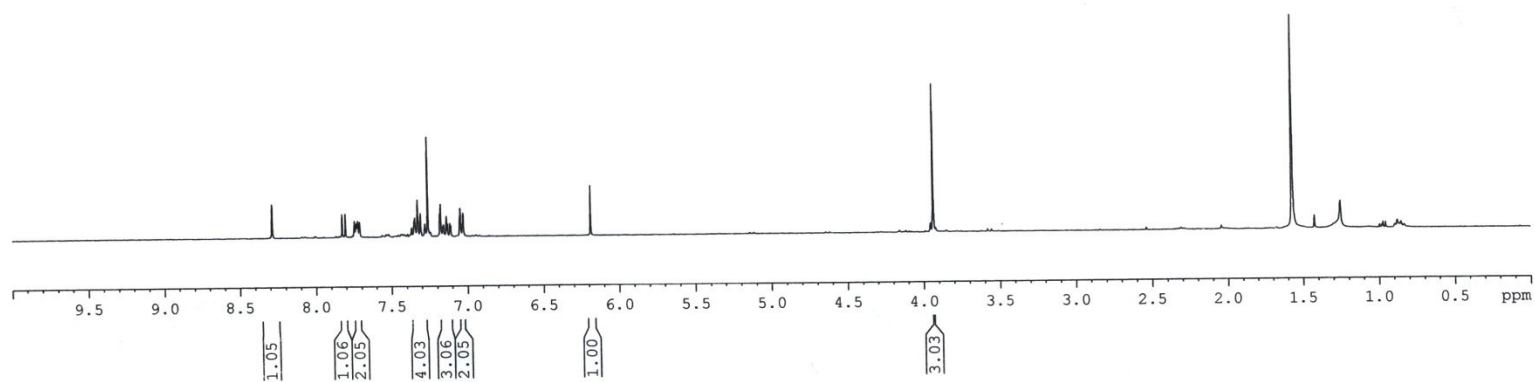
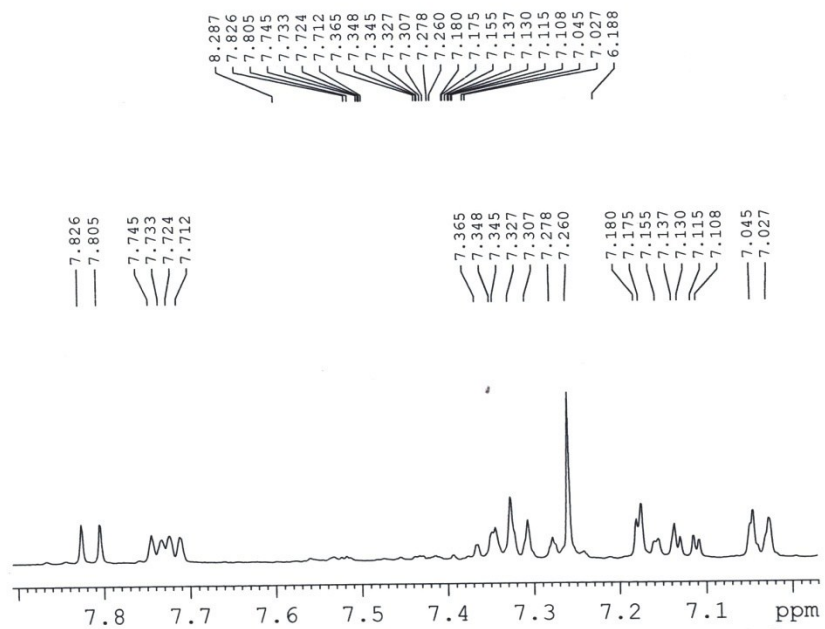
Current Data Parameters
NAME Dr. A HAJRA-2019-13C
EXPNO 445
PROCNO 1

F2 - Acquisition Parameters
Date_ 20191022
Time_ 20.57
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDC13
NS 620
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 186.42
DW 20.800 usec
DE 6.50 usec
TE 299.6 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

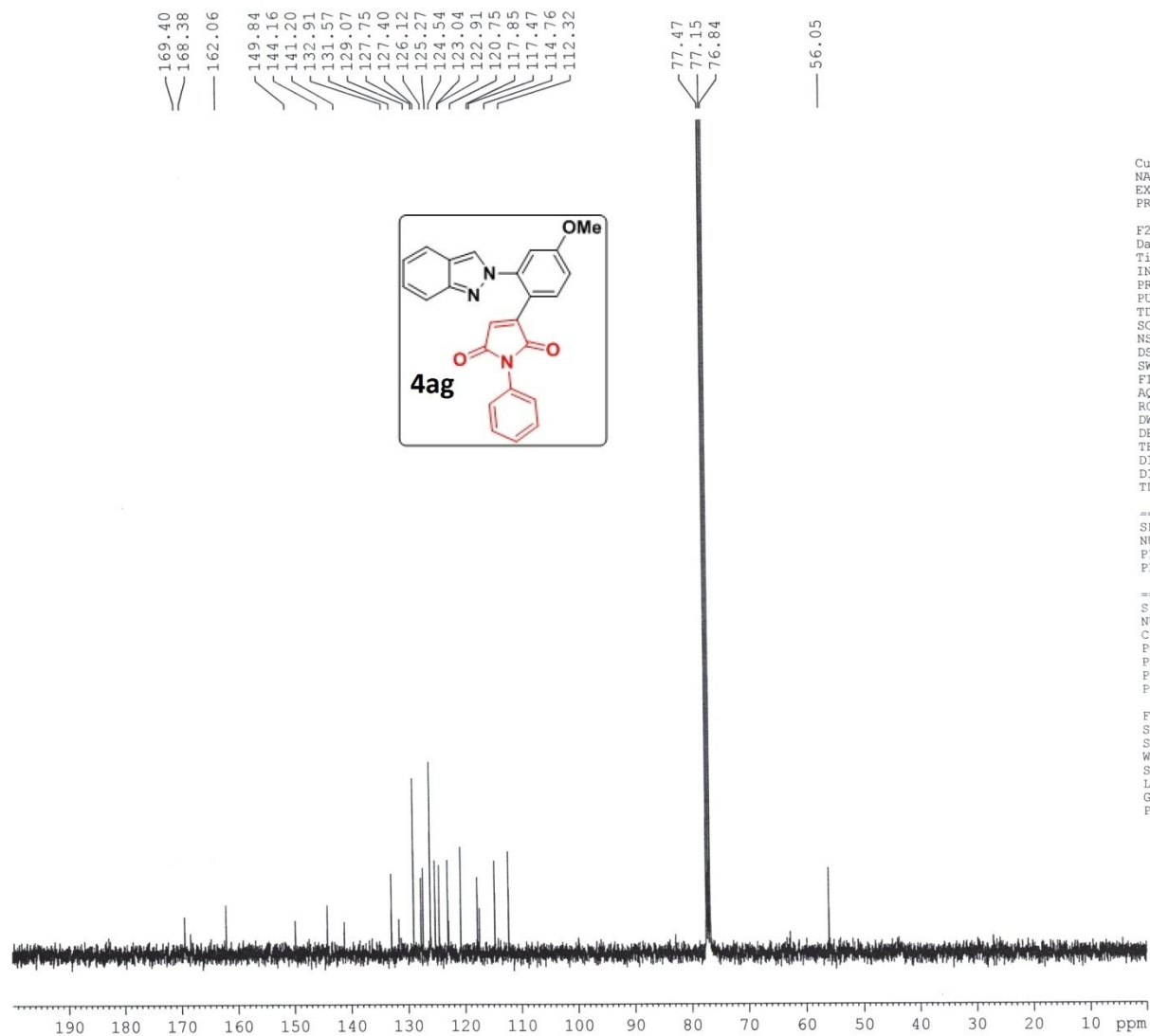
===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG2 waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177846 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40



— 3.937



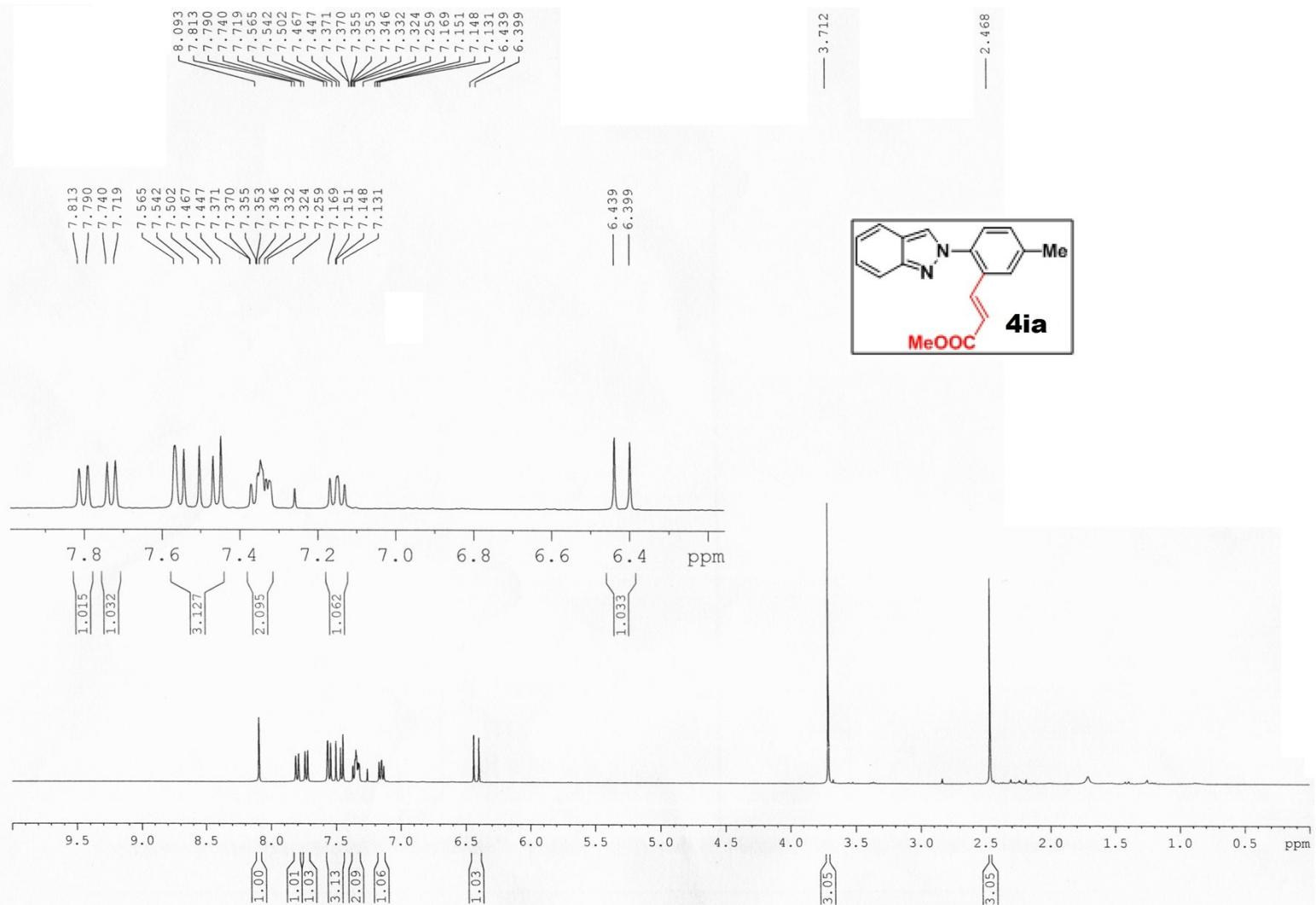
Current Data Parameters
 NAME Dr. A HAJRA-2019-13C
 EXPNO 491
 PROCNO 1

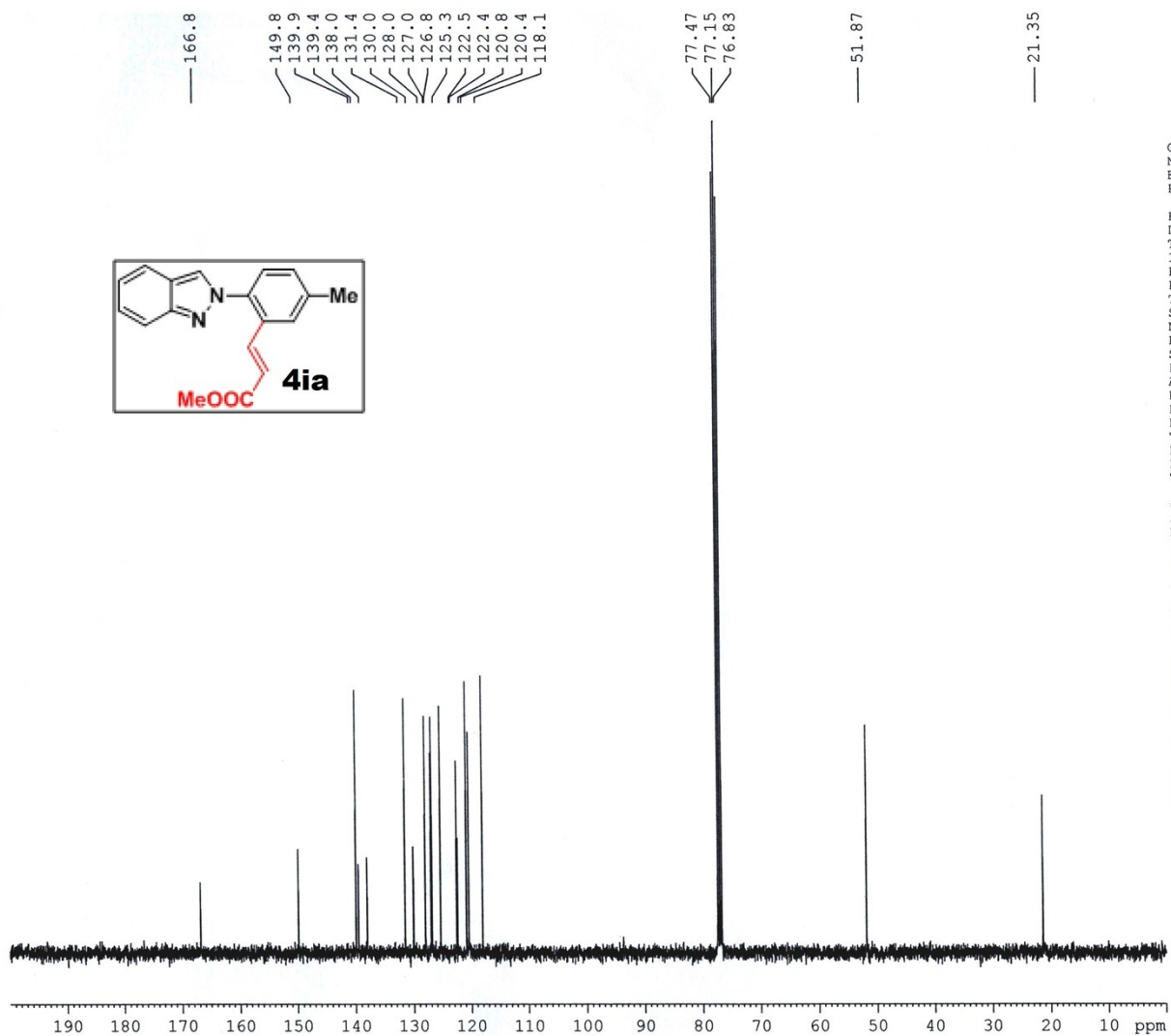
F2 - Acquisition Parameters
 Date_ 20191103
 Time 11.55
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDCl3
 NS 512
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 297.8 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TDO 1

===== CHANNEL f1 =====
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

===== CHANNEL f2 =====
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG[2] waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177843 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40





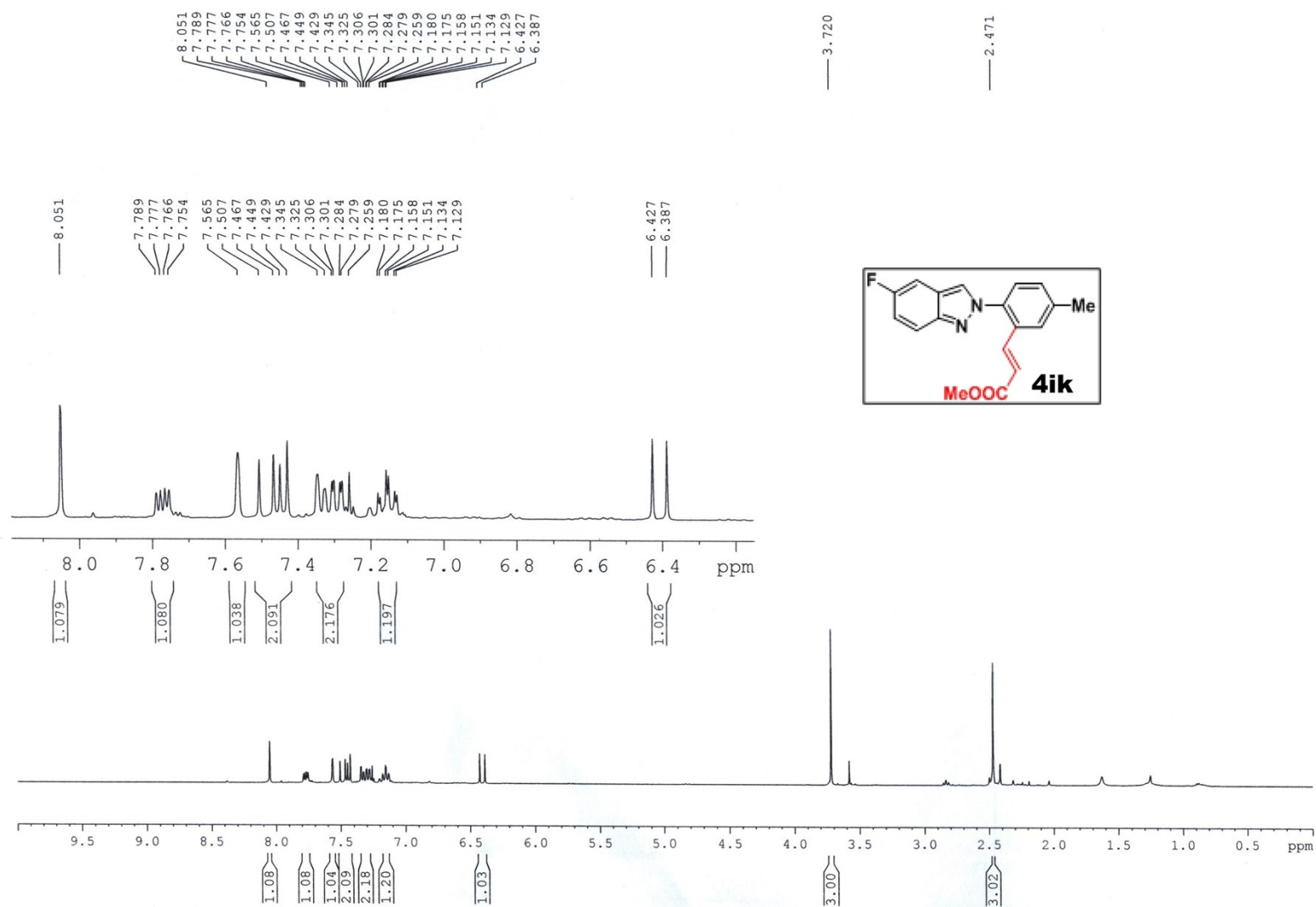
Current Data Parameters
 NAME Dr. A HAJRA-2019-13C
 EXPNO 591
 PROCNO 1

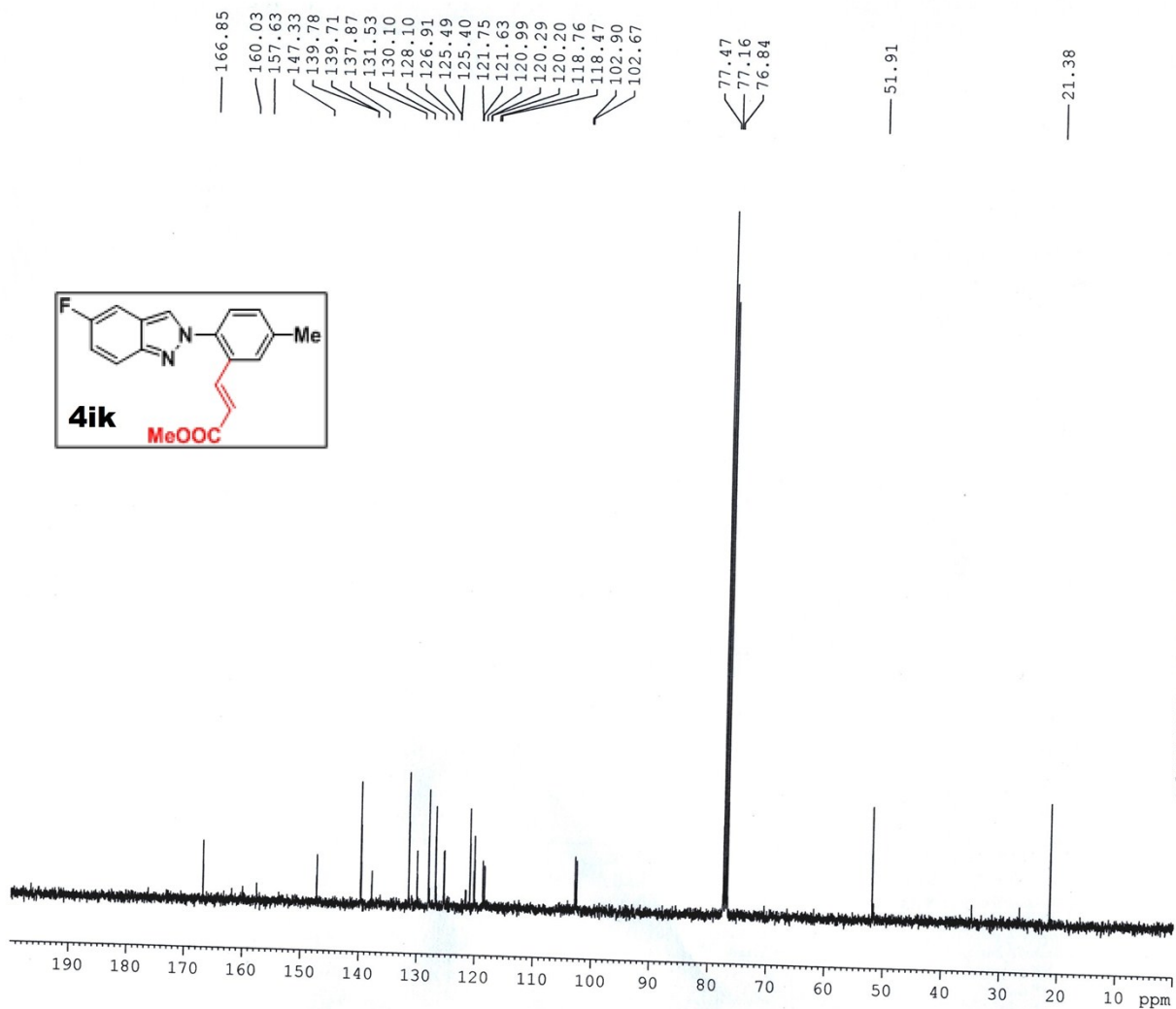
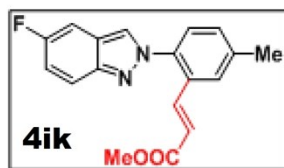
F2 - Acquisition Parameters
 Date_ 20191223
 Time 20.15
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDCl3
 NS 220
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 297.0 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TD0 1

===== CHANNEL f1 =====
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

===== CHANNEL f2 =====
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG[2] waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177872 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40





Current Data Parameters
 NAME Dr. A HAJRA-2019-13C
 EXPNO 592
 PROCNO 1

F2 - Acquisition Parameters
 Date 20191225
 Time 15.47
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDC13
 NS 350
 DS 2
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 298.5 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TDO 1

===== CHANNEL f1 =====
 SFO1 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

===== CHANNEL f2 =====
 SFO2 400.1516006 MHz
 NUC2 1H
 CPDPRG[2] waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16212000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177844 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40