

Supporting Information

Synthesis of 3-Azabicyclo[3.1.0]hexane Derivatives via Palladium-Catalyzed Cyclopropanation of Maleimides with *N*-Tosylhydrazones: Practical and Facile Access to CP-

866,087

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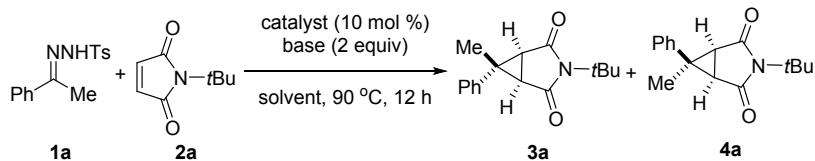
A. General Information

Melting points were measured using a melting point instrument and are uncorrected. Chemical shifts were reported in ppm from the solvent resonance as the internal standard (CDCl_3 $\delta_{\text{H}} = 7.26$ ppm, $\delta_{\text{C}} = 77.16$ ppm). Multiplicity was indicated as follows: s (singlet), d (doublet), t (triplet), q (quartet), m (multiplet). Coupling constants were reported in Hertz (Hz). IR spectra were obtained with an infrared spectrometer on either potassium bromide pellets or liquid films between two potassium bromide pellets. GC–MS data were obtained using electron ionization. HRMS was carried out on a high-resolution mass spectrometer (LCMS-IT-TOF). TLC was performed using commercially available 100–400 mesh silica gel plates (GF₂₅₄). X-ray structural analyses were conducted on an X-ray analysis instrument. The 25 mL Schlenk tube was purchased commercially and used directly to optimize the reaction conditions.

Materials. *N*-Tosylhydrazones were synthesized according to literature procedure. Other commercially available reagents and solvents were purchased and used without further purification. Analytical thin-layer chromatography was performed on 0.20 mm silica gel plates (GF₂₅₄) using UV light as a visualizing agent. Flash column chromatography was carried out using silica gel (200–300 mesh) with the indicated solvent system. All reactions were conducted in oven-dried Schlenk tubes. All the reaction temperatures reported are oil bath temperatures.

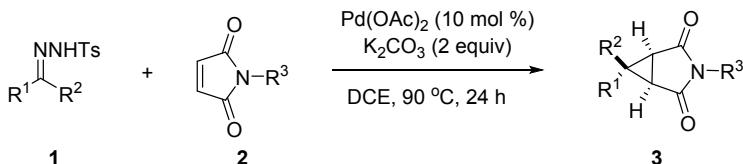
Caution: Normally, *N*-tosylhydrazones are potentially explosive at high temperature. All these grams-scale experiments must be handled carefully. Although no accident happened during our investigation, we strongly recommend all operations involving *N*-tosylhydrazones should carry out in a well-ventilated hood behind a blast shield.

B. General Condition Optimizing Process for the Reaction of Non-Terminal Olefin **2a** and *N*-Tosylhydrazones **1**



A 25 mL round-bottom flask equipped with a magnetic stirring bar, a reflux condenser, **1a** (0.4 mmol), **2** (0.2 mmol), catalyst (10 mol %), base (2 equiv), solvent (2 mL). The mixture was stirred at 90 °C in an oil bath under N₂. After 12 h, the resulting solution was cooled to room temperature, added water (10 mL), and then extracted with EtOAc (3 × 10 mL). The combined organic phases were dried over anhydrous Na₂SO₄, filtered and concentrated *in vacuo*. The ratio of **3a** and **4a** were determined by GC-MS analysis of the obtained crude product. Further purification by flash column chromatography on silica gel (eluting with petroleum ether/ethyl acetate) provided the pure product **3a** and **4a**.

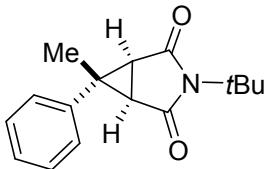
C. General Process for Substrate Scope of *N*-Tosylhydrazones and Maleimides



A 250 mL round-bottom flask equipped with a magnetic stirring bar, a reflux condenser, **1** (20 mmol), **2** (10 mmol), Pd(OAc)₂ (10 mol %), K₂CO₃ (2 equiv), solvent (100 mL). The mixture was stirred at 90 °C in an oil bath under N₂. After 24 h, the resulting solution was cooled to room temperature, added water (100 mL), and then extracted with EtOAc (3 × 100 mL). The combined organic phases were dried over anhydrous Na₂SO₄, filtered and concentrated *in vacuo*. Further purification by flash column chromatography on silica gel (eluting with petroleum ether/ethyl acetate) provided the pure product **3**.

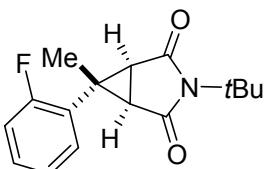
D. Analysis Data for the Products

trans-3-(*tert*-Butyl)-6-Methyl-6-Phenyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3a)



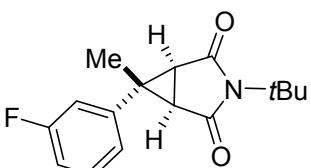
1.82 g, 71% yield; white solid, mp: 149–150 °C ; TLC (petroleum ether/ethyl acetate, 10:1 v/v): R_f = 0.60; ^1H NMR (400 MHz, CDCl_3) δ 7.53–7.25 (m, 5H), 2.65 (s, 2H), 1.58 (s, 9H), 1.57 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.6, 143.0, 128.9, 127.6, 127.3, 58.1, 41.2, 32.5, 28.5, 17.0; IR (KBr): 2973, 2931, 1765, 1702, 1531, 1454, 1339, 1263, 1166, 1083, 1010 cm^{-1} ; HRMS (ESI, m/z): [M+H]⁺ Calcd. for $\text{C}_{16}\text{H}_{20}\text{NO}_2$, 258.1489; found, 258.1488.

trans-3-(*tert*-Butyl)-6-(2-Fluorophenyl)-6-Methyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3b)



2.09 g, 76% yield; white solid, mp: 96–97 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.48; ^1H NMR (400 MHz, CDCl_3) δ 7.31–7.25 (m, 2H), 7.13–7.04 (m, 2H), 2.63 (s, 2H), 1.59 (s, 9H), 1.54 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.2, 160.8 (d, $^1J_{\text{F-C}} = 248.0$ Hz), 130.0 (d, $^2J_{\text{F-C}} = 13.6$ Hz), 129.7 (d, $^3J_{\text{F-C}} = 8.1$ Hz), 129.6 (d, $^4J_{\text{F-C}} = 3.5$ Hz), 124.5 (d, $^3J_{\text{F-C}} = 3.7$ Hz), 116.2 (d, $^2J_{\text{F-C}} = 21.1$ Hz), 58.1, 36.6, 31.7, 28.5, 16.1; ^{19}F NMR (376 MHz, CDCl_3) δ -114.6 – -114.7 (m, 1F); IR (KBr): 3074, 2976, 1768, 1704, 1490, 1455, 1341, 1262, 1216, 1166, 1080, 1012 cm^{-1} ; HRMS (ESI, m/z): [M+H]⁺ Calcd. For $\text{C}_{16}\text{H}_{19}\text{FNO}_2$, 276.1394; found, 276.1395.

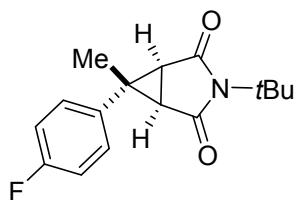
trans-3-(*tert*-Butyl)-6-(3-Fluorophenyl)-6-Methyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3c)



2.28 g, 83% yield; white solid, mp: 92–93 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f =

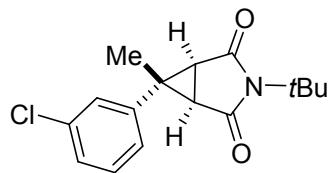
0.50; ^1H NMR (400 MHz, CDCl_3) δ 7.30 (q, $J = 7.2$ Hz, 1H), 7.08 (d, $J = 7.6$ Hz, 1H), 6.98 (m, 2H), 2.63 (s, 2H), 1.58 (s, 9H), 1.57 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.4, 162.9 (d, $^1J_{\text{F-C}} = 245.9$ Hz), 145.2 (d, $^3J_{\text{F-C}} = 7.4$ Hz), 130.5 (d, $^3J_{\text{F-C}} = 8.3$ Hz), 122.8 (d, $^4J_{\text{F-C}} = 2.8$ Hz), 114.7 (d, $^2J_{\text{F-C}} = 16.6$ Hz), 114.5 (d, $^2J_{\text{F-C}} = 17.9$ Hz), 58.2, 40.3, 32.6, 28.5, 16.6; ^{19}F NMR (376 MHz, CDCl_3) δ -111.8 – -111.9 (m, 1F); IR (KBr): 3061, 2971, 2927, 1691, 1584, 1444, 1355, 1261, 1172, 1090, 1021 cm^{-1} ; HRMS (ESI, m/z): $[\text{M}+\text{H}]^+$ Calcd. for $\text{C}_{16}\text{H}_{19}\text{FNO}_2$, 276.1394; found, 276.1393.

***trans*-3-(*tert*-Butyl)-6-(4-Fluorophenyl)-6-Methyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3d)**



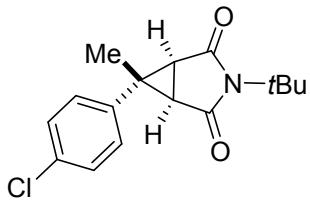
2.42 g, 88% yield; white solid, mp: 95–96 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): $R_f = 0.48$; ^1H NMR (400 MHz, CDCl_3) δ 7.28 (dd, $J_1 = 7.6$ Hz, $J_2 = 5.2$ Hz, 2H), 7.02 (t, $J = 8.4$ Hz, 2H), 2.61 (s, 2H), 1.58 (s, 9H), 1.55 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.4, 161.9 (d, $^1J_{\text{F-C}} = 245.7$ Hz), 138.8 (d, $^4J_{\text{F-C}} = 3.4$ Hz), 129.1 (d, $^3J_{\text{F-C}} = 8.1$ Hz), 115.8 (d, $^2J_{\text{F-C}} = 21.5$ Hz), 58.1, 40.5, 32.6, 28.5, 17.1; ^{19}F NMR (376 MHz, CDCl_3) δ -114.1 – -114.2 (m, 1F); IR (KBr): 2973, 2930, 2868, 1767, 1703, 1513, 1443, 1343, 1224, 1166, 1085, 1011 cm^{-1} ; HRMS (ESI, m/z): $[\text{M}+\text{H}]^+$ Calcd. for $\text{C}_{16}\text{H}_{19}\text{FNO}_2$, 276.1394; found, 276.1399.

***trans*-3-(*tert*-Butyl)-6-(3-Chlorophenyl)-6-Methyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3e)**



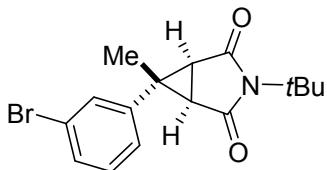
2.36 g, 81% yield; white solid, mp: 102–103 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): $R_f = 0.37$; ^1H NMR (400 MHz, CDCl_3) δ 7.31 (s, 1H), 7.27–7.18 (m, 3H), 2.63 (s, 2H), 1.58 (s, 9H), 1.56 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.1, 144.7, 134.7, 130.2, 127.9, 125.5, 58.2, 40.4, 32.3, 28.5, 16.8; IR (KBr): 3263, 2974, 2923, 1764, 1691, 1464, 1357, 1260, 1174, 1095, 1022, cm^{-1} ; HRMS (ESI, m/z): $[\text{M}+\text{H}]^+$ Calcd. for $\text{C}_{16}\text{H}_{19}\text{ClNO}_2$, 292.1099; found, 292.1098.

***trans*-3-(*tert*-Butyl)-6-(4-Chlorophenyl)-6-Methyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3f)**



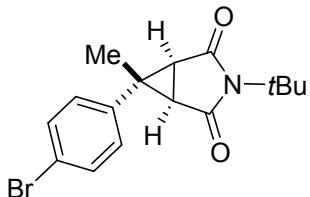
2.56 g, 88% yield; white solid, mp: 102–103 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.33; ^1H NMR (400 MHz, CDCl_3) δ 7.31 (d, J = 8.4 Hz, 2H), 7.25 (d, J = 8.4 Hz, 2H), 2.61 (s, 2H), 1.58 (s, 9H), 1.55 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.2, 141.4, 133.5, 129.1, 128.7, 58.1, 40.3, 32.5, 28.5, 16.9; IR (KBr): 2970, 2933, 2872, 1767, 1702, 1483, 1340, 1165, 1093, 1011 cm^{-1} ; HRMS (ESI, m/z): [M+Na] $^+$ Calcd. for $\text{C}_{16}\text{H}_{18}\text{ClNNaO}_2$, 314.0918; found, 314.0920.

***trans*-6-(3-Bromophenyl)-3-(*tert*-Butyl)-6-Methyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3g)**



2.68 g, 80% yield; white solid, mp: 108–109 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.44; ^1H NMR (400 MHz, CDCl_3) δ 7.47 (s, 1H), 7.40 (d, J = 7.2 Hz, 1H), 7.27–7.19 (m, 2H), 2.63 (s, 2H), 1.58 (s, 9H), 1.56 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.1, 145.0, 130.8, 130.7, 130.5, 126.0, 122.8, 58.2, 40.4, 32.3, 28.5, 16.9; IR (KBr): 3064, 2972, 2930, 1765, 1698, 1564, 1465, 1260, 1168, 1077, 1007 cm^{-1} ; HRMS (ESI, m/z): [M+H] $^+$ Calcd. for $\text{C}_{16}\text{H}_{19}\text{BrNO}_2$, 336.0594; found, 336.0594.

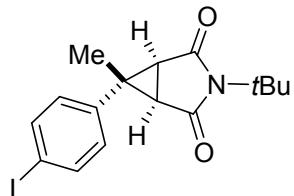
***trans*-6-(4-Bromophenyl)-3-(*tert*-Butyl)-6-Methyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3h)**



2.78 g, 83% yield; white solid, mp: 94–95 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.35; ^1H NMR (400 MHz, CDCl_3) δ 7.46 (d, J = 8.4 Hz, 2H), 7.19 (d, J = 8.4 Hz, 2H), 2.61 (s, 2H), 1.58 (s, 9H), 1.55 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.2, 141.9, 132.1, 129.1, 121.6,

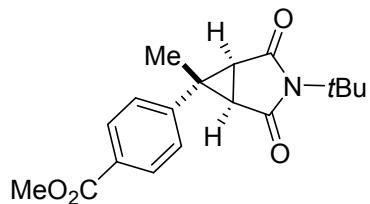
58.2, 40.4, 32.4, 28.5, 16.8; IR (KBr): 3074, 2973, 2930, 1765, 1702, 1485, 1341, 1263, 1166, 1085, 1010 cm^{-1} ; HRMS (ESI, m/z): $[\text{M}+\text{Na}]^+$ Calcd. for $\text{C}_{16}\text{H}_{18}\text{BrNNaO}_2$, 358.0413; found, 358.0410.

trans-3-(*tert*-Butyl)-6-(4-Iodophenyl)-6-Methyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3i)



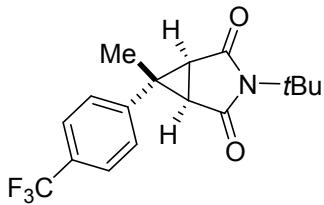
2.07 g, 67% yield; white solid, mp: 123–124 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.47; ^1H NMR (400 MHz, CDCl_3) δ 7.67 (d, J = 8.0 Hz, 2H), 7.06 (d, J = 7.6 Hz, 2H), 2.60 (s, 2H), 1.57 (s, 9H), 1.54 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.2, 142.6, 138.0, 129.3, 93.1, 58.2, 40.5, 32.4, 28.5, 16.8; IR (KBr): 2972, 2928, 2864, 1766, 1701, 1482, 1339, 1264, 1166, 1087, 1007 cm^{-1} ; HRMS (ESI, m/z): $[\text{M}+\text{Na}]^+$ Calcd. for $\text{C}_{16}\text{H}_{18}\text{INNaO}_2$, 406.0274; found, 406.0272.

Methyl 4-(*trans*-3-(*tert*-butyl)-6-Methyl-2,4-Dioxo-3-Azabicyclo[3.1.0]hexan-6-yl)benzoate (3j)



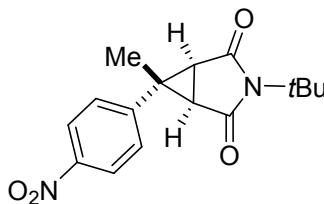
2.43 g, 77% yield; white solid, mp: 137–138 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.22; ^1H NMR (400 MHz, CDCl_3) δ 8.01 (d, J = 8.4 Hz, 2H), 7.38 (d, J = 8.0 Hz, 2H), 3.91 (s, 3H), 2.66 (s, 2H), 1.59 (s, 12H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.1, 166.4, 147.6, 130.2, 129.5, 127.3, 58.2, 52.2, 40.5, 32.5, 28.5, 16.5; IR (KBr): 3069, 2941, 1701, 1610, 1450, 1343, 1278, 1170, 1106, 1012 cm^{-1} ; HRMS (ESI, m/z): $[\text{M}+\text{H}]^+$ Calcd. for $\text{C}_{18}\text{H}_{22}\text{NO}_4$, 316.1543; found, 316.1538.

trans-3-(*tert*-Butyl)-6-Methyl-6-(4-(trifluoromethyl)phenyl)-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3k)



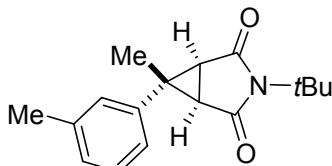
2.60 g, 80% yield; white solid, mp: 177–178 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.30; ¹H NMR (400 MHz, CDCl₃) δ 7.61 (d, J = 8.0 Hz, 2H), 7.44 (d, J = 8.0 Hz, 2H), 2.65 (s, 2H), 1.59 (s, 12H); ¹³C NMR (100 MHz, CDCl₃) δ 174.1, 146.7, 130.0 (q, $^2J_{F-C}$ = 32.5 Hz), 127.8, 126.0 (d, $^3J_{F-C}$ = 3.6 Hz), 123.8, (q, $^1J_{F-C}$ = 270.1 Hz), 58.3, 30.4, 32.3, 28.5, 16.7; ¹⁹F NMR (376 MHz, CDCl₃) δ -62.7 (s, 3F); IR (KBr): 3062, 2980, 2927, 2856, 1689, 1463, 1342, 1263, 1163, 1083, 1027 cm⁻¹; HRMS (ESI, m/z): [M+Na]⁺ Calcd. for C₁₇H₁₈F₃NNaO₂, 348.1182; found, 348.1180.

***trans*-3-(tert-Butyl)-6-Methyl-6-(4-Nitrophenyl)-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3l)**



2.20 g, 73% yield; white solid, mp: 167–168 °C; TLC (petroleum ether/ethyl acetate, 10:1 v/v): R_f = 0.32; ¹H NMR (400 MHz, CDCl₃) δ 8.21 (d, J = 8.8 Hz, 2H), 7.50 (d, J = 8.4 Hz, 2H), 2.68 (s, 2H), 1.61 (s, 3H), 1.59 (s, 9H); ¹³C NMR (100 MHz, CDCl₃) δ 173.6, 149.6, 147.2, 128.3, 124.2, 58.4, 39.9, 32.4, 28.5, 16.5; IR (KBr): 3116, 3059, 2978, 2931, 2867, 1764, 1693, 1515, 1354, 1265, 1177, 1087, 1030 cm⁻¹; HRMS (ESI, m/z): [M+Na]⁺ Calcd. for C₁₆H₁₈N₂NaO₄, 325.1159; found, 325.1162.

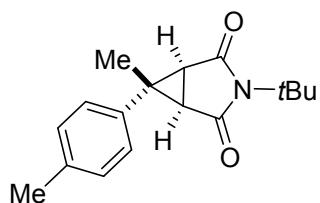
***trans*-3-(tert-Butyl)-6-Methyl-6-(*m*-Tolyl)-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3m)**



2.09 g, 77% yield; white solid, mp: 96–97 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f =

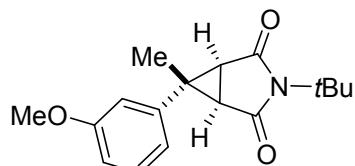
0.56; ^1H NMR (400 MHz, CDCl_3) δ 7.20 (t, $J = 7.2$ Hz, 1H), 7.10 (m, 3H), 2.63 (s, 2H), 2.34 (s, 3H), 1.58 (s, 9H), 1.56 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.7, 143.0, 138.7, 129.8, 128.3, 128.2, 124.4, 58.0, 41.4, 32.5, 28.5, 21.3, 17.1; IR (KBr): 3067, 2971, 2926, 2868, 1763, 1698, 1460, 1335, 1266, 1165, 1082, 1007 cm^{-1} ; HRMS (ESI, m/z): [M+Na] $^+$ Calcd. for $\text{C}_{17}\text{H}_{21}\text{NNaO}_2$, 294.1464; found, 294.1469.

***trans*-3-(*tert*-Butyl)-6-Methyl-6-(*p*-Yolyl)-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3n)**



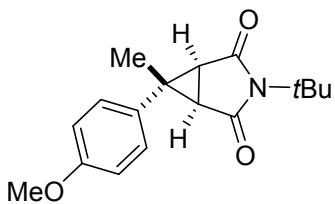
2.01 g, 74% yield; white solid, mp: 93–94 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): $R_f = 0.55$; ^1H NMR (400 MHz, CDCl_3) δ 7.20–7.14 (m, 4H), 2.62 (s, 2H), 2.32 (s, 3H), 1.58 (s, 9H), 1.56 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.7, 140.1, 137.4, 129.5, 127.2, 58.0, 41.0, 32.7, 28.5, 21.0, 17.0; IR (KBr): 2972, 2926, 2865, 1765, 1703, 1462, 1338, 1264, 1164, 1085, 1009 cm^{-1} ; HRMS (ESI, m/z): [M+Na] $^+$ Calcd. for $\text{C}_{17}\text{H}_{21}\text{NNaO}_2$, 294.1464 found, 294.1468.

***trans*-3-(*tert*-Butyl)-6-(3-Methoxyphenyl)-6-Methyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3o)**



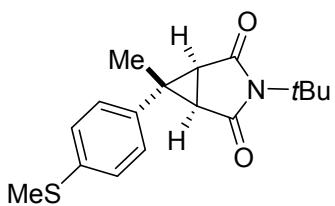
1.89 g, 66% yield; white solid, mp: 92–93 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): $R_f = 0.43$; ^1H NMR (400 MHz, CDCl_3) δ 7.24 (t, $J = 8.0$ Hz, 1H), 6.90–6.79 (m, 3H), 3.80 (s, 3H), 2.64 (s, 2H), 1.58 (s, 9H), 1.57 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.6, 159.9, 144.5, 130.0, 119.5, 113.2, 113.1, 58.0, 55.3, 41.2, 32.6, 28.5, 16.9; IR (KBr): 2972, 2928, 2836, 1765, 1702, 1464, 1339, 1270, 1164, 1045, 1008 cm^{-1} ; HRMS (ESI, m/z): [M+Na] $^+$ Calcd. for $\text{C}_{17}\text{H}_{21}\text{NNaO}_3$, 310.1414; found, 310.1415.

***trans*-3-(*tert*-Butyl)-6-(4-Methoxyphenyl)-6-Methyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3p)**



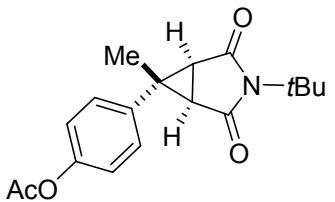
1.61 g, 56% yield; white solid, mp: 94–95 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): $R_f = 0.41$; ^1H NMR (400 MHz, CDCl_3) δ 7.22 (d, $J = 8.8$ Hz, 2H), 6.85 (d, $J = 8.8$ Hz, 2H), 3.79 (s, 3H), 2.61 (s, 2H), 1.58 (s, 9H), 1.55 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.7, 158.9, 135.1, 128.3, 114.2, 58.0, 55.3, 40.8, 32.8, 28.5, 17.1; IR (KBr): 2970, 2926, 2838, 1763, 1699, 1462, 1340, 1252, 1167, 1084, 1027 cm^{-1} ; HRMS (ESI, m/z): $[\text{M}+\text{Na}]^+$ Calcd. for $\text{C}_{17}\text{H}_{21}\text{NNaO}_3$, 310.1414; found, 310.1415.

**trans-3-(tert-Butyl)-6-Methyl-6-(4-(Methylthio)phenyl)-3-Azabicyclo[3.1.0]hexane-2,4-Dione
(3q)**



1.76 g, 58% yield; white solid, mp: 116–117 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): $R_f = 0.30$; ^1H NMR (400 MHz, CDCl_3) δ 7.21 (s, 4H), 2.61 (s, 2H), 2.47 (s, 3H), 1.58 (s, 9H), 1.55 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.5, 139.7, 138.2, 127.8, 126.8, 58.1, 40.7, 32.6, 28.5, 16.9, 15.7; IR (KBr): 3061, 2972, 2926, 1762, 1691, 1456, 1351, 1263, 1164, 1086, 1015 cm^{-1} ; HRMS (ESI, m/z): $[\text{M}+\text{Na}]^+$ Calcd. for $\text{C}_{17}\text{H}_{21}\text{NNaO}_2\text{S}$, 326.1185; found, 326.1183.

4-(*trans*-3-(tert-Butyl)-6-Methyl-2,4-Dioxo-3-Azabicyclo[3.1.0]hexan-6-yl)phenyl acetate (3r)

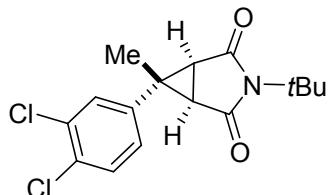


2.24 g, 71% yield; white solid, mp: 151–152 °C; TLC (petroleum ether/ethyl acetate, 10:1 v/v): $R_f = 0.30$; ^1H NMR (400 MHz, CDCl_3) δ 7.32 (d, $J = 8.4$ Hz, 2H), 7.05 (d, $J = 8.0$ Hz, 2H), 2.63 (s, 2H), 2.29 (s, 3H), 1.58 (s, 9H), 1.56 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.4, 169.3, 149.9, 140.5, 128.6, 122.1, 58.1, 40.6, 32.5, 28.5, 21.1, 17.1; IR (KBr): 3075, 2975, 2934, 1762, 1699,

1459, 1364, 1264, 1198, 1083, 1011 cm⁻¹; HRMS (ESI, m/z): [M+H]⁺ Calcd. for C₁₈H₂₂NO₄, 316.1543; found, 316.1545.

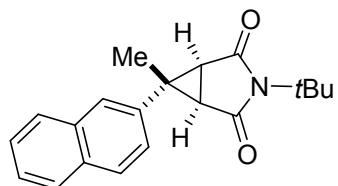
***trans*-3-(*tert*-Butyl)-6-(3,4-Dichlorophenyl)-6-Methyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione**

(3s)



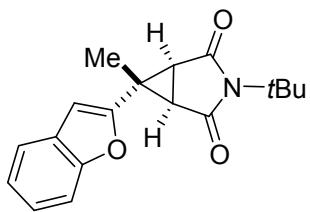
2.41 g, 74% yield; white solid, mp: 158–159 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.40; ¹H NMR (400 MHz, CDCl₃) δ 7.42–7.39 (m, 2H), 7.15 (d, J = 8.4 Hz, 1H), 2.61 (s, 2H), 1.58 (s, 9H), 1.55 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 173.9, 142.9, 133.0, 130.9, 129.5, 126.7, 58.3, 39.6, 32.3, 28.4, 16.7; IR (KBr): 3070, 2975, 2935, 1767, 1702, 1470, 1340, 1265, 1168, 1087, 1020 cm⁻¹; HRMS (ESI, m/z): [M+H]⁺ Calcd. for C₁₆H₁₈Cl₂NO₂, 326.0709; found, 326.0704.

***trans*-3-(*tert*-Butyl)-6-Methyl-6-(Naphthalen-2-yl)-3-Azabicyclo[3.1.0]hexane-2,4-Dione(3t)**



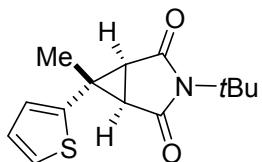
2.21 g, 72% yield; white solid, mp: 161–162 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.56; ¹H NMR (400 MHz, CDCl₃) δ 7.8 (t, J = 8.0 Hz, 3H), 7.76 (s, 1H), 7.51–7.46 (m, 2H), 7.42 (d, J = 8.4 Hz, 1H), 2.75 (s, 2H), 1.66 (s, 3H), 1.61 (s, 9H); ¹³C NMR (100 MHz, CDCl₃) δ 174.6, 140.2, 133.2, 132.6, 128.9, 127.8, 127.6, 126.6, 126.4, 126.1, 125.4, 58.1, 41.4, 32.6, 28.5, 17.0; IR (KBr): 2976, 2921, 2853, 1761, 1698, 1467, 1337, 1268, 1161 cm⁻¹; HRMS (ESI, m/z): [M+Na]⁺ Calcd. for C₂₀H₂₁NNaO₂, 330.1464; found, 330.1467.

***trans*-6-(Benzofuran-2-yl)-3-(*tert*-Butyl)-6-Methyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3u)**



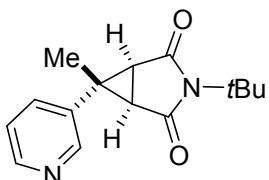
1.90 g, 64% yield; white solid, mp: 133–134 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.48; ^1H NMR (400 MHz, CDCl_3) δ 7.50 (d, J = 7.2 Hz 1H), 7.37 ((d, J = 8.0 Hz 1H), 7.27–7.19 (m, 2H), 6.59 (s, 1H), 2.96 (s, 2H), 1.69 (s, 3H), 1.59 (s, 9H); ^{13}C NMR (100 MHz, CDCl_3) δ 173.5, 156.4, 154.3, 128.1, 124.5, 123.1, 120.9, 111.0, 103.3, 58.3, 33.6, 32.6, 28.5, 11.8; IR (KBr): 3108, 2976, 2928, 1764, 1702, 1459, 1349, 1265, 1163, 1083, 1014 cm^{-1} ; HRMS (ESI, m/z): [M+Na] $^+$ Calcd. for $\text{C}_{18}\text{H}_{19}\text{NNaO}_3$, 320.1257; found, 320.1260.

trans-3-(tert-Butyl)-6-Methyl-6-(thiophen-2-yl)-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3v)



1.32 g, 50% yield; white solid, mp: 74–75 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.57; ^1H NMR (400 MHz, CDCl_3) δ 7.17 (d, J = 4.8 Hz, 1H), 6.91 (t, J = 4.0 Hz, 1H), 6.86 (d, J = 3.6 Hz, 1H), 2.71 (s, 2H), 1.68 (s, 3H), 1.57 (s, 9H); ^{13}C NMR (100 MHz, CDCl_3) δ 173.5, 146.9, 126.9, 124.3, 123.8, 58.2, 35.9, 35.2, 28.5, 15.9; IR (KBr): 2973, 2929, 1766, 1702, 1463, 1338, 1264, 1165, 1075, 1007 cm^{-1} ; HRMS (ESI, m/z): [M+Na] $^+$ Calcd. for $\text{C}_{14}\text{H}_{17}\text{NNaO}_2\text{S}$, 286.0872; found, 286.0872.

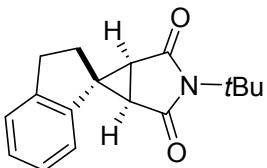
trans-3-(tert-Butyl)-6-Methyl-6-(Pyridin-3-yl)-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3w)



1.47 g, 57% yield; orange solid, mp: 150–151 °C; TLC (petroleum ether/ethyl acetate, 1:1 v/v): R_f = 0.45; ^1H NMR (400 MHz, CDCl_3) δ 8.57 (d, J = 27.6 Hz, 2H), 7.62 (s, 1H), 7.27 (s, 1H), 2.66 (s, 2H), 1.59 (s, 9H), 1.58 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 173.9, 156.9, 148.9, 138.4, 135.0,

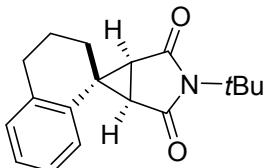
123.6, 58.3, 38.4, 31.9, 28.5, 16.7; IR (KBr): 3042, 2981, 2937, 1762, 1700, 1467, 1364, 1262, 1171, 1083, 1022 cm^{-1} ; HRMS (ESI, m/z): [M+H]⁺ Calcd. for C₁₅H₁₉N₂O₂, 259.1441; found, 259.1443.

***trans*-3-(*tert*-Butyl)-2',3'-Dihydro-3-Azaspido[bicyclo[3.1.0]hexane-6,1'-indene]-2,4-Dione (3x)**



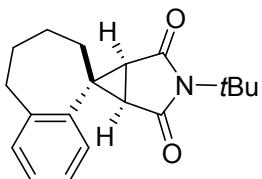
0.81 g, 30% yield; white solid, mp: 119–120 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.50; ¹H NMR (400 MHz, CDCl₃) δ 7.22–7.14 (m, 3H), 6.67 (d, J = 7.2 Hz, 1H), 3.12 (t, J = 7.6 Hz, 2H), 2.64 (s, 2H), 2.28 (t, J = 7.6 Hz, 2H), 1.57 (s, 9H); ¹³C NMR (100 MHz, CDCl₃) δ 174.3, 143.7, 141.9, 128.1, 127.1, 124.9, 119.1, 57.9, 46.5, 34.2, 29.9, 28.5, 26.2; IR (KBr): 2970, 2928, 2855, 1764, 1701, 1464, 1344, 1265, 1163, 1008 cm^{-1} ; HRMS (ESI, m/z): [M+Na]⁺ Calcd. for C₁₇H₁₉NNaO₂, 292.1308; found, 292.1309.

***trans*-3-(*tert*-Butyl)-3',4'-Dihydro-2'H-3-Azaspido[bicyclo[3.1.0]hexane-6,1'-naphthalene]-2,4-Dione (3y)**



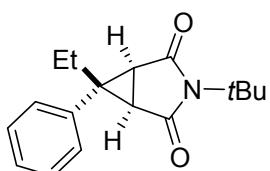
1.13 g, 40% yield; white solid, mp: 72–73 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.58; ¹H NMR (400 MHz, CDCl₃) δ 7.16–7.09 (m, 3H), 6.65 (d, J = 6.8 Hz, 1H), 2.88 (t, J = 5.2 Hz, 2H), 2.69 (s, 2H), 1.96–1.91 (m, 4H), 1.58 (s, 9H); ¹³C NMR (100 MHz, CDCl₃) δ 173.9, 137.9, 136.6, 129.3, 126.9, 126.6, 121.8, 58.1, 38.2, 36.4, 30.2, 28.5, 24.2, 21.9; IR (KBr): 2932, 2865, 1763, 1700, 1458, 1343, 1263, 1166, 1009 cm^{-1} ; HRMS (ESI, m/z): [M+Na]⁺ Calcd. for C₁₈H₂₁NNaO₂, 306.1464; found, 306.1466.

***trans*-3'-(*tert*-Butyl)-6,7,8,9-Tetrahydro-3'-Azaspido[benzo[7]annulene-5,6'-bicyclo[3.1.0]hexane]-2',4'-Dione (3z)**



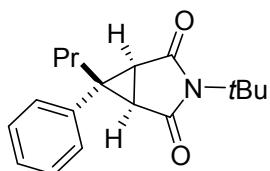
0.92 g, 31% yield; white solid, mp: 134–135 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.61; ^1H NMR (400 MHz, CDCl_3) δ 7.19–7.15 (m, 4H), 2.98 (d, J = 10.4 Hz, 2H), 2.65 (s, 2H), 1.87–1.80 (m, 4H), 1.68 (s, 2H), 1.58 (s, 9H); ^{13}C NMR (100 MHz, CDCl_3) δ 175.0, 143.4, 141.7, 129.7, 127.9, 126.4, 126.2, 57.9, 45.0, 34.9, 32.8, 28.5, 28.4, 28.3, 27.3; IR (KBr): 2930, 2857, 1765, 1701, 1454, 1341, 1266, 1209, 1165, 1015 cm^{-1} ; HRMS (ESI, m/z): [M+Na]⁺ Calcd. for $\text{C}_{19}\text{H}_{23}\text{NNaO}_2$, 320.1621; found, 320.1620.

trans-3-(*tert*-Butyl)-6-Ethyl-6-Phenyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3aa)



1.71 g, 63% yield; white solid, mp: 75–76 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.40; ^1H NMR (400 MHz, CDCl_3) δ 7.32 (s, 5H), 2.64 (s, 2H), 1.86 (d, J = 7.2 Hz, 2H), 1.58 (s, 9H), 0.91 (t, J = 7.2 Hz, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.6, 141.3, 128.8, 128.7, 127.7, 57.9, 47.6, 32.2, 28.5, 24.0, 10.6; IR (KBr): 3069, 2973, 2877, 1766, 1689, 1458, 1340, 1265, 1166, 1077, 1016 cm^{-1} ; HRMS (ESI, m/z): [M+Na]⁺ Calcd. for $\text{C}_{17}\text{H}_{21}\text{NNaO}_2$, 294.1464; found, 294.1467.

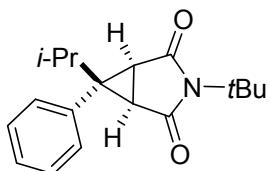
trans-3-(*tert*-Butyl)-6-Phenyl-6-Propyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3ab)



1.91 g, 67% yield; white solid, mp: 73–74 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.47; ^1H NMR (400 MHz, CDCl_3) δ 7.31 (s, 5H), 2.63 (s, 2H), 1.80 (t, J = 8.4 Hz, 2H), 1.59 (s, 9H), 1.39–1.28 (m, 2H), 0.82 (t, J = 7.2 Hz, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.6, 141.6, 128.7, 128.6, 127.6, 57.9, 46.6, 33.0, 32.2, 28.5, 19.7, 13.9; IR (KBr): 2964, 2872, 1766, 1703,

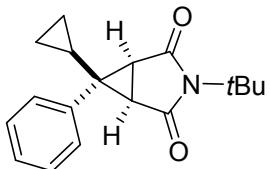
1457, 1339, 1264, 1166, 1014 cm⁻¹; HRMS (ESI, m/z): [M+Na]⁺ Calcd. for C₁₈H₂₃NNaO₂, 308.1621; found, 308.1624.

***trans*-3-(*tert*-Butyl)-6-Isopropyl-6-Phenyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3ac)**



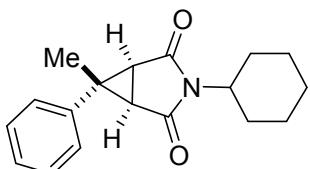
1.37 g, 48% yield; white solid, mp: 142–143 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.40; ¹H NMR (400 MHz, CDCl₃) δ 7.30 (s, 5H), 2.64 (s, 2H), 1.95–1.88 (m, 1H), 1.58 (s, 9H), 0.93 (d, J = 6.8 Hz, 6H); ¹³C NMR (100 MHz, CDCl₃) δ 174.5, 138.2, 130.8, 128.1, 127.8, 58.0, 52.0, 33.2, 28.5, 27.4, 20.0; IR (KBr): 3072, 2967, 2927, 1759, 1698, 1454, 1370, 1259, 1172, 1079, 1017 cm⁻¹; HRMS (ESI, m/z): [M+Na]⁺ Calcd. for C₁₈H₂₃NNaO₂, 308.1621; found, 308.1624.

***trans*-3-(*tert*-Butyl)-6-Cyclopropyl-6-Phenyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3ad)**



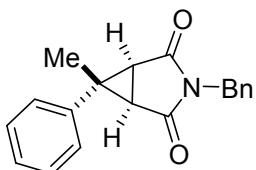
0.99 g, 35% yield; white solid, mp: 90–91 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.43; ¹H NMR (400 MHz, CDCl₃) δ 7.31–7.21 (m, 5H), 2.68 (s, 2H), 1.59 (s, 9H), 1.15–1.12 (m, 1H), 0.50 (d, J = 8.0 Hz, 2H), 0.22 (d, J = 5.2 Hz, 2H); ¹³C NMR (100 MHz, CDCl₃) δ 174.7, 140.2, 129.2, 128.5, 127.7, 57.9, 46.3, 32.3, 28.4, 11.3, 4.9; IR (KBr): 2974, 2929, 1766, 1702, 1458, 1344, 1266, 1210, 1166, 1017 cm⁻¹; HRMS (ESI, m/z): [M+Na]⁺ Calcd. for C₁₈H₂₁NNaO₂, 306.1464; found, 306.1469.

***trans*-3-Cyclohexyl-6-Methyl-6-Phenyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3ae)**



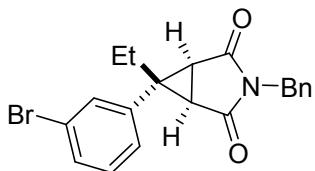
1.90 g, 67% yield; white solid, mp: 94–95 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.56; ^1H NMR (400 MHz, CDCl_3) δ 7.32 (s, 5H), 3.93 (t, J = 12.0 Hz, 1H), 2.20–2.11 (m, 2H), 1.83 (d, J = 12.4 Hz, 2H), 1.67–1.60 (m, 3H), 1.53 (s, 3H), 1.36–1.21 (m, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 173.6, 142.9, 128.9, 127.7, 127.3, 51.5, 42.0, 32.4, 29.0, 25.9, 25.0, 17.2; IR (KBr): 3076, 2930, 2857, 1765, 1698, 1449, 1371, 1266, 1188, 1052 cm^{-1} ; HRMS (ESI, m/z): $[\text{M}+\text{Na}]^+$ Calcd. for $\text{C}_{18}\text{H}_{21}\text{NNaO}_2$, 306.1464; found, 306.1467.

***trans*-3-Benzyl-6-Methyl-6-Phenyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3af)**



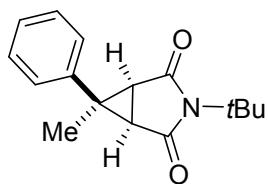
1.78 g, 61% yield; white solid, mp: 151–152 °C; TLC (petroleum ether/ethyl acetate, 10:1 v/v): R_f = 0.43; ^1H NMR (400 MHz, CDCl_3) δ 7.43 (d, J = 7.2 Hz, 2H), 7.34–7.25 (m, 8H), 4.61 (s, 2H), 2.76 (s, 2H), 1.27 (s, 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 173.2, 142.8, 135.7, 129.3, 129.0, 128.6, 128.1, 127.7, 127.2, 42.8, 42.1, 32.5, 17.0; IR (KBr): 3077, 2967, 2858, 1760, 1695, 1440, 1392, 1336, 1167, 1058, 1000 cm^{-1} ; HRMS (ESI, m/z): $[\text{M}+\text{Na}]^+$ Calcd. for $\text{C}_{19}\text{H}_{17}\text{NNaO}_2$, 314.1151; found, 314.1152.

***trans*-3-Benzyl-6-(3-Bromophenyl)-6-Ethyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (3ag)**



2.80 g, 73% yield; white solid, mp: 126–127 °C; TLC (petroleum ether/ethyl acetate, 15:1 v/v): R_f = 0.25; ^1H NMR (400 MHz, CDCl_3) δ 7.45–7.40 (m, 4H), 7.33–7.17 (m, 5H), 4.61 (s, 2H), 2.73 (s, 2H), 1.43 (q, J = 7.6 Hz 2H), 0.70 (t, J = 7.6 Hz 3H); ^{13}C NMR (100 MHz, CDCl_3) δ 172.8, 143.2, 135.7, 131.8, 131.1, 130.3, 129.3, 128.6, 128.1, 127.4, 122.7, 48.3, 42.2, 31.9, 23.9, 10.3; IR (KBr): 3066, 2926, 2858, 1768, 1703, 1567, 1394, 1342, 1164, 1064, 1002 cm^{-1} ; HRMS (ESI, m/z): $[\text{M}+\text{Na}]^+$ Calcd. for $\text{C}_{20}\text{H}_{18}\text{NNaO}_2$, 406.0413; found, 406.0404.

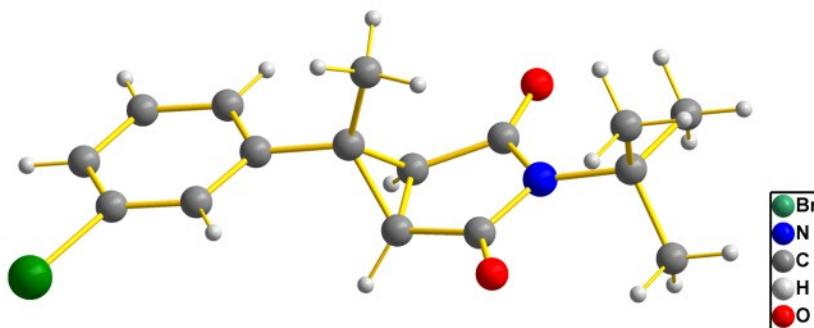
cis-3-(*tert*-Butyl)-6-Methyl-6-Phenyl-3-Azabicyclo[3.1.0]hexane-2,4-Dione (4a)



0.44 g, 16% yield; white solid, mp: 120–121 °C; TLC (petroleum ether/ethyl acetate, 10:1 v/v): R_f = 0.30; ^1H NMR (400 MHz, CDCl_3) δ 7.36–7.27 (m, 5H), 2.55 (s, 2H), 1.44 (s, 3H), 0.98 (s, 9H); ^{13}C NMR (100 MHz, CDCl_3) δ 174.4, 138.1, 129.0, 128.9, 127.9, 56.9, 42.7, 33.2, 27.6, 27.5; IR (KBr): 3059, 3019, 2968, 1766, 1701, 1451, 1358, 1266, 1170, 1071, 1005 cm^{-1} ; HRMS (ESI, m/z): [M+Na]⁺ Calcd. for $\text{C}_{16}\text{H}_{19}\text{NNaO}_2$, 280.1308; found, 280.1308.

E. X-ray Crystallographic Data for 3g

The X-ray crystallographic structures for **3g**. ORTEP representation with 50% probability thermal ellipsoids. Solvent and hydrogen are omitted for clarity. Crystal data have been deposited to CCDC, number 1489545.

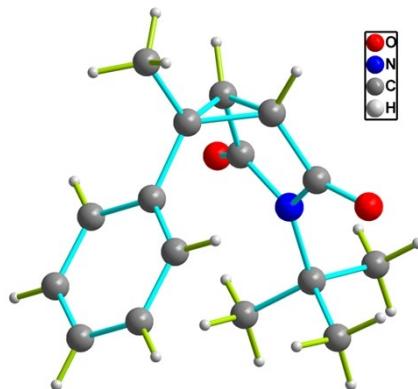


Empirical formula	$\text{C}_{16}\text{H}_{18}\text{BrNO}_2$
Formula weight	336.22
Temperature	293(2) K
Wavelength	0.71073 Å
Crystal system, space group	MONOCLINIC, P2(1)/c
Unit cell dimensions	$a = 6.9645(14)$ Å $\alpha = 90$ deg. $b = 14.893(3)$ Å $\beta = 101.19(3)$ deg. $c = 14.570(3)$ Å $\gamma = 90$ deg.

Volume	1482.5(5) Å ³
Z, Calculated density	4, 1.506 Mg/m ³
Absorption coefficient	2.774 mm ⁻¹
F(000)	688
Crystal size	0.10×0.10×0.10 mm
Theta range for data collection	5.50 to 73.85 deg.
Limiting indices	-8 ≤ h ≤ 8, -17 ≤ k ≤ 17, -17 ≤ l ≤ 15
Reflections collected / unique	10300 / 3491 [R(int) = 0.0420]
Completeness to theta = 25.00	99.8%
Refinement method	Full-matrix least-squares on F ²
Data / restraints / parameters	2607 / 0 / 182
Goodness-of-fit on F ²	1.153
Final R indices [I>2sigma(I)]	R1 = 0.0396, wR2 = 0.0774
R indices (all data)	R1 = 0.0845, wR2 = 0.1299
Extinction coefficient	0.0032(9)

F. X-ray Crystallographic Data for 4a

The X-ray crystallographic structures for **4a**. ORTEP representation with 50% probability thermal ellipsoids. Solvent and hydrogen are omitted for clarity. Crystal data have been deposited to CCDC, number 1489546.

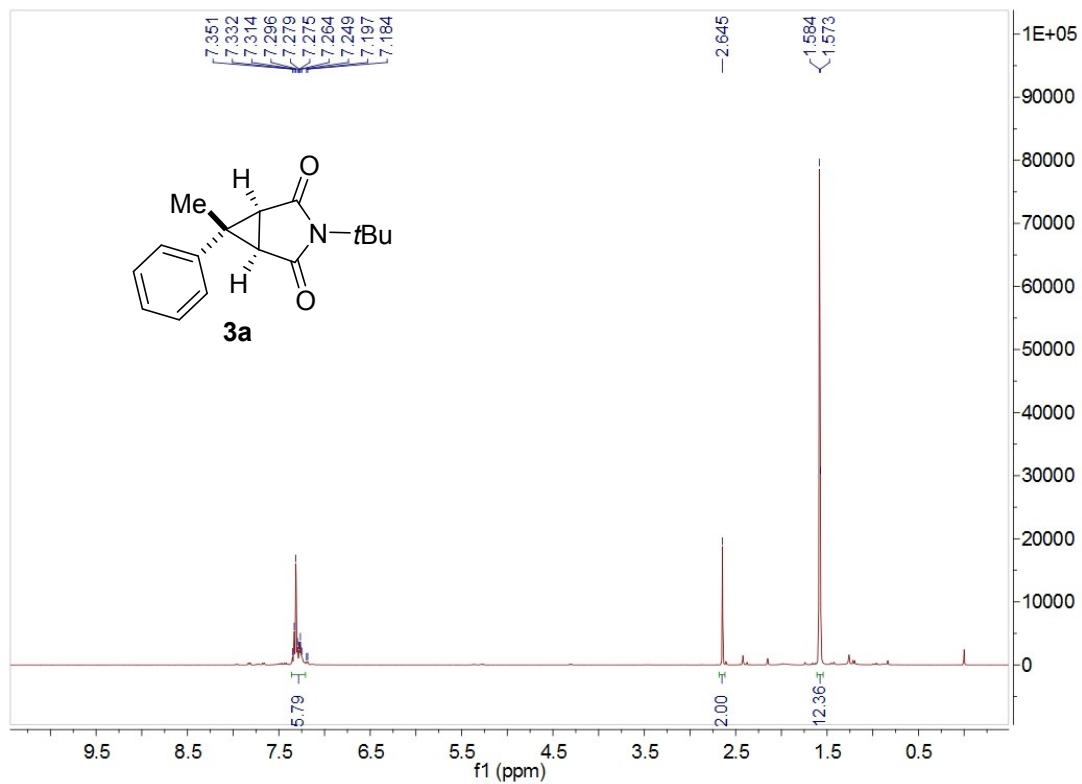


Empirical formula	C ₁₆ H ₁₉ NO ₂
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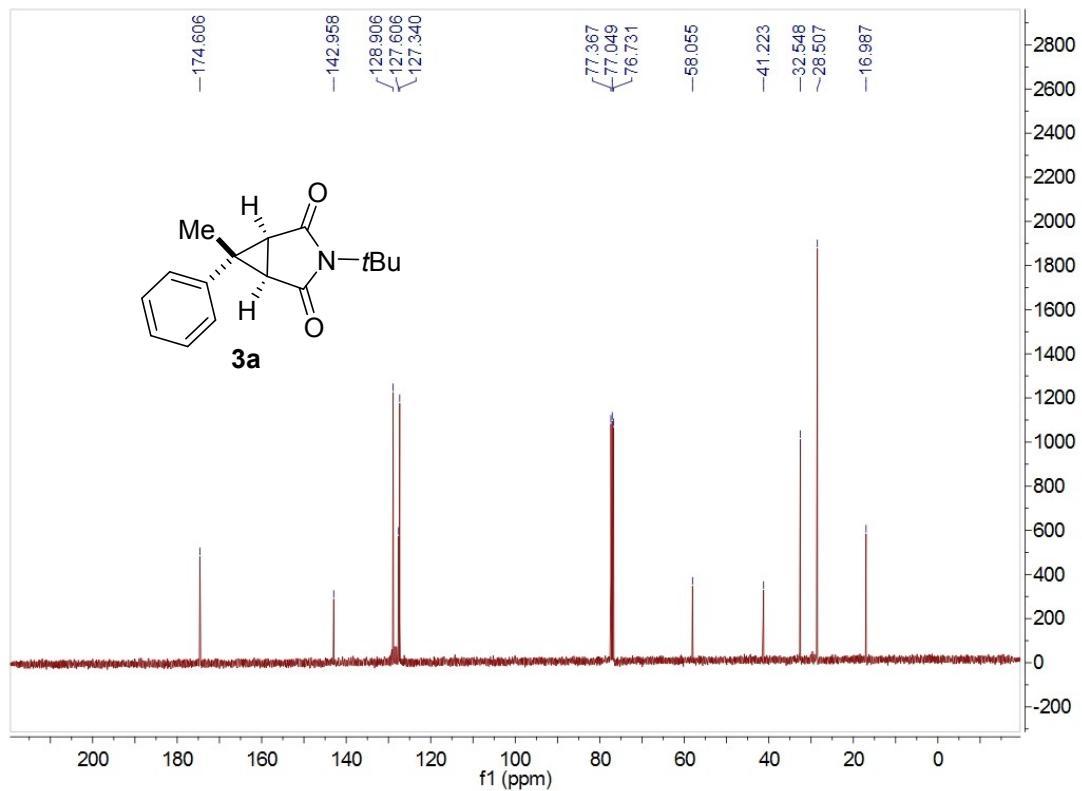
Formula weight	257.32
Temperature	293(2) K
Wavelength	0.71073 Å
Crystal system, space group	MONOCLINIC, P2(1)/c
Unit cell dimensions	a = 8.8657(18) Å alpha = 90 deg. b = 11.708(2) Å beta = 90 deg. c = 13.548(3) Å gamma = 90 deg.
Volume	1406.3(5) Å ³
Z, Calculated density	4, 1.215 Mg/m ³
Absorption coefficient	0.080 mm ⁻¹
F(000)	552
Crystal size	0.10×0.10×0.10 mm
Theta range for data collection	3.01 to 27.48 deg.
Limiting indices	-10 ≤ h ≤ 11, -15 ≤ k ≤ 15, -17 ≤ l ≤ 17
Reflections collected / unique	13757 / 3199 [R(int) = 0.0962]
Completeness to theta = 25.00	99.7%
Refinement method	Full-matrix least-squares on F ²
Data / restraints / parameters	3199 / 0 / 173
Goodness-of-fit on F ²	1.040
Final R indices [I>2sigma(I)]	R1 = 0.0585, wR2 = 0.1521
R indices (all data)	R1 = 0.0972, wR2 = 0.2084
Extinction coefficient	0.31(3)

G. NMR Spectra of New Compounds

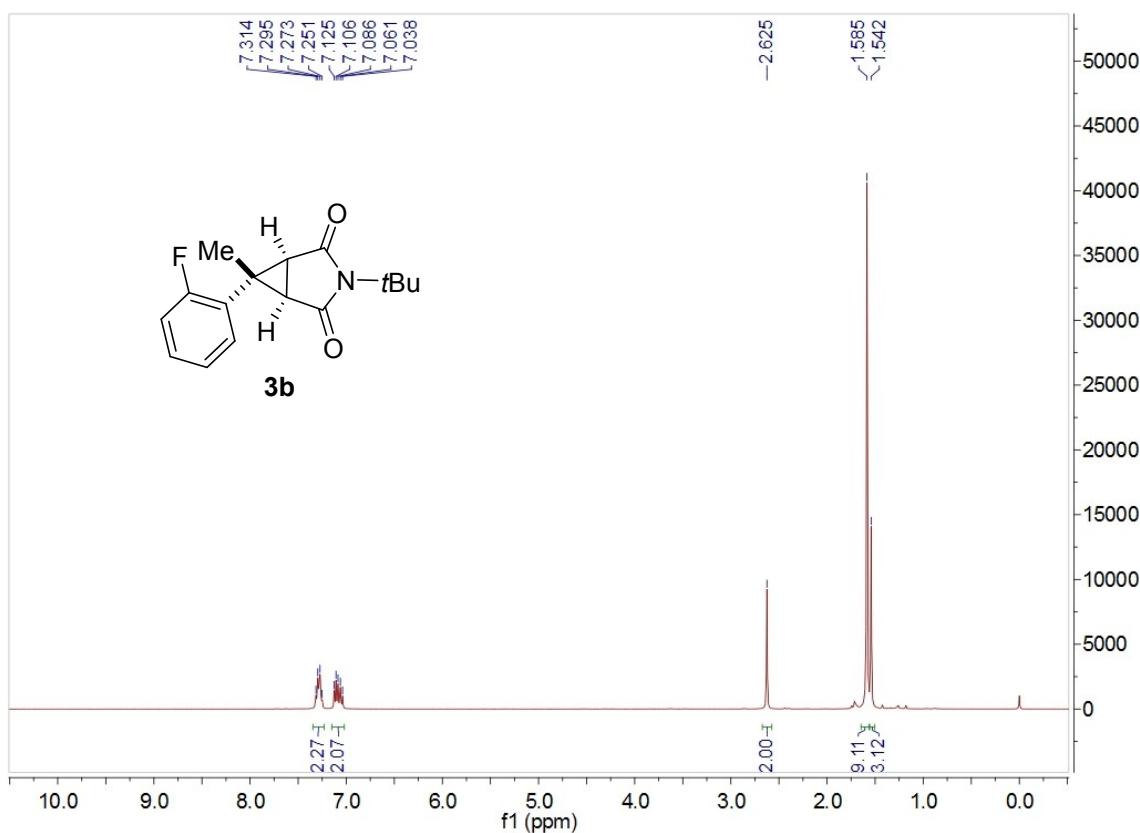
¹H NMR (400 MHz, CDCl₃) spectrum for 3a



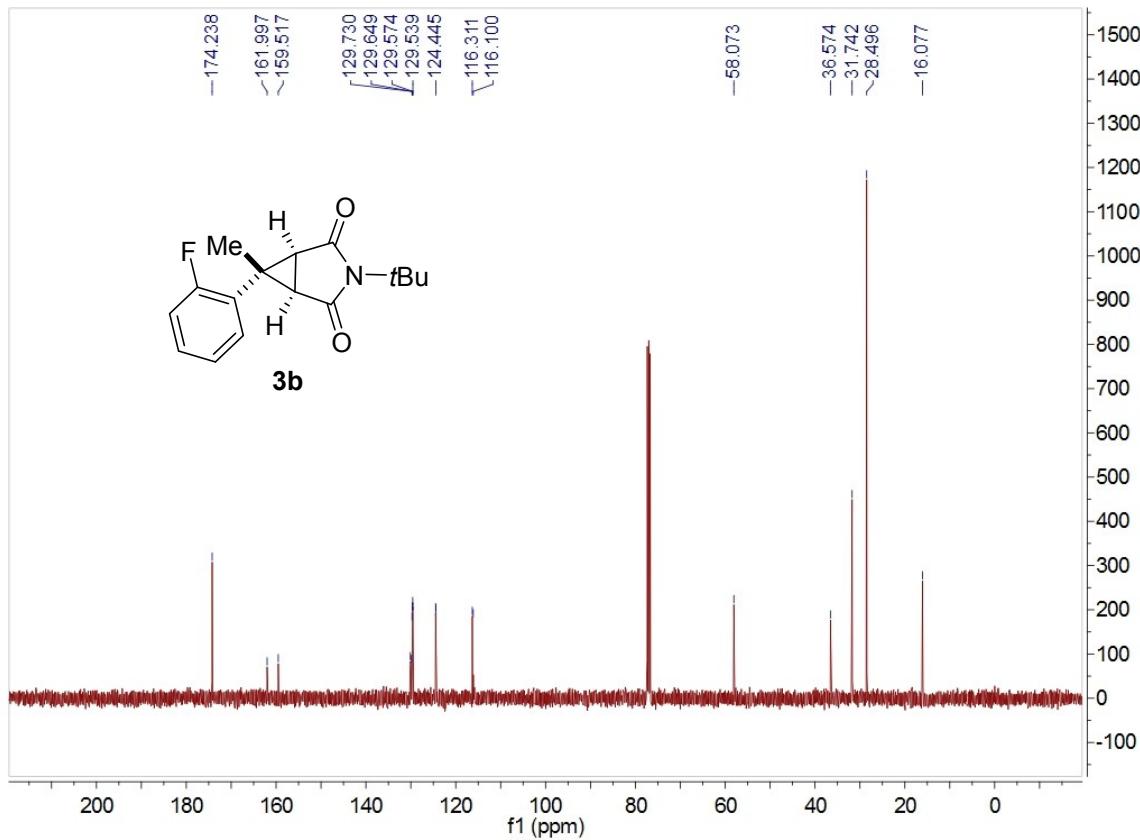
¹³C NMR (100 MHz, CDCl₃) spectrum for 3a



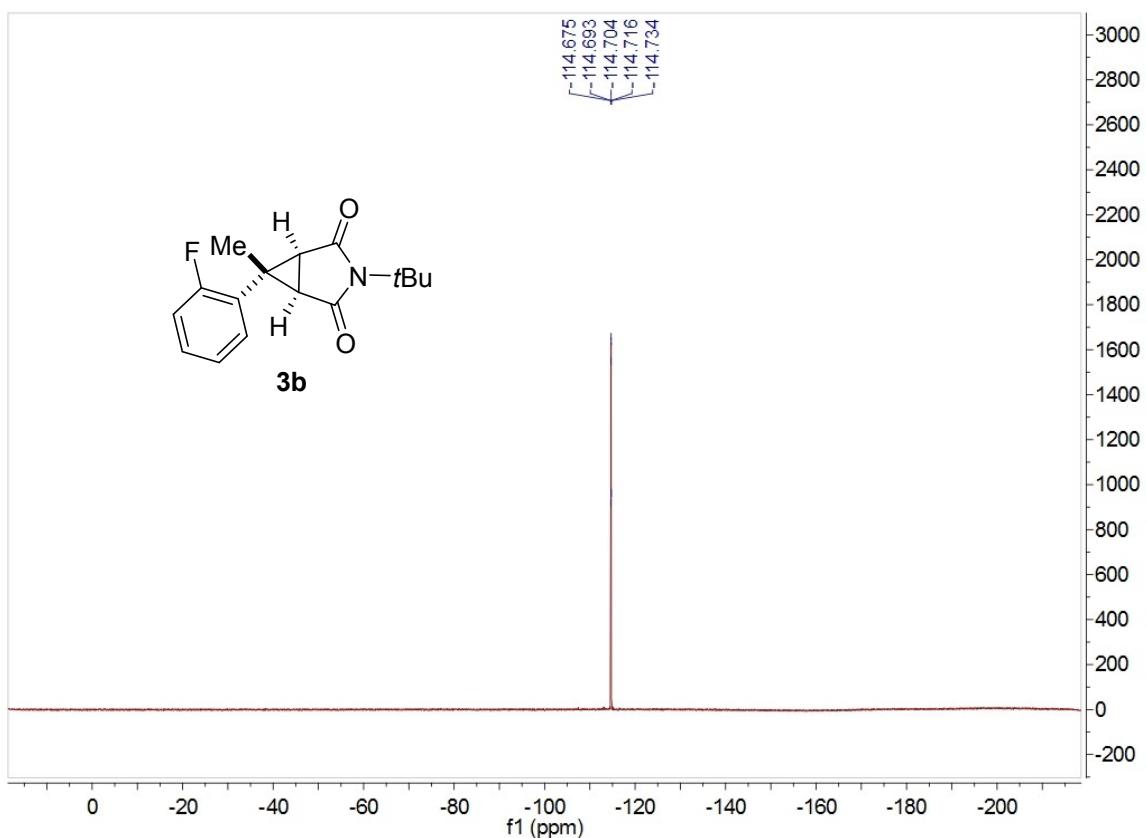
¹H NMR (400 MHz, CDCl₃) spectrum for 3b



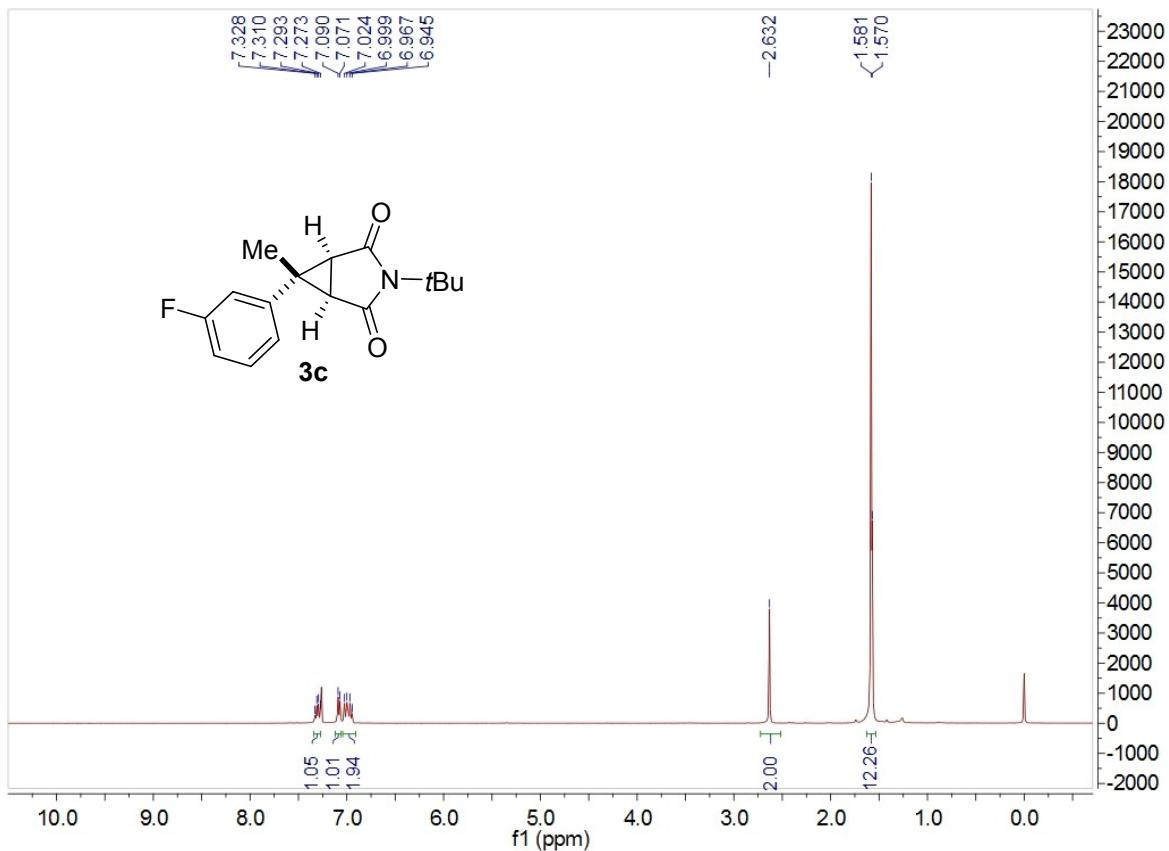
¹³C NMR (100 MHz, CDCl₃) spectrum for 3b



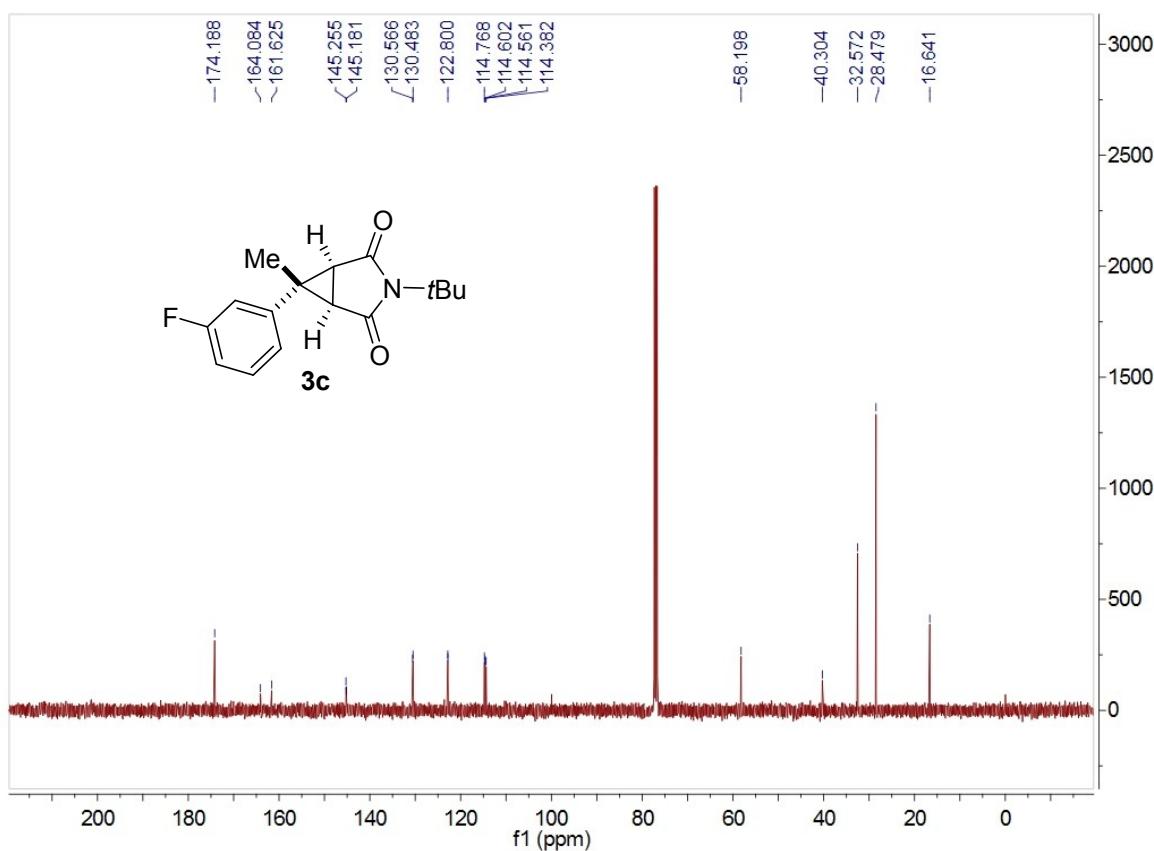
¹⁹F NMR (376 MHz, CDCl₃) spectrum for 3b



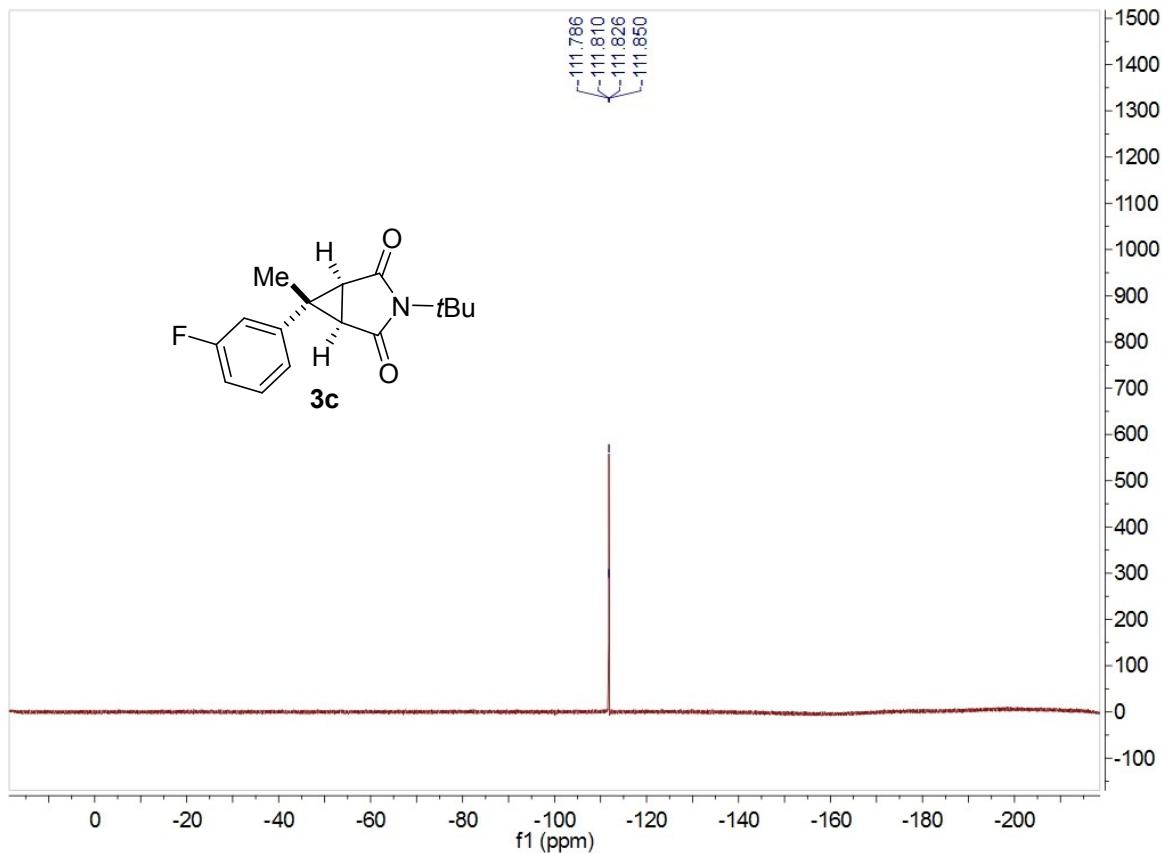
¹H NMR (400 MHz, CDCl₃) spectrum for 3c



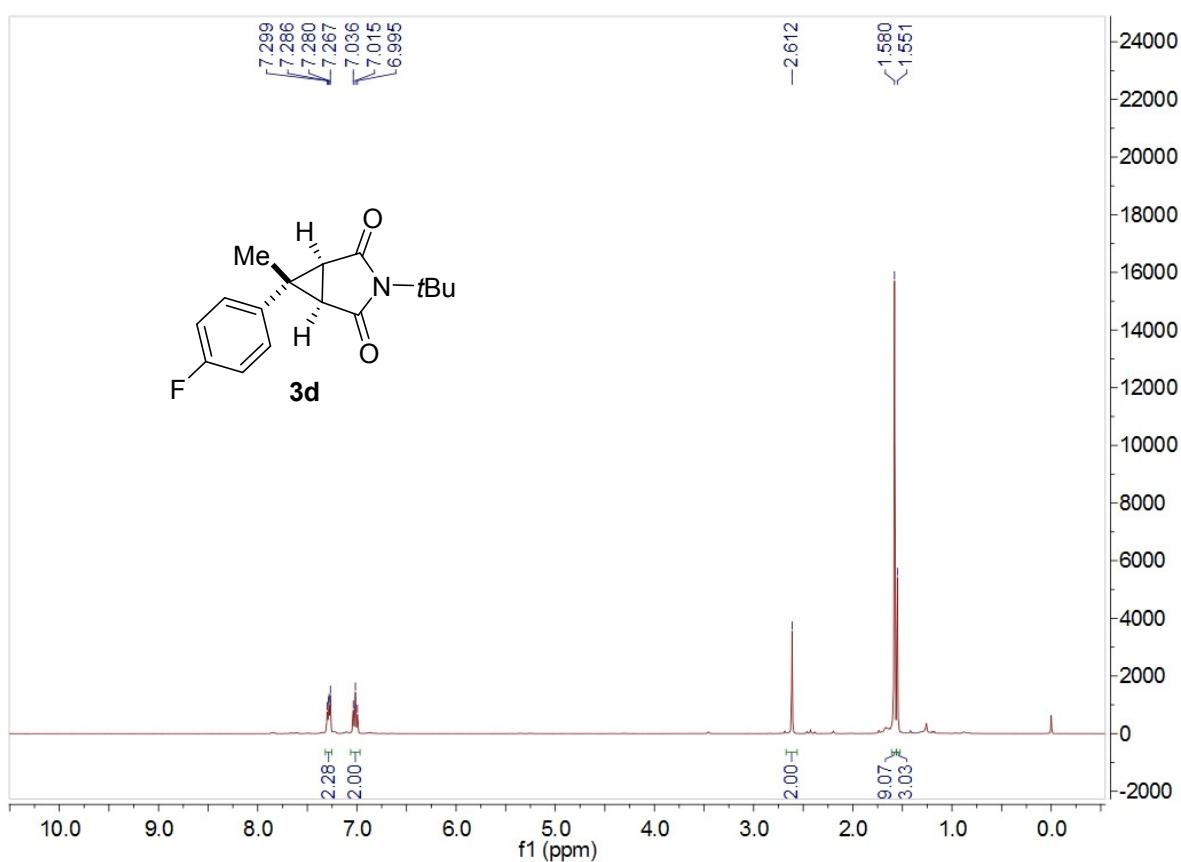
¹³C NMR (100 MHz, CDCl₃) spectrum for 3c



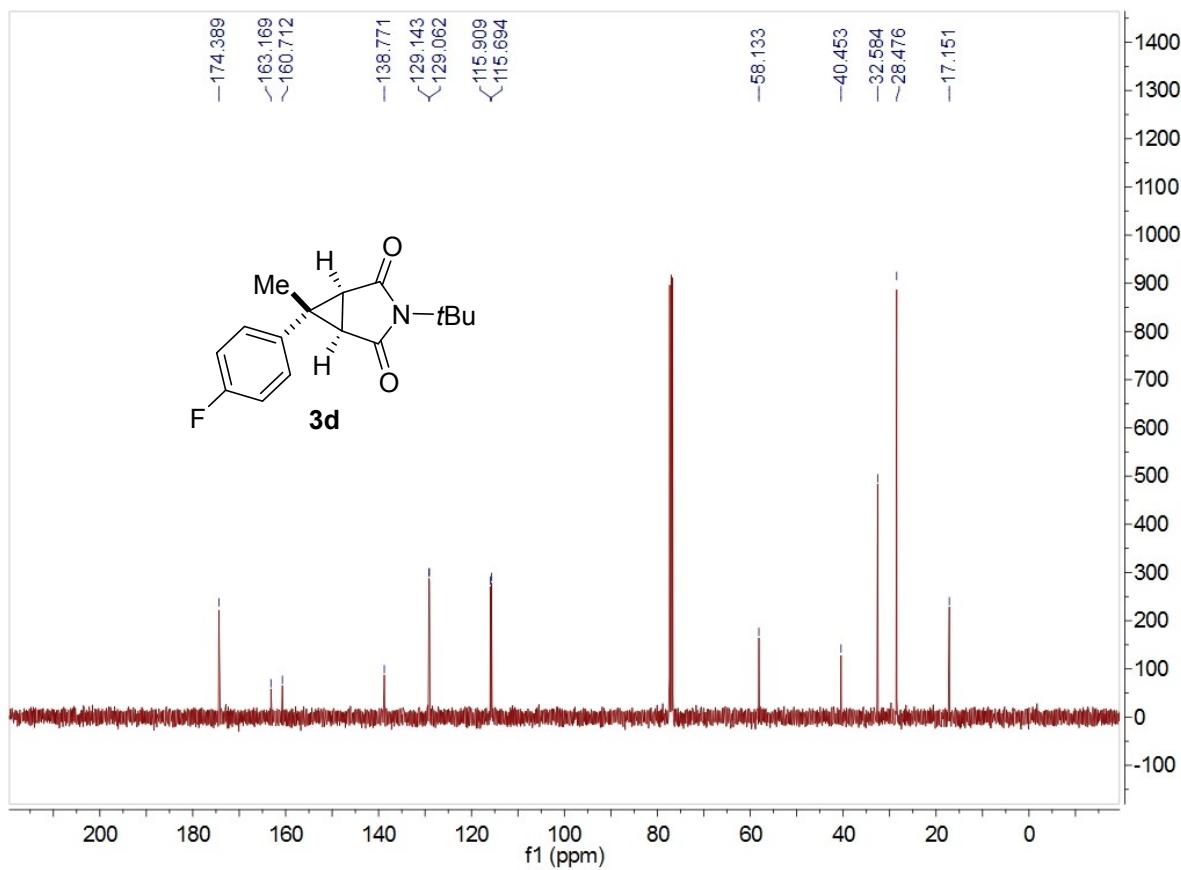
¹⁹F NMR (376 MHz, CDCl₃) spectrum for 3c



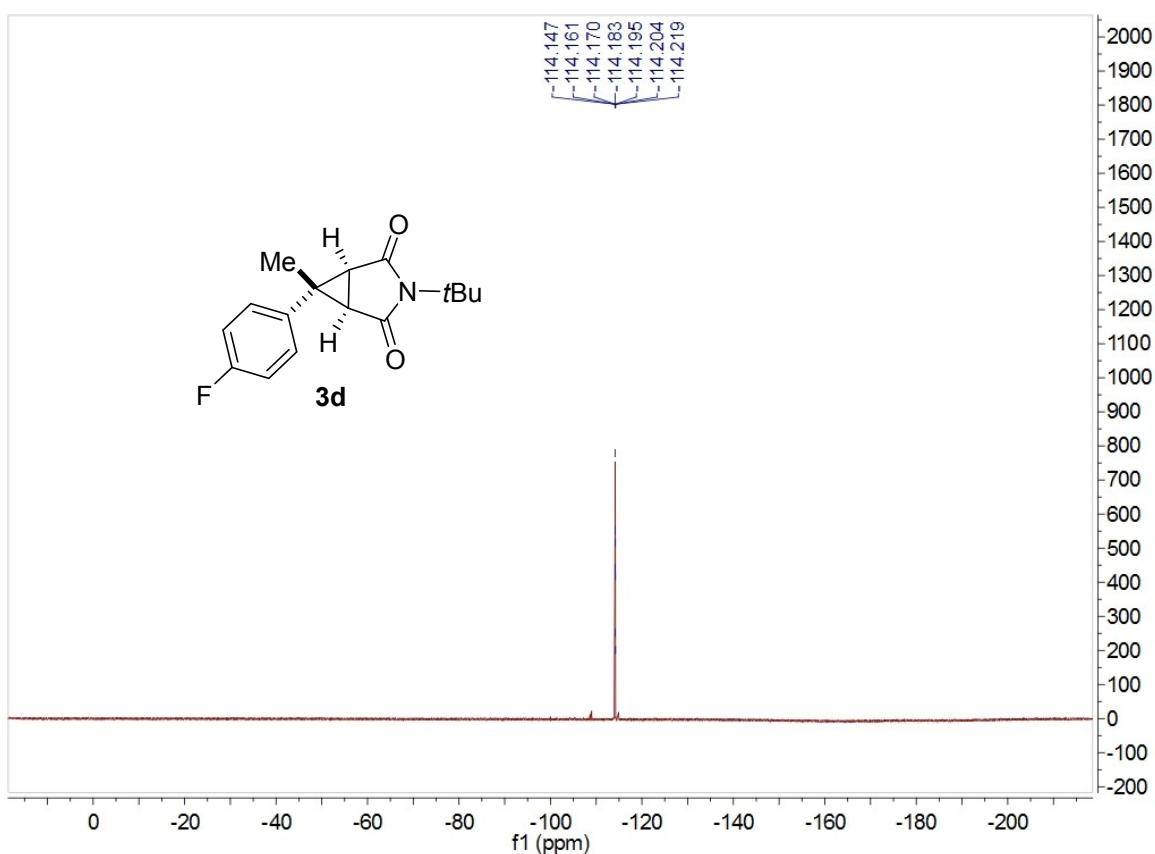
¹H NMR (400 MHz, CDCl₃) spectrum for 3d



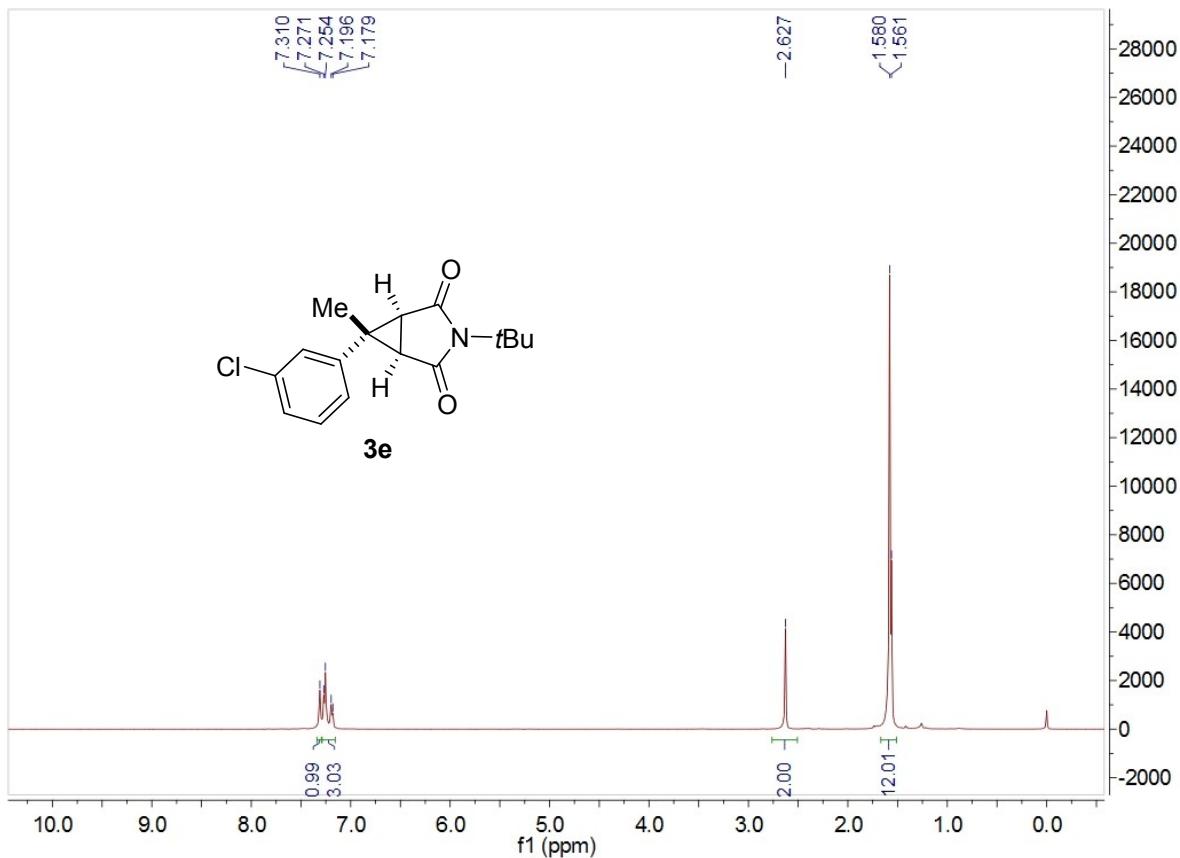
¹³C NMR (100 MHz, CDCl₃) spectrum for 3d



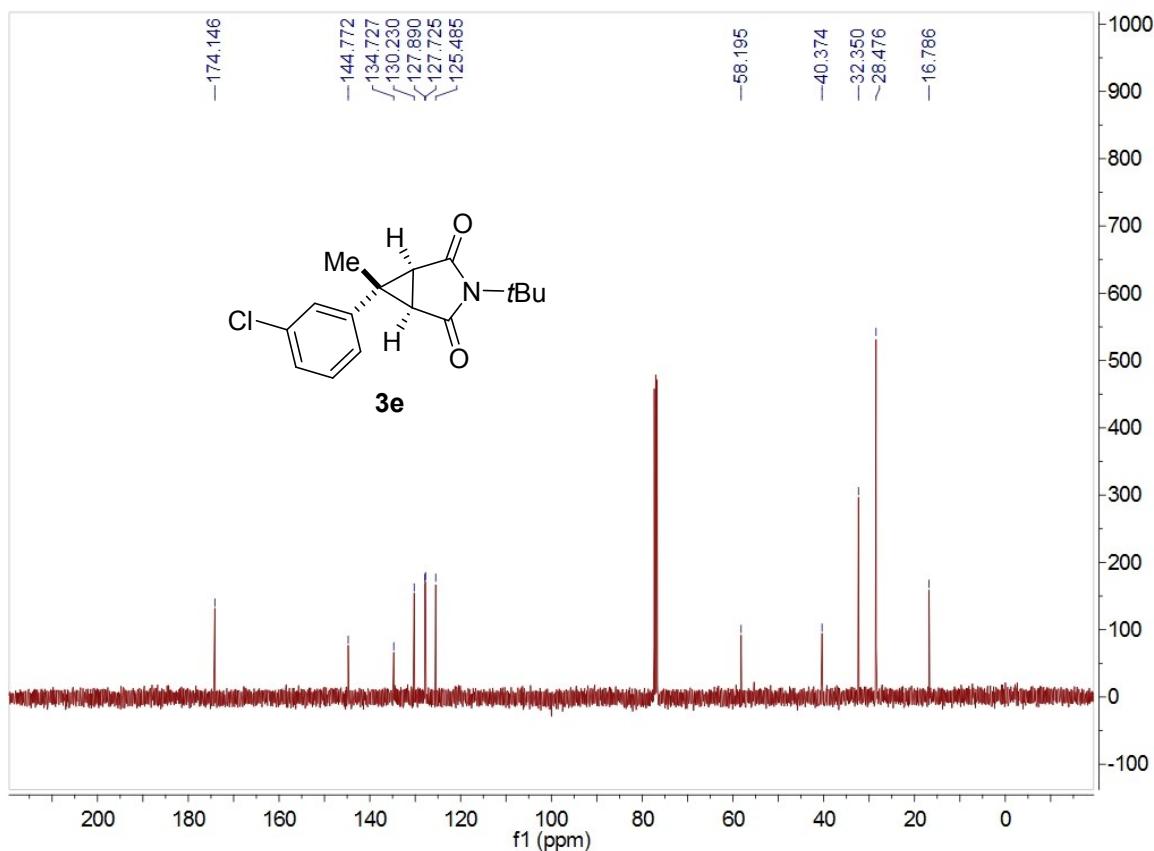
¹⁹F NMR (376 MHz, CDCl₃) spectrum for 3d



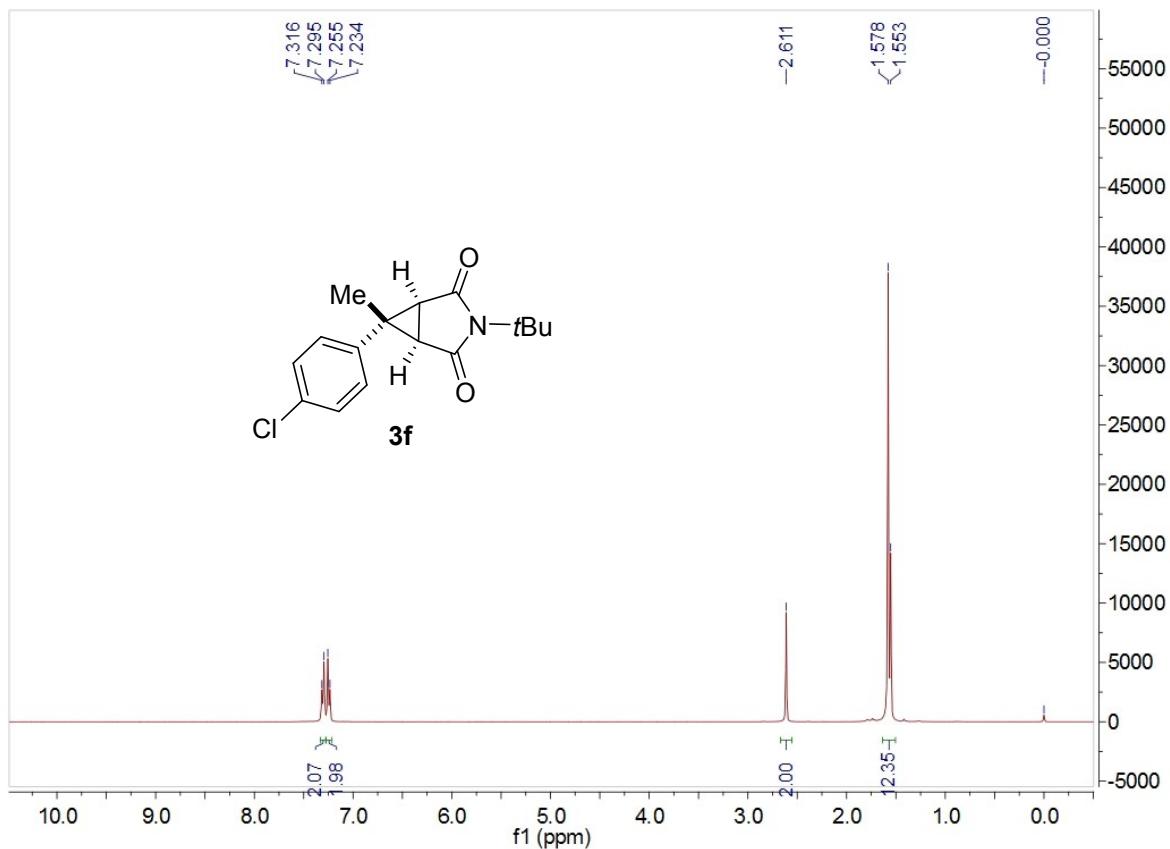
¹H NMR (400 MHz, CDCl₃) spectrum for 3e



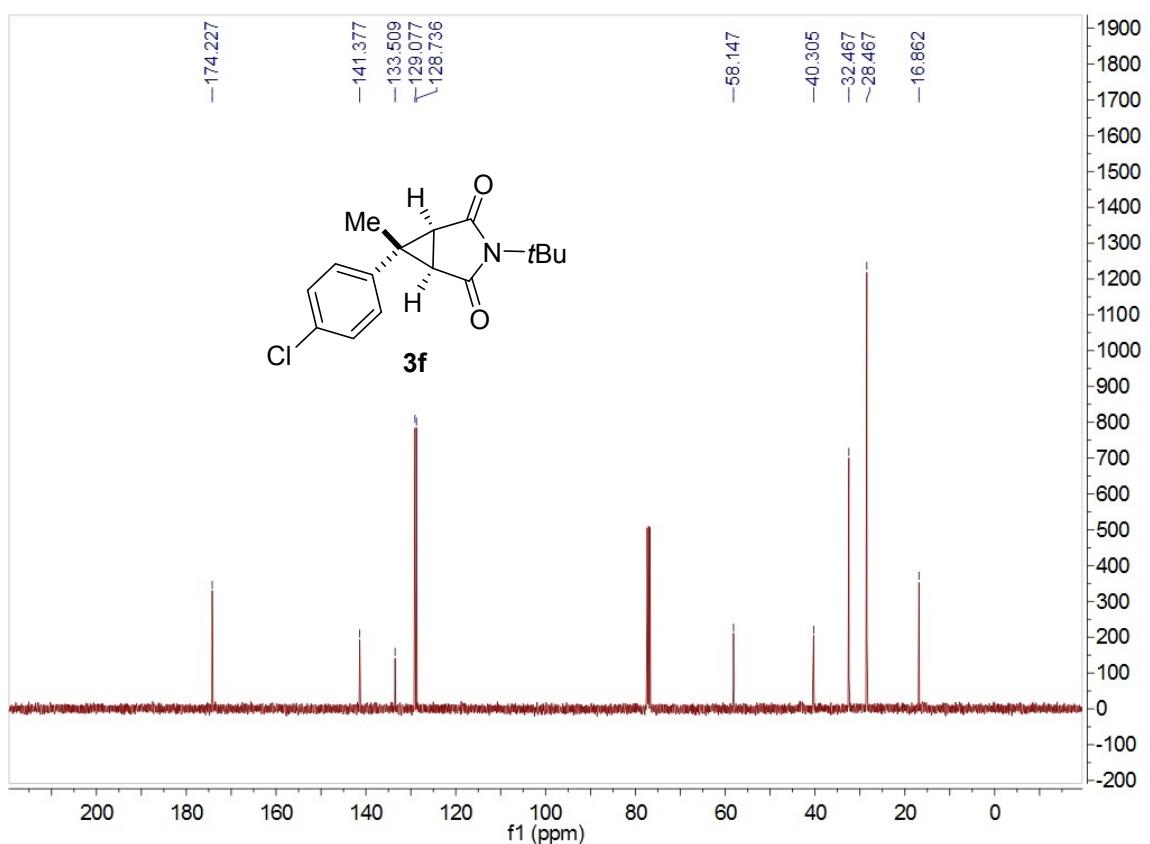
¹³C NMR (100 MHz, CDCl₃) spectrum for 3e



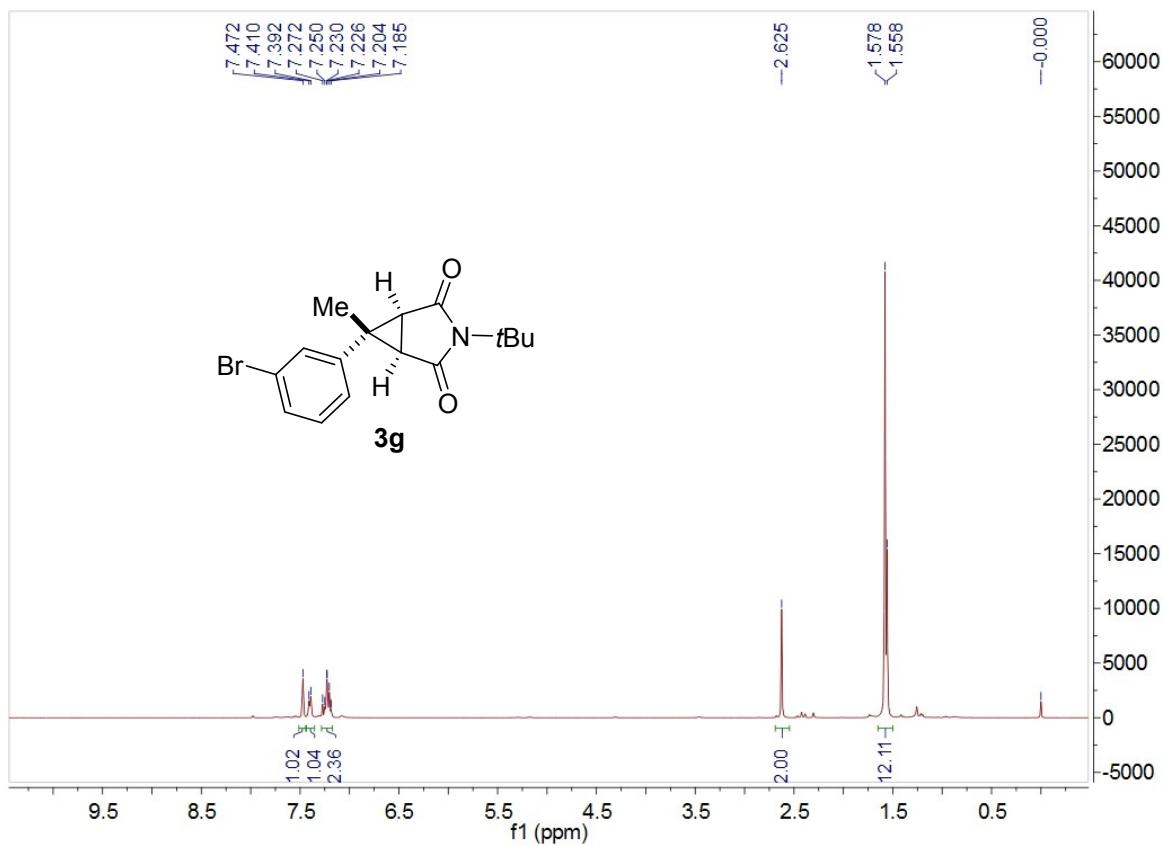
¹H NMR (400 MHz, CDCl₃) spectrum for 3f



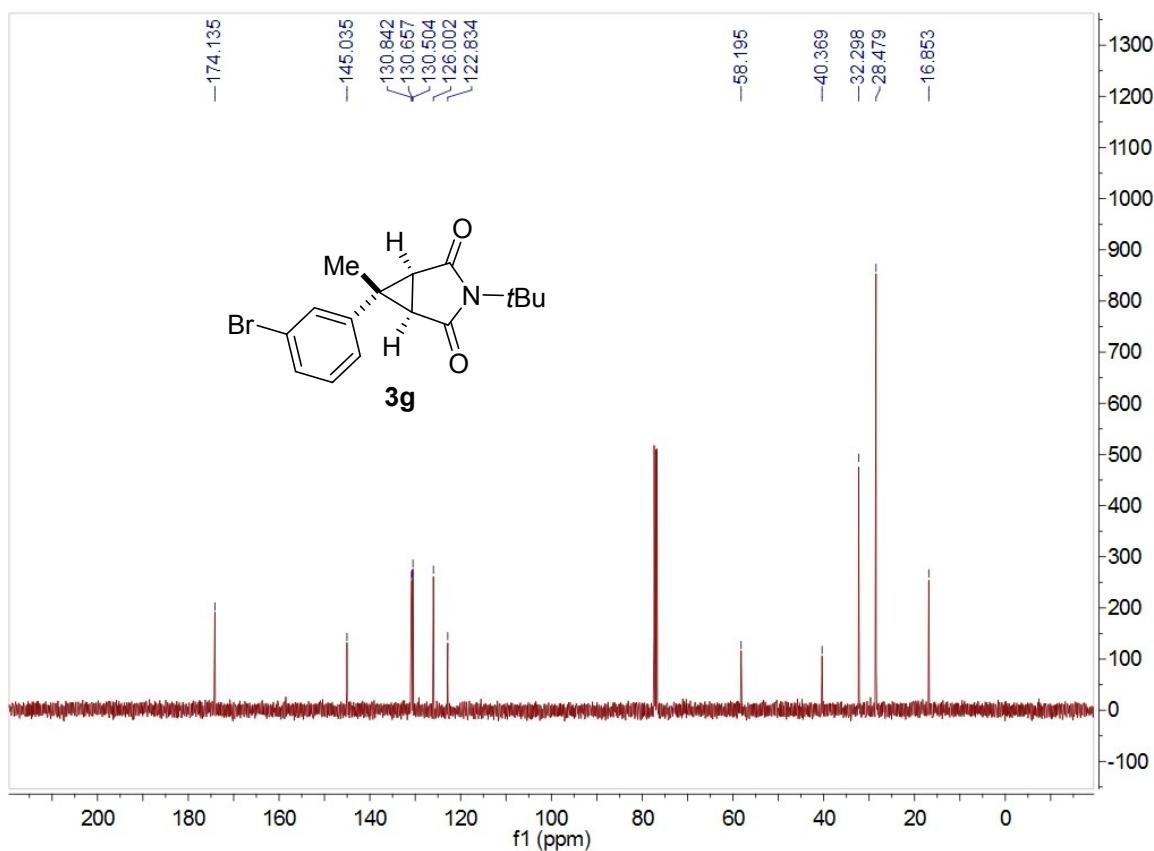
¹³C NMR (100 MHz, CDCl₃) spectrum for 3f



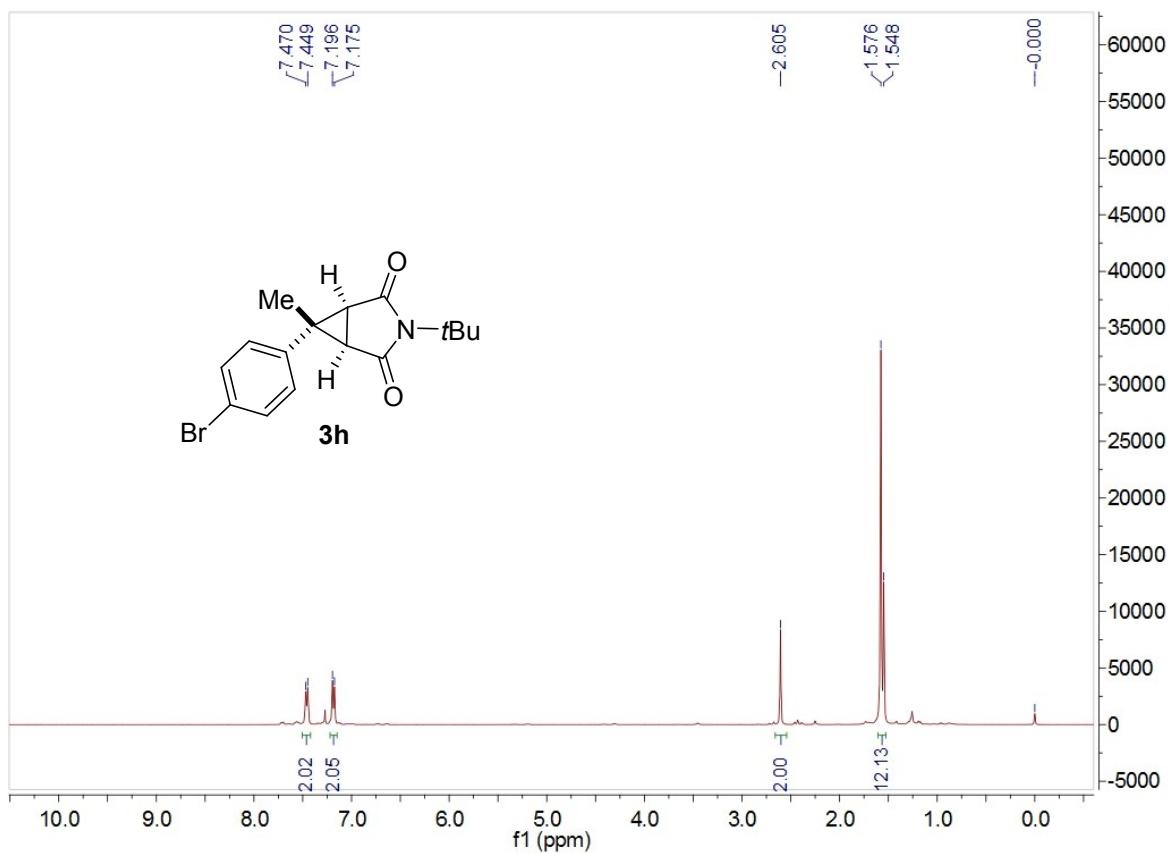
¹H NMR (400 MHz, CDCl₃) spectrum for 3g



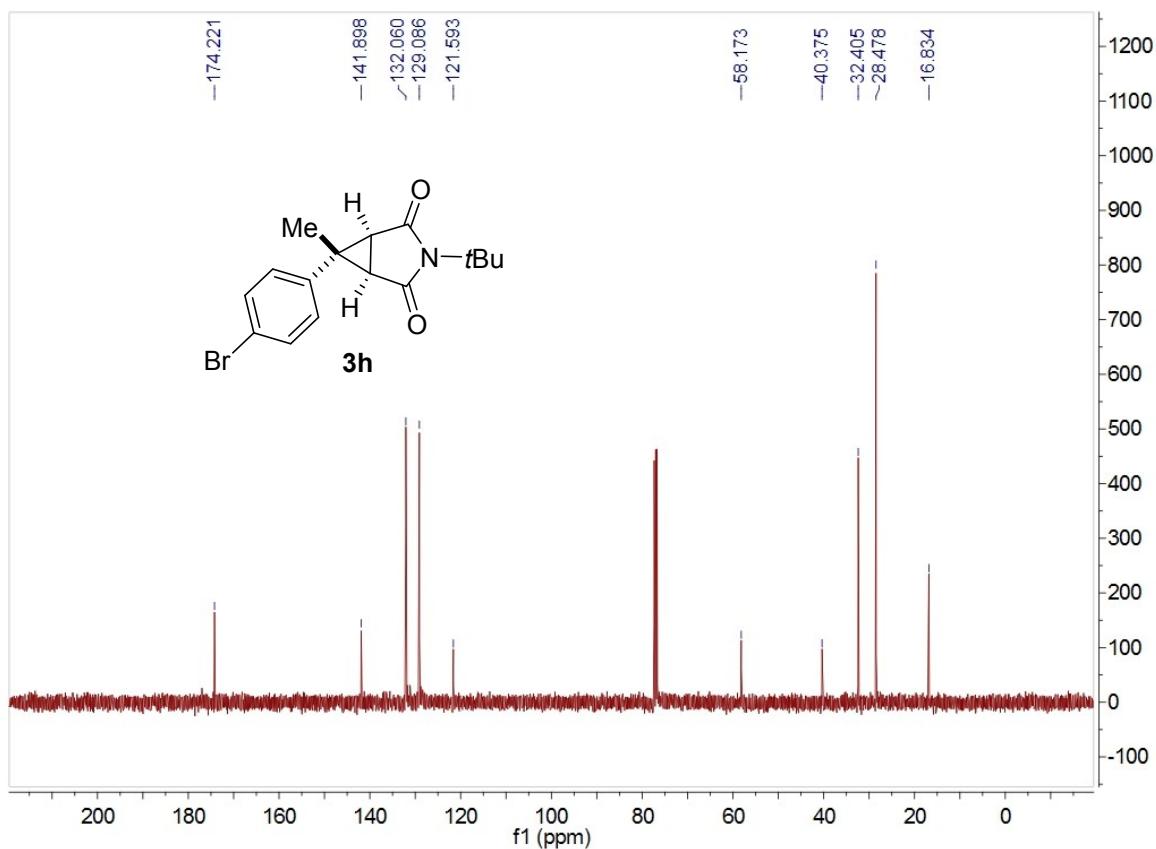
¹³C NMR (100 MHz, CDCl₃) spectrum for 3g



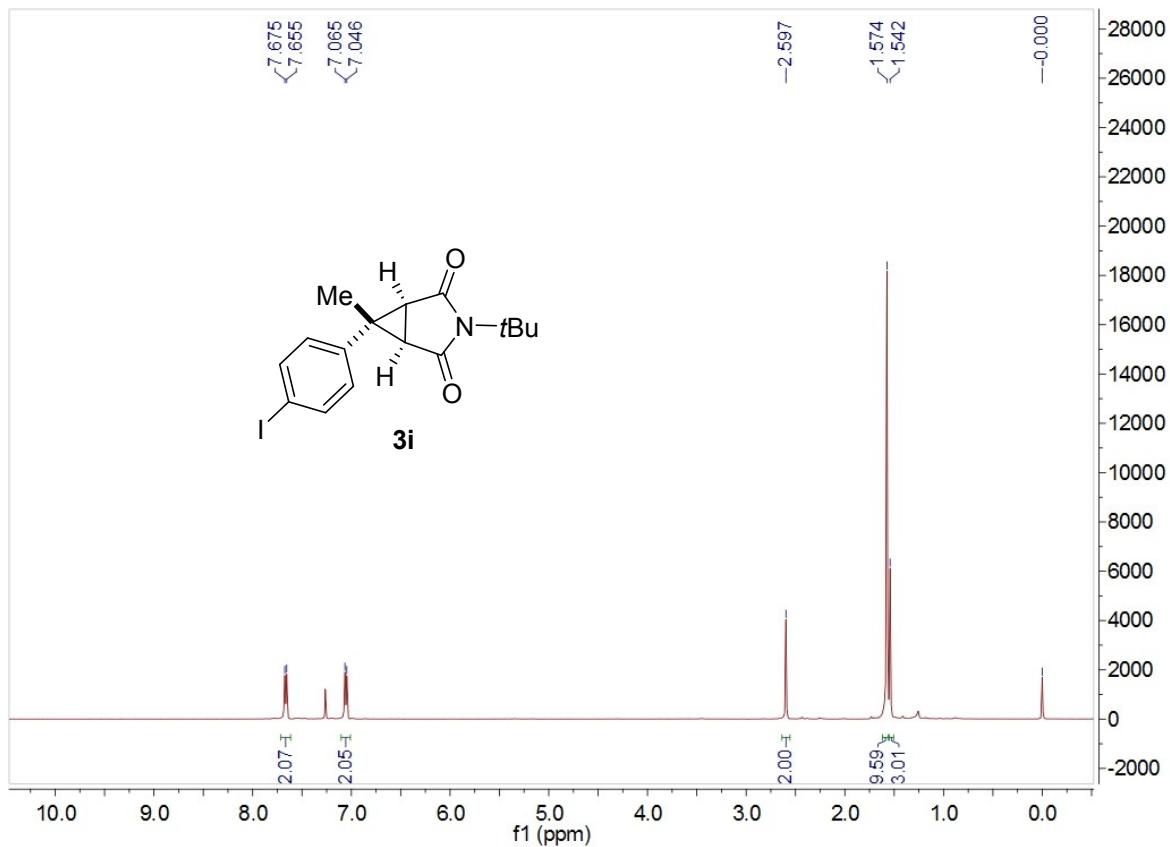
¹H NMR (400 MHz, CDCl₃) spectrum for 3h



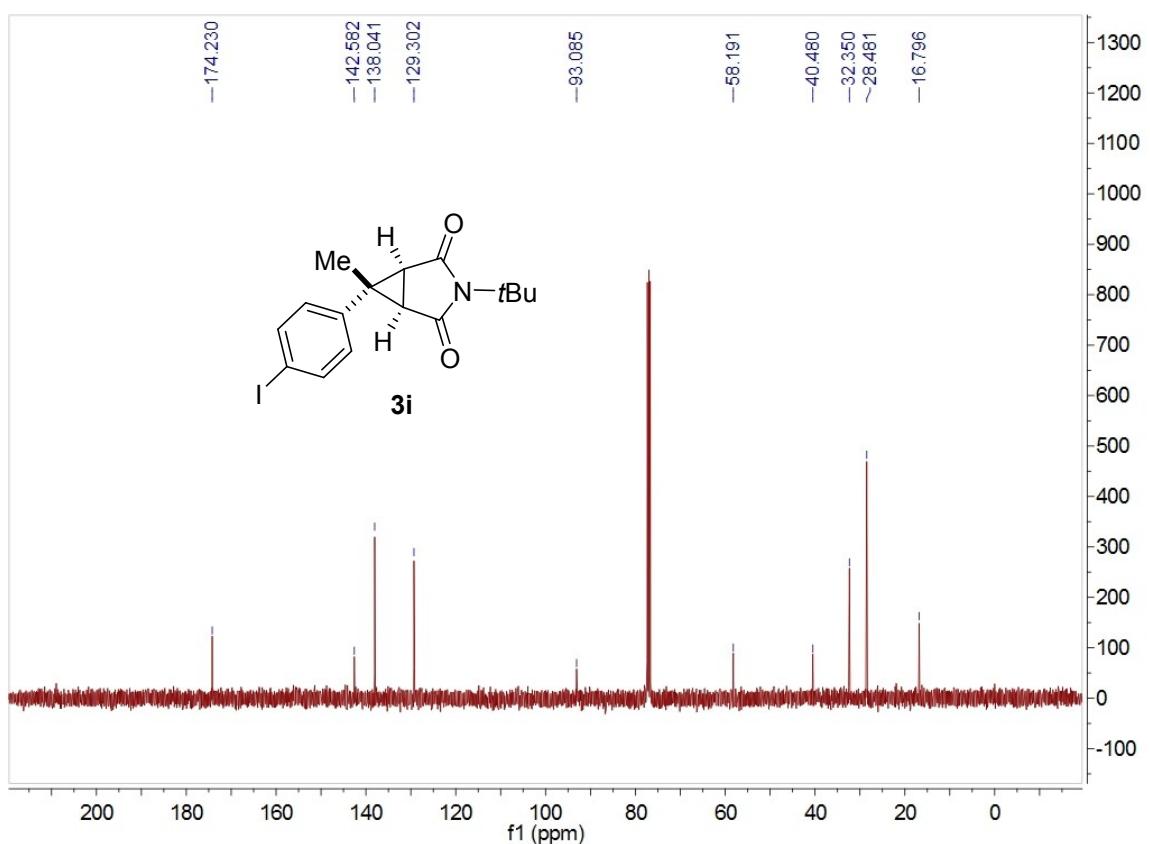
¹³C NMR (100 MHz, CDCl₃) spectrum for 3h



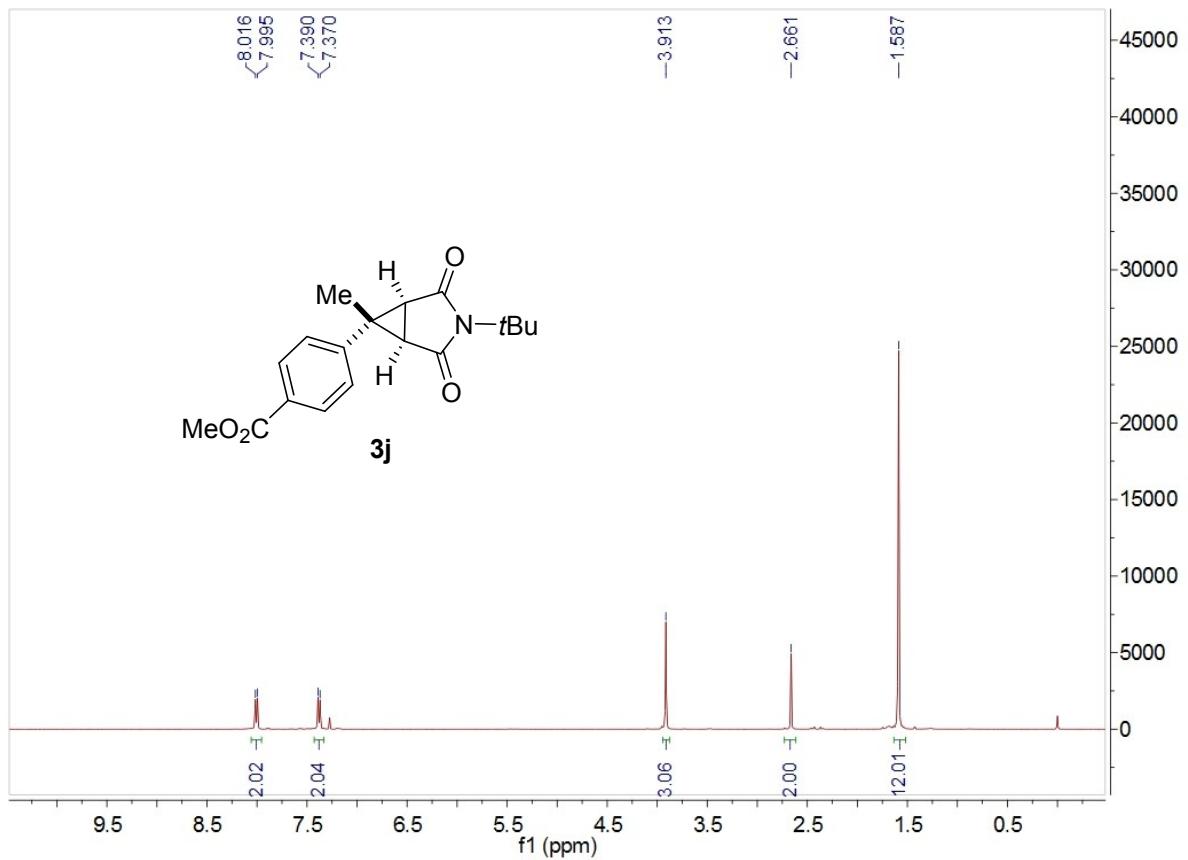
¹H NMR (400 MHz, CDCl₃) spectrum for 3i



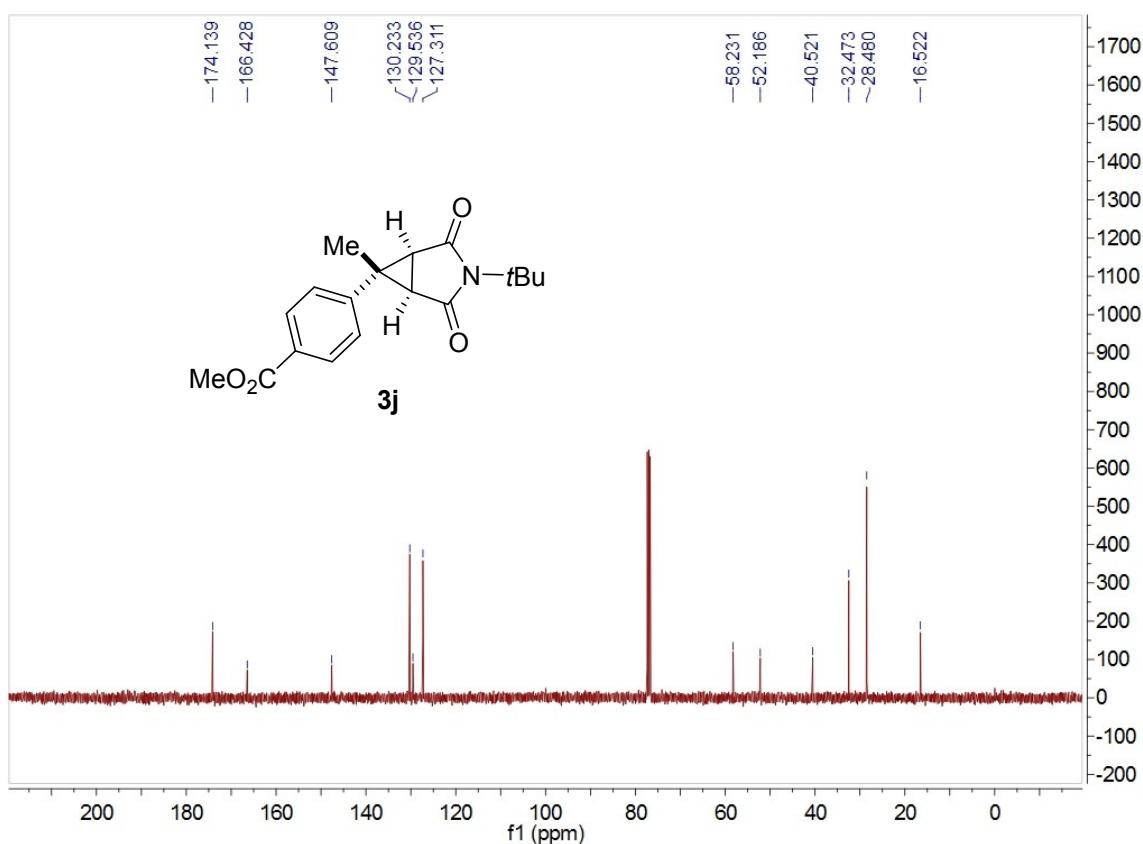
¹³C NMR (100 MHz, CDCl₃) spectrum for 3i



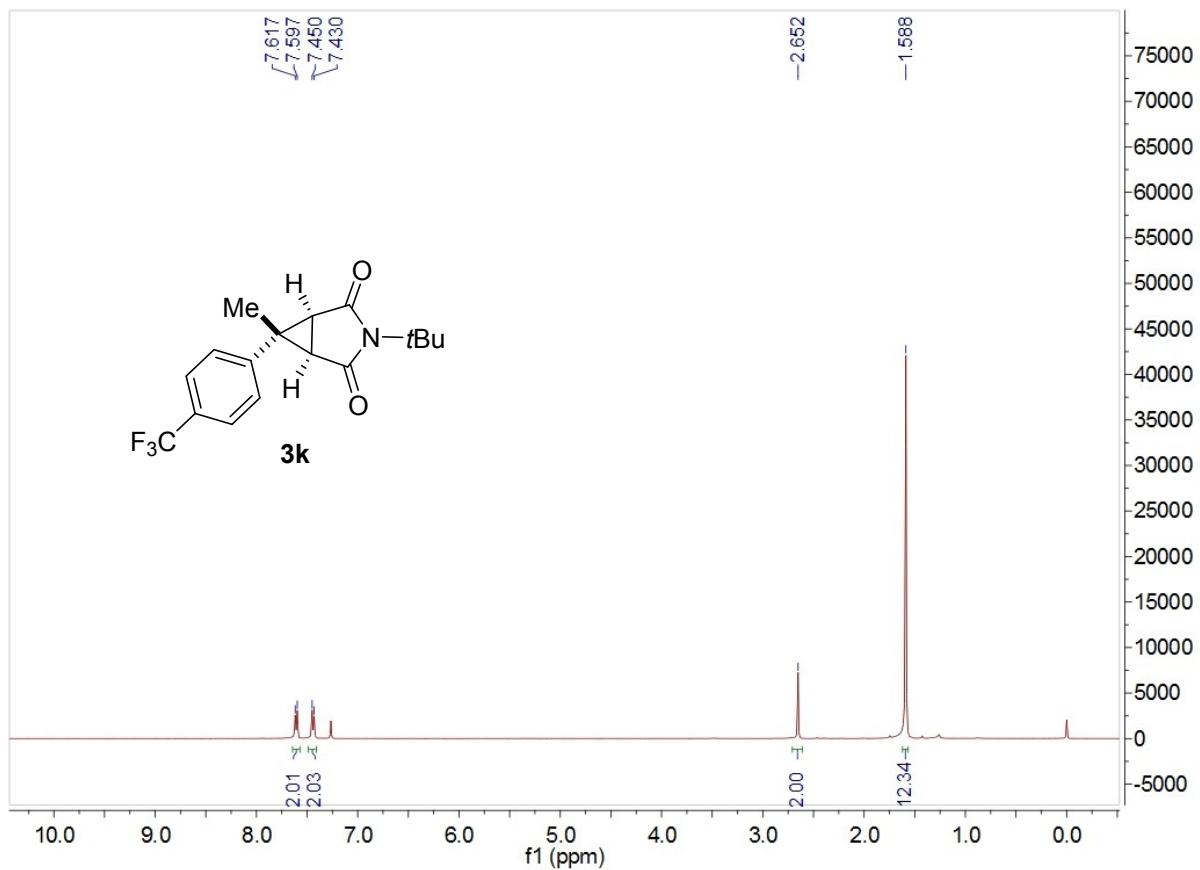
¹H NMR (400 MHz, CDCl₃) spectrum for 3j



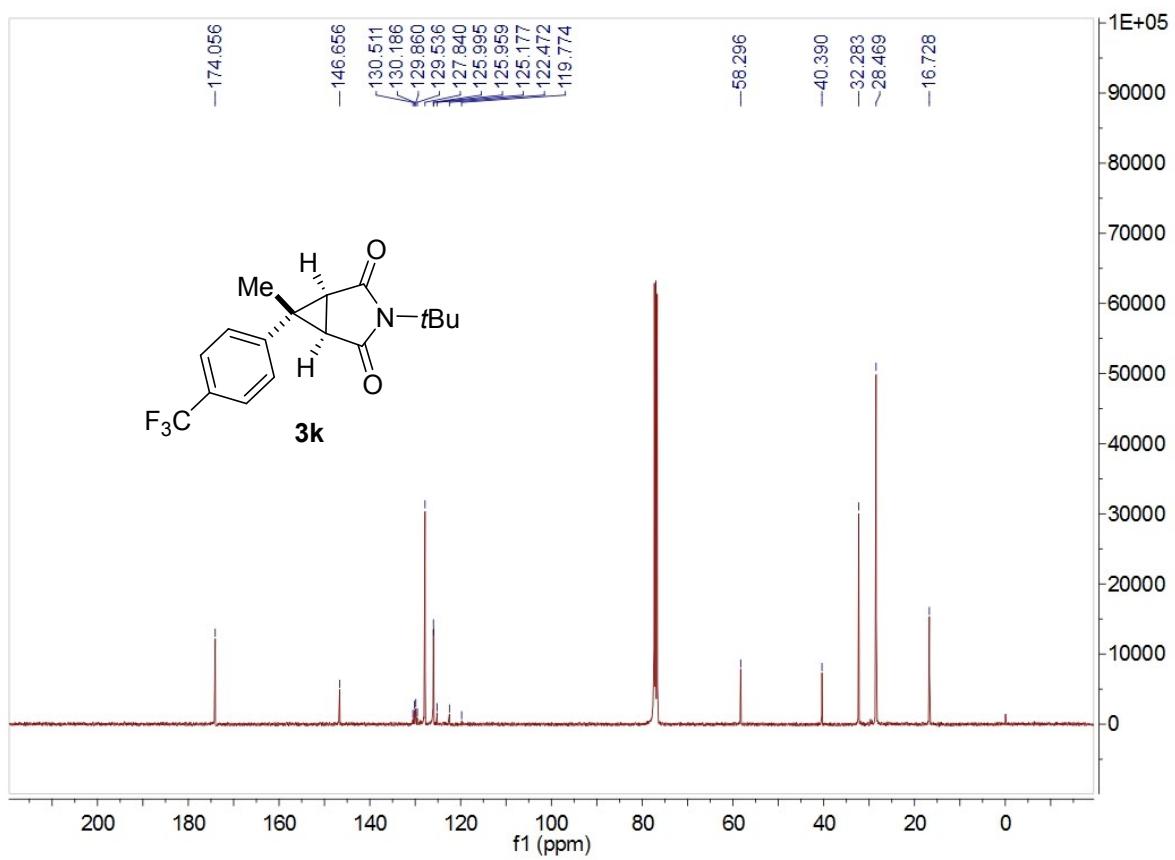
¹³C NMR (100 MHz, CDCl₃) spectrum for 3j



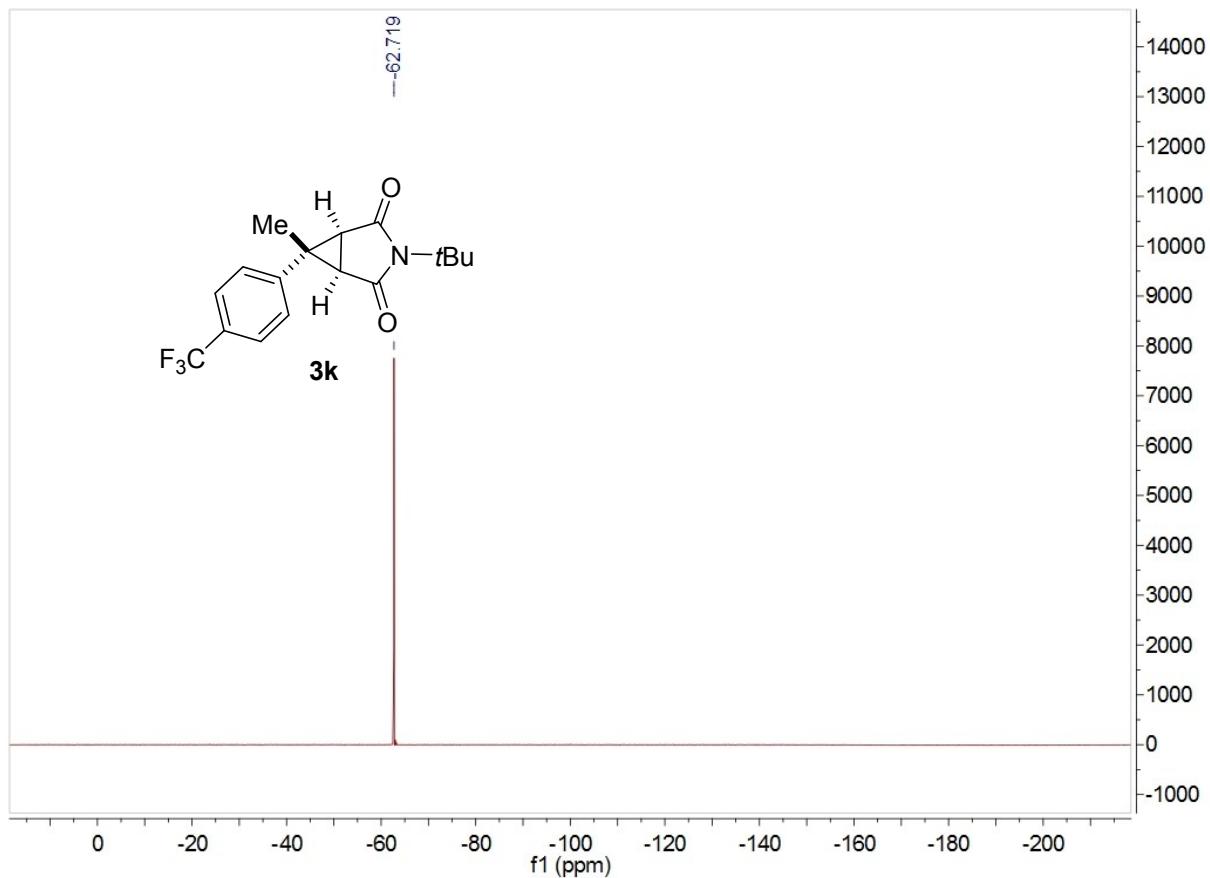
¹H NMR (400 MHz, CDCl₃) spectrum for 3k



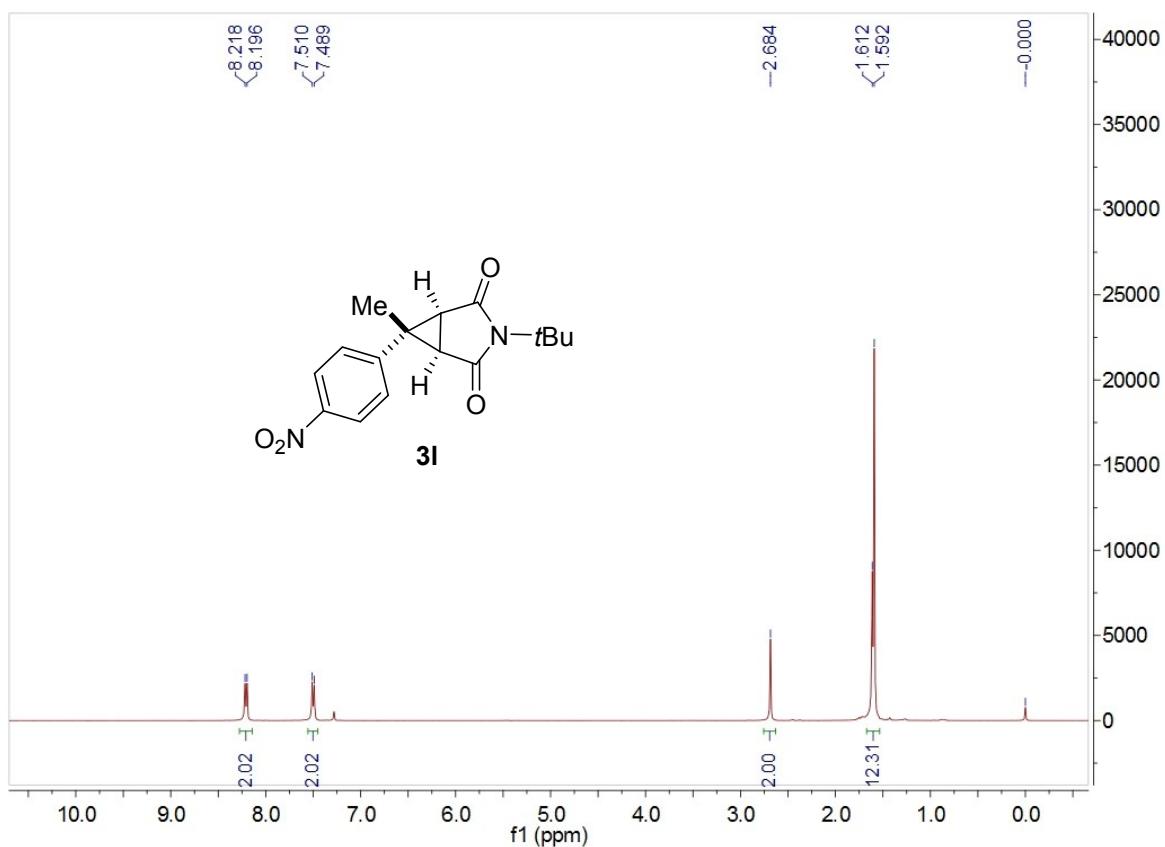
¹³C NMR (100 MHz, CDCl₃) spectrum for 3k



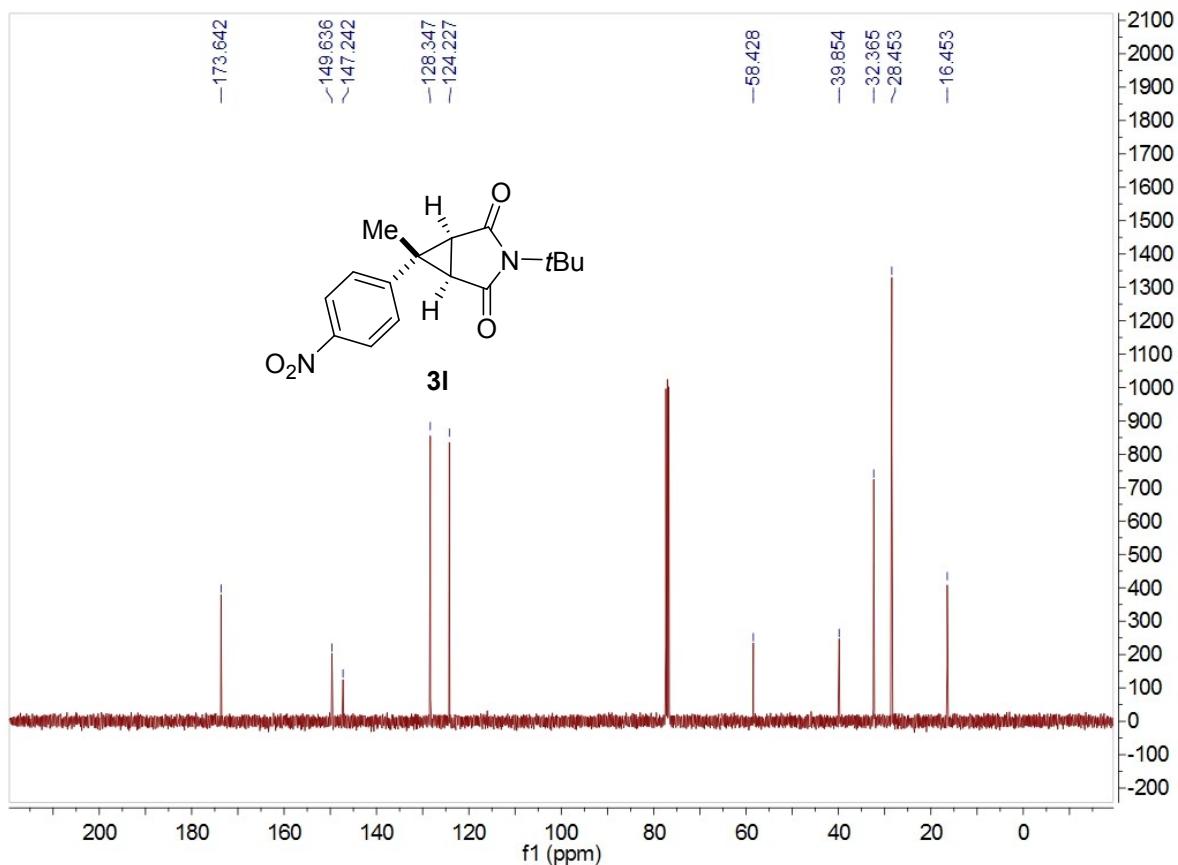
¹⁹F NMR (376 MHz, CDCl₃) spectrum for 3k



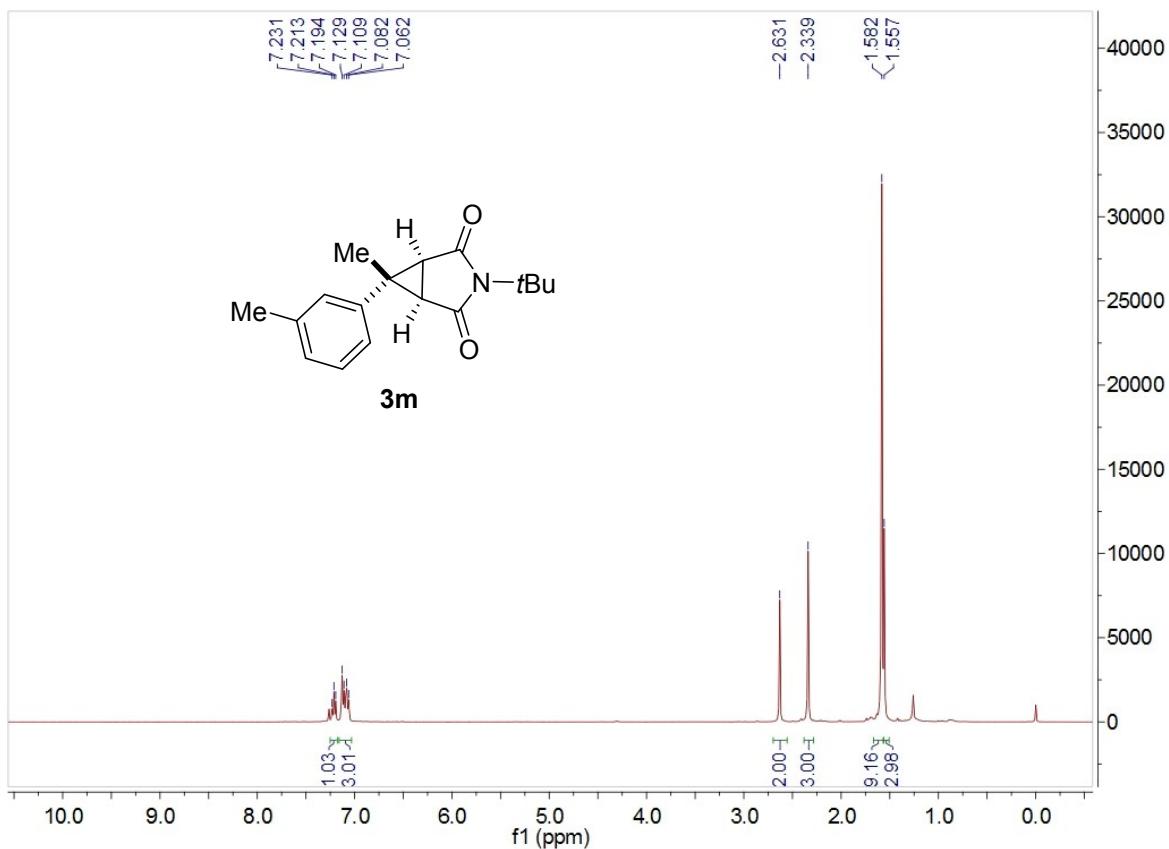
¹H NMR (400 MHz, CDCl₃) spectrum for 3l



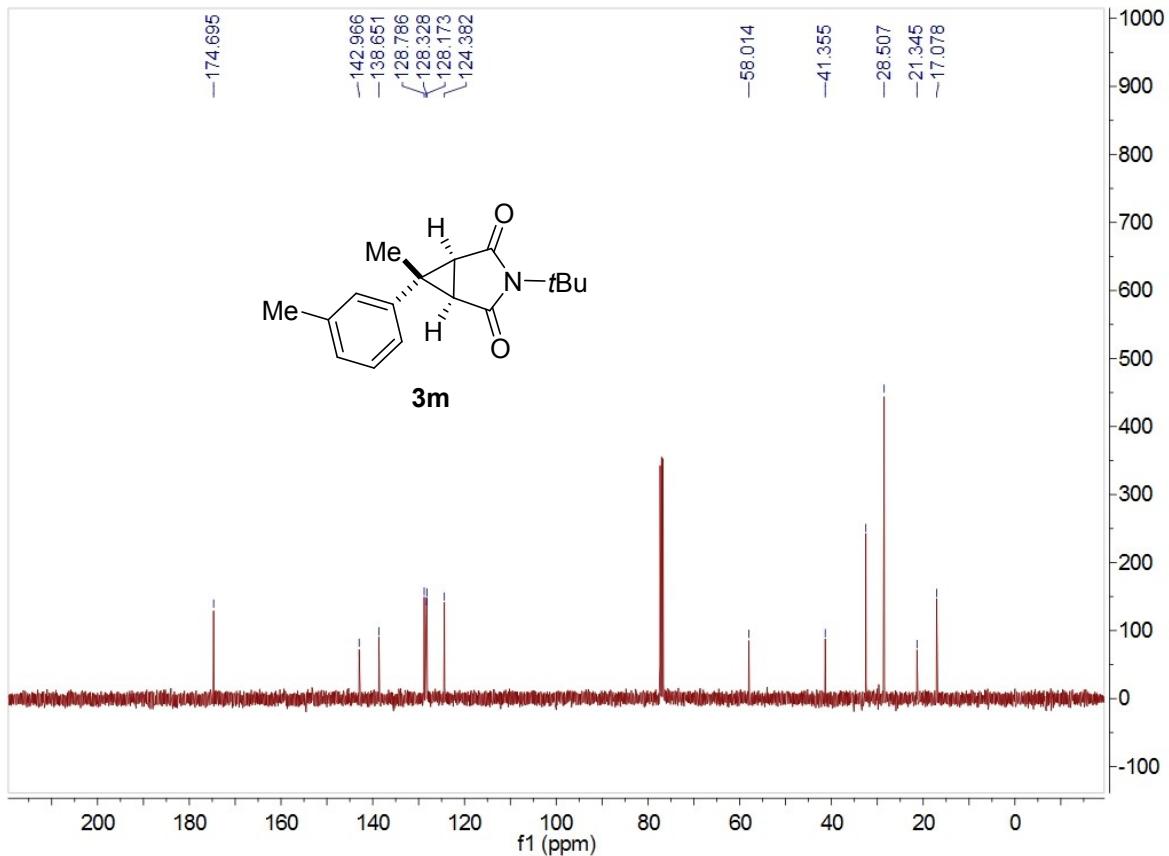
¹³C NMR (100 MHz, CDCl₃) spectrum for 3l



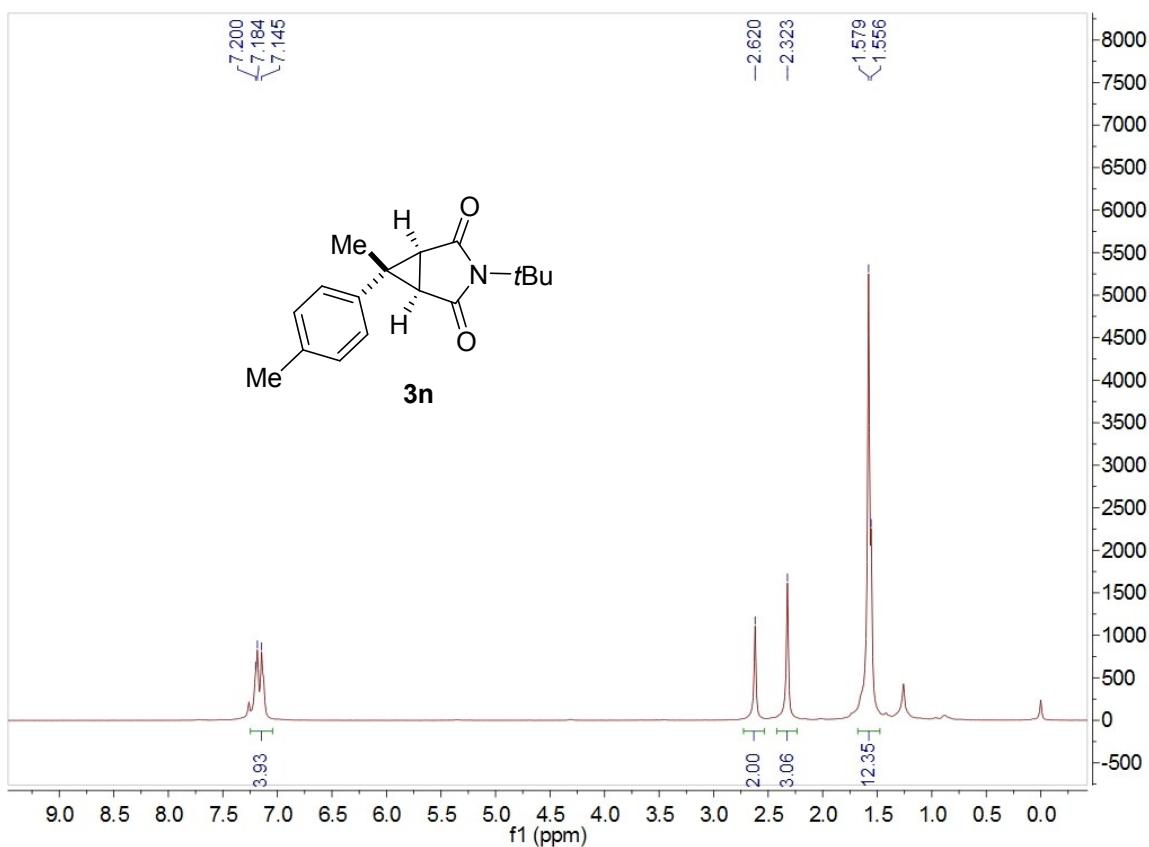
¹H NMR (400 MHz, CDCl₃) spectrum for 3m



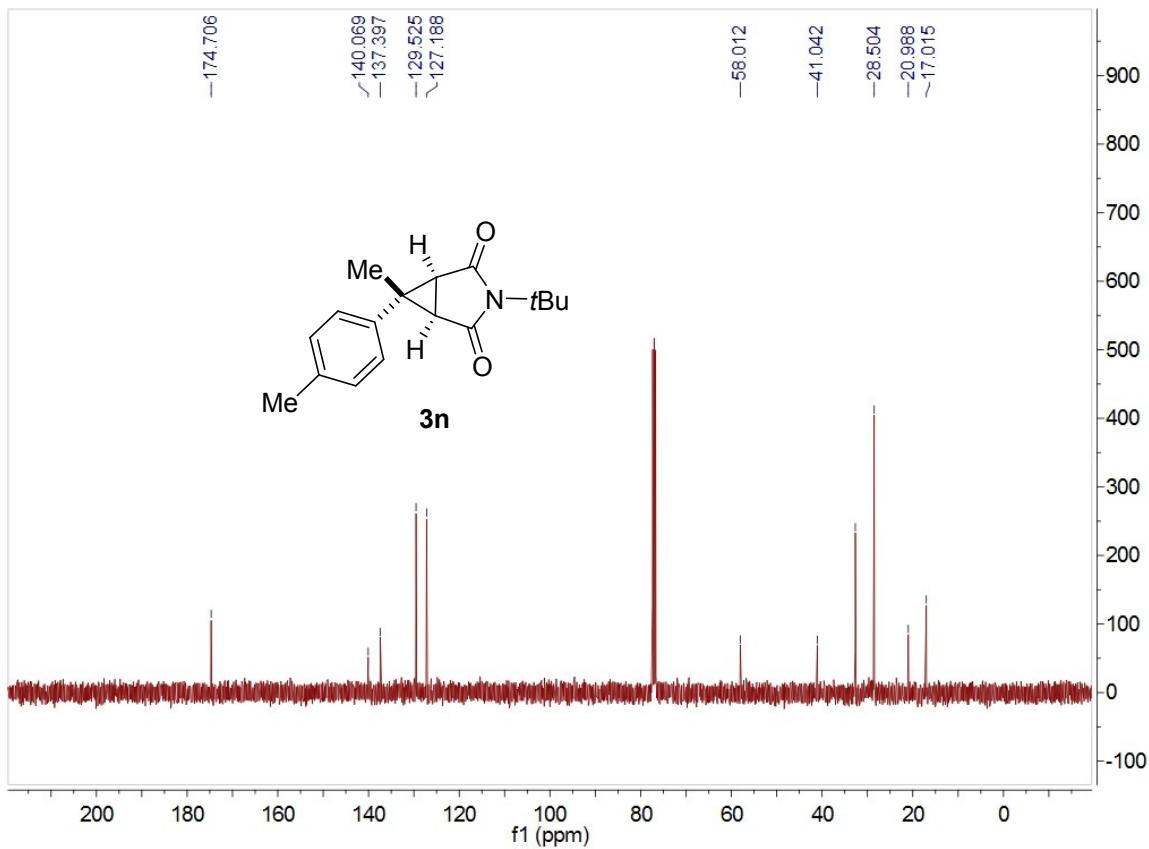
¹³C NMR (100 MHz, CDCl₃) spectrum for 3m



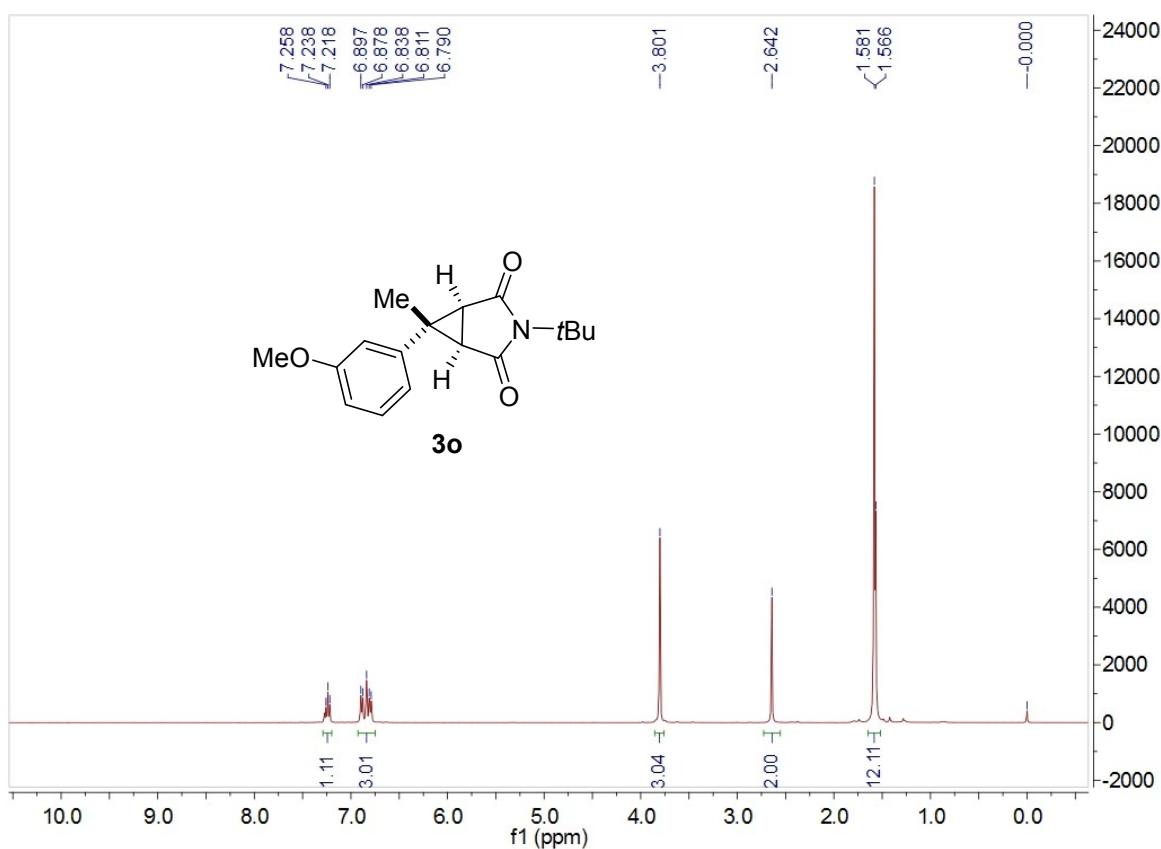
¹H NMR (400 MHz, CDCl₃) spectrum for 3n



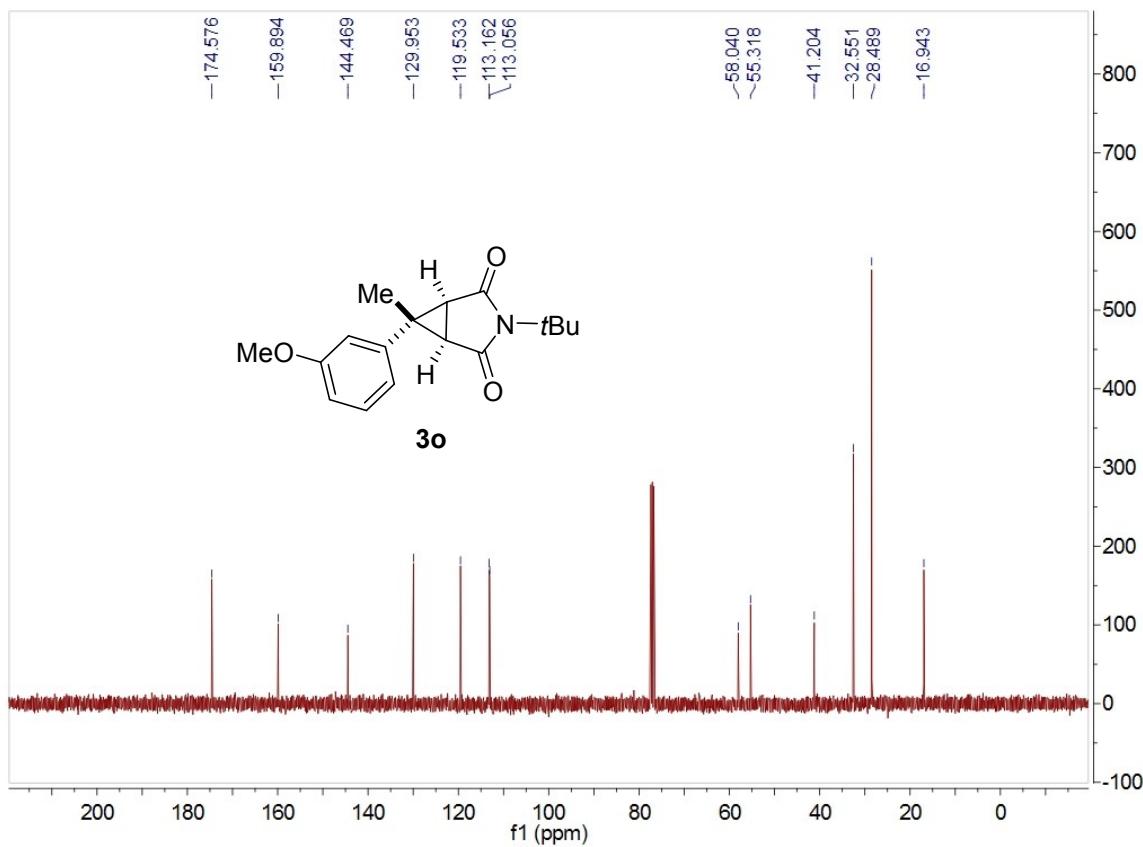
¹³C NMR (100 MHz, CDCl₃) spectrum for 3n



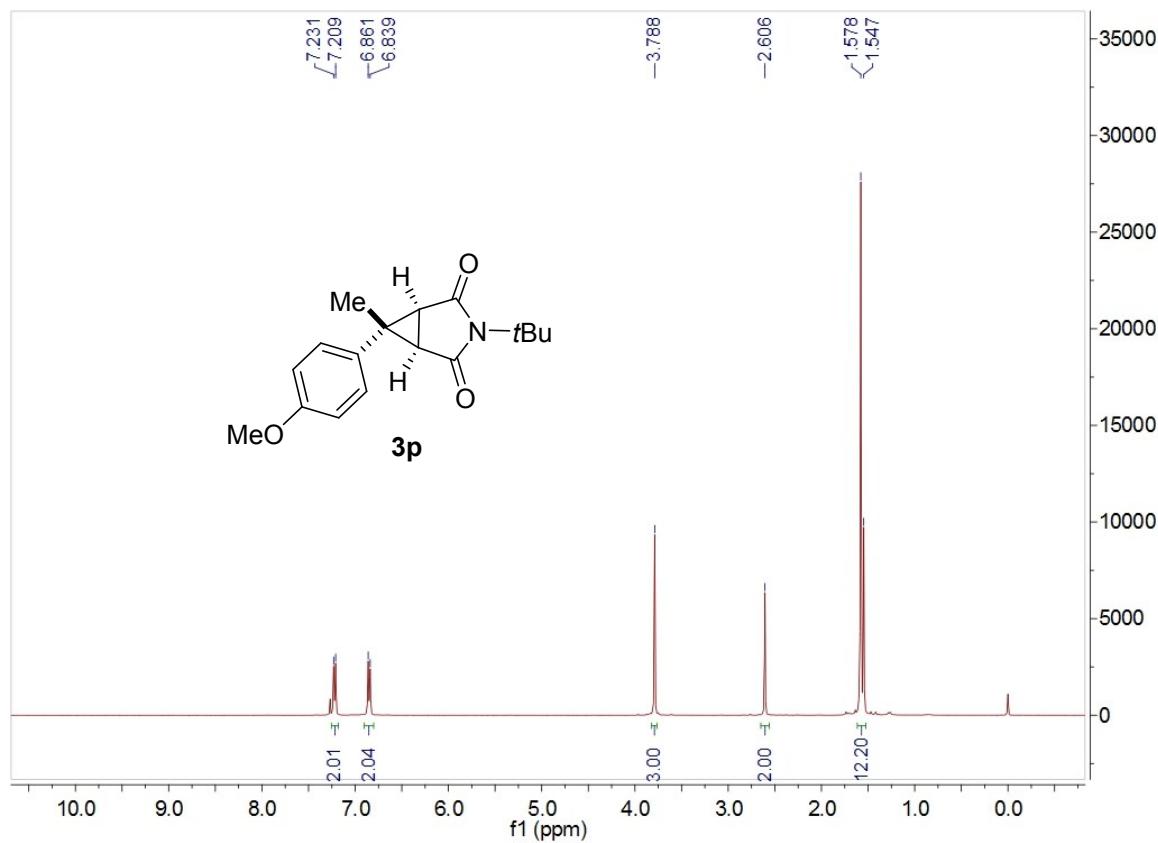
¹H NMR (400 MHz, CDCl₃) spectrum for 3o



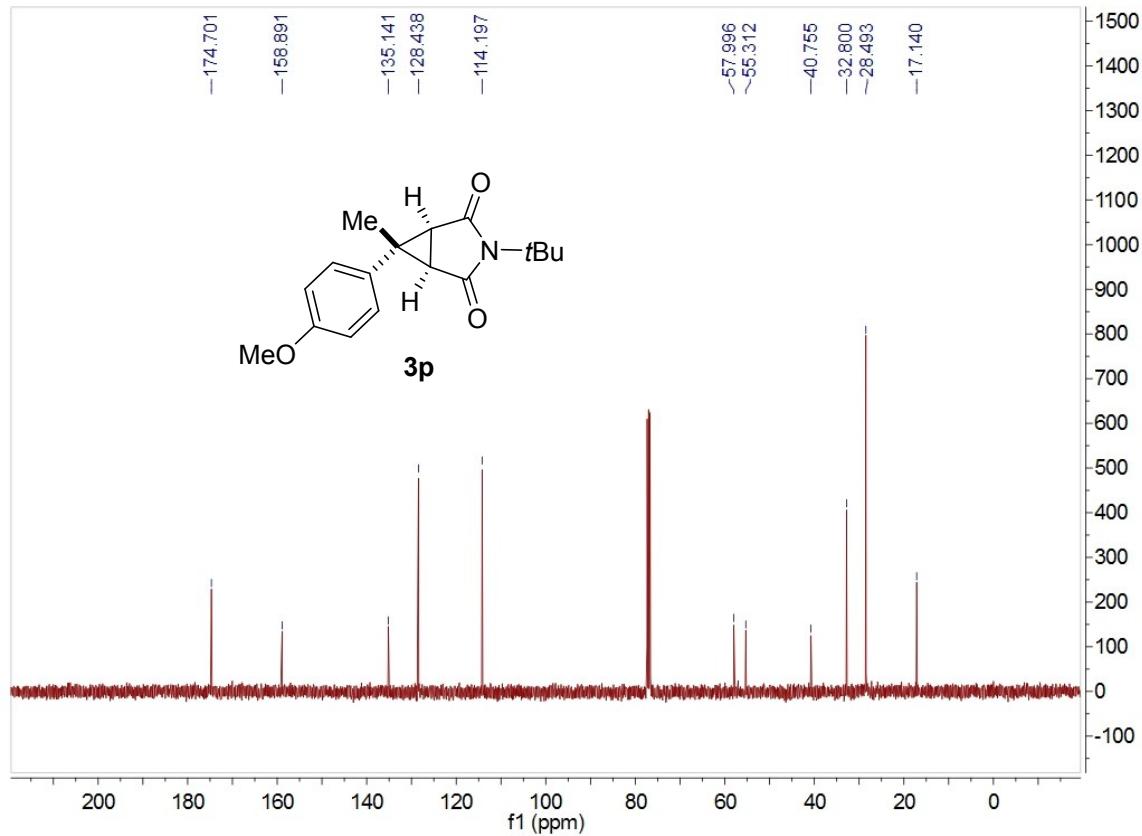
¹³C NMR (100 MHz, CDCl₃) spectrum for 3o



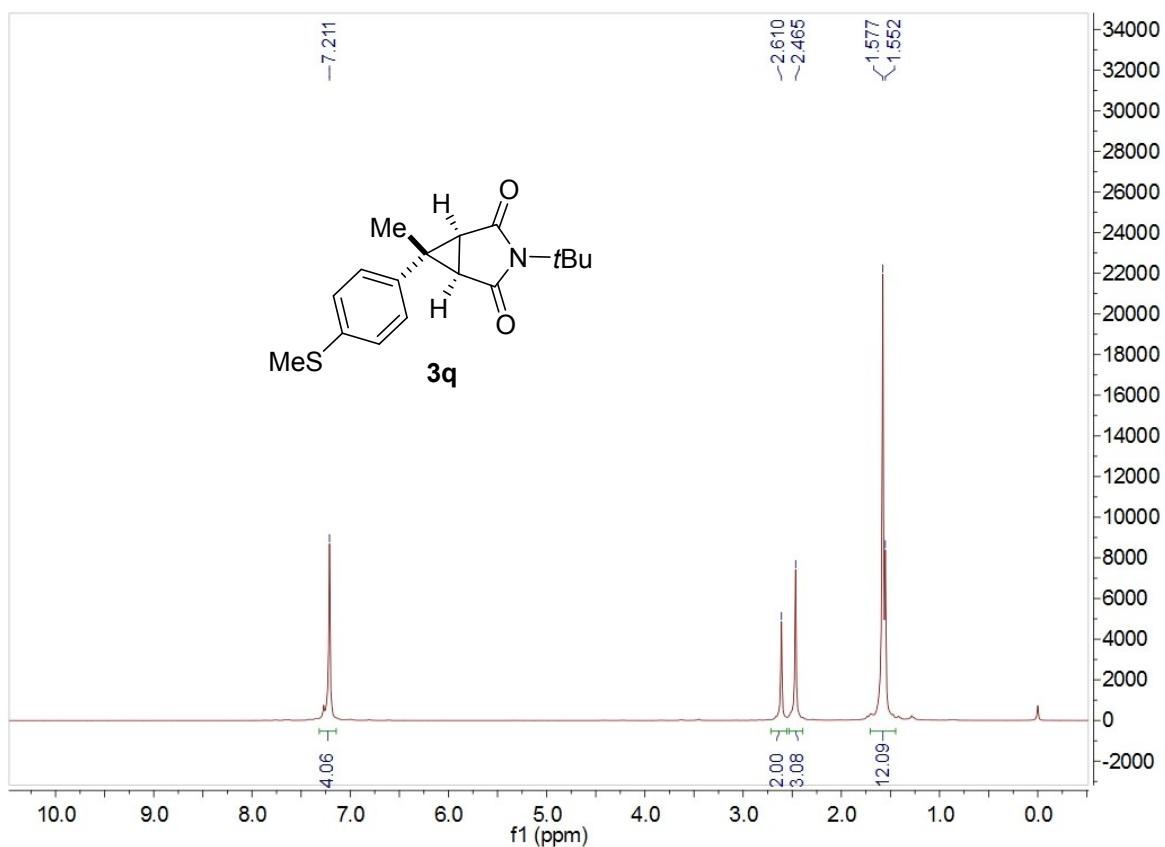
¹H NMR (400 MHz, CDCl₃) spectrum for 3p



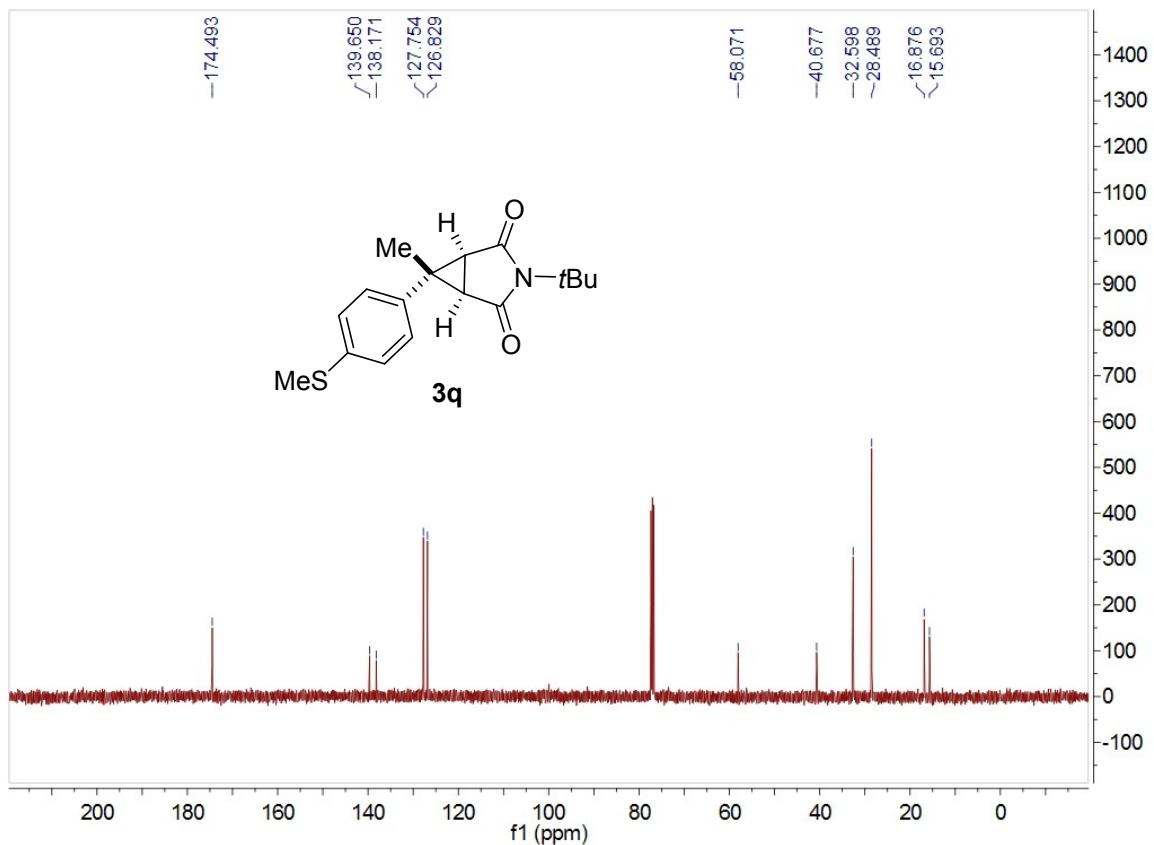
¹³C NMR (100 MHz, CDCl₃) spectrum for 3p



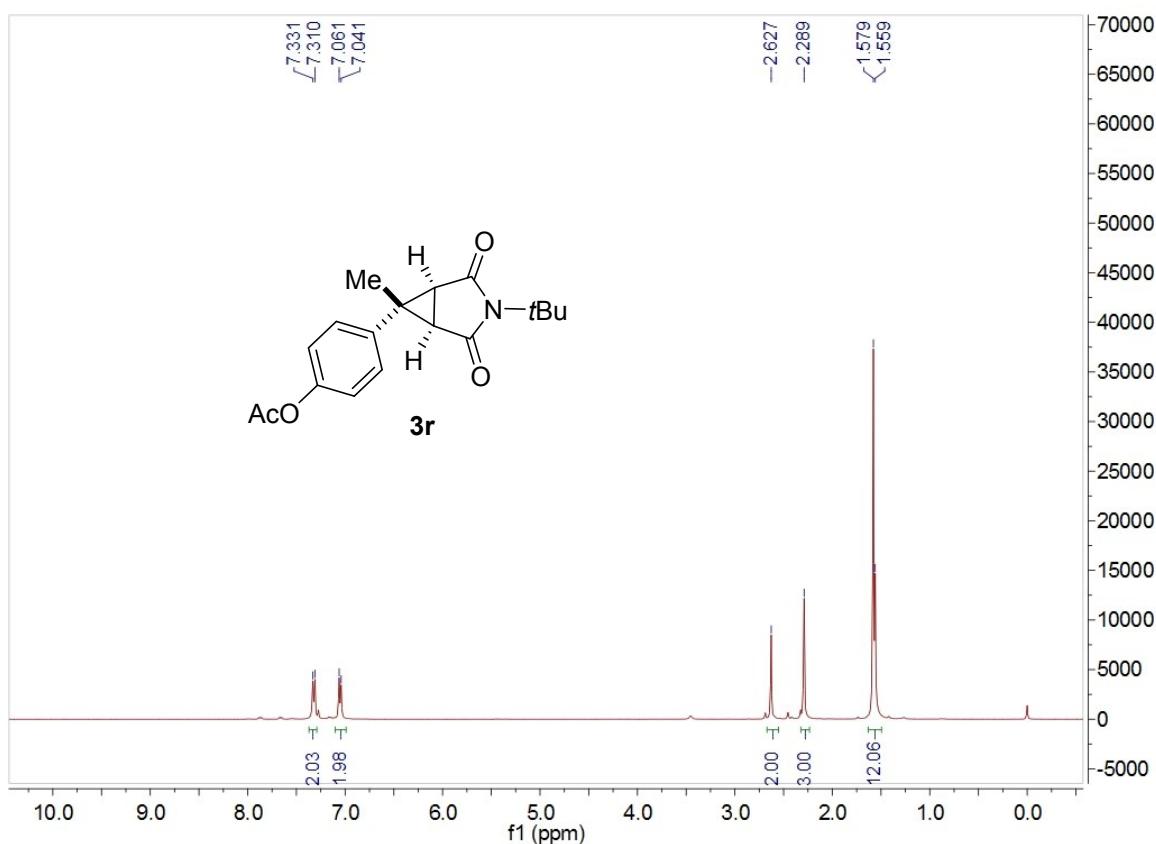
¹H NMR (400 MHz, CDCl₃) spectrum for 3q



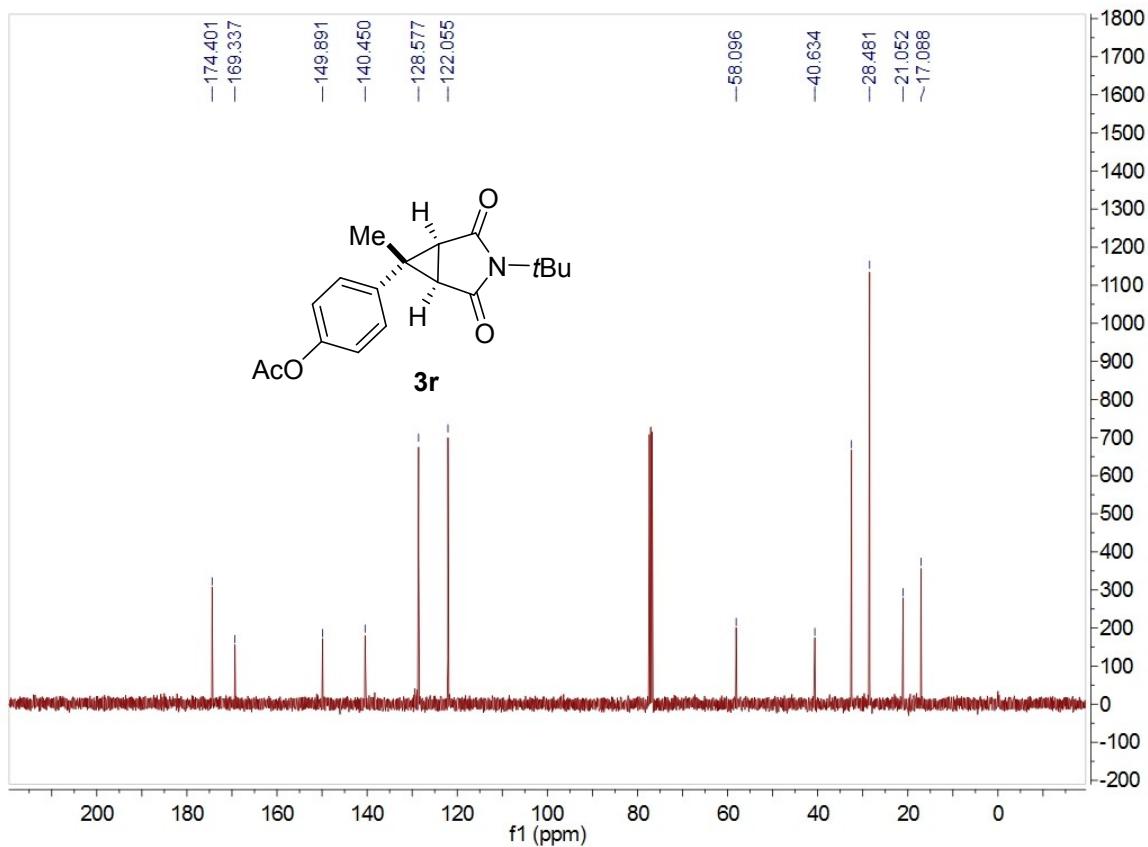
¹³C NMR (100 MHz, CDCl₃) spectrum for 3q



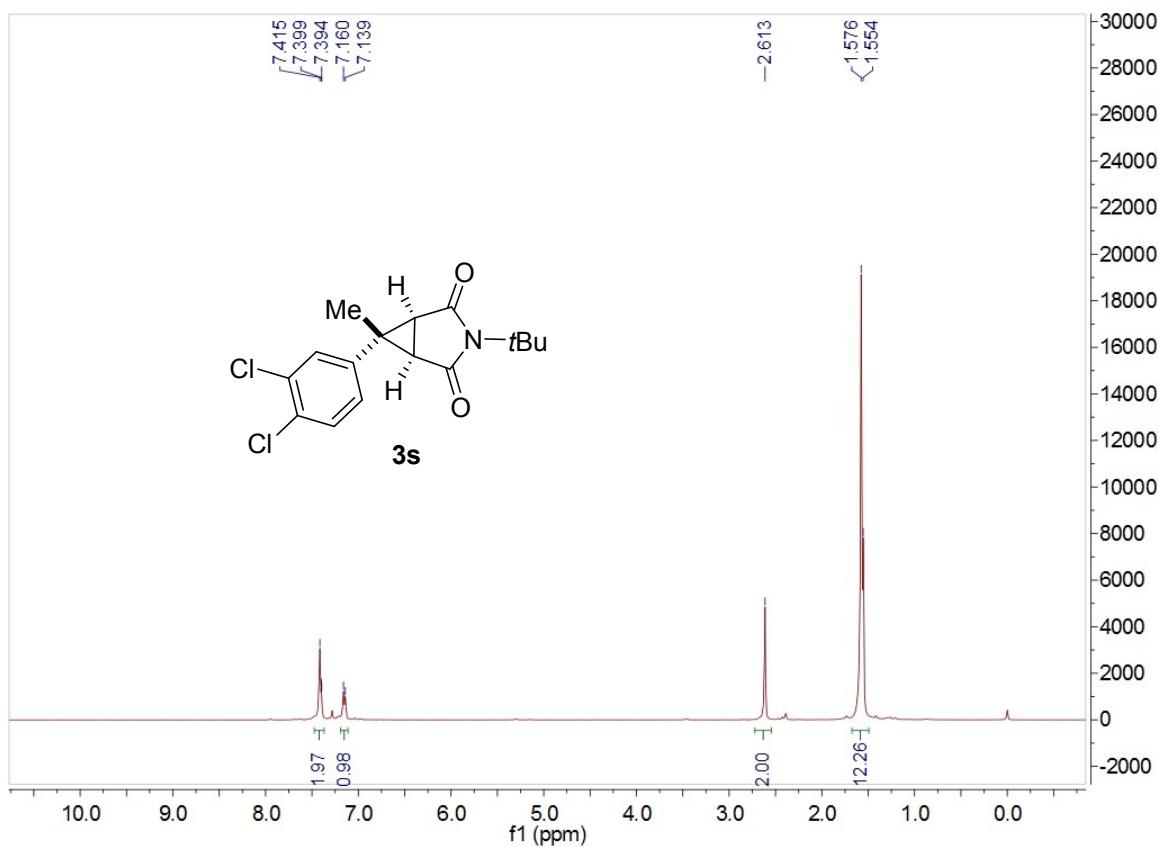
¹H NMR (400 MHz, CDCl₃) spectrum for 3r



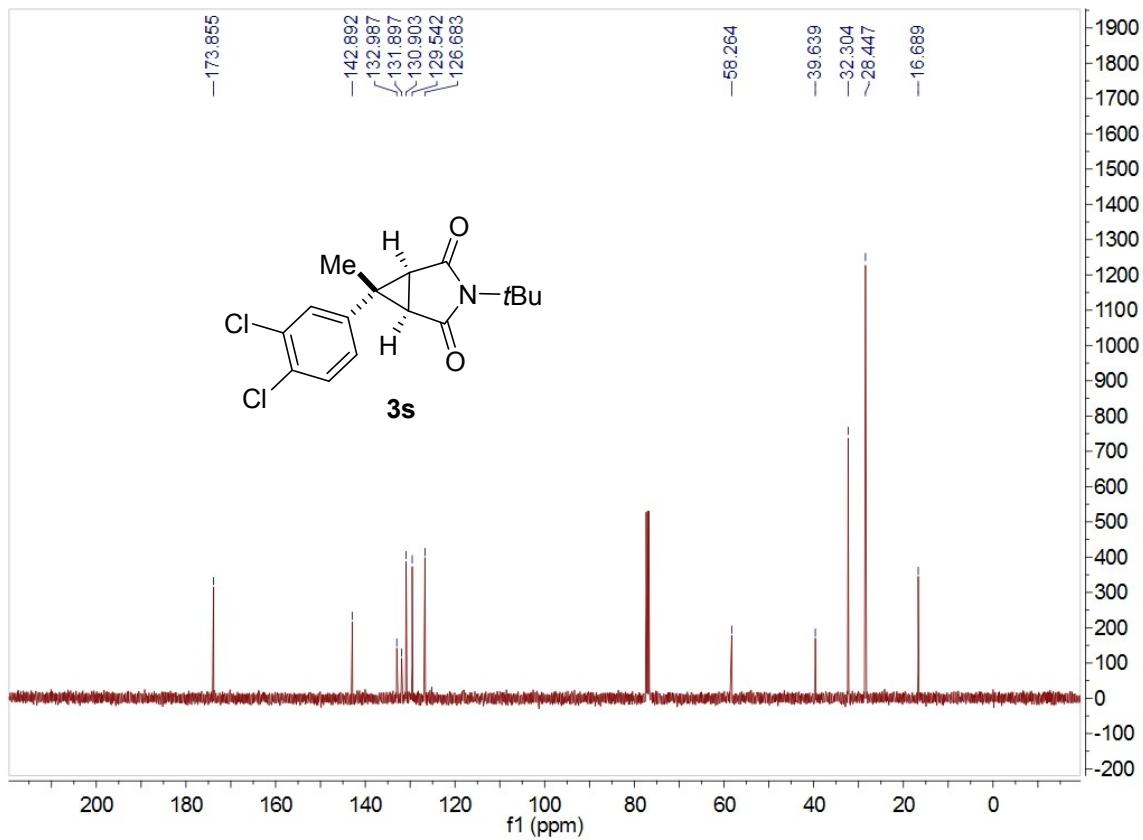
¹³C NMR (100 MHz, CDCl₃) spectrum for 3r



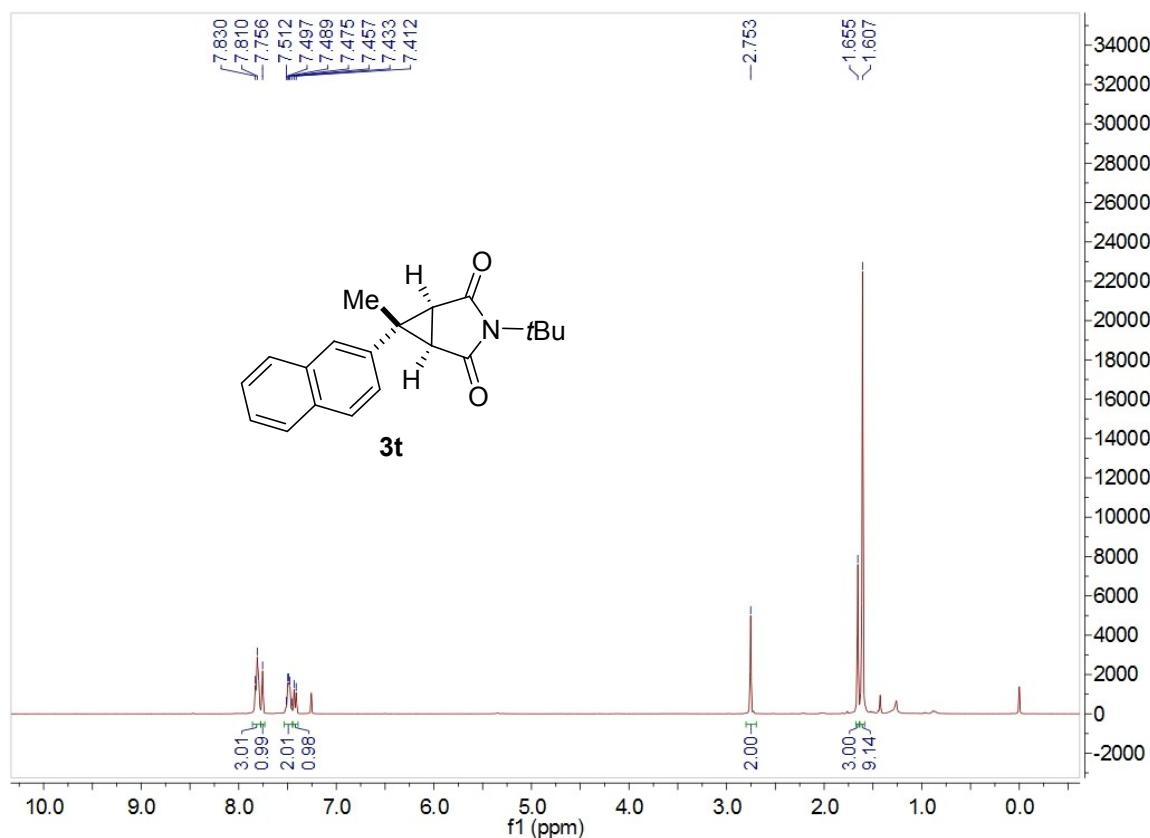
¹H NMR (400 MHz, CDCl₃) spectrum for 3s



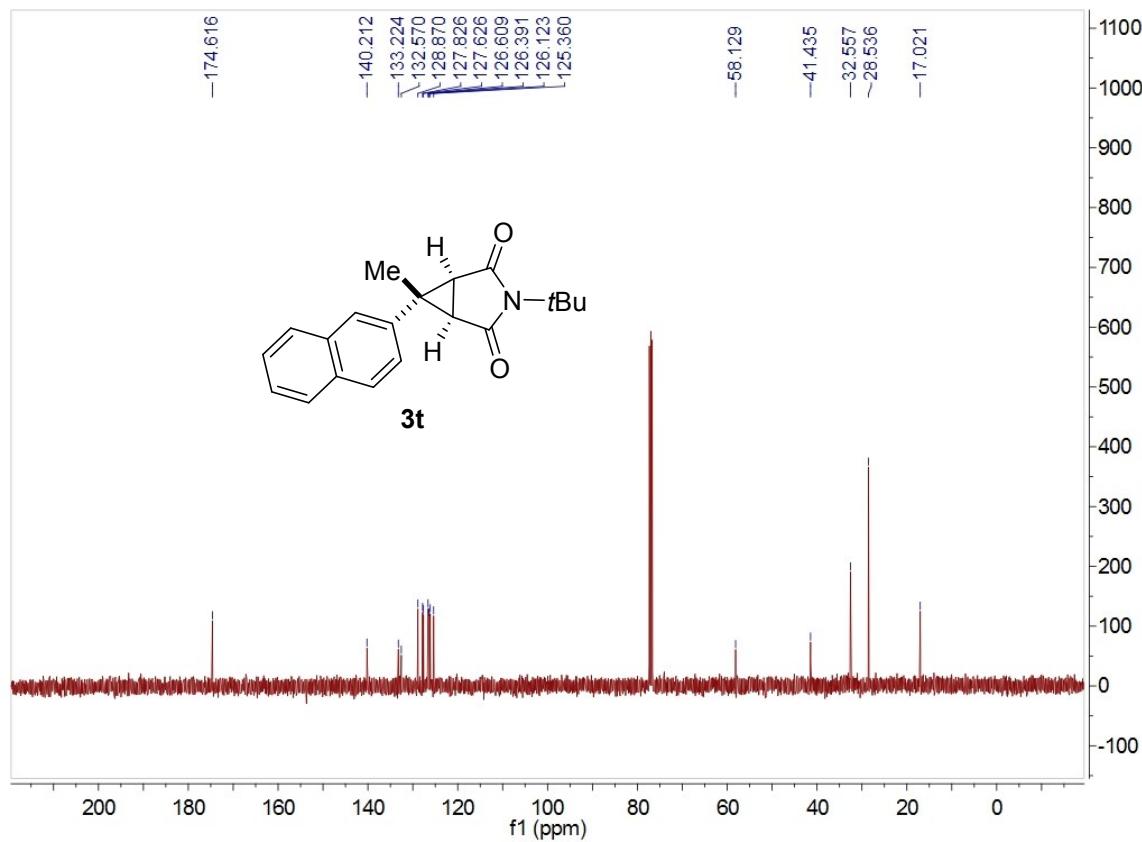
¹³C NMR (100 MHz, CDCl₃) spectrum for 3s



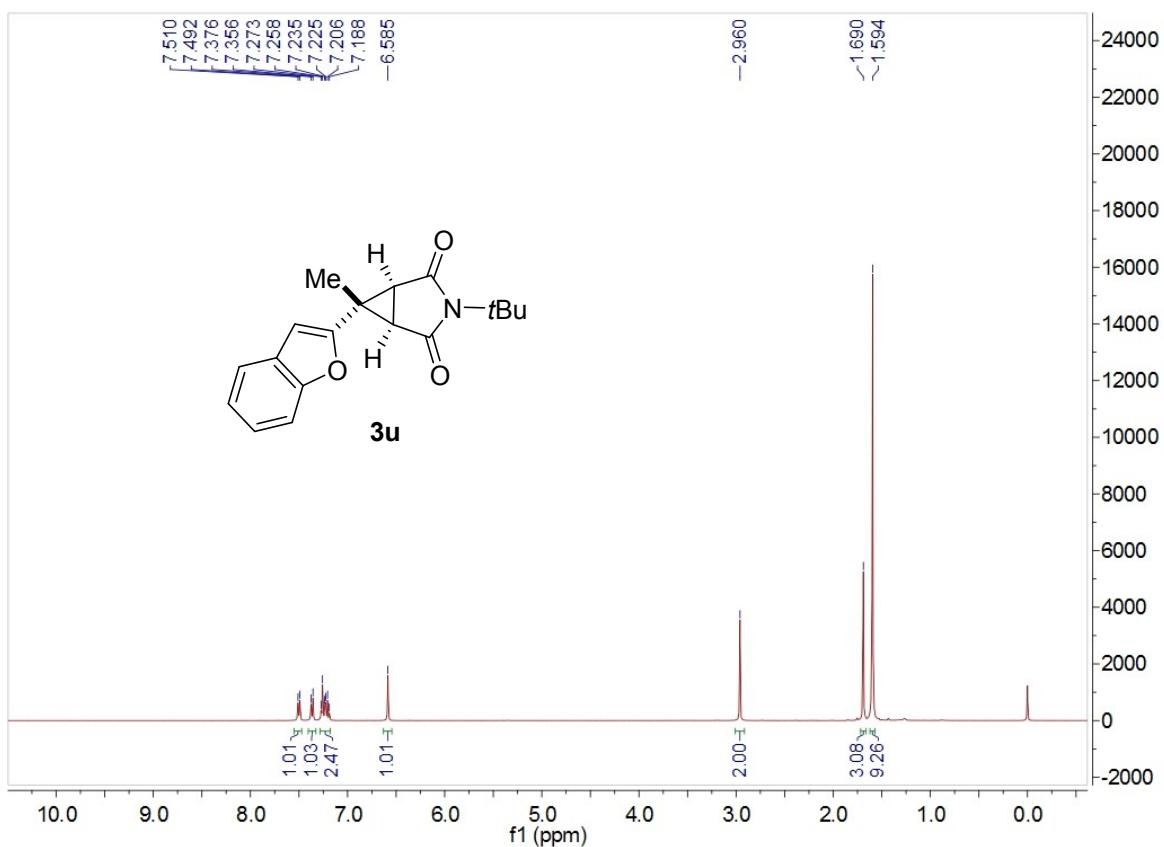
¹H NMR (400 MHz, CDCl₃) spectrum for 3t



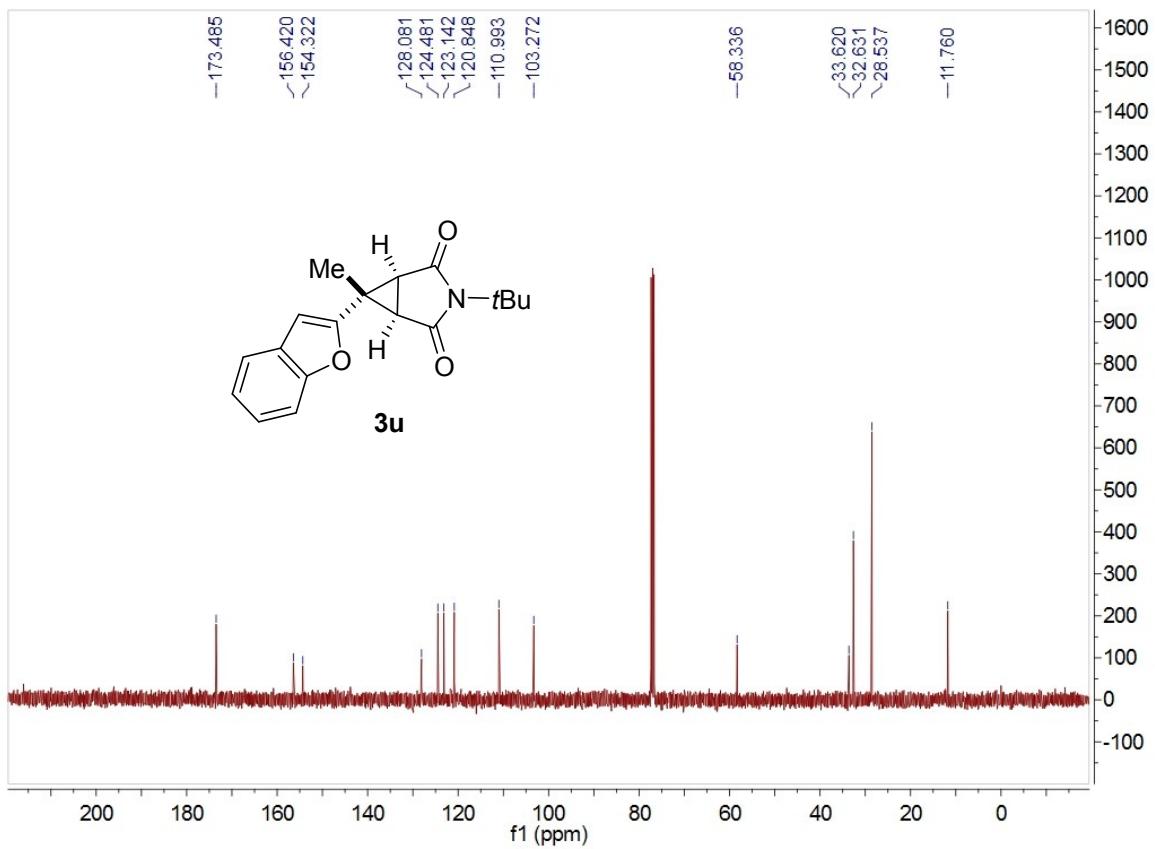
¹³C NMR (100 MHz, CDCl₃) spectrum for 3t



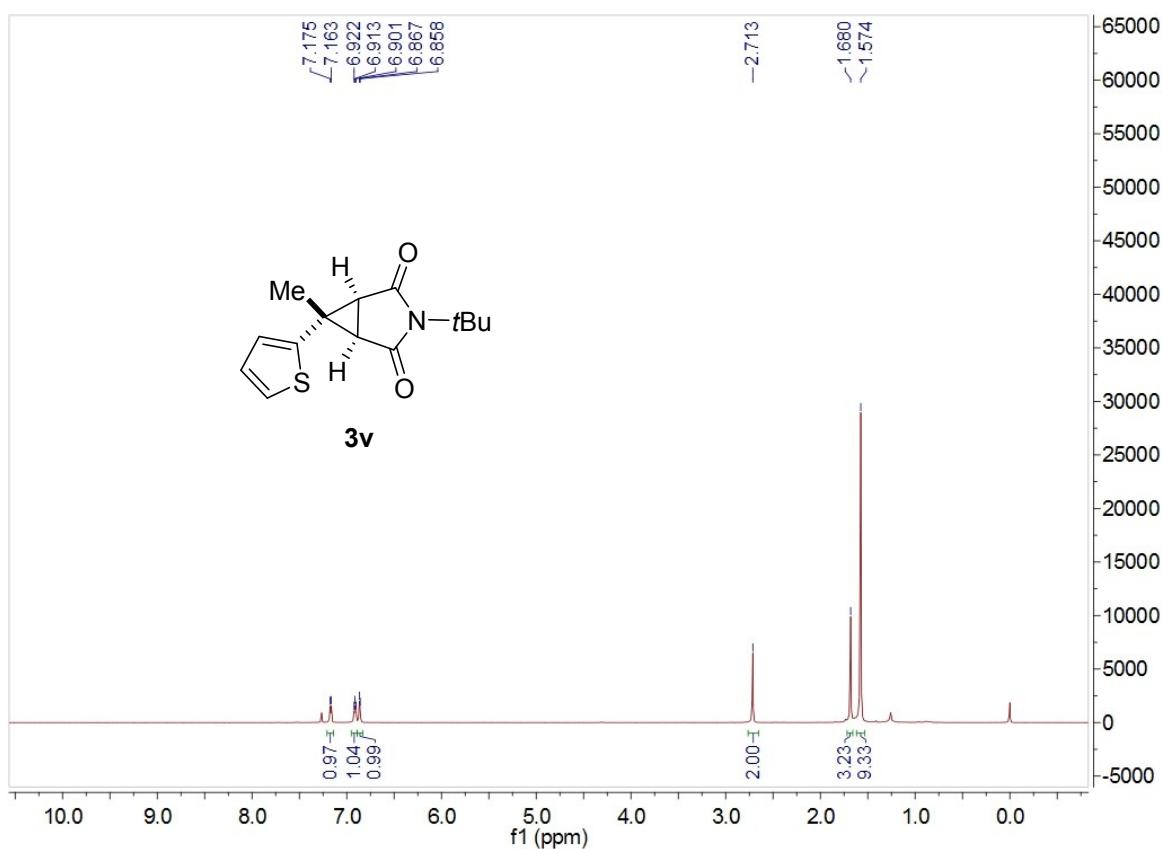
¹H NMR (400 MHz, CDCl₃) spectrum for 3u



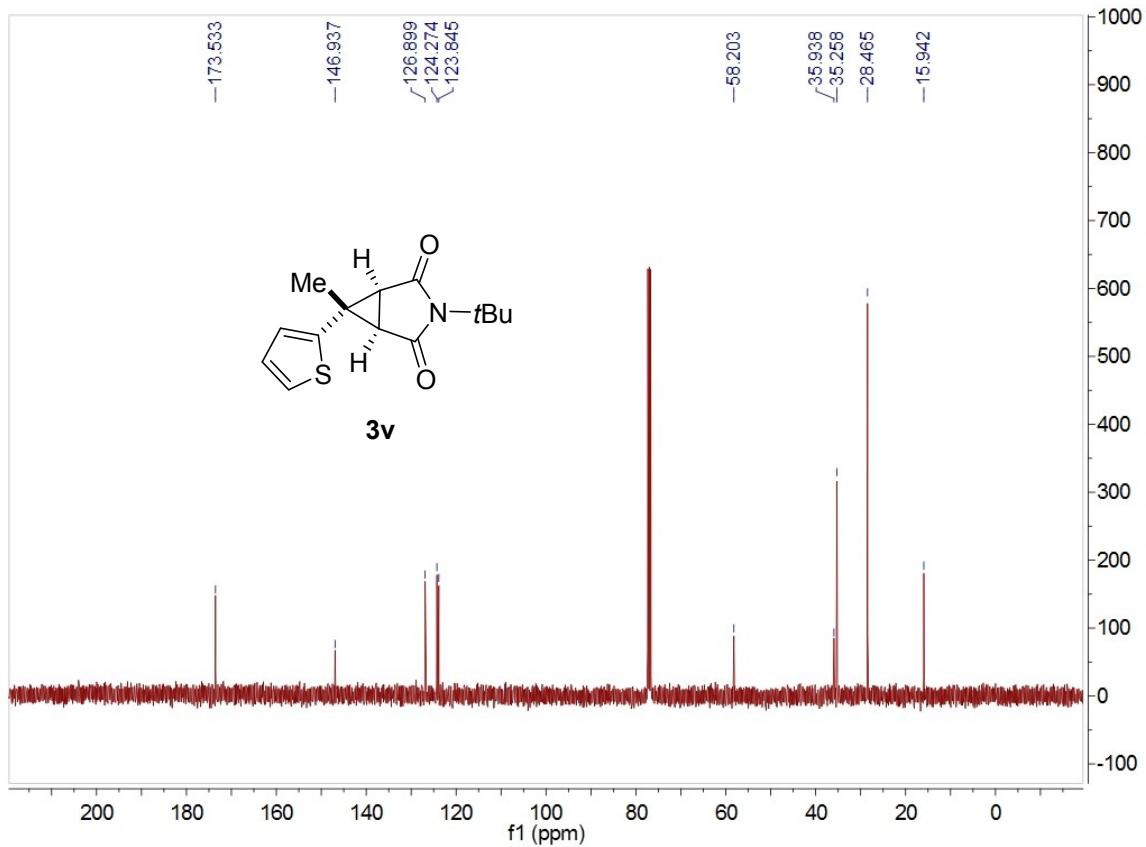
¹³C NMR (100 MHz, CDCl₃) spectrum for 3u



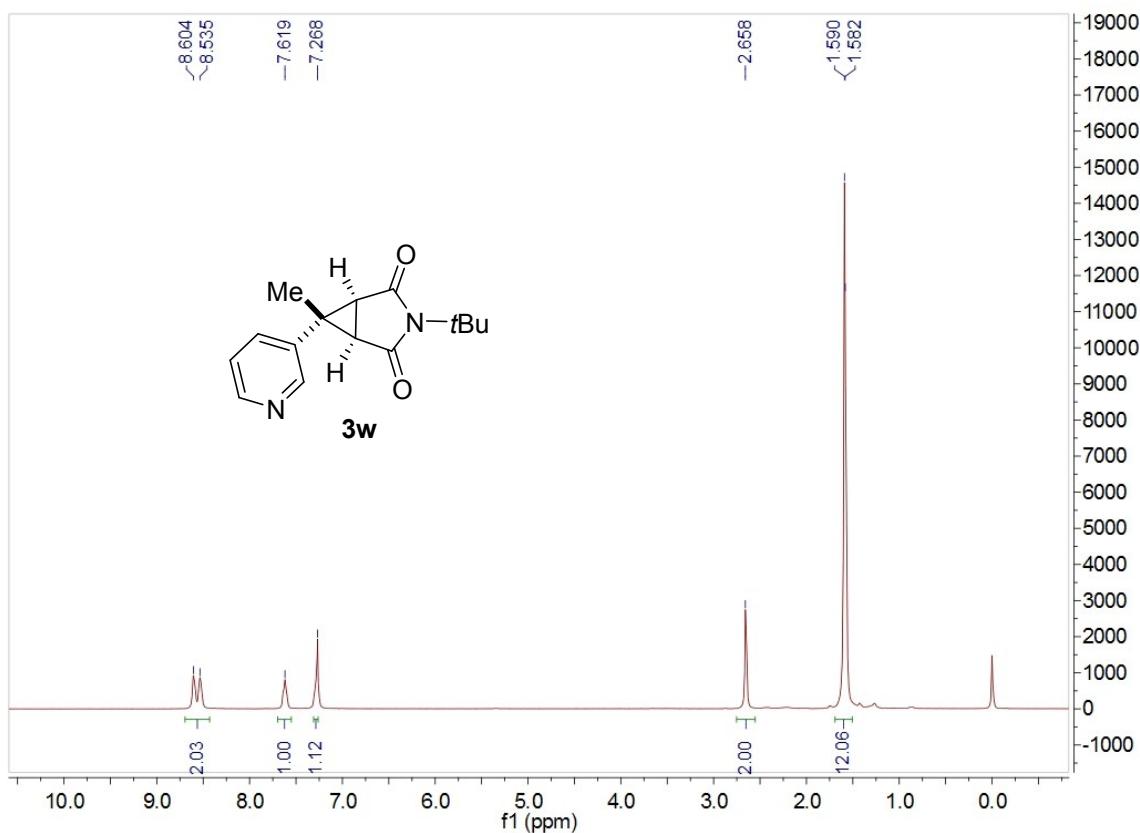
¹H NMR (400 MHz, CDCl₃) spectrum for 3v



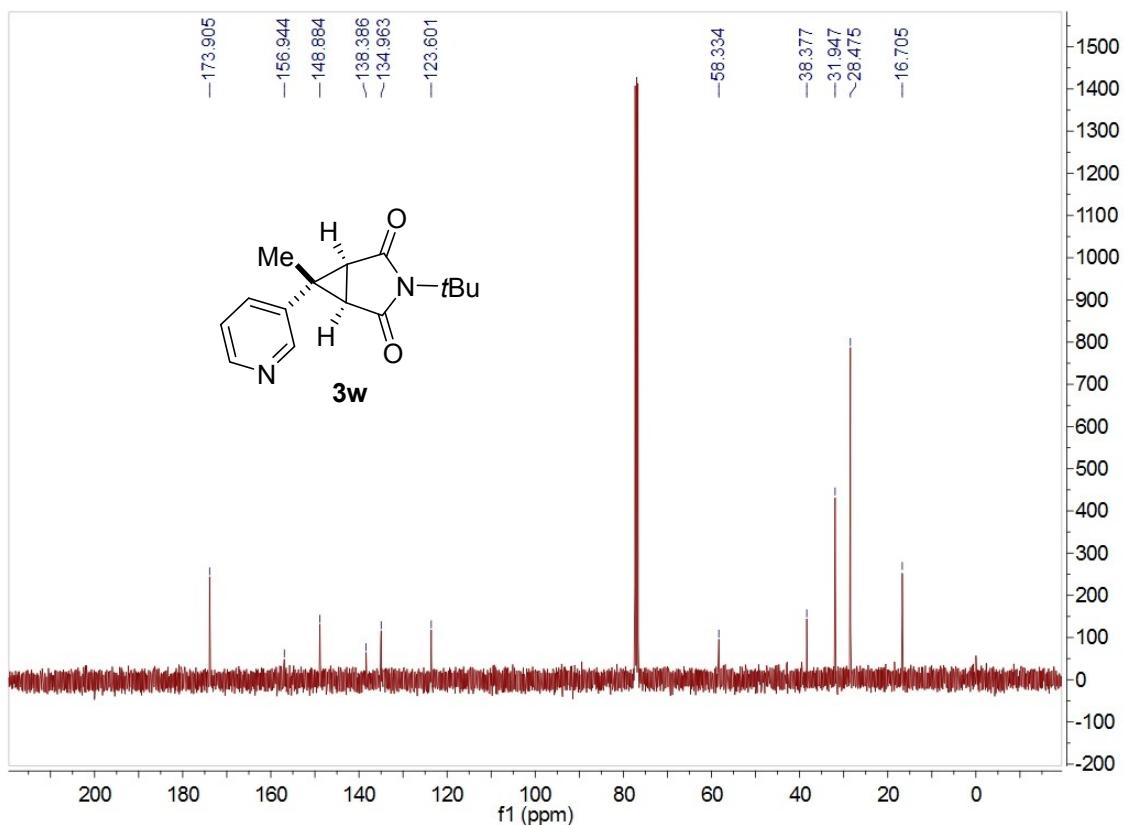
¹³C NMR (100 MHz, CDCl₃) spectrum for 3v



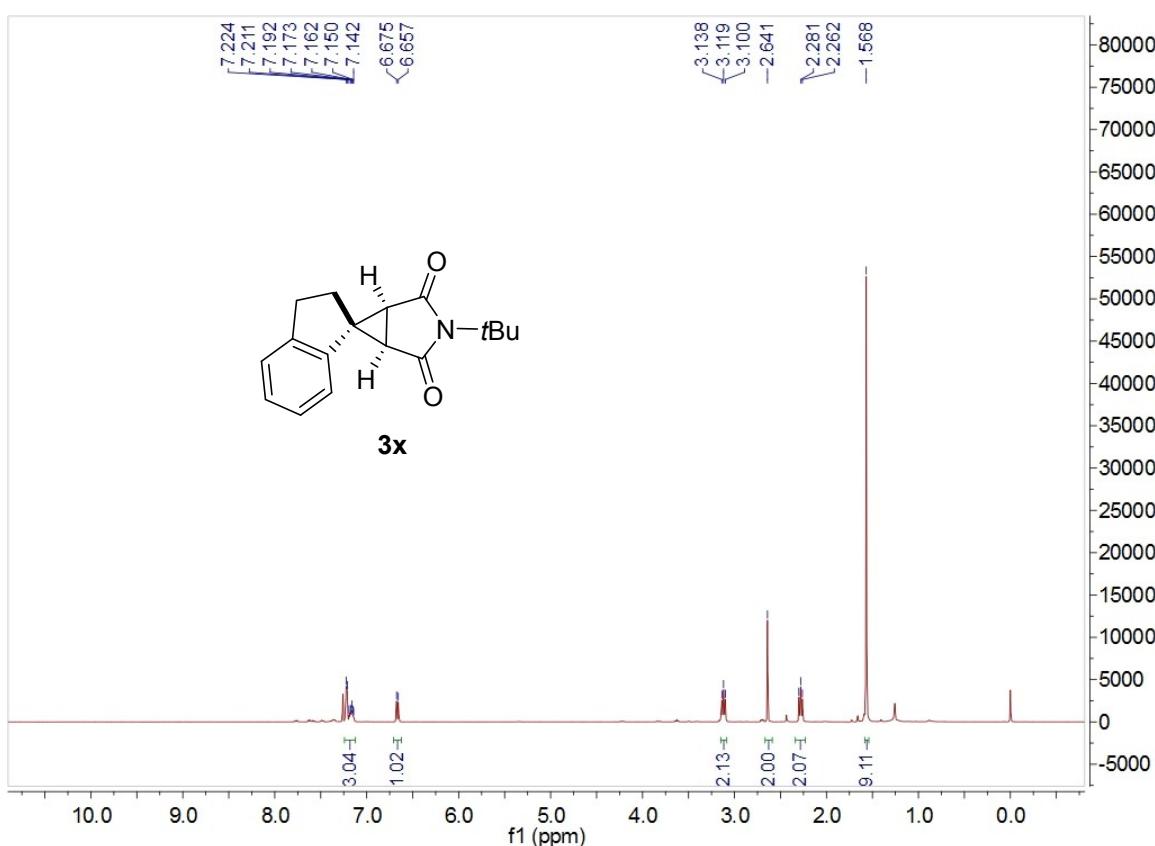
¹H NMR (400 MHz, CDCl₃) spectrum for 3w



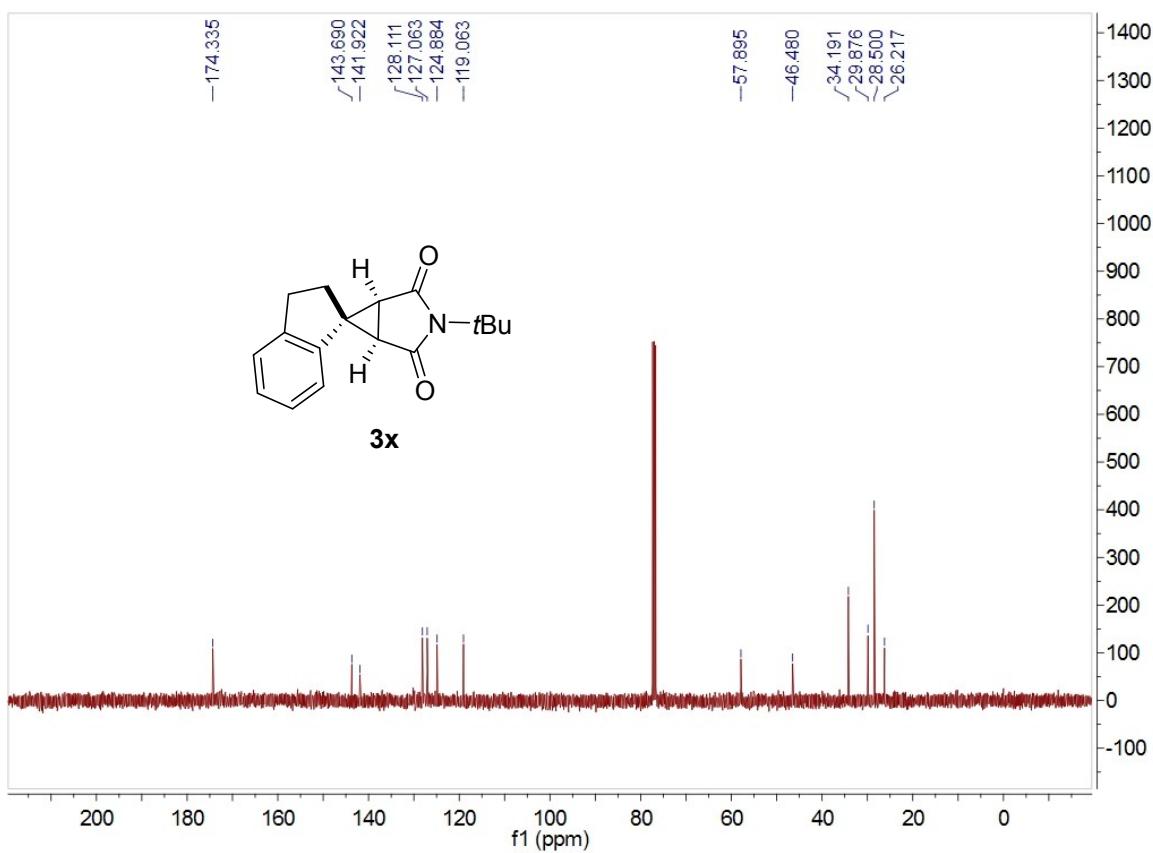
¹³C NMR (100 MHz, CDCl₃) spectrum for 3w



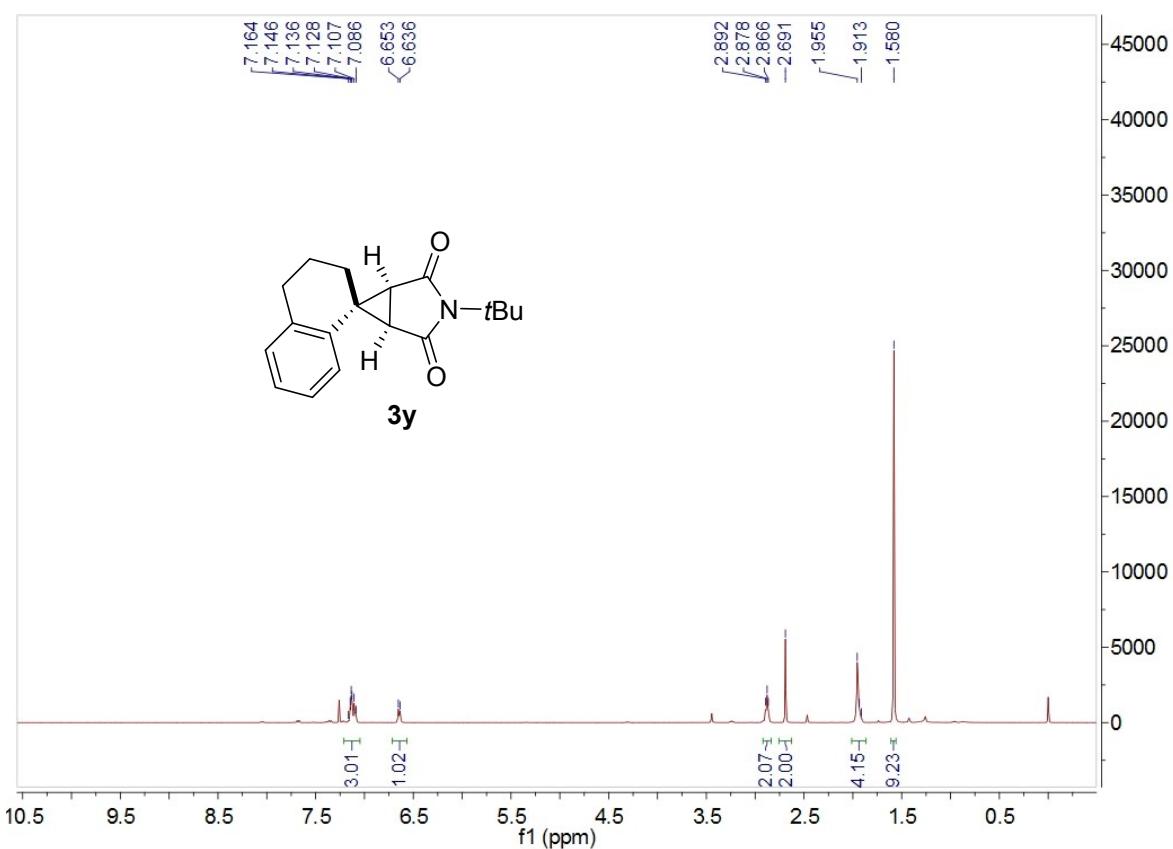
¹H NMR (400 MHz, CDCl₃) spectrum for 3x



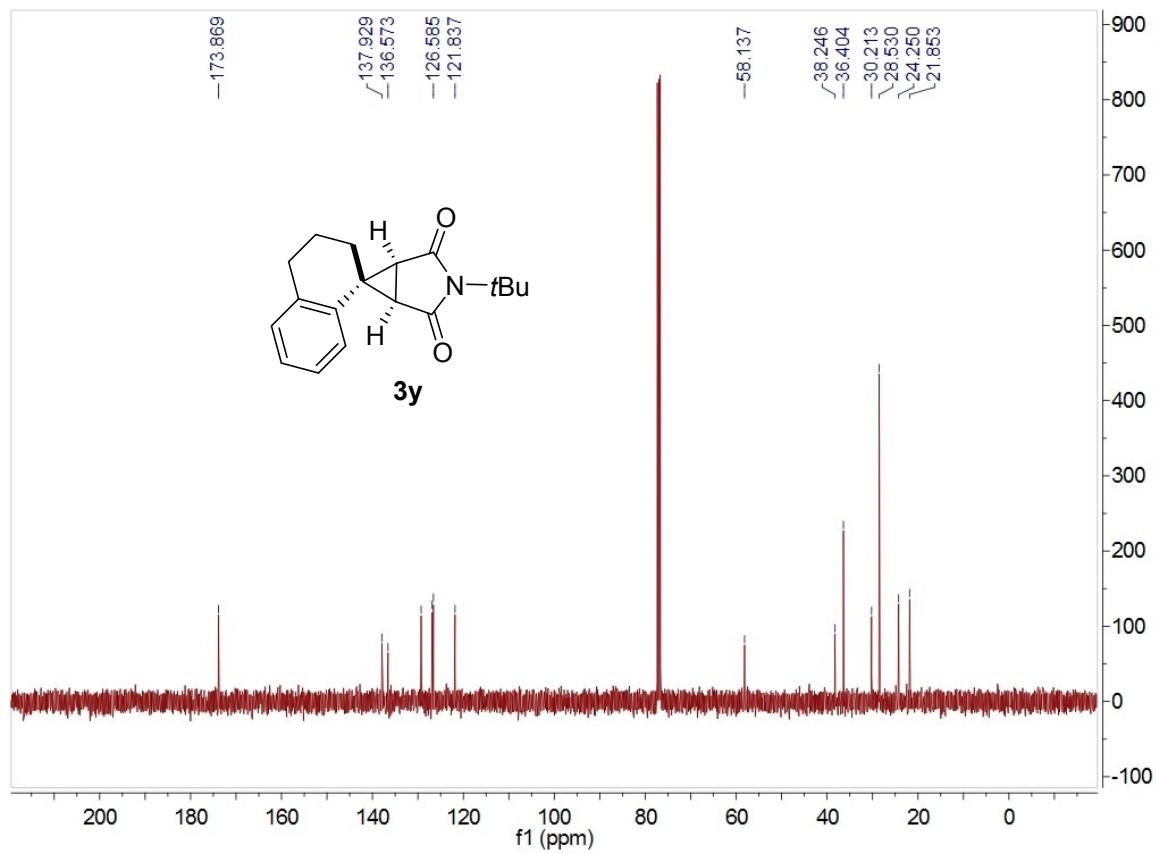
¹³C NMR (100 MHz, CDCl₃) spectrum for 3x



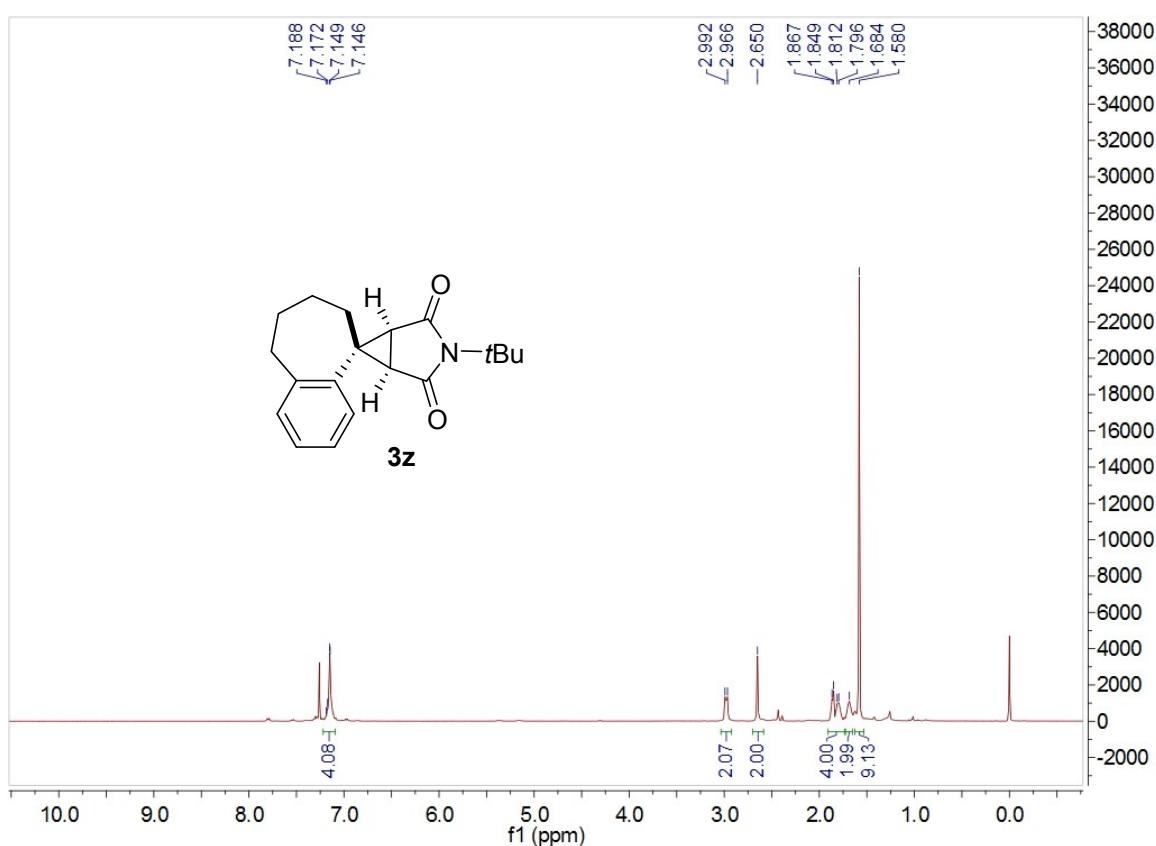
¹H NMR (400 MHz, CDCl₃) spectrum for 3y



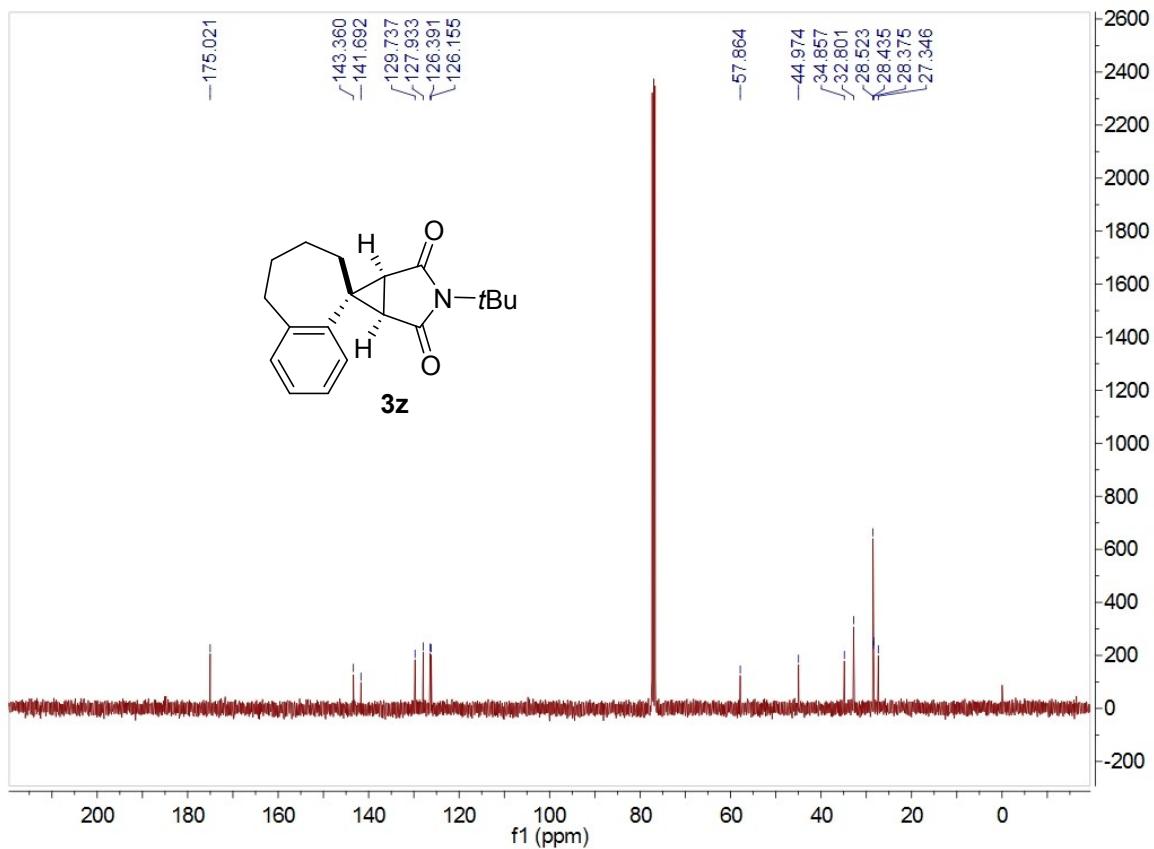
¹³C NMR (100 MHz, CDCl₃) spectrum for 3y



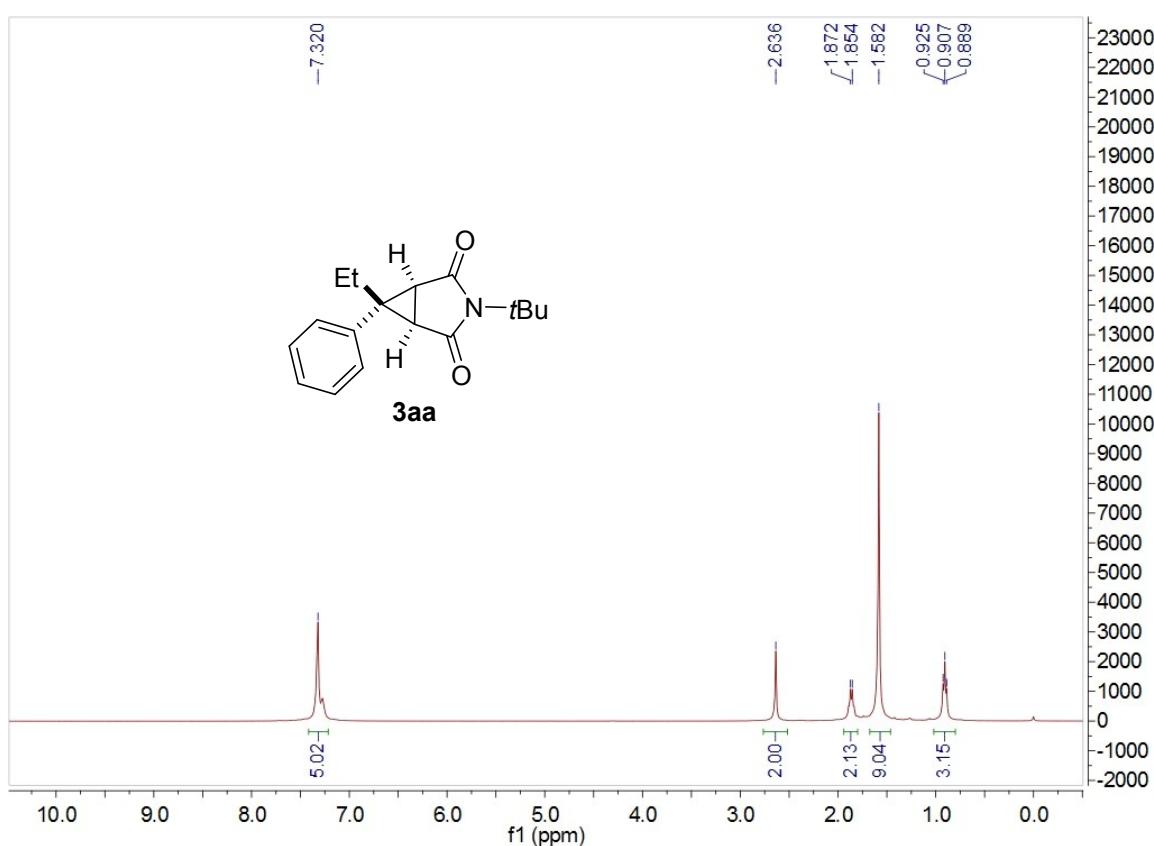
¹H NMR (400 MHz, CDCl₃) spectrum for 3z



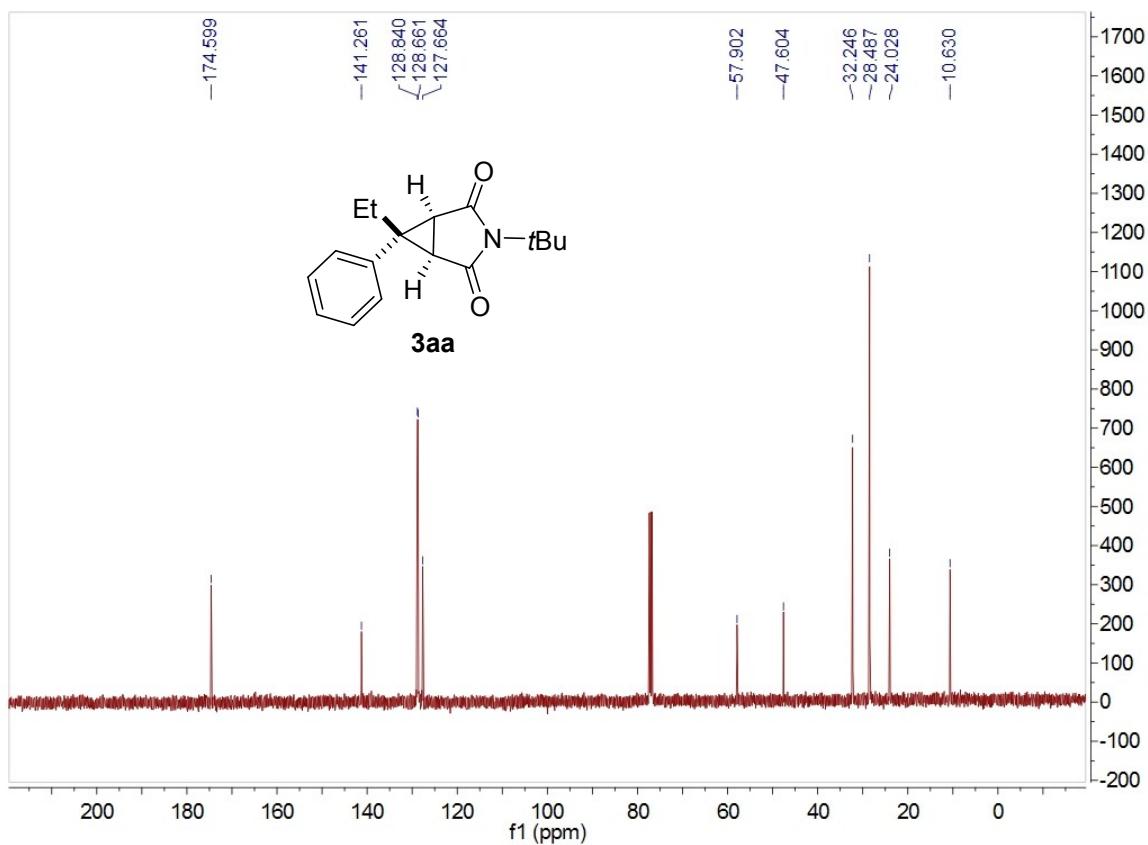
¹³C NMR (100 MHz, CDCl₃) spectrum for 3z



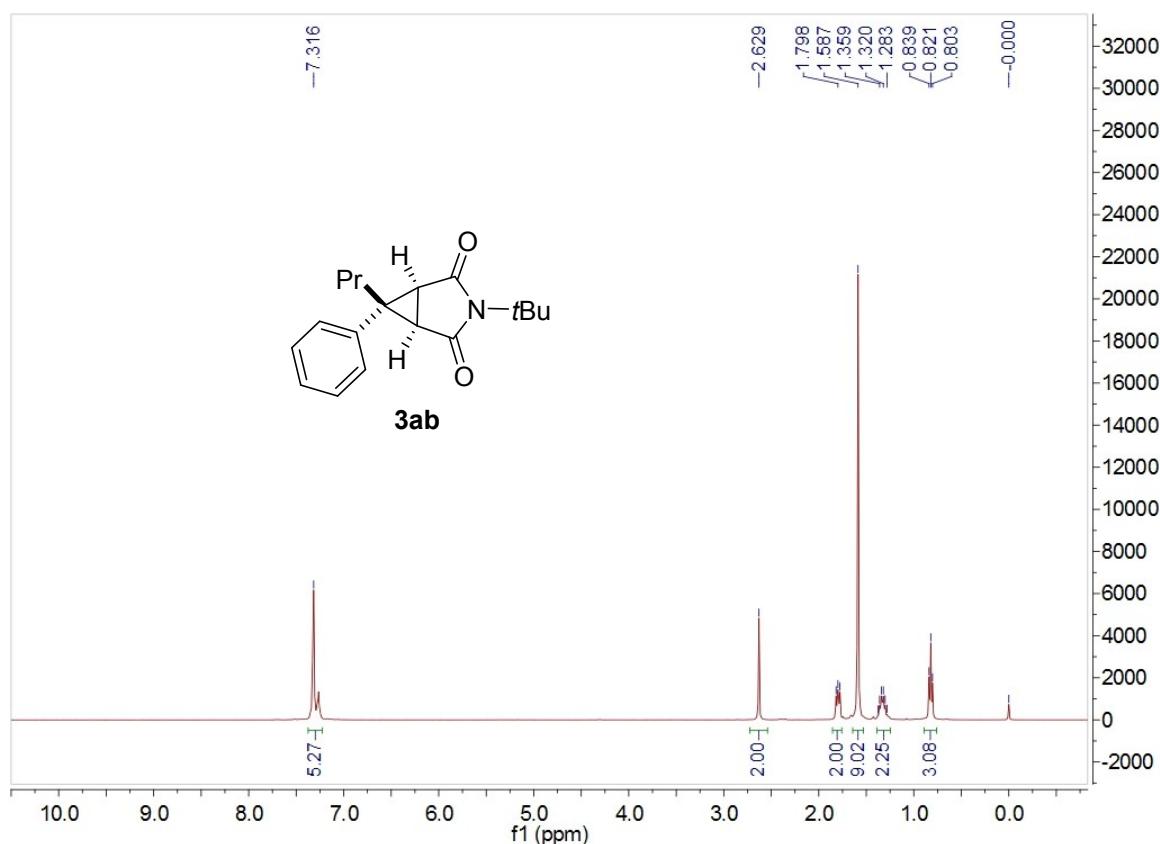
¹H NMR (400 MHz, CDCl₃) spectrum for 3aa



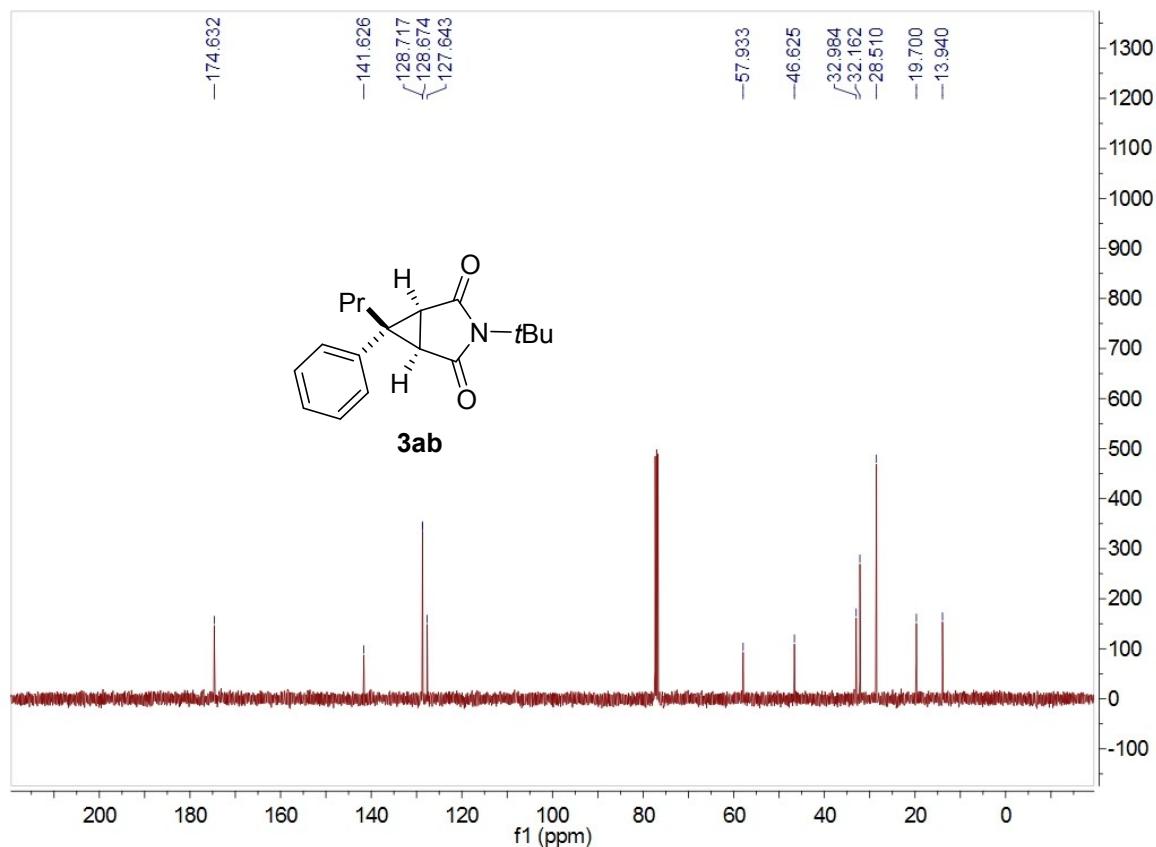
¹³C NMR (100 MHz, CDCl₃) spectrum for 3aa



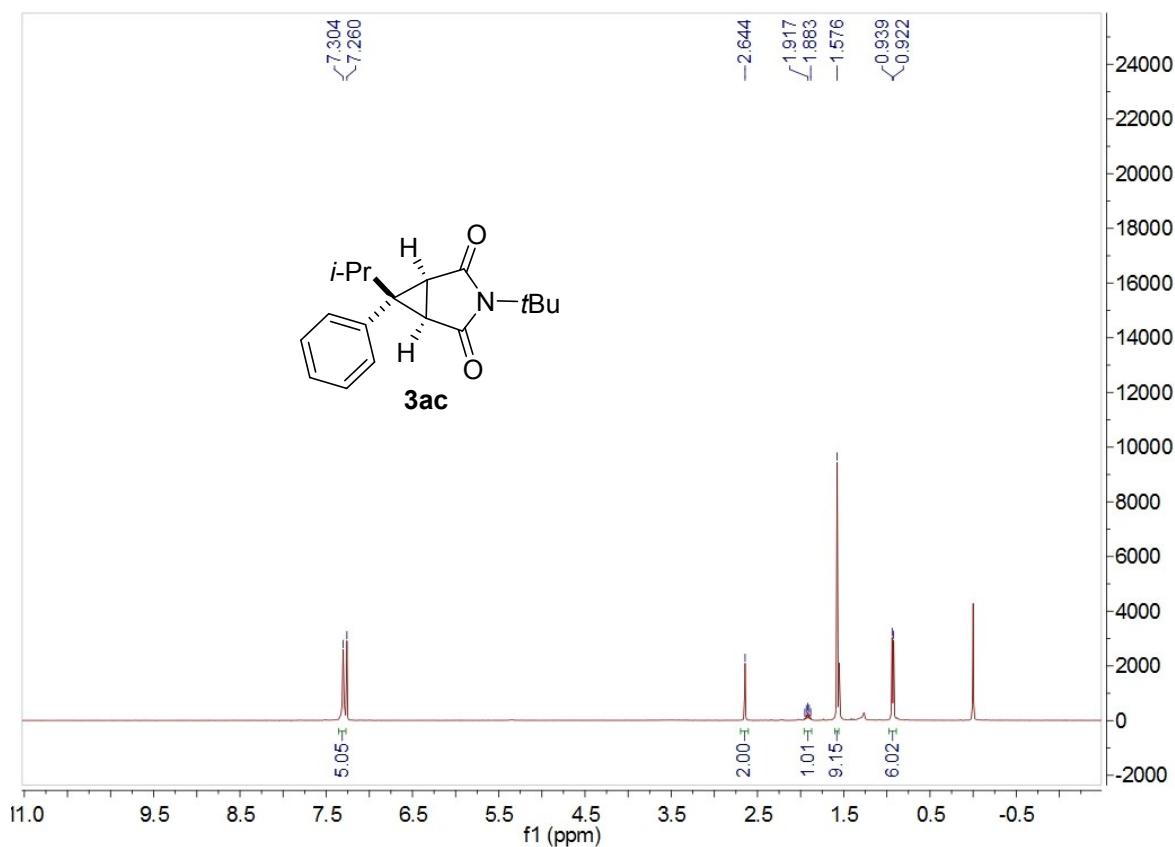
¹H NMR (400 MHz, CDCl₃) spectrum for 3ab



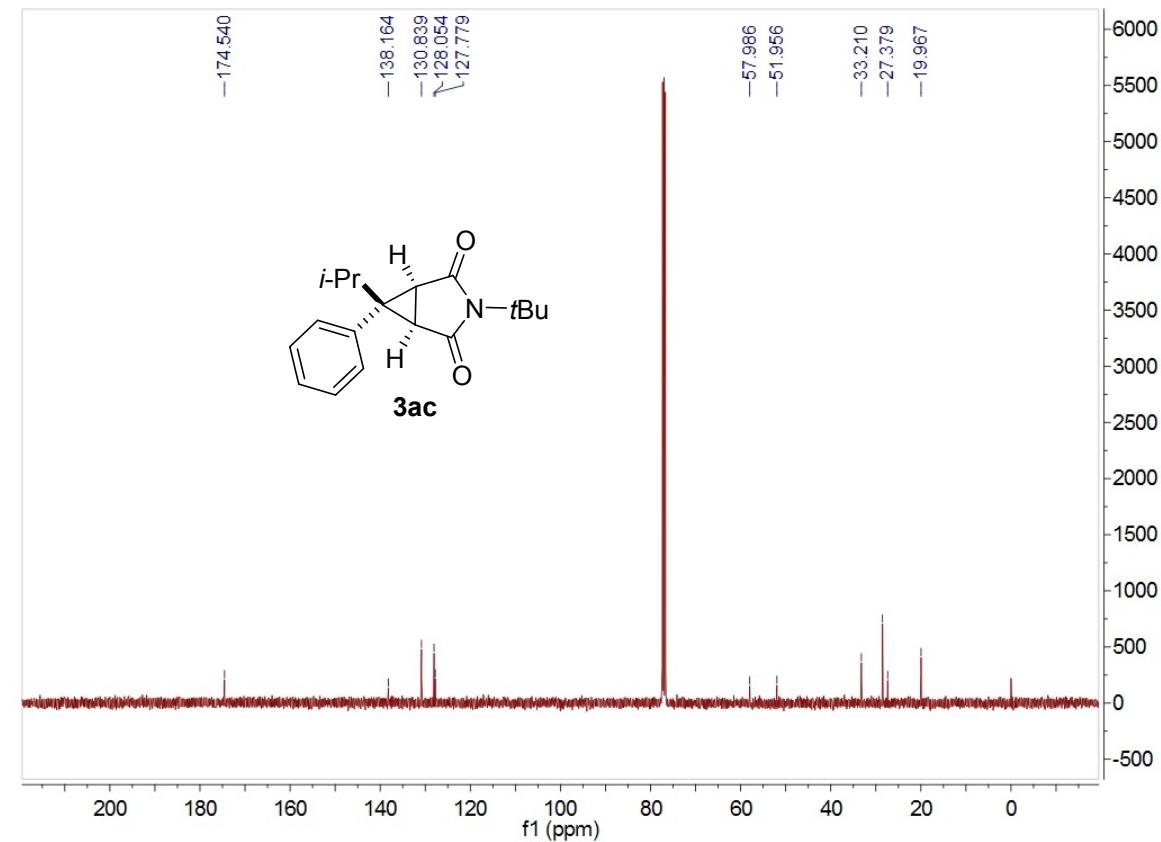
¹³C NMR (100 MHz, CDCl₃) spectrum for 3ab



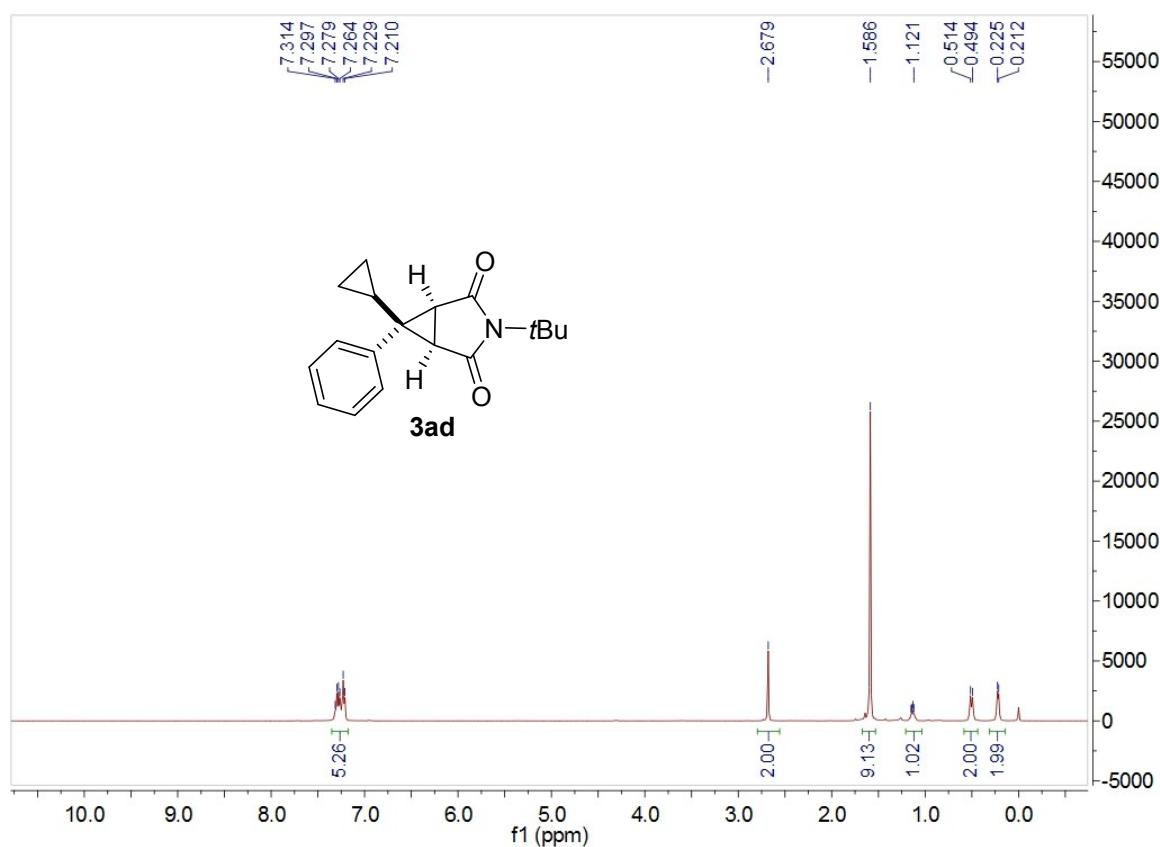
¹H NMR (400 MHz, CDCl₃) spectrum for 3ac



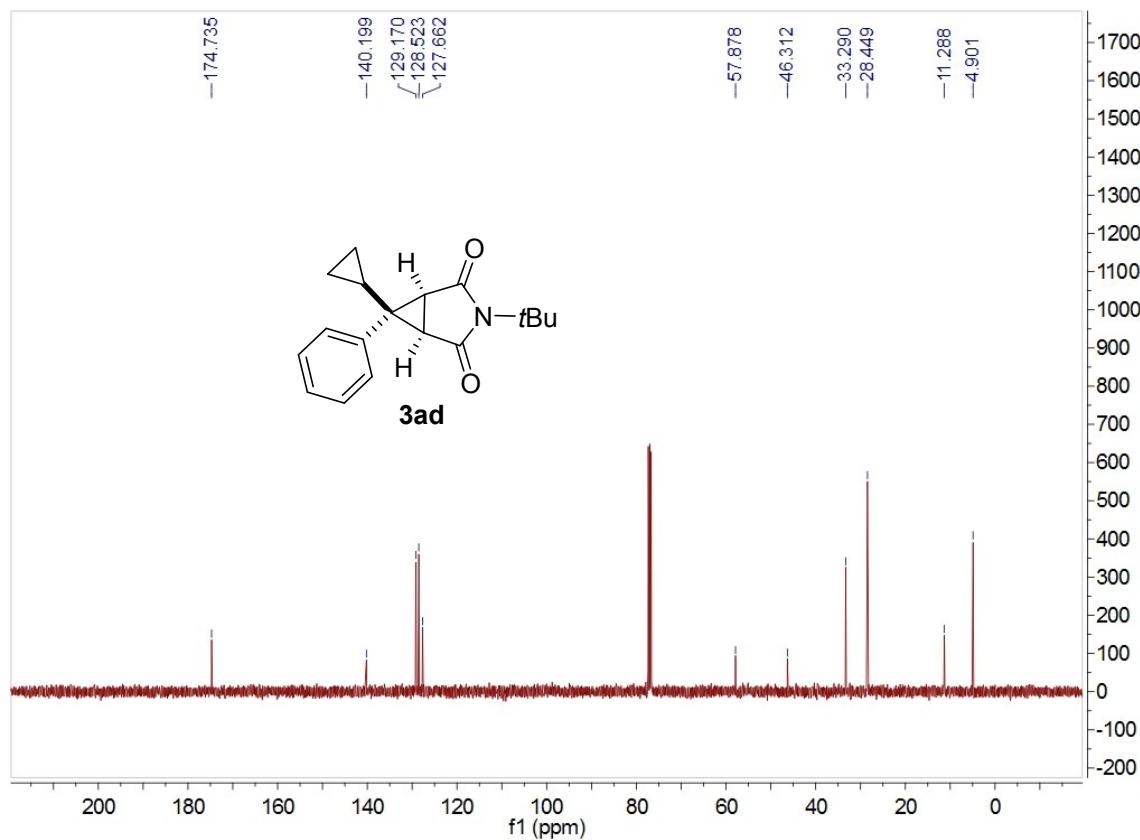
¹³C NMR (100 MHz, CDCl₃) spectrum for 3ac



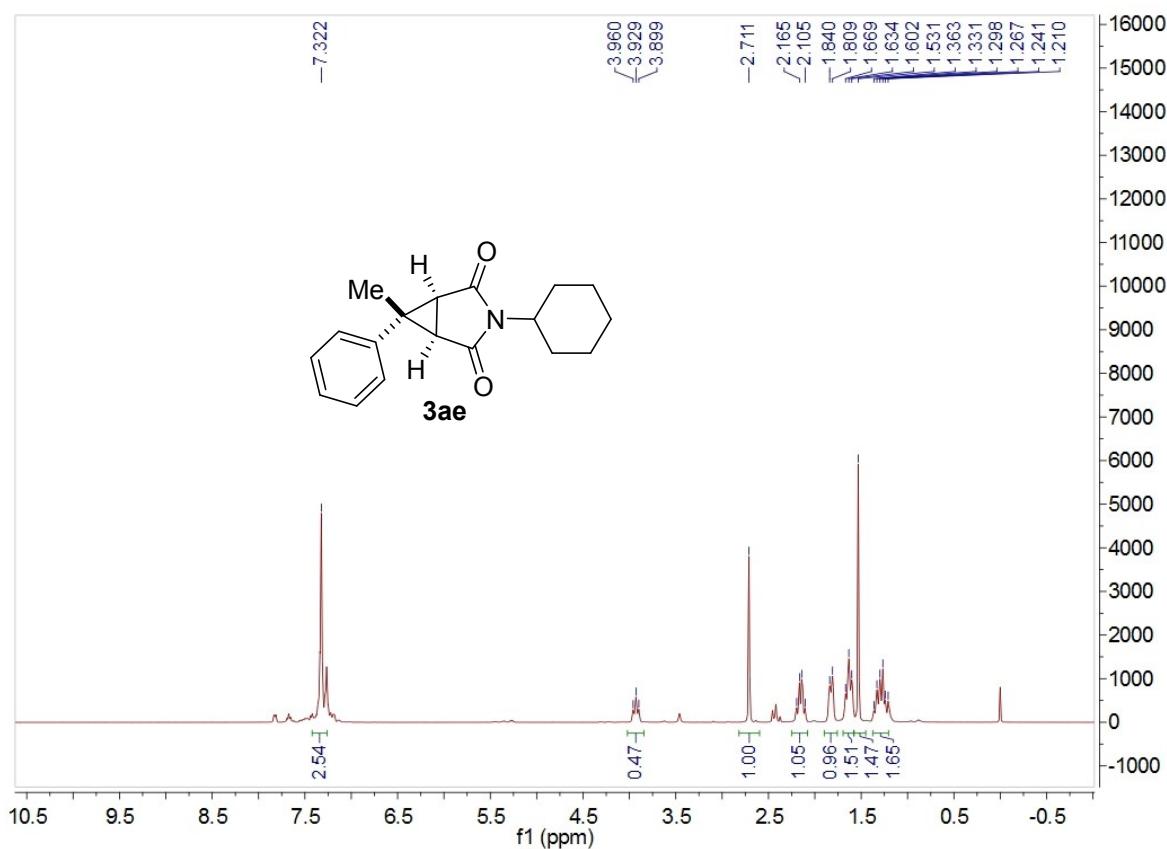
¹H NMR (400 MHz, CDCl₃) spectrum for 3ad



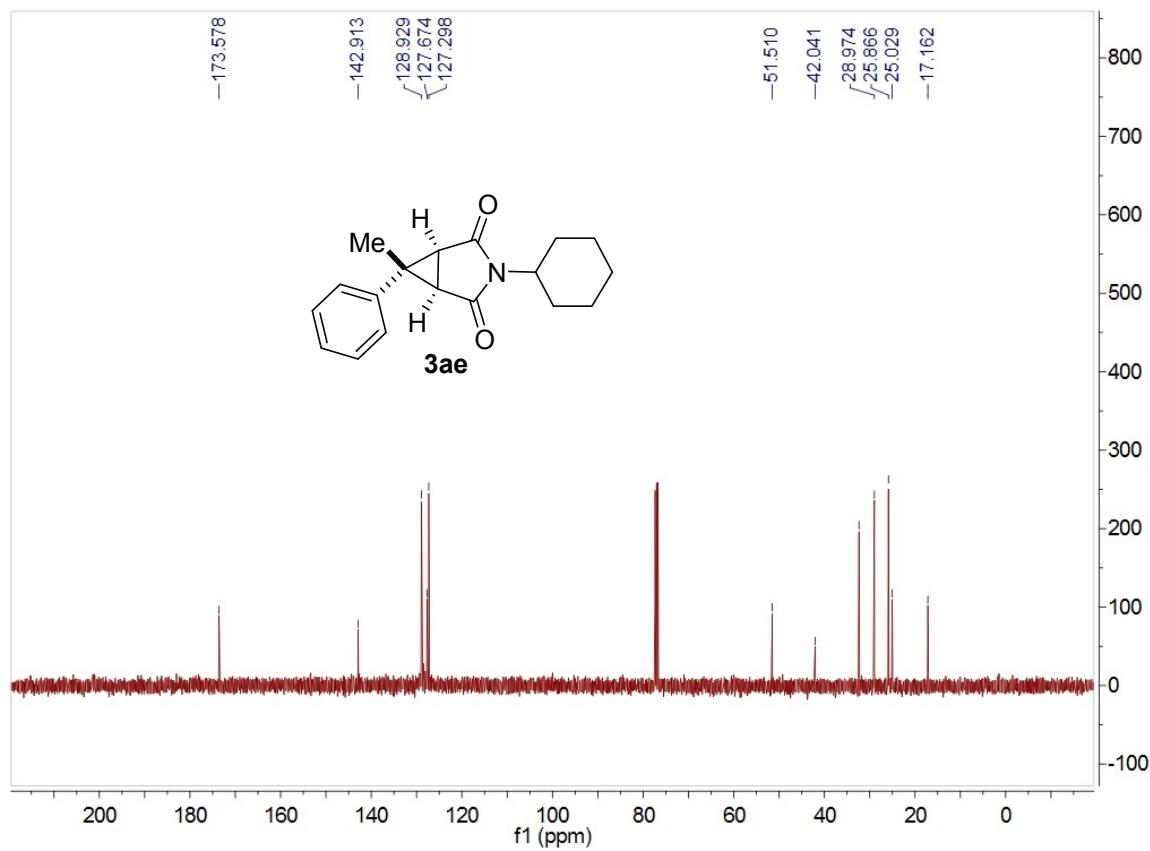
¹³C NMR (100 MHz, CDCl₃) spectrum for 3ad



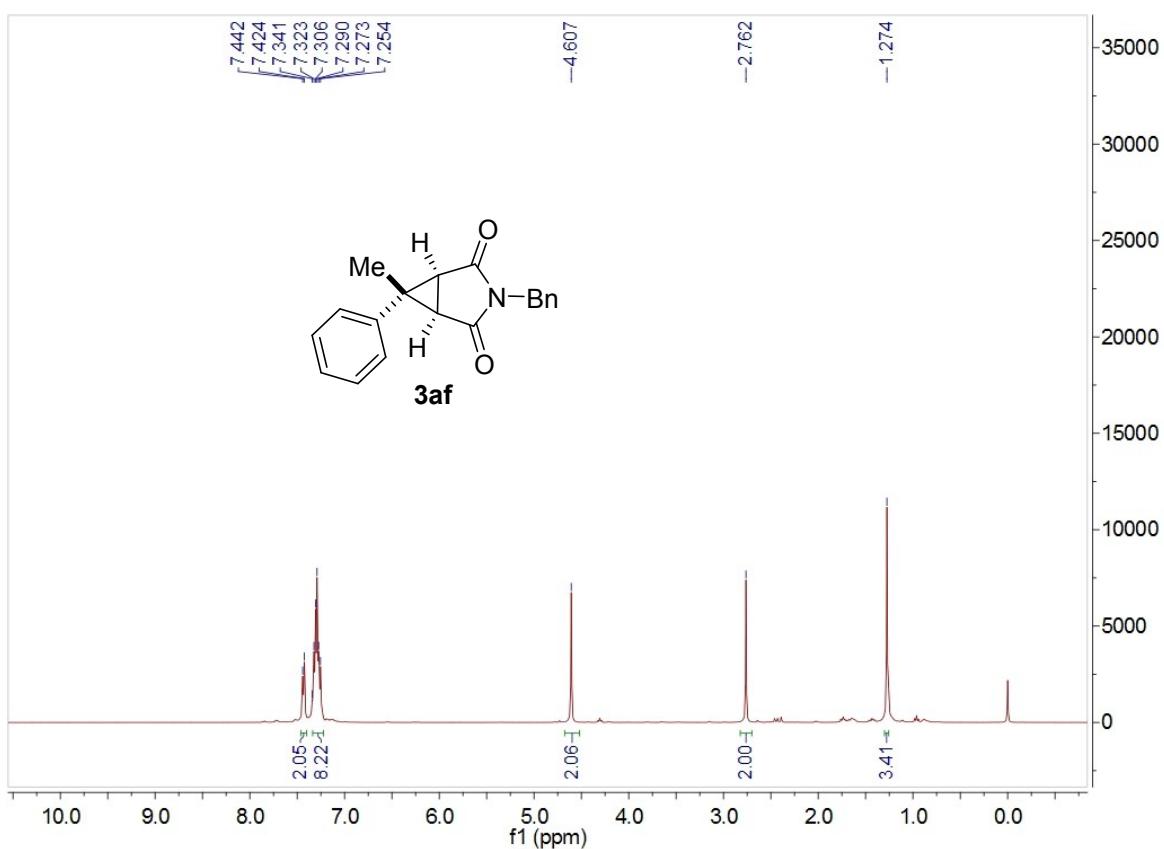
¹H NMR (400 MHz, CDCl₃) spectrum for 3ae



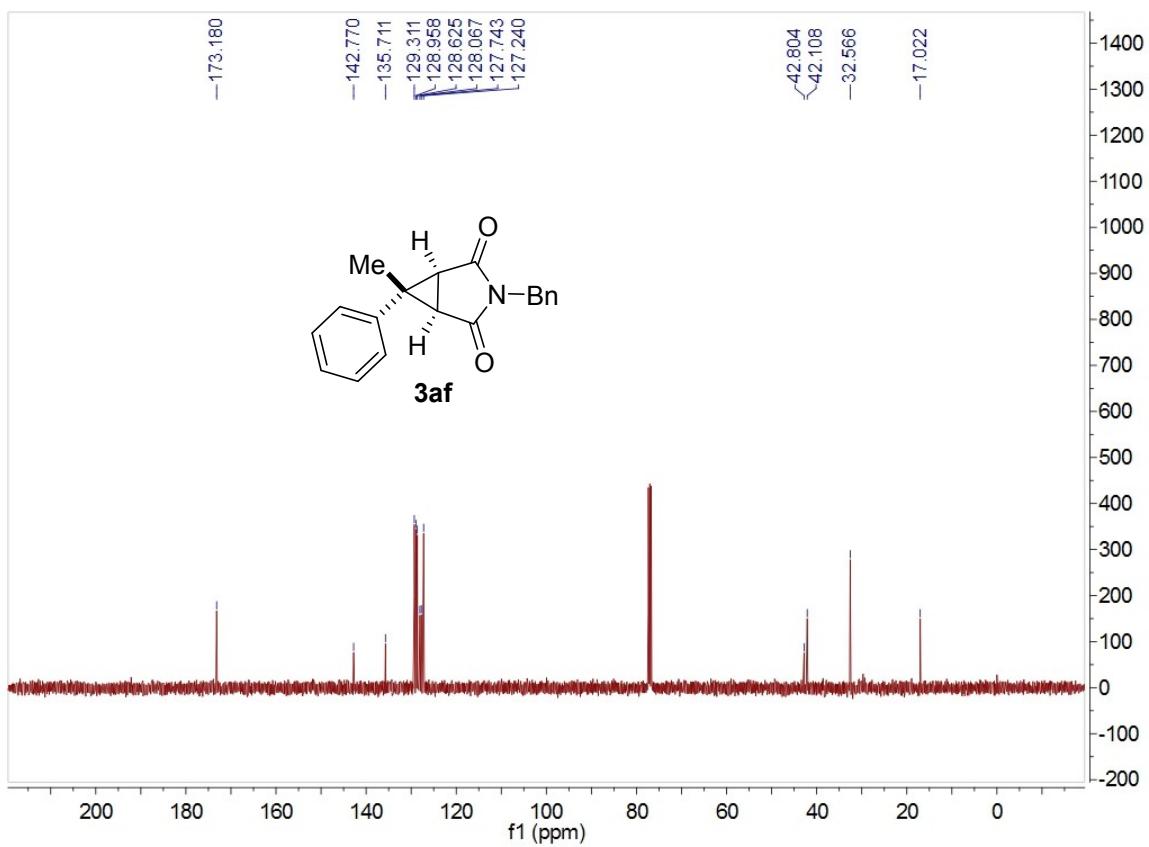
¹³C NMR (100 MHz, CDCl₃) spectrum for 3ae



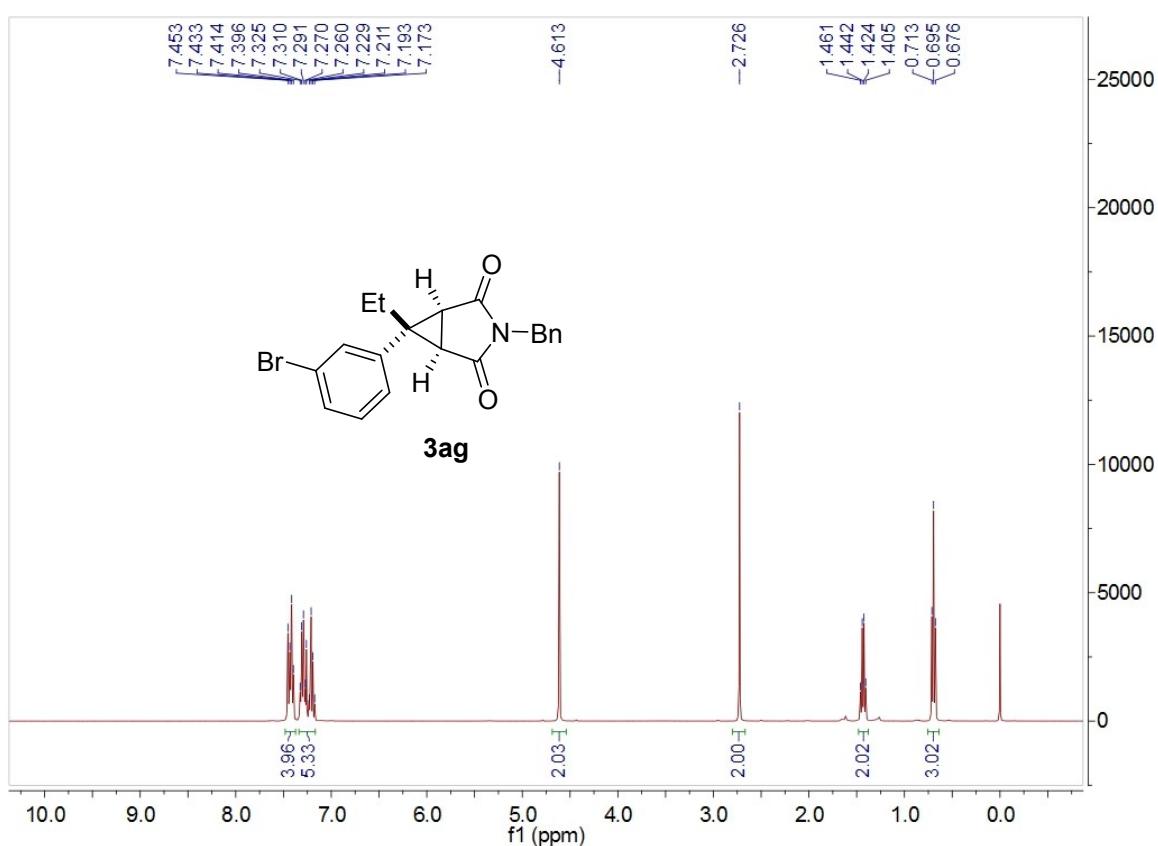
¹H NMR (400 MHz, CDCl₃) spectrum for 3af



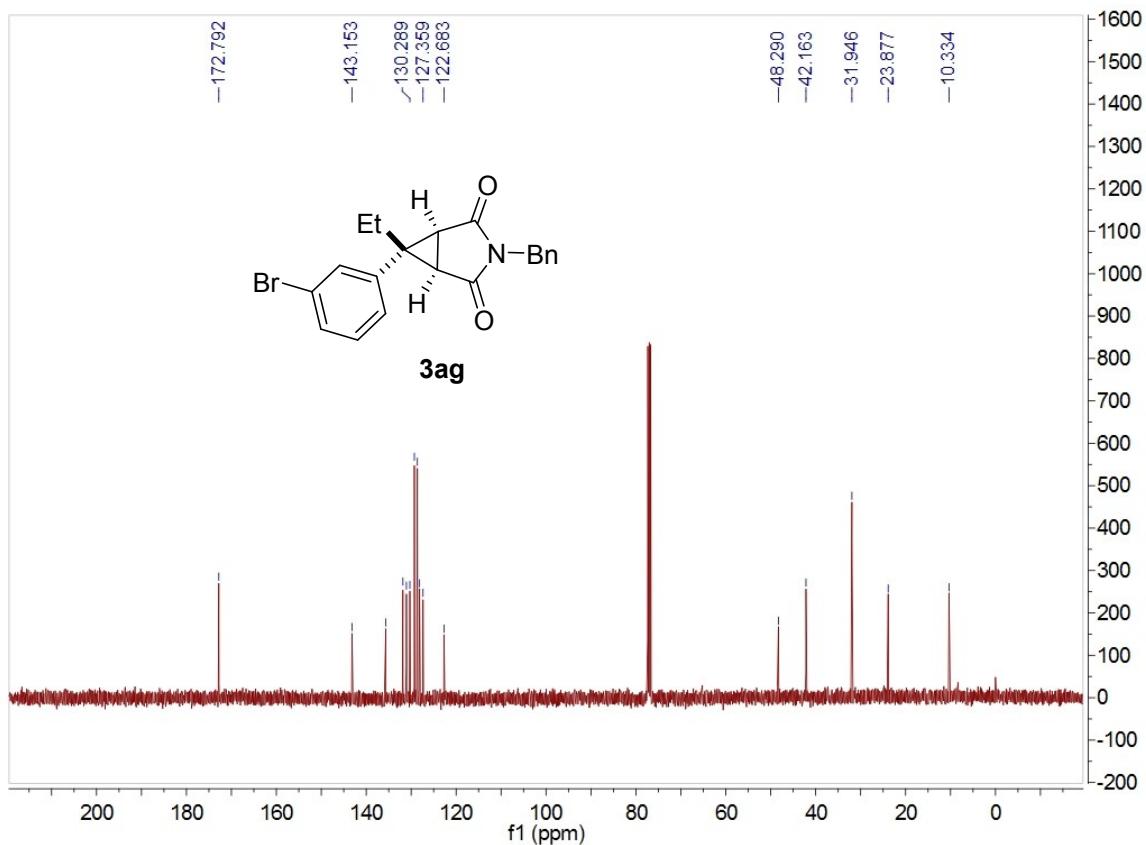
¹³C NMR (100 MHz, CDCl₃) spectrum for 3af



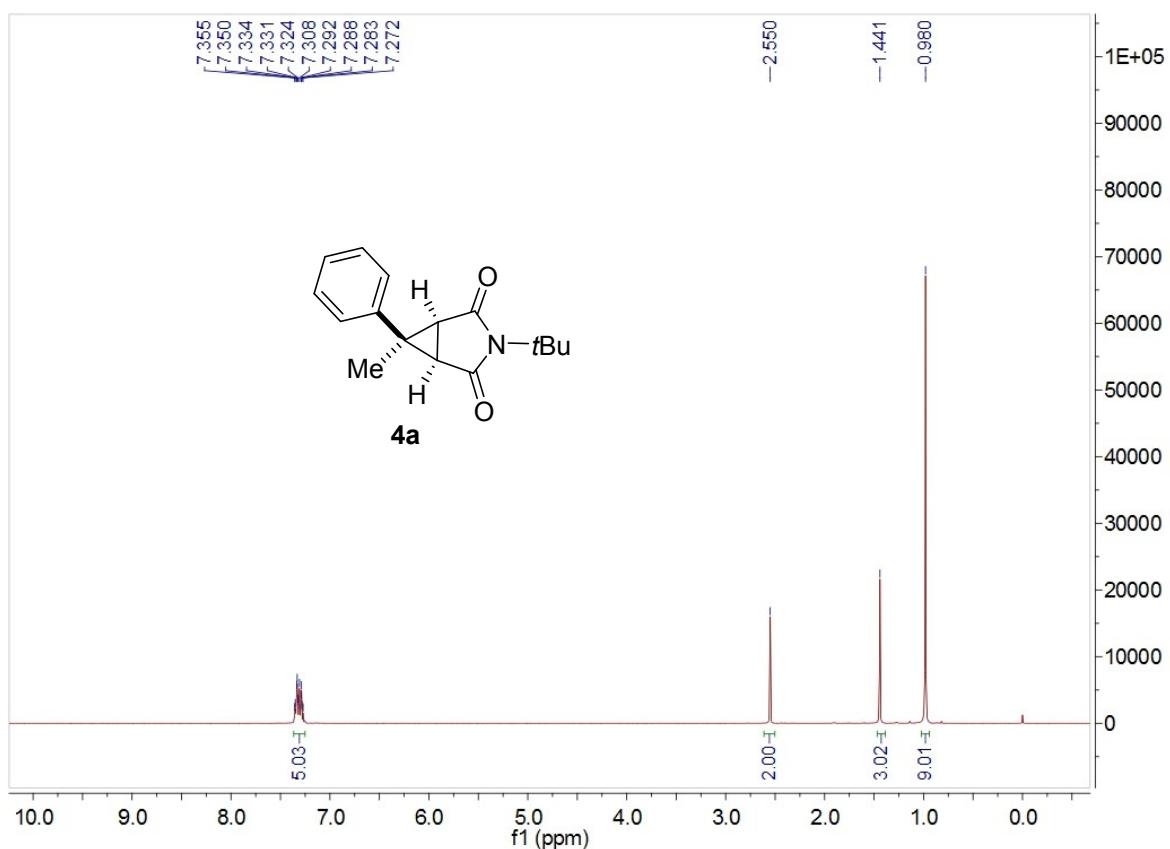
¹H NMR (400 MHz, CDCl₃) spectrum for 3ag



¹³C NMR (100 MHz, CDCl₃) spectrum for 3ag



¹H NMR (400 MHz, CDCl₃) spectrum for 4a



¹³C NMR (100 MHz, CDCl₃) spectrum for 4a

