

## Supporting Information

# Lewis acid-catalyzed [3+2] annulations of oxindole based spirocyclic donor–acceptor cyclopropanes with ynamides

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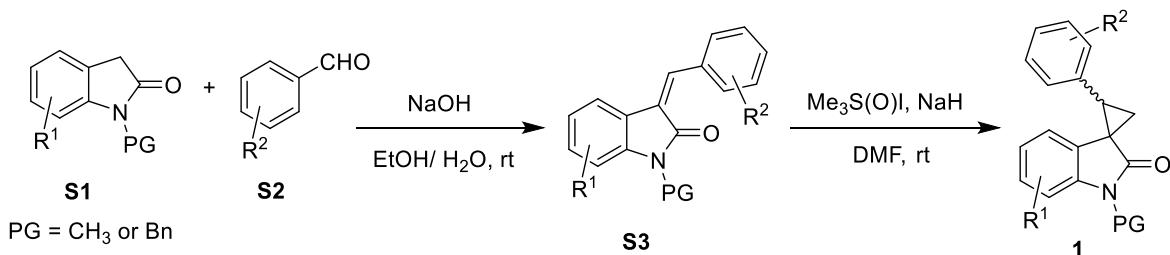
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## 1. General

<sup>1</sup>H NMR and <sup>13</sup>C NMR were recorded on a Bruker-400 MHz spectrometer. Proton chemical shifts are reported in ppm downfield from tetramethylsilane or from the residual solvent as internal standard in CDCl<sub>3</sub> ( $\delta$  7.26 ppm) and in (CD<sub>3</sub>)<sub>2</sub>SO (2.50 ppm). Carbon chemical shifts were internally referenced to the deuterated solvent signals in CDCl<sub>3</sub> ( $\delta$  77.0 ppm) and in (CD<sub>3</sub>)<sub>2</sub>SO (39.5 ppm). High-resolution mass spectra were recorded on a Thermo Scientific LTQ Orbitrap ESI ion trap mass spectrometer. Reagents obtained from commercial sources are used without further purification and all solvents were purified and dried according to standard methods prior to use, unless stated otherwise. Spirocyclopropyl oxindoles **1**<sup>1</sup> and ynamides **2**<sup>2</sup> were synthesized according to reported procedures.

## 2. Procedures for synthesis of spirocyclopropyl oxindoles **1**

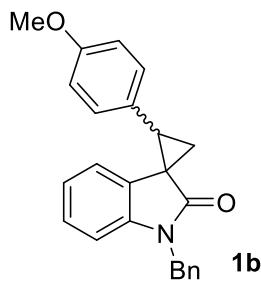


**General Procedure:** Spirocyclopropyl oxindoles **1** were synthesized in two steps starting from *N*-protected oxindoles.<sup>1</sup> **1a**, **1d**, **1f** are known compounds and the characterization data all corresponded to the reported values.<sup>1</sup>

**Step 1:** To a solution of *N*-protected oxindole **S1** (1.0 equiv) in ethanol (1 M) at room temperature was added aromatic aldehyde **S2** (1.1 equiv) under vigorous stirring, and then the solution was added a solution of NaOH (2.0 equiv) in ethanol-water (1:2, 0.5 M) dropwise. The resulting mixture stirred for a specified time. Then water was added to the resulting mixture. After extraction with EtOAc, the organic layer was dried over anhydrous MgSO<sub>4</sub>, and then concentrated under reduced pressure, the crude product was purified by flash silica gel column chromatography.

**Step 2:** A suspension of sodium hydride (2.2 equiv) in anhydrous DMF (50 mg/ mL) was added trimethylsulfoxonium iodide (2.0 equiv) and the suspension was stirred for 30 min at room temperature, then solution of **S3** (1.0 equiv) in DMF was added dropwise. After completion of the addition (TLC controlled), the resulting mixture was quenched with distilled water, after extraction with EtOAc, the organic layer was washed with brine three times, the organic layer was dried over anhydrous MgSO<sub>4</sub>, and then concentrated under reduced pressure, the crude product was purified by flash silica gel column chromatography.

### 1'-benzyl-2-(4-methoxyphenyl)spiro[cyclopropane-1,3'-indolin]-2'-one (**1b**)

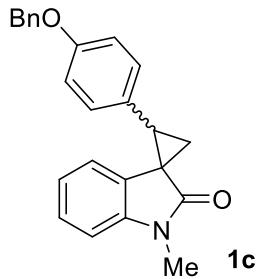


Synthesized by the general procedure using 1-benzylindolin-2-one (500 mg, 2.24 mmol), 4-methoxybenzaldehyde (335.5 mg, 2.46 mmol), and trimethylsulfoxonium iodide (887.3 mg, 4.04 mmol), flash chromatography (petroleum ether:ethyl acetate = 8:1 to 4:1), separated diastereomers.

**1b-1:** 315.8 mg (40%),  $R_f$  = 0.42 (petroleum ether:ethyl acetate = 6:1). Light yellow amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.35-7.42 (m, 4H), 7.28-7.31 (m, 1H), 7.18 (d,  $J$  = 8.5 Hz, 2H), 7.07 (t,  $J$  = 7.7 Hz, 1H), 6.89 (d,  $J$  = 8.6 Hz, 2H), 6.82 (d,  $J$  = 7.8 Hz, 1H), 6.72 (t,  $J$  = 7.5 Hz, 1H), 6.10 (d,  $J$  = 7.4 Hz, 1H), 5.12 (d,  $J$  = 15.7 Hz, 1H), 5.05 (d,  $J$  = 15.7 Hz, 1H), 3.79 (s, 3H), 3.45 (t,  $J$  = 8.5 Hz, 1H), 2.33 (dd,  $J$  = 4.4 Hz, 9.1 Hz, 1H), 2.06 (dd,  $J$  = 4.5 Hz, 7.9 Hz, 1H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 176.3, 158.5, 142.6, 136.0, 130.7, 128.5, 127.3, 127.2, 127.0, 126.8, 126.1, 121.2, 120.6, 113.5, 108.5, 54.8, 43.8, 35.4, 33.0, 22.6 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{24}\text{H}_{22}\text{NO}_2$  356.1645; Found 356.1647.

**1b-2:** 280 mg (35%),  $R_f$  = 0.35 (petroleum ether:ethyl acetate = 6:1). Light yellow amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.28-7.35 (m, 7H), 7.19-7.23 (m, 1H), 7.07-7.11 (m, 1H), 7.0 (dd,  $J$  = 0.6 Hz, 7.3 Hz, 1H), 6.94 (d,  $J$  = 8.7 Hz, 2H), 6.85 (d,  $J$  = 7.7 Hz, 1H), 4.96 (d,  $J$  = 15.6 Hz, 1H), 4.91 (d,  $J$  = 15.6 Hz, 1H), 3.83 (s, 3H), 3.21 (t,  $J$  = 8.8 Hz, 1H), 2.51 (dd,  $J$  = 4.9 Hz, 8.6 Hz, 1H), 2.16 (dd,  $J$  = 4.9 Hz, 9.0 Hz, 1H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 173.7, 158.5, 142.2, 136.2, 130.7, 130.1, 128.4, 127.2, 126.4, 126.1, 121.7, 117.9, 113.2, 108.5, 54.9, 43.6, 38.1, 33.7, 22.3 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{24}\text{H}_{22}\text{NO}_2$  356.1645; Found 356.1648.

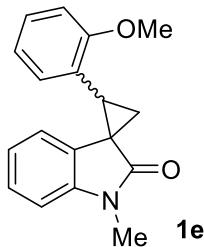
### 2-(4-(benzyloxy)phenyl)-1'-methylspiro[cyclopropane-1,3'-indolin]-2'-one (1c)



Synthesized by the general procedure using 1-methylindolin-2-one (500 mg, 3.40 mmol), 4-methoxybenzaldehyde (335.5 mg, 3.74 mmol), and trimethylsulfoxonium iodide (1.34 g, 6.05 mmol), flash chromatography (petroleum ether:ethyl acetate = 8:1 to 4:1),  $dr$  = 1:1, 1.07 g (77%).

Yellow amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.47-7.49 (m, 4H), 7.41-7.45 (m, 4H), 7.31-7.39 (m, 5H), 7.21 (t,  $J$  = 7.7 Hz, 1H), 7.12-7.17 (m, 3H), 6.96-7.01 (m, 5H), 6.88 (d,  $J$  = 7.8 Hz, 2H), 6.79 (t,  $J$  = 7.6 Hz, 1H), 6.14 (d,  $J$  = 7.4 Hz, 1H), 5.00-5.05 (m, 2H + 2H), 3.39 (t,  $J$  = 8.4 Hz, 1H), 3.33 (s, 3H), 3.17 (s, 3H), 3.14 (t,  $J$  = 8.9 Hz, 1H), 2.43 (dd,  $J$  = 4.8 Hz, 7.9 Hz, 1H), 2.25 (dd,  $J$  = 4.4 Hz, 9.1 Hz, 1H), 2.07 (dd,  $J$  = 4.9 Hz, 9.1 Hz, 1H), 2.01 (dd,  $J$  = 4.5 Hz, 7.9 Hz, 1H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 175.7, 173.1, 157.5, 157.4, 143.3, 142.8, 136.6, 136.4, 130.5, 130.3, 129.8, 128.0, 127.43, 127.36, 127.02, 126.98, 126.9, 126.2, 126.1, 126.0, 121.3, 120.9, 120.2, 117.5, 114.2, 113.6, 107.3, 69.3, 69.2, 37.3, 34.8, 33.4, 32.9, 26.0, 25.8, 22.0 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{24}\text{H}_{22}\text{NO}_2$  356.1645; Found 356.1648.

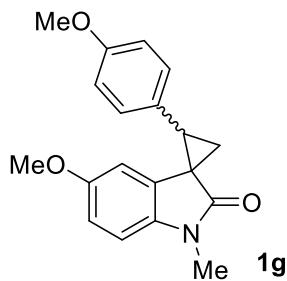
### 2-(2-methoxyphenyl)-1'-methylspiro[cyclopropane-1,3'-indolin]-2'-one (1e)



Synthesized by the general procedure using 1-methylindolin-2-one (500 mg, 3.40 mmol), 2-methoxybenzaldehyde (509.2 mg, 3.74 mmol), and trimethylsulfoxonium iodide (1.27 g, 5.78 mmol), flash chromatography (petroleum ether:ethyl acetate = 8:1 to 4:1), *dr* = 3:1, 646 mg (68%).

Light yellow amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.39 (d,  $J$  = 7.2 Hz, 1H), 7.33 (d,  $J$  = 7.4 Hz, 1H × 3), 7.22-7.28 (m, 2H + 1H × 3), 7.11 (td,  $J$  = 7.7 Hz, 1.0 Hz, 1H × 3), 7.06 (t,  $J$  = 7.5 Hz, 1H), 6.96-7.00 (m, 2H + 1H × 3), 6.87 (d,  $J$  = 7.8 Hz, 1H), 6.84 (d,  $J$  = 7.7 Hz, 1H × 3), 6.77 (d,  $J$  = 8.2 Hz, 1H), 6.62-6.67 (m, 2H × 3), 5.89 (d,  $J$  = 7.4 Hz, 1H × 3), 3.55 (s, 3H), 3.33 (s, 3H × 3), 3.27 (s, 3H × 3), 3.16 (s, 3H), 3.15 (t,  $J$  = 8.5 Hz, 1H × 3), 3.06 (t,  $J$  = 8.7 Hz, 1H), 2.28 (dd,  $J$  = 4.8 Hz, 8.9 Hz, 1H), 2.22 (dd,  $J$  = 4.6 Hz, 8.1 Hz, 1H × 3), 2.03 (dd,  $J$  = 4.8 Hz, 8.9 Hz, 1H), 1.95 (dd,  $J$  = 4.6 Hz, 8.1 Hz, 1H × 3) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ): MAJOR,  $\delta$  = 176.5, 158.8, 143.4, 129.3, 128.5, 128.0, 125.9, 124.1, 120.8, 119.84, 119.1, 110.3, 107.1, 54.8, 32.5, 32.0, 26.3, 21.5 ppm; MINOR,  $\delta$  = 173.9, 158.2, 143.1, 130.9, 129.8, 128.2, 126.2, 123.3, 121.4, 119.76, 117.8, 109.7, 107.2, 54.9, 32.9, 32.6, 26.0, 22.0 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{18}\text{H}_{18}\text{NO}_2$  280.1332; Found 280.1335.

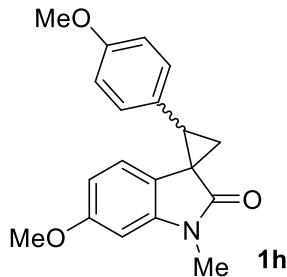
### 5'-methoxy-2-(4-methoxyphenyl)-1'-methylspiro[cyclopropane-1,3'-indolin]-2'-one (1g)



Synthesized by the general procedure using 5-methoxy-1-methylindolin-2-one (500 mg, 2.82 mmol), 4-methoxybenzaldehyde (422.3 mg, 3.10 mmol), and trimethylsulfoxonium iodide (1.04 g, 4.74 mmol), flash chromatography (petroleum ether:ethyl acetate = 8:1 to 4:1), major diastereomer, *dr* = 6.2:1, 601 mg (69%).

Yellow amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ): MAJOR  $\delta$  = 7.15 (d,  $J$  = 8.7 Hz, 2H), 6.87 (d,  $J$  = 8.6 Hz, 2H), 6.78-6.79 (m, 1H), 6.72 (dd,  $J$  = 2.4 Hz, 8.5 Hz, 1H), 5.68 (d,  $J$  = 2.4 Hz, 1H), 3.80 (s, 3H), 3.52 (s, 3H), 3.32-3.36 (m, 1H), 3.33 (s, 3H), 2.22 (dd,  $J$  = 4.5 Hz, 9.1 Hz, 1H), 1.98 (dd,  $J$  = 4.5 Hz, 7.9 Hz, 1H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 176.1, 158.8, 154.9, 137.4, 130.9, 130.2, 128.9, 126.99, 113.7, 113.3, 111.2, 107.9, 55.4, 55.2, 35.2, 34.2, 33.6, 26.6, 22.6 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{19}\text{H}_{20}\text{NO}_3$  310.1438; Found 310.1441.

#### **6'-methoxy-2-(4-methoxyphenyl)-1'-methylspiro[cyclopropane-1,3'-indolin]-2'-one (1h)**

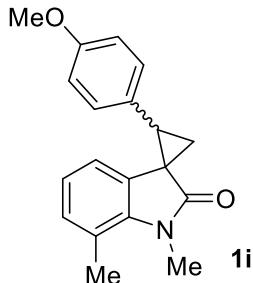


Synthesized by the general procedure using 6-methoxy-1-methylindolin-2-one (500 mg, 2.82 mmol), 4-methoxybenzaldehyde (422.3 mg, 3.10 mmol), and trimethylsulfoxonium iodide (1.05 g, 4.79 mmol), flash chromatography (petroleum ether:ethyl acetate = 8:1 to 4:1), *dr* = 3.7:1, 593 mg (68%).

Yellow amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.22 (d,  $J$  = 8.6 Hz, 2H), 7.08 (d,  $J$  = 8.4 Hz, 2H  $\times$  3.7), 6.79-6.84 (m, 3H + 2H  $\times$  3.7), 6.56 (dd,  $J$  = 2.2 Hz, 8.1 Hz, 1H), 6.46-6.47 (m, 1H  $\times$  4.7), 6.22 (dd,  $J$  = 2.3 Hz, 8.2 Hz, 1H  $\times$  3.7), 5.88 (d,  $J$  = 8.3 Hz, 1H  $\times$  3.7), 3.82 (s, 3H), 3.76 (s, 3H  $\times$  3.7), 3.74 (s, 3H), 3.71 (s, 3H  $\times$  3.7), 3.27 (s, 3H  $\times$  3.7), 3.19 (t,  $J$  = 8.4 Hz, 1H  $\times$  3.7), 3.12 (s, 3H), 3.02 (t,  $J$  = 8.8 Hz, 1H), 2.78 (dd,  $J$  = 4.9 Hz, 8.5 Hz, 1H), 2.09 (dd,  $J$  = 4.4 Hz, 9.1 Hz, 1H  $\times$  3.7), 1.96 (dd,  $J$  = 4.9 Hz, 9.0 Hz, 1H), 1.87 (dd,  $J$  = 4.5 Hz, 7.8 Hz, 1H  $\times$  3.7) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ): MAJOR,  $\delta$  = 177.1, 159.2, 158.5, 144.9, 130.9, 127.4, 121.2, 119.4, 113.7, 105.4, 96.1, 55.4, 55.2, 34.5, 32.9, 26.6, 22.1 ppm; MINOR,  $\delta$

= 174.5, 159.4, 158.7, 144.5, 130.2, 126.4, 122.9, 118.5, 113.3, 105.8, 95.9, 55.6, 55.1, 37.2, 33.4, 26.4, 22.2 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for C<sub>19</sub>H<sub>20</sub>NO<sub>3</sub> 310.1438; Found 310.1442.

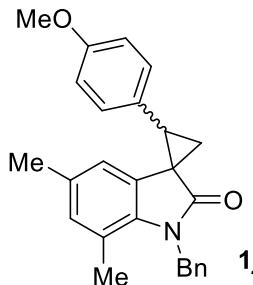
### **2-(4-methoxyphenyl)-1',7'-dimethylspiro[cyclopropane-1,3'-indolin]-2'-one (1i)**



Synthesized by the general procedure using 1,7-dimethylindolin-2-one (500 mg, 3.10 mmol), 4-methoxybenzaldehyde (464.3 mg, 3.41 mmol), and trimethylsulfoxonium iodide (1.09 g, 4.96 mmol), flash chromatography (petroleum ether:ethyl acetate = 8:1 to 4:1), major diastereomer, *dr* = 6:1, 582 mg (64%).

Light yellow amorphous solid. New compound. <sup>1</sup>H NMR (400 MHz, CDCl<sub>3</sub>): MAJOR  $\delta$  = 7.10 (d, *J* = 8.6 Hz, 2H), 6.87 (d, *J* = 7.8 Hz, 1H), 6.82 (d, *J* = 8.6 Hz, 2H), 6.59 (t, *J* = 7.6 Hz, 1H), 5.84 (d, *J* = 7.4 Hz, 1H), 3.77 (s, 3H), 3.61 (s, 3H), 3.28 (t, *J* = 8.5 Hz, 1H), 2.59 (s, 3H), 2.17 (dd, *J* = 4.4 Hz, 9.1 Hz, 1H), 1.91 (dd, *J* = 4.4 Hz, 8.0 Hz, 1H) ppm; <sup>13</sup>C{<sup>1</sup>H} NMR (100 MHz, CDCl<sub>3</sub>):  $\delta$  = 177.2, 158.7, 141.6, 131.1, 130.3, 128.1, 127.1, 121.3, 119.4, 118.7, 113.7, 113.3, 55.2, 35.8, 32.9, 29.9, 23.2, 19.7, 19.2 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for C<sub>19</sub>H<sub>20</sub>NO<sub>2</sub> 294.1489; Found 294.1492.

### **1'-benzyl-2-(4-methoxyphenyl)-5',7'-dimethylspiro[cyclopropane-1,3'-indolin]-2'-one (1j)**

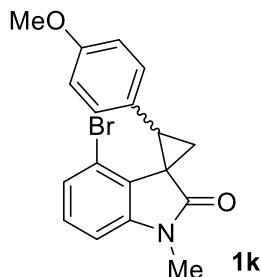


Synthesized by the general procedure using 1-benzyl-5,7-dimethylindolin-2-one (500 mg, 1.99 mmol), 4-methoxybenzaldehyde (298.2 mg, 2.19 mmol), and trimethylsulfoxonium iodide (744.5 mg, 3.38 mmol), flash chromatography (petroleum ether:ethyl acetate = 8:1 to 4:1), *dr* = 1:1, 261 mg (68%).

Light yellow amorphous solid. New compound. <sup>1</sup>H NMR (400 MHz, CDCl<sub>3</sub>):  $\delta$  = 7.26 (d, *J* = 8.0 Hz, 4H), 7.13-7.20 (m, 6H), 7.10 (d, *J* = 8.4 Hz, 2H), 7.01 (d, *J* = 7.2 Hz, 2H), 6.81 (dd, *J* = 2.1 Hz, 8.7 Hz, 4H), 6.69 (s, 1H), 6.61 (s, 1H), 6.57 (s, 1H), 5.69 (s, 1H), 5.26 (s, 2H), 5.02-5.15 (m, 2H), 3.72 (s, 3H), 3.70 (s, 3H), 3.33 (t, *J* = 8.4 Hz, 1H), 3.07 (t, *J* = 8.7 Hz, 1H), 2.41 (dd, *J* = 4.8 Hz, 8.5 Hz, 1H), 2.27 (s, 3H), 2.24-2.25 (m, 1H), 2.22 (s, 3H), 2.20 (s, 3H), 2.04 (dd, *J* = 4.8 Hz,

9.0 Hz, 1H), 2.00 (s, 3H), 1.92-1.94 (m, 1H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ): MAJOR,  $\delta$  = 177.6, 158.6, 138.6, 138.2, 131.5, 131.1, 130.92, 130.2, 128.6, 127.1, 126.8, 125.7, 119.6, 119.1, 113.7, 55.2, 45.2, 38.9, 33.6, 23.2, 20.8, 18.61 ppm; MINOR,  $\delta$  = 174.7, 158.8, 138.2, 138.1, 131.4, 130.94, 130.7, 128.7, 128.2, 127.0, 126.4, 125.6, 119.3, 116.6, 113.2, 55.0, 44.9, 36.2, 32.8, 22.6, 20.7, 18.56 ppm. HRMS (ESI) m/z:  $[\text{M} + \text{H}]^+$  Calcd. for  $\text{C}_{26}\text{H}_{26}\text{NO}_2$  384.1958; Found 384.1961.

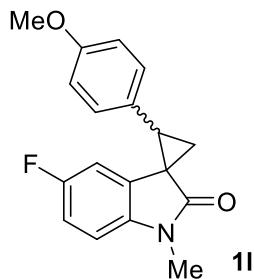
#### **4'-bromo-2-(4-methoxyphenyl)-1'-methylspiro[cyclopropane-1,3'-indolin]-2'-one (1k)**



Synthesized by the general procedure using 4-bromo-1-methylindolin-2-one (500 mg, 2.21 mmol), 4-methoxybenzaldehyde (331.0 mg, 2.43 mmol), and trimethylsulfoxonium iodide (779.0 mg, 3.54 mmol), flash chromatography (petroleum ether:ethyl acetate = 8:1 to 4:1), single diastereomer, 475 mg (60%).

Light yellow amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.25 (d,  $J$  = 8.7 Hz, 2H), 7.16 (dd,  $J$  = 0.7 Hz, 8.1 Hz, 1H), 7.08-7.12 (m, 1H), 6.81-6.87 (m, 3H), 4.05 (t,  $J$  = 9.0 Hz, 1H), 3.79 (s, 3H), 3.15 (s, 3H), 2.87 (dd,  $J$  = 4.8 Hz, 9.2 Hz, 1H), 2.21 (dd,  $J$  = 4.8 Hz, 8.8 Hz, 1H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 172.8, 158.4, 145.3, 130.5, 127.6, 126.7, 126.0, 125.8, 114.6, 113.1, 106.9, 54.9, 35.4, 32.5, 26.4, 19.0 ppm. HRMS (ESI) m/z:  $[\text{M} + \text{H}]^+$  Calcd. for  $\text{C}_{18}\text{H}_{17}\text{BrNO}_2$  358.0437; Found 358.0440.

#### **5'-fluoro-2-(4-methoxyphenyl)-1'-methylspiro[cyclopropane-1,3'-indolin]-2'-one (1l)**

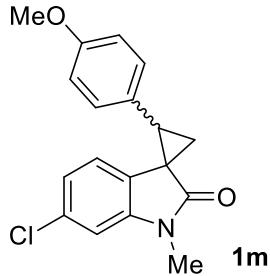


Synthesized by the general procedure using 5-fluoro-1-methylindolin-2-one (500 mg, 3.03 mmol), 4-methoxybenzaldehyde (453.8 mg, 3.33 mmol), and trimethylsulfoxonium iodide (1.12 g, 5.09 mmol), flash chromatography (petroleum ether:ethyl acetate = 8:1 to 4:1), major diastereomer,  $dr$  = 4.5:1, 558.6 mg (62%).

Light yellow amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ): MAJOR  $\delta$  = 7.05 (d,  $J$  = 8.6 Hz, 2H), 6.76-6.81 (m, 3H), 6.70 (dd,  $J$  = 4.3 Hz, 8.4 Hz, 1H), 5.73 (dd,  $J$  = 2.5 Hz,

8.6 Hz, 1H), 3.72 (s, 3H), 3.27 (t,  $J$  = 8.6 Hz, 1H), 3.25 (s, 3H), 2.15 (dd,  $J$  = 4.5 Hz, 9.1 Hz, 1H), 1.91 (dd,  $J$  = 4.6 Hz, 8.0 Hz, 1H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 175.8, 158.7, 158.1 (d,  $J_{\text{C}-\text{F}}$  = 237.3 Hz), 139.4 (d,  $J_{\text{C}-\text{F}}$  = 2.0 Hz), 130.5, 129.1 (d,  $J_{\text{C}-\text{F}}$  = 8.8 Hz), 126.2, 113.6, 112.3 (d,  $J_{\text{C}-\text{F}}$  = 23.6 Hz), 108.4 (d,  $J_{\text{C}-\text{F}}$  = 26.0 Hz), 107.7 (d,  $J_{\text{C}-\text{F}}$  = 8.5 Hz), 54.8, 35.4, 33.4 (d,  $J_{\text{C}-\text{F}}$  = 2.2 Hz), 26.4, 22.7 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{18}\text{H}_{17}\text{FNO}_2$  298.1238; Found 298.1242.

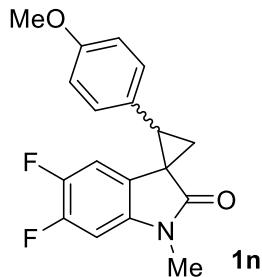
### **6'-chloro-2-(4-methoxyphenyl)-1'-methylspiro[cyclopropane-1,3'-indolin]-2'-one (1m)**



Synthesized by the general procedure using 6-chloro-1-methylindolin-2-one (500 mg, 2.75 mmol), 4-methoxybenzaldehyde (411.2 mg, 3.02 mmol), and trimethylsulfoxonium iodide (992.5 mg, 4.51 mmol), flash chromatography (petroleum ether:ethyl acetate = 8:1 to 4:1), major diastereomer,  $dr$  = 14:1, 569.5 mg (66%).

Yellow amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ): MAJOR  $\delta$  = 7.10 (d,  $J$  = 8.6 Hz, 2H), 6.87 (d,  $J$  = 1.8 Hz, 1H), 6.84 (d,  $J$  = 8.7 Hz, 2H), 6.69 (dd,  $J$  = 1.8 Hz, 8.0 Hz, 1H), 5.89 (d,  $J$  = 8.0 Hz, 1H), 3.80 (s, 3H), 3.28-3.32 (m, 1H), 3.31 (s, 3H), 2.20 (dd,  $J$  = 4.5 Hz, 9.1 Hz, 1H), 1.98 (dd,  $J$  = 4.5 Hz, 7.9 Hz, 1H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 176.4, 158.9, 144.9, 132.3, 130.9, 126.7, 125.9, 121.4, 121.3, 121.2, 113.8, 108.4, 55.2, 35.6, 33.1, 26.7, 22.8 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{18}\text{H}_{17}\text{ClNO}_2$  314.0942; Found 314.0945.

### **5',6'-difluoro-2-(4-methoxyphenyl)-1'-methylspiro[cyclopropane-1,3'-indolin]-2'-one (1n)**

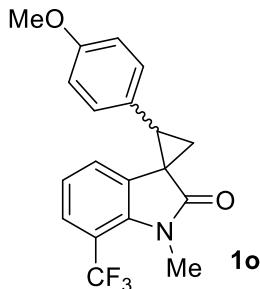


Synthesized by the general procedure using 5,6-difluoro-1-methylindolin-2-one (500 mg, 2.73 mmol), 4-methoxybenzaldehyde (408.4 mg, 3.00 mmol), and trimethylsulfoxonium iodide (913.3 mg, 4.15 mmol), flash chromatography (petroleum ether:ethyl acetate = 8:1 to 4:1),  $dr$  = 2.6:1, 482 mg (56%).

Light yellow amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.19-7.23 (m, 2H), 7.05-7.08 (m, 2H  $\times$  2.6), 6.90-6.95 (m, 1H), 6.79-6.83 (m, 1H + 2H  $\times$  2.6), 6.72-6.76 (m, 1H

$\times 2.6)$ , 6.64-6.69 (m, 2H), 5.73-5.82 (m, 1H  $\times$  2.6), 3.75 (s, 3H  $\times$  2.6), 3.74 (s, 3H), 3.28 (s, 3H  $\times$  2.6), 3.25 (s, 3H), 3.04-3.13 (m, 1H  $\times$  3.6), 2.33-2.38 (m, 1H), 2.14-2.19 (m, 1H  $\times$  2.6), 1.98-2.03 (m, 1H), 1.90-1.95 (m, 1H  $\times$  2.6) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 176.2, 176.0, 149.5, (dd,  $J_{\text{C}-\text{F}} = 14.1$  Hz, 243.52 Hz), 145.6 (dd,  $J_{\text{C}-\text{F}} = 13.1$  Hz, 239.0 Hz), 139.71, 139.69, 130.8, 130.7, 130.19, 130.18, 129.45, 129.36, 126.4, 126.2, 125.8, 125.5, 123.0 (dd,  $J_{\text{C}-\text{F}} = 3.6$  Hz, 7.3 Hz), 113.9, 113.8, 112.5 (d,  $J_{\text{C}-\text{F}} = 23.5$  Hz), 110.1 (d,  $J_{\text{C}-\text{F}} = 21.3$  Hz), 108.8, 108.6, 107.9, 107.8, 55.1, 55.0, 35.7, 35.5, 33.7, 33.68, 26.7, 26.6, 22.9, 22.8 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{18}\text{H}_{16}\text{F}_2\text{NO}_2$  316.1144; Found 316.1147.

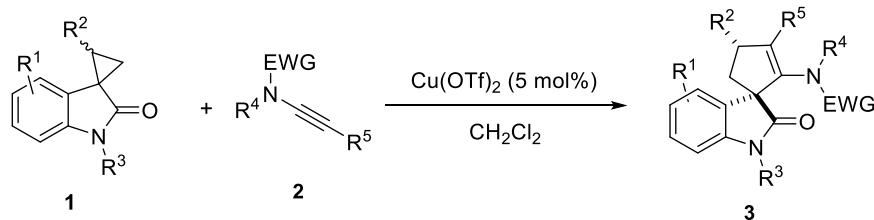
### 2-(4-methoxyphenyl)-1'-methyl-7'-(trifluoromethyl)spiro[cyclopropane-1,3'-indolin]-2'-one (1o)



Synthesized by the general procedure using 1-methyl-7-(trifluoromethyl)indolin-2-one (500 mg, 2.32 mmol), 4-methoxybenzaldehyde (347.4 mg, 2.55 mmol), and trimethylsulfoxonium iodide (796.5 mg, 3.62 mmol), flash chromatography (petroleum ether:ethyl acetate = 8:1 to 4:1), major diastereomer, *dr* = 25:1, 483 mg (60%).

Light yellow amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ): MAJOR  $\delta$  = 7.45 (d,  $J = 8.0$  Hz, 1H), 7.12 (d,  $J = 8.6$  Hz, 2H), 6.86 (d,  $J = 8.6$  Hz, 2H), 6.77 (t,  $J = 7.8$  Hz, 1H), 6.16 (d,  $J = 7.4$  Hz, 1H), 3.81 (s, 3H), 3.57 (d,  $J = 2.3$  Hz, 3H), 3.39 (t,  $J = 8.6$  Hz, 1H), 2.29 (dd,  $J = 4.5$  Hz, 9.2 Hz, 1H), 2.05 (dd,  $J = 4.5$  Hz, 8.1 Hz, 1H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 177.2, 158.9, 141.39, 141.38, 131.0, 129.9, 126.4, 124.3 (q,  $J_{\text{C}-\text{F}} = 5.9$  Hz), 123.7 (q,  $J_{\text{C}-\text{F}} = 269.7$  Hz), 123.8, 120.6, 113.8, 112.0 (q,  $J_{\text{C}-\text{F}} = 32.5$  Hz), 55.1, 36.9, 32.2, 29.1 (q,  $J_{\text{C}-\text{F}} = 6.4$  Hz), 23.8 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{19}\text{H}_{17}\text{F}_3\text{NO}_2$  348.1206; Found 348.1210.

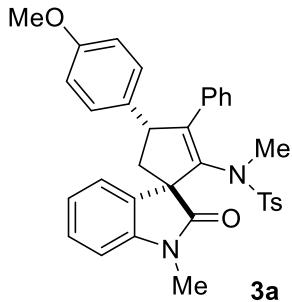
### 3. Procedures for the synthesis of spirocyclopenteneoxindoles 3



**General Procedure:** To a mixture of spirocyclopropyl oxindole **1** (0.10 mmol, 1.0 equiv, diastereomeric mixture) and ynamide **2** (0.12 mmol, 1.2 equiv) in  $\text{CH}_2\text{Cl}_2$  (1.0 mL),  $\text{Cu}(\text{OTf})_2$

(1.8 mg, 0.005 mmol, 0.05 equiv) was added. The mixture was stirred at room temperature (25 °C) in air for 4 h. Then, the solvent was removed in vacuo and the residue was purified by silica gel flash chromatography (petroleum ether:ethyl acetate = 8:1 to 4:1) to provide product **3**. Ynamide **2** are known compounds and the characterization data all corresponded to the reported values.<sup>2</sup>

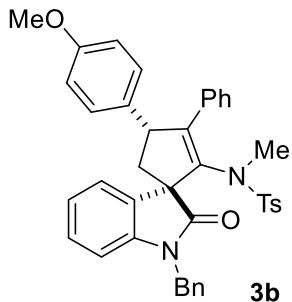
**N-(4-(4-methoxyphenyl)-1'-methyl-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N,N-dimethylbenzenesulfonamide (3a)**



Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer, *dr* = 9:1, 48.6 mg (86%).

Colorless solid. mp 253-255 °C. New compound. <sup>1</sup>H NMR (400 MHz, CDCl<sub>3</sub>): δ = 7.61 (d, *J* = 5.7 Hz, 1H), 7.27-7.33 (m, 5H), 7.12-7.14 (m, 4H), 6.80-6.93 (m, 6H), 6.74 (d, *J* = 7.7 Hz, 1H), 4.71 (t, *J* = 7.9 Hz, 1H), 3.67 (s, 3H), 3.13 (s, 3H), 2.96 (dd, *J* = 8.4 Hz, 13.4 Hz, 1H), 2.65 (s, 3H), 2.32 (dd, *J* = 7.8 Hz, 13.4 Hz, 1H), 2.26 (s, 3H) ppm; <sup>13</sup>C{<sup>1</sup>H} NMR (100 MHz, CDCl<sub>3</sub>): δ = 178.7, 158.0, 149.5, 143.8, 142.2, 136.8, 136.2, 135.4, 134.5, 130.7, 129.0, 128.7, 128.3, 128.1, 127.6, 127.4, 127.3, 126.9, 122.4, 113.7, 107.6, 61.5, 54.8, 50.1, 45.4, 37.0, 26.1, 21.0 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for C<sub>34</sub>H<sub>33</sub>N<sub>2</sub>O<sub>4</sub>S 565.2156; Found 565.2158.

**N-(1'-benzyl-4-(4-methoxyphenyl)-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N,N-dimethylbenzenesulfonamide (3b)**

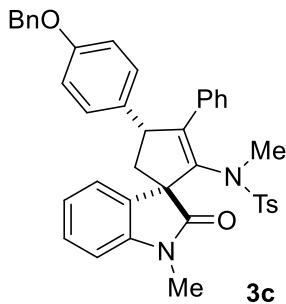


Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer, *dr* > 20:1, 41.0 mg (64%).

Colorless amorphous solid. New compound. <sup>1</sup>H NMR (400 MHz, CDCl<sub>3</sub>): δ = 7.71-7.72 (m, 1H), 7.15-7.36 (m, 11H), 7.08-7.11 (m, 4H), 6.81-6.87 (m, 4H), 6.71-6.75 (m, 2H), 5.02 (d, *J* = 15.5

Hz, 1H), 4.70-4.76 (m, 2H), 3.75 (s, 3H), 2.98 (dd,  $J$  = 8.4 Hz, 13.4 Hz, 1H), 2.49 (s, 3H), 2.36 (dd,  $J$  = 8.1 Hz, 13.4 Hz, 1H), 2.26 (s, 3H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 179.2, 158.2, 151.0, 143.2, 142.6, 137.1, 136.2, 135.6, 134.5, 131.3, 131.0, 129.3, 129.2, 129.0, 128.8, 128.5, 128.4, 128.2, 127.8, 127.4, 127.3, 126.9, 122.9, 113.9, 108.8, 62.1, 55.1, 52.6, 50.5, 45.6, 43.8, 21.3 ppm. HRMS (ESI) m/z:  $[\text{M} + \text{H}]^+$  Calcd. for  $\text{C}_{40}\text{H}_{37}\text{N}_2\text{O}_4\text{S}$  641.2469; Found 641.2472.

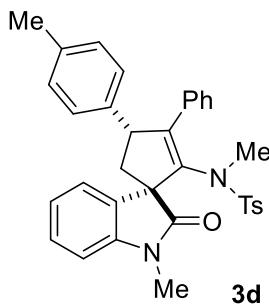
**N-(4-(4-(benzyloxy)phenyl)-1'-methyl-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N,4-dimethylbenzenesulfonamide (3c)**



Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer,  $dr > 20:1$ , 51.3 mg (80%).

Colorless amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.63-7.65 (m, 1H), 7.24-7.44 (m, 10H), 7.16-7.20 (m, 5H), 6.92-6.95 (m, 5H), 6.78 (d,  $J$  = 7.7 Hz, 1H), 5.01 (s, 2H), 4.72 (t,  $J$  = 8.0 Hz, 1H), 3.18 (s, 3H), 2.98 (dd,  $J$  = 8.4 Hz, 13.4 Hz, 1H), 2.66 (s, 3H), 2.35 (dd,  $J$  = 7.8 Hz, 13.4 Hz, 1H), 2.32 (s, 3H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 179.0, 157.4, 149.7, 143.9, 142.4, 137.0, 136.8, 136.4, 135.9, 134.7, 130.9, 129.2, 128.8, 128.5, 128.4, 128.3, 127.8, 127.6, 127.3, 127.1, 122.6, 114.8, 107.8, 69.8, 61.7, 50.3, 45.5, 37.4, 26.2, 21.2 ppm. HRMS (ESI) m/z:  $[\text{M} + \text{H}]^+$  Calcd. for  $\text{C}_{40}\text{H}_{37}\text{N}_2\text{O}_4\text{S}$  641.2469; Found 641.2471.

**N,4-dimethyl-N-(1'-methyl-2'-oxo-3-phenyl-4-(*p*-tolyl)spiro[cyclopentane-1,3'-indolin]-2-en-2-yl)benzenesulfonamide (3d)**

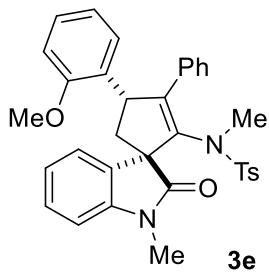


Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1),  $dr = 2:1$ , 39.0 mg (71%).

Colorless amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.61-7.66 (m, 3H), 7.12-7.40 (m, 28H), 7.09 (d,  $J$  = 7.8 Hz, 4H), 7.02 (d,  $J$  = 7.9 Hz, 2H), 6.80-6.86 (m, 12H), 6.76

(d,  $J = 7.7$  Hz, 2H), 4.70 (t,  $J = 8.2$  Hz, 1H  $\times$  3), 3.19 (s, 3H), 3.16 (s, 3H  $\times$  2), 2.95 (dd,  $J = 8.5$  Hz, 13.4 Hz, 1H  $\times$  2), 2.78 (s, 3H), 2.72-2.76 (m, 1H), 2.62-2.67 (m, 1H), 2.62 (s, 3H  $\times$  2), 2.30-2.35 (m, 2H), 2.31 (s, 3H), 2.30 (s, 3H  $\times$  2), 2.28 (s, 3H  $\times$  2), 2.23 (s, 3H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta = 179.1, 178.8, 150.0, 149.8, 144.0, 143.5, 142.6, 142.5, 140.6, 139.7, 137.2, 136.5, 136.1, 136.0, 135.9, 135.8, 135.1, 134.8, 131.7, 131.0, 129.3, 129.0, 128.9, 128.6, 128.4, 128.2, 128.1, 128.0, 127.8, 127.6, 127.5, 127.2, 125.5, 122.7, 107.9, 107.8, 61.8, 61.7, 53.4, 52.6, 51.0, 50.8, 45.2, 44.6, 37.8, 37.2, 26.5, 26.3, 21.3, 21.0$  ppm. HRMS (ESI) m/z:  $[\text{M} + \text{H}]^+$  Calcd. for  $\text{C}_{34}\text{H}_{33}\text{N}_2\text{O}_3\text{S}$  549.2206; Found 549.2207.

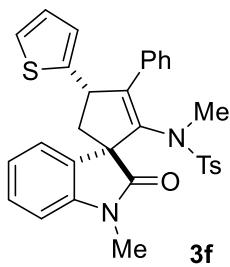
**N-(4-(2-methoxyphenyl)-1'-methyl-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N,4-dimethylbenzenesulfonamide (3e)**



Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer,  $dr = 3.5:1$ , 42.9 mg (76%).

Colorless amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta = 7.54-7.56$  (m, 1H), 7.35-7.39 (m, 2H), 7.29 (td,  $J = 7.7$  Hz, 0.8 Hz, 1H), 7.12-7.20 (m, 6H), 6.87-6.95 (m, 5H), 6.78 (d,  $J = 8.1$  Hz, 1H), 6.69 (d,  $J = 7.5$  Hz, 1H), 5.10 (t,  $J = 8.0$  Hz, 1H), 3.80 (s, 3H), 3.10 (s, 3H), 2.97 (dd,  $J = 8.8$  Hz, 13.1 Hz, 1H), 2.71 (s, 3H), 2.33-2.41 (m, 1H), 2.30 (s, 3H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta = 179.0, 157.2, 149.9, 143.4, 142.5, 142.2, 136.1, 136.0, 135.0, 134.6, 131.6, 129.4, 128.8, 128.3, 128.0, 127.8, 127.7, 127.6, 126.9, 122.6, 120.4, 110.3, 107.6, 61.8, 55.0, 45.4, 43.4, 37.2, 26.2, 21.2$  ppm. HRMS (ESI) m/z:  $[\text{M} + \text{H}]^+$  Calcd. for  $\text{C}_{34}\text{H}_{33}\text{N}_2\text{O}_4\text{S}$  565.2156; Found 565.2159.

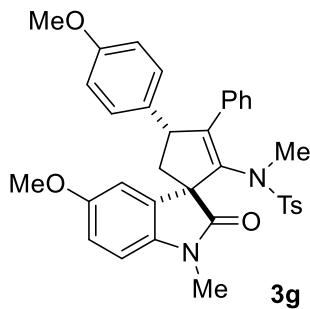
**N,4-dimethyl-N-(1'-methyl-2'-oxo-3-phenyl-4-(thiophen-2-yl)spiro[cyclopentane-1,3'-indolin]-2-en-2-yl)benzenesulfonamide (3f)**



Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer,  $dr = 15:1$ , 44.3 mg (82%).

Colorless amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.55-7.56 (m, 1H), 7.34 (td,  $J$  = 7.7 Hz, 0.9 Hz, 1H), 7.13-7.26 (m, 7H), 6.96-6.99 (m, 4H), 6.86-6.87 (m, 2H), 6.77 (d,  $J$  = 7.7 Hz, 1H), 5.00 (t,  $J$  = 7.8 Hz, 1H), 3.16 (s, 3H), 3.04 (dd,  $J$  = 8.2 Hz, 13.3 Hz, 1H), 2.62 (s, 3H), 2.48 (dd,  $J$  = 7.5 Hz, 13.3 Hz, 1H), 2.32 (s, 3H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 178.7, 149.0, 147.1, 144.0, 142.6, 137.1, 136.3, 134.3, 130.8, 129.0, 128.6, 128.4, 127.8, 127.3, 126.7, 125.5, 124.1, 122.7, 107.9, 61.6, 52.6, 46.3, 45.9, 37.9, 26.4, 21.3 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{31}\text{H}_{29}\text{N}_2\text{O}_3\text{S}_2$  541.1614; Found 541.1619.

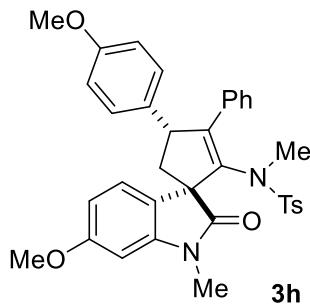
**N-(5'-methoxy-4-(4-methoxyphenyl)-1'-methyl-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N,4-dimethylbenzenesulfonamide (3g)**



Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer, *dr* = 17:1, 51.1 mg (86%).

Colorless amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.31-7.36 (m, 3H), 7.23-7.27 (m, 5H), 6.97-6.99 (m, 4H), 6.88-6.91 (m, 3H), 6.69 (d,  $J$  = 8.4 Hz, 1H), 4.73 (t,  $J$  = 7.8 Hz, 1H), 3.86 (s, 3H), 3.79 (s, 3H), 3.18 (s, 3H), 3.02 (dd,  $J$  = 8.4 Hz, 13.3 Hz, 1H), 2.71 (s, 3H), 2.35 (s, 3H), 2.34 (dd,  $J$  = 7.4 Hz, 13.3 Hz, 1H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 178.6, 158.1, 155.97, 149.5, 142.4, 137.2, 136.8, 136.5, 135.8, 135.7, 134.8, 134.7, 131.98, 129.3, 129.2, 128.8, 128.3, 127.97, 127.79, 127.6, 127.0, 113.8, 113.5, 112.98, 108.0, 62.1, 55.96, 55.89, 55.84, 55.79, 54.97, 50.2, 45.7, 26.3, 21.2 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{35}\text{H}_{35}\text{N}_2\text{O}_5\text{S}$  595.2261; Found 591.2265.

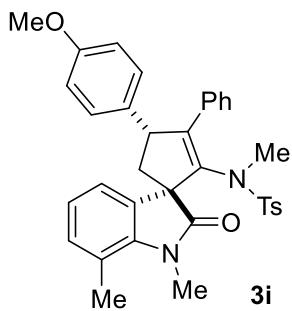
**N-(6'-methoxy-4-(4-methoxyphenyl)-1'-methyl-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N,4-dimethylbenzenesulfonamide (3h)**



Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer, *dr* > 40:1, 49.9 mg (84%).

Colorless amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.45 (d,  $J$  = 7.6 Hz, 1H), 7.13-7.24 (m, 7H), 6.94 (br, 4H), 6.82 (d,  $J$  = 8.5 Hz, 2H), 6.63 (dd,  $J$  = 2.2 Hz, 8.2 Hz, 1H), 6.29 (s, 1H), 4.66 (t,  $J$  = 7.9 Hz, 1H), 3.85 (s, 3H), 3.75 (s, 3H), 3.12 (s, 3H), 2.92 (dd,  $J$  = 8.5 Hz, 13.4 Hz, 1H), 2.65 (s, 3H), 2.31 (s, 3H), 2.24 (dd,  $J$  = 7.6 Hz, 13.4 Hz, 1H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 179.8, 160.7, 158.2, 149.5, 145.2, 142.4, 137.3, 136.9, 135.9, 134.9, 129.3, 128.9, 128.4, 127.9, 127.7, 127.0, 122.8, 113.9, 106.3, 96.1, 61.3, 55.5, 55.1, 50.3, 45.7, 29.7, 26.4, 21.4 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{35}\text{H}_{35}\text{N}_2\text{O}_5\text{S}$  595.2261; Found 591.2263.

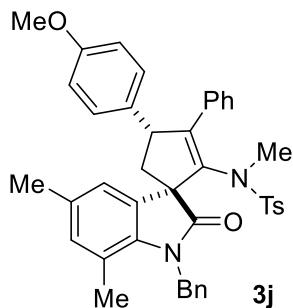
**N-(4-(4-methoxyphenyl)-1',7'-dimethyl-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N,4-dimethylbenzenesulfonamide (3i)**



Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer, *dr* = 20:1, 46.3 mg (80%).

Colorless amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.40-7.51 (m, 2H), 7.32-7.35 (m, 3H), 7.20-7.24 (m, 3H), 7.08-7.11 (m, 2H), 7.01-7.05 (m, 4H), 6.88 (d,  $J$  = 8.6 Hz, 2H), 4.78 (t,  $J$  = 8.0 Hz, 1H), 3.78 (s, 3H), 3.47 (s, 3H), 3.01 (dd,  $J$  = 8.5 Hz, 13.4 Hz, 1H), 2.74 (s, 3H), 2.56 (s, 3H), 2.38 (s, 3H), 2.33 (dd,  $J$  = 7.8 Hz, 13.4 Hz, 1H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 179.7, 158.2, 149.6, 142.3, 141.8, 139.8, 137.4, 136.9, 135.8, 134.8, 132.3, 131.7, 129.4, 129.3, 129.2, 128.9, 128.8, 128.7, 128.6, 128.4, 128.0, 127.9, 127.7, 127.5, 127.0, 122.6, 119.3, 113.9, 61.2, 55.1, 50.3, 46.0, 29.7, 21.3, 18.96 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{35}\text{H}_{35}\text{N}_2\text{O}_4\text{S}$  579.2312; Found 579.2313.

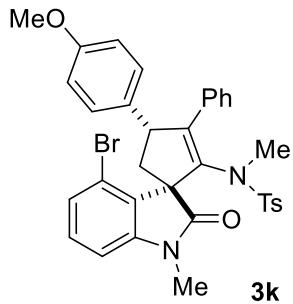
**N-(1'-benzyl-4-(4-methoxyphenyl)-5',7'-dimethyl-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N,4-dimethylbenzenesulfonamide (3j)**



Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer, *dr* = 15:1, 46.8 mg (70%).

Colorless amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.38-7.42 (m, 2H), 7.21-7.28 (m, 6H), 7.12-7.16 (m, 1H), 7.02-7.06 (m, 2H), 6.81-6.93 (m, 9H), 5.16 (d,  $J$  = 17.2 Hz, 1H), 5.03 (d,  $J$  = 17.2 Hz, 1H), 4.75 (t,  $J$  = 8.4 Hz, 1H), 3.75 (s, 3H), 2.96 (dd,  $J$  = 8.4 Hz, 13.4 Hz, 1H), 2.51 (s, 3H), 2.38 (s, 3H), 2.33-2.37 (m, 1H), 2.28 (s, 3H), 2.20 (s, 3H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 180.1, 158.2, 151.1, 146.2, 144.7, 142.6, 139.2, 137.7, 137.6, 136.8, 135.7, 134.6, 132.8, 132.4, 131.9, 129.3, 129.0, 128.5, 128.4, 127.8, 127.5, 126.8, 125.4, 119.0, 113.9, 61.6, 55.1, 50.4, 46.4, 45.1, 37.1, 21.3, 20.8, 18.5 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{42}\text{H}_{41}\text{N}_2\text{O}_4\text{S}$  669.2782; Found 669.2785.

***N*-(4'-bromo-4-(4-methoxyphenyl)-1'-methyl-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-*N*,4-dimethylbenzenesulfonamide (3k)**

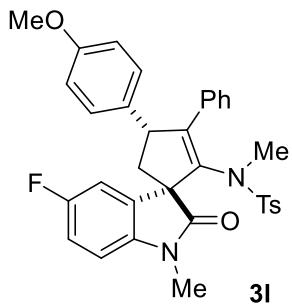


Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), *dr* = 1.3:1, 42.5 mg (66%).

Colorless amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.43-7.47 (m, 3H  $\times$  2.3), 7.17-7.25 (m, 5H  $\times$  2.3), 7.08 (t,  $J$  = 7.4 Hz, 1H), 6.95-7.01 (m, 1H + 2H  $\times$  1.3), 6.77-6.83 (m, 4H  $\times$  2.3), 6.58 (d,  $J$  = 7.3 Hz, 2H), 6.41 (d,  $J$  = 7.6 Hz, 2H  $\times$  1.3), 4.61 (dd,  $J$  = 7.0 Hz, 10.0 Hz, 1H  $\times$  1.3), 4.55 (dd,  $J$  = 4.8 Hz, 10.5 Hz, 1H), 3.73 (s, 3H  $\times$  1.3), 3.72 (s, 3H), 3.28 (s, 3H  $\times$  1.3), 3.22 (s, 3H), 2.97 (m, 1H  $\times$  2.3), 2.72 (s, 3H  $\times$  1.3), 2.60 (s, 3H), 2.50-2.62 (m, 1H  $\times$  2.3), 2.43 (s, 3H  $\times$  1.3), 2.41 (s, 3H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 178.6, 178.2, 158.3, 158.0, 147.0, 146.3, 145.8, 143.23, 143.20, 137.2, 135.6, 135.3, 135.2, 135.1, 134.8, 134.7, 133.2, 130.6, 129.9, 129.7, 129.3, 128.5, 128.4, 128.0, 127.9, 127.8, 127.6, 127.5, 127.2, 127.1, 126.3, 126.1, 119.5, 119.4, 113.7, 113.6, 107.2, 107.1, 63.5, 62.7, 55.1, 55.0, 52.4, 50.9, 42.4, 39.8,

39.0, 26.6, 26.5, 21.5 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for C<sub>34</sub>H<sub>32</sub>BrN<sub>2</sub>O<sub>4</sub>S 643.1261; Found 643.1263.

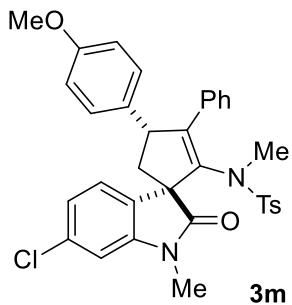
**N-(5'-fluoro-4-(4-methoxyphenyl)-1'-methyl-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N,4-dimethylbenzenesulfonamide (3l)**



Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer, *dr* = 3:1, 41.4 mg (71%).

Colorless amorphous solid. New compound. <sup>1</sup>H NMR (400 MHz, CDCl<sub>3</sub>): δ = 7.31 (d, *J* = 8.6 Hz, 1H), 7.23-7.26 (m, 2H), 7.13-7.17 (m, 4H), 6.95-7.01 (m, 5H), 6.83 (d, *J* = 8.5 Hz, 2H), 6.74-6.76 (m, 1H), 6.63-6.66 (m, 1H), 4.66 (t, *J* = 8.0 Hz, 1H), 3.75 (s, 3H), 3.15 (s, 3H), 2.94 (dd, *J* = 8.4 Hz, 13.4 Hz, 1H), 2.66 (s, 3H), 2.31 (s, 3H), 2.26 (dd, *J* = 7.8 Hz, 13.4 Hz, 1H) ppm; <sup>13</sup>C{<sup>1</sup>H} NMR (100 MHz, CDCl<sub>3</sub>): δ = 178.8, 159.3 (d, *J*<sub>C-F</sub> = 237.2 Hz), 158.3, 149.8, 142.6, 139.9, 136.4 (d, *J*<sub>C-F</sub> = 9.6 Hz), 135.4, 135.0 (d, *J*<sub>C-F</sub> = 24.9 Hz), 134.6, 132.6 (d, *J*<sub>C-F</sub> = 7.9 Hz), 129.2, 129.0, 128.3, 127.9, 127.6, 127.0, 114.7 (d, *J*<sub>C-F</sub> = 23.4 Hz), 114.0, 108.1 (d, *J*<sub>C-F</sub> = 8.0 Hz), 62.1, 55.1, 50.3, 45.7, 37.2, 26.5, 21.3 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for C<sub>34</sub>H<sub>32</sub>FN<sub>2</sub>O<sub>4</sub>S 583.2061; Found 583.2058.

**N-(6'-chloro-4-(4-methoxyphenyl)-1'-methyl-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N,4-dimethylbenzenesulfonamide (3m)**

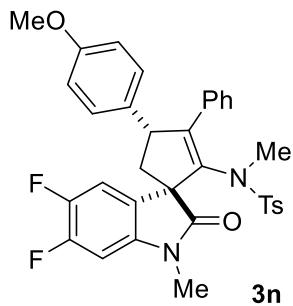


Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), *dr* = 2:1, 40.7 mg (68%).

Colorless amorphous solid. New compound. <sup>1</sup>H NMR (400 MHz, CDCl<sub>3</sub>): δ = 7.51 (d, *J* = 7.9 Hz, 1H), 7.44 (d, *J* = 7.9 Hz, 2H), 7.29-7.32 (m, 3H), 7.22-7.26 (m, 6H), 7.12-7.17 (m, 10H), 7.06-7.09 (m, 5H), 6.99-7.01 (m, 10H), 6.89 (d, *J* = 8.0 Hz, 2H), 6.8 (d, *J* = 8.6 Hz, 4H),

6.74-6.78 (m, 3H), 6.61 (br, 2H), 4.63-4.70 (m, 1H × 3), 3.73 (s, 3H × 2), 3.69 (s, 3H), 3.15 (s, 3H), 3.10 (s, 3H × 2), 2.92 (dd,  $J$  = 8.5 Hz, 13.4 Hz, 1H × 2), 2.80 (s, 3H), 2.70 (s, 3H × 2), 2.68-2.73 (m, 1H), 2.59 (dd,  $J$  = 7.1 Hz, 13.5 Hz, 1H), 2.34 (s, 3H × 2), 2.33 (s, 3H), 2.22 (dd,  $J$  = 7.5 Hz, 13.4 Hz, 1H × 2) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 178.9, 178.8, 158.2, 158.1, 150.1, 149.6, 145.0, 144.6, 142.8, 142.6, 136.5, 135.9, 135.4, 135.1, 134.8, 134.6, 134.39, 134.37, 134.3, 130.0, 129.3, 129.2, 129.1, 129.0, 128.9, 128.22, 128.19, 128.1, 127.9, 127.8, 127.6, 127.3, 126.5, 126.2, 122.44, 122.39, 113.9, 113.7, 108.7, 108.4, 61.5, 61.4, 55.03, 54.96, 50.5, 50.1, 45.4, 44.5, 37.5, 26.5, 26.4, 21.3 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{34}\text{H}_{32}\text{ClN}_2\text{O}_4\text{S}$  599.1766; Found 599.1771.

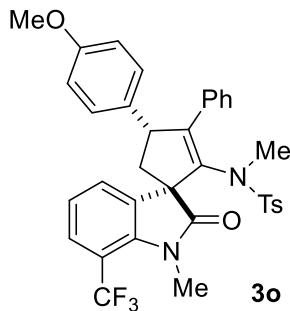
**N-(5',6'-difluoro-4-(4-methoxyphenyl)-1'-methyl-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N,4-dimethylbenzenesulfonamide (3n)**



Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1),  $dr$  = 1.5:1, 43.2 mg (72%).

Colorless amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.38 (dd,  $J$  = 2.5 Hz, 8.0 Hz, 1H), 7.33 (d,  $J$  = 8.6 Hz, 2H × 1.5), 7.24-7.28 (m, 3H), 7.14-7.21 (m, 4H × 2.5), 7.01-7.07 (m, 1H + 4H × 1.5), 6.97-6.99 (m, 2H × 1.5 + 1H), 6.92 (d,  $J$  = 8.0 Hz, 2H), 6.85 (d,  $J$  = 8.6 Hz, 2H × 1.5), 6.76-6.78 (m, 3H), 6.66 (dd,  $J$  = 4.0 Hz, 8.6 Hz, 1H × 1.5), 4.64-4.69 (m, 1H × 2.5), 3.77 (s, 3H × 1.5), 3.72 (s, 3H), 3.22 (s, 3H), 3.18 (s, 3H × 1.5), 2.95 (dd,  $J$  = 8.3 Hz, 13.4 Hz, 1H × 1.5), 2.78 (s, 3H), 2.72-2.76 (m, 1H), 2.67 (s, 3H × 1.5), 2.63 (dd,  $J$  = 7.1 Hz, 13.4 Hz, 1H), 2.33 (s, 3H × 2.5), 2.28 (dd,  $J$  = 7.7 Hz, 13.4 Hz, 1H × 1.5) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 178.8, 178.6, 159.4 (d,  $J_{\text{C}-\text{F}} = 221.8$  Hz), 159.3 (d,  $J_{\text{C}-\text{F}} = 237.7$  Hz), 150.2, 149.8, 142.7 (d,  $J_{\text{C}-\text{F}} = 21.2$  Hz), 139.9, 139.4, 135.7, 135.4, 135.2, 134.9, 134.6 (d,  $J_{\text{C}-\text{F}} = 8.0$  Hz), 133.3 (d,  $J_{\text{C}-\text{F}} = 7.7$  Hz), 132.6 (d,  $J_{\text{C}-\text{F}} = 8.0$  Hz), 129.4, 129.0, 128.3, 128.2 (d,  $J_{\text{C}-\text{F}} = 18.2$  Hz), 127.8 (d,  $J_{\text{C}-\text{F}} = 27.2$  Hz), 127.1, 114.8 (d,  $J_{\text{C}-\text{F}} = 23.5$  Hz), 114.0, 113.7, 113.5 (d,  $J_{\text{C}-\text{F}} = 32.5$  Hz), 108.2 (dd,  $J_{\text{C}-\text{F}} = 8.1$  Hz, 10.0 Hz), 62.1, 55.1, 55.0, 50.6, 50.4, 45.7, 44.5, 37.7, 26.7, 26.5, 21.3 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{34}\text{H}_{31}\text{F}_2\text{N}_2\text{O}_4\text{S}$  601.1967; Found 601.1969.

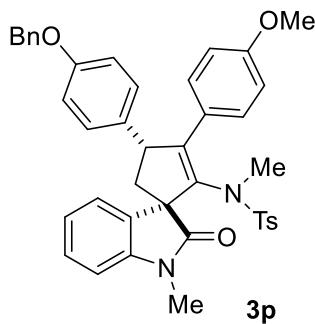
**N-(4-(4-methoxyphenyl)-1'-methyl-2'-oxo-3-phenyl-7'-(trifluoromethyl)spiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N,4-dimethylbenzenesulfonamide (3o)**



Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer, *dr* = 2.3:1, 41.1 mg (65%).

Colorless amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.78-7.82 (m, 1H  $\times$  3.3), 7.65 (d,  $J$  = 8.0 Hz, 1H), 7.59 (d,  $J$  = 8.0 Hz, 1H  $\times$  2.3), 7.32 (d,  $J$  = 8.6 Hz, 2H  $\times$  2.3), 7.23-7.27 (m, 4H + 2H  $\times$  2.3), 7.18-7.21 (m, 3H  $\times$  3.3), 7.11-7.14 (m, 1H  $\times$  3.3), 6.92-6.98 (m, 3H  $\times$  3.3), 6.83 (d,  $J$  = 8.6 Hz, 3H + 2H  $\times$  2.3), 6.77 (d,  $J$  = 8.6 Hz, 1H  $\times$  2.3), 4.67-4.73 (m, 1H  $\times$  3.3), 3.75 (s, 3H  $\times$  2.3), 3.71 (s, 3H), 3.38-3.39 (m, 3H), 3.35-3.36 (m, 3H  $\times$  2.3), 3.20-3.23 (m, 1H), 2.94 (dd,  $J$  = 8.4 Hz, 13.4 Hz, 1H  $\times$  2.3), 2.82 (s, 3H), 2.64 (s, 3H  $\times$  2.3), 2.36-2.39 (m, 1H), 2.31 (s, 3H  $\times$  3.3), 2.23-2.27 (m, 1H  $\times$  2.3) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 179.9, 179.7, 158.4, 158.3, 143.0, 142.9, 135.9, 135.4, 134.9, 134.6, 134.5, 133.6, 129.3, 129.1, 128.4, 128.3 (q,  $J_{\text{C}-\text{F}}$  = 214.7 Hz), 128.2, 128.1, 127.9 (q,  $J_{\text{C}-\text{F}}$  = 243.9 Hz), 126.5 (q,  $J_{\text{C}-\text{F}}$  = 6.0 Hz), 122.1 (q,  $J_{\text{C}-\text{F}}$  = 3.8 Hz), 114.1, 113.8, 112.3 (q,  $J_{\text{C}-\text{F}}$  = 32.9 Hz), 112.1 (q,  $J_{\text{C}-\text{F}}$  = 32.3 Hz), 60.5, 60.3, 55.2, 55.1, 50.7, 50.3, 46.0, 45.1, 37.6, 29.1, 29.0, 21.3 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for  $\text{C}_{35}\text{H}_{32}\text{F}_3\text{N}_2\text{O}_4\text{S}$  633.2029; Found 633.2027.

***N*-(4-(4-(benzyloxy)phenyl)-3-(4-methoxyphenyl)-1'-methyl-2'-oxospiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-*N*,4-dimethylbenzenesulfonamide (3p)**

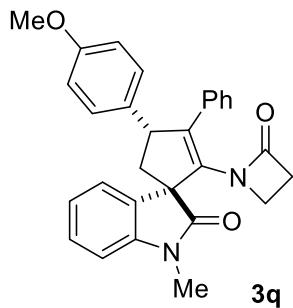


Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer, *dr* > 20:1, 55.7 mg (83%).

Colorless amorphous solid. New compound.  $^1\text{H}$  NMR (400 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 7.25-7.61 (m, 1H), 7.21-7.43 (m, 10H), 7.12-7.18 (m, 2H), 6.90-6.98 (m, 5H), 6.67-6.74 (m, 3H), 5.01 (s, 2H), 4.66 (t,  $J$  = 7.8 Hz, 1H), 3.74 (s, 3H), 3.14 (s, 3H), 2.93 (dd,  $J$  = 8.5 Hz, 13.4 Hz, 1H), 2.67 (s, 3H), 2.31 (s, 3H), 2.26 (dd,  $J$  = 7.6 Hz, 13.4 Hz, 1H) ppm;  $^{13}\text{C}\{\text{H}\}$  NMR (100 MHz,  $\text{CDCl}_3$ ):  $\delta$  = 179.3, 159.2, 157.5, 149.1, 144.0, 142.4, 137.0, 136.9, 136.5, 136.2, 131.2, 129.7, 129.3, 128.9,

128.5, 127.9, 127.5, 127.2, 127.0, 122.8, 115.0, 113.4, 107.8, 70.0, 61.9, 55.0, 50.2, 45.5, 37.0, 26.4, 21.4 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for C<sub>41</sub>H<sub>39</sub>N<sub>2</sub>O<sub>5</sub>S 671.2574; Found 671.2577.

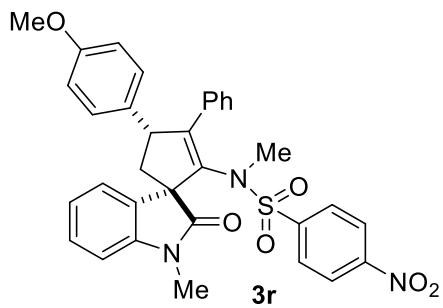
**4-(4-methoxyphenyl)-1'-methyl-2-(2-oxoazetidin-1-yl)-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2'-one (3q)**



Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer, *dr* = 8:1, 36.0 mg (80%).

Colorless amorphous solid. New compound. <sup>1</sup>H NMR (400 MHz, CDCl<sub>3</sub>): δ = 7.30 (d, *J* = 7.5 Hz, 2H), 7.27-7.28 (m, 1H), 7.26 (d, *J* = 2.9 Hz, 1H), 7.14-7.21 (m, 5H), 7.05-7.09 (m, 1H), 6.88 (d, *J* = 7.5 Hz, 1H), 6.75 (d, *J* = 8.7 Hz, 2H), 4.47 (t, *J* = 8.2 Hz, 1H), 3.73 (s, 3H), 3.28 (s, 3H), 2.96-3.01 (m, 1H), 2.49-2.69 (m, 5H) ppm; <sup>13</sup>C{<sup>1</sup>H} NMR (100 MHz, CDCl<sub>3</sub>): δ = 178.9, 163.8, 158.2, 144.3, 134.99, 134.9, 132.7, 132.6, 132.5, 129.5, 129.4, 129.0, 128.5, 127.8, 127.7, 127.4, 122.2, 121.8, 113.9, 113.7, 108.4, 60.1, 55.1, 51.5, 46.3, 41.1, 36.2, 29.7, 26.7 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for C<sub>29</sub>H<sub>27</sub>N<sub>2</sub>O<sub>3</sub> 451.2016; Found 451.2020.

**N-(4-(4-methoxyphenyl)-1'-methyl-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N-methyl-4-nitrobenzenesulfonamide (3r)**

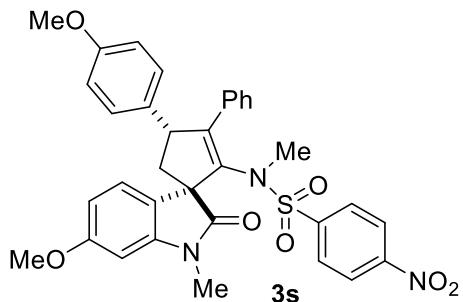


Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer, *dr* > 40:1, 52.5 mg (88%).

Light yellow amorphous solid. New compound. <sup>1</sup>H NMR (400 MHz, CDCl<sub>3</sub>): δ = 7.91 (d, *J* = 8.9 Hz, 2H), 7.56 (br, 1H), 7.34 (td, *J* = 7.7 Hz, 1.0 Hz, 2H), 7.14-7.19 (m, 9H), 6.75-6.81 (m, 3H), 4.69-4.71 (m, 1H), 3.74 (s, 3H), 3.17 (s, 3H), 2.93 (dd, *J* = 8.3 Hz, 13.4 Hz, 1H), 2.79 (s, 3H), 2.29 (dd, *J* = 8.0 Hz, 13.4 Hz, 1H) ppm; <sup>13</sup>C{<sup>1</sup>H} NMR (100 MHz, CDCl<sub>3</sub>): δ = 178.8, 158.3, 151.1, 149.3, 145.3, 144.2, 136.3, 135.3, 134.4, 130.5, 129.2, 129.0, 128.1, 123.5, 122.8,

114.1, 108.1, 61.6, 55.1, 50.2, 45.7, 26.5 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for C<sub>33</sub>H<sub>30</sub>N<sub>3</sub>O<sub>6</sub>S 596.1850; Found 596.1855.

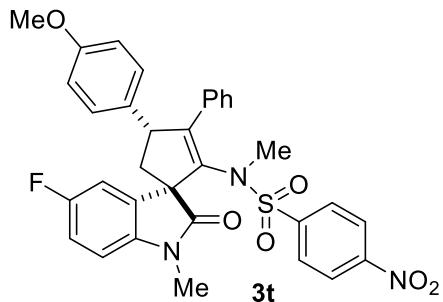
**N-(6'-methoxy-4-(4-methoxyphenyl)-1'-methyl-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N-methyl-4-nitrobenzenesulfonamide (3s)**



Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer, *dr* > 40:1, 56.3 mg (90%).

Light yellow amorphous solid. New compound. <sup>1</sup>H NMR (400 MHz, CDCl<sub>3</sub>): δ = 7.93 (d, *J* = 8.9 Hz, 2H), 7.15-7.22 (m, 10H), 6.80-6.83 (m, 3H), 6.61-6.63 (m, 1H), 4.69 (t, *J* = 7.8 Hz, 1H), 3.81 (s, 3H), 3.73 (s, 3H), 3.14 (s, 3H), 2.94 (dd, *J* = 8.3 Hz, 13.3 Hz, 1H), 2.81 (s, 3H), 2.26 (dd, *J* = 7.7 Hz, 13.3 Hz, 1H) ppm; <sup>13</sup>C{<sup>1</sup>H} NMR (100 MHz, CDCl<sub>3</sub>): δ = 177.8, 157.7, 155.5, 150.1, 148.6, 144.8, 136.8, 135.5, 134.7, 133.7, 131.0, 128.6, 127.5, 127.4, 122.8, 113.4, 112.9, 112.5, 107.7, 61.4, 55.3, 54.5, 49.6, 45.3, 29.0, 25.9 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for C<sub>34</sub>H<sub>32</sub>N<sub>3</sub>O<sub>7</sub>S 626.1955; Found 626.1959.

**N-(5'-fluoro-4-(4-methoxyphenyl)-1'-methyl-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N-methyl-4-nitrobenzenesulfonamide (3t)**

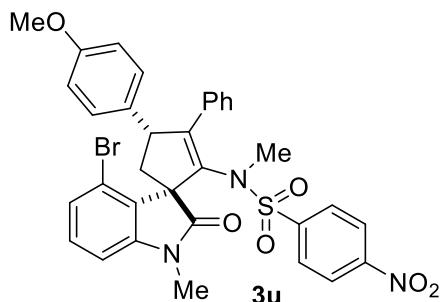


Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer, *dr* > 40:1, 51.5 mg (84%).

Light yellow amorphous solid. New compound. <sup>1</sup>H NMR (400 MHz, CDCl<sub>3</sub>): δ = 7.95 (d, *J* = 8.8 Hz, 2H), 7.31-7.35 (m, 3H), 7.09-7.17 (m, 6H), 7.02 (td, *J* = 2.4 Hz, 8.8 Hz, 2H), 6.82 (d, *J* = 7.9 Hz, 2H), 6.67-6.72 (m, 1H), 4.69 (t, *J* = 7.6 Hz, 1H), 3.75 (s, 3H), 3.19 (s, 3H), 2.94 (dd, *J* = 8.2 Hz, 13.5 Hz, 1H), 2.86 (s, 3H), 2.27 (dd, *J* = 8.1 Hz, 13.5 Hz, 1H) ppm; <sup>13</sup>C{<sup>1</sup>H} NMR (100 MHz, CDCl<sub>3</sub>): δ = 178.6, 159.4 (d, *J*<sub>C-F</sub> = 239.9 Hz), 158.6, 151.3, 149.5, 145.3, 140.2, 135.9,

135.1, 134.3, 132.1 (d,  $J_{C-F} = 8$  Hz), 129.3, 128.2, 123.7, 115.4 (d,  $J_{C-F} = 23.4$  Hz), 114.2, 108.6 (d,  $J_{C-F} = 8$  Hz), 62.1 (d,  $J_{C-F} = 1.5$  Hz), 55.3, 53.6, 50.3, 46.1, 29.8, 26.8 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for C<sub>33</sub>H<sub>29</sub>FN<sub>3</sub>O<sub>6</sub>S 614.1756; Found 614.1759.

**N-(4'-bromo-4-(4-methoxyphenyl)-1'-methyl-2'-oxo-3-phenylspiro[cyclopentane-1,3'-indolin]-2-en-2-yl)-N-methyl-4-nitrobenzenesulfonamide (3u)**



Synthesized by the general procedure, flash chromatography (petroleum ether:ethyl acetate = 10:1 to 5:1), major diastereomer,  $dr > 40:1$ , 54.5 mg (81%).

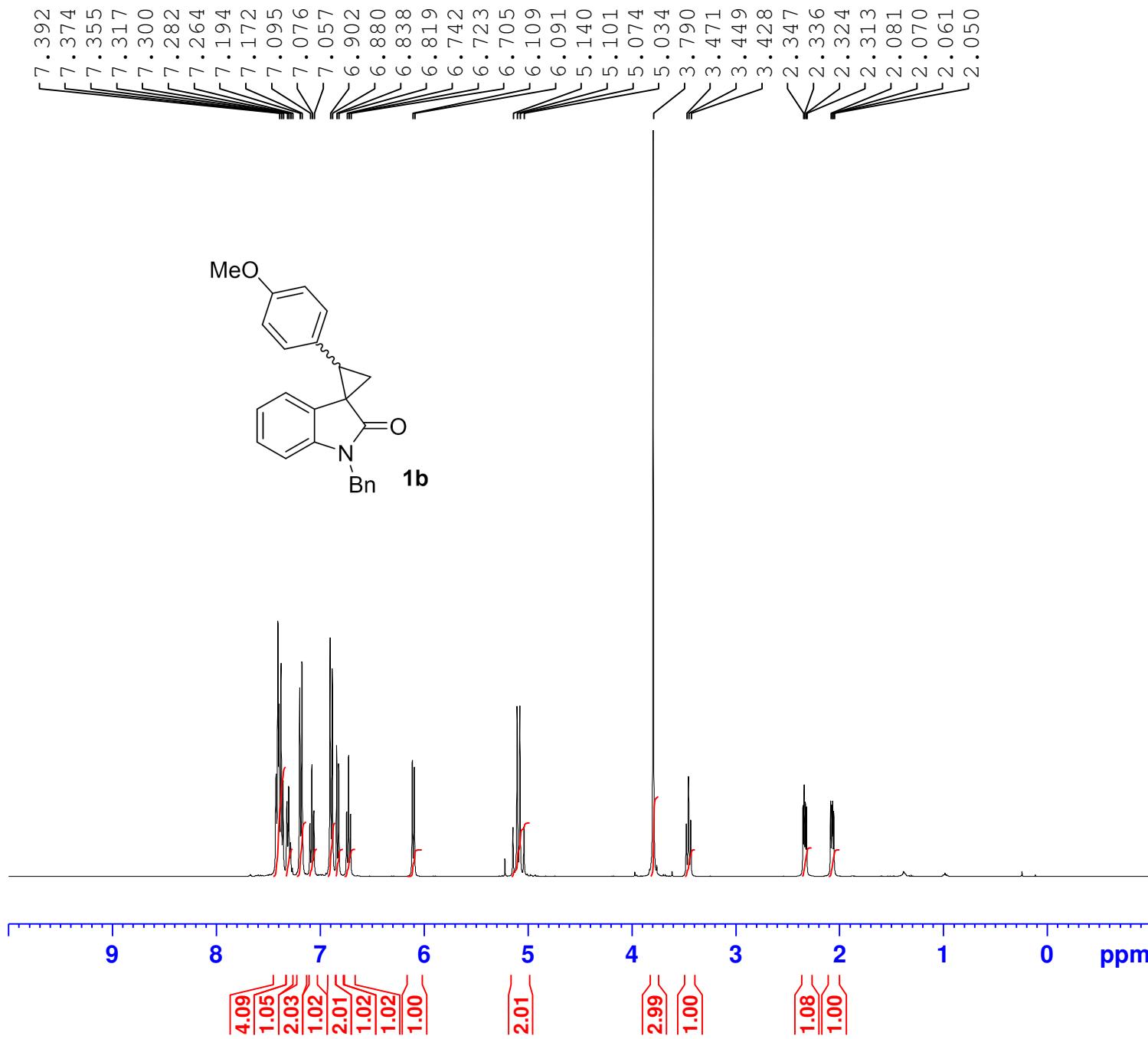
Light yellow amorphous solid. New compound. <sup>1</sup>H NMR (400 MHz, CDCl<sub>3</sub>):  $\delta$  = 8.07 (d,  $J$  = 8.8 Hz, 2H), 7.70 (d,  $J$  = 8.8 Hz, 2H), 7.24-7.26 (m, 2H), 7.12 (d,  $J$  = 8.6 Hz, 2H), 6.98 (t,  $J$  = 7.5 Hz, 1H), 6.84-6.87 (m, 1H), 6.80 (t,  $J$  = 7.7 Hz, 2H), 6.7 (d,  $J$  = 8.6 Hz, 2H), 6.52 (d,  $J$  = 7.5 Hz, 2H), 4.69 (dd,  $J$  = 7.3 Hz, 9.9 Hz, 1H), 3.71 (s, 3H), 3.29 (s, 3H), 3.05 (dd,  $J$  = 10.0 Hz, 12.8 Hz, 1H), 3.01 (s, 3H), 2.62 (dd,  $J$  = 7.3 Hz, 12.8 Hz, 1H) ppm; <sup>13</sup>C{<sup>1</sup>H} NMR (100 MHz, CDCl<sub>3</sub>):  $\delta$  = 178.1, 158.3, 149.9, 149.6, 147.0, 143.5, 135.8, 134.5, 134.4, 130.4, 129.7, 129.0, 127.96, 127.6, 127.5, 126.5, 123.7, 119.7, 113.8, 107.6, 62.9, 55.1, 50.8, 41.97, 40.7, 26.6 ppm. HRMS (ESI) m/z: [M + H]<sup>+</sup> Calcd. for C<sub>33</sub>H<sub>29</sub>BrN<sub>3</sub>O<sub>6</sub>S 674.0955; Found 674.0958.

#### 4. A gram-scale synthesis of spirocyclopenteneoxindole 3a

To a mixture of spirocyclopropyl oxindole **1a** (1.50 g, 5.40 mmol, 1.0 equiv, diastereomeric mixture) and ynamide **2a** (1.84 g, 6.44 mmol, 1.2 equiv) in CH<sub>2</sub>Cl<sub>2</sub> (20 mL), Cu(OTf)<sub>2</sub> (115.5 mg, 0.32 mmol, 0.05 equiv) was added. The mixture was stirred at room temperature (25 °C) in air for 8 h. Then, the mixture was diluted with CH<sub>2</sub>Cl<sub>2</sub> (25.0 mL) and washed with saturated NaHCO<sub>3</sub>. The aqueous phase was extracted with CH<sub>2</sub>Cl<sub>2</sub> (25 mL × 3). The organic layers were combined, washed with brine, dried over Na<sub>2</sub>SO<sub>4</sub>, and concentrated. The residue was purified by silica gel flash chromatography (petroleum ether:ethyl acetate = 8:1 to 4:1) to provide product **3a** (2.41 g, 79%,  $dr$  = 9:1).

#### 6. References

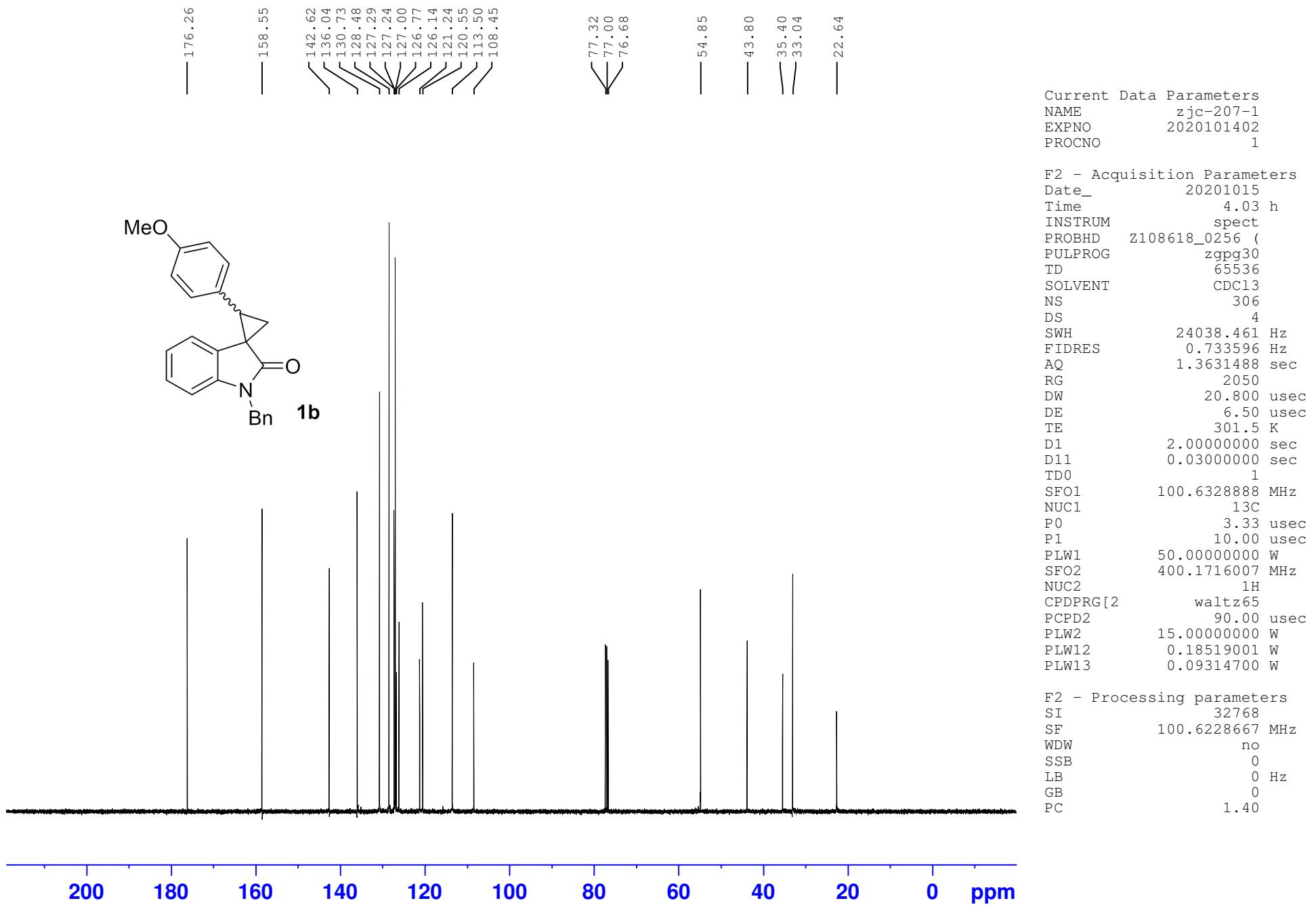
1. A. A. Akaev, E. V. Villemson, N. S. Vorobyeva, A. G. Majouga, E. M. Budynina and M. Y. Melnikov, *J. Org. Chem.*, 2017, **82**, 5689-5701.
2. D.-X. Zhang, J.-J. Man, Y. Chen, L. Yin, J.-C. Zhong and Q.-F. Zhang, *RSC Adv.*, 2019, **9**, 12567-12571.

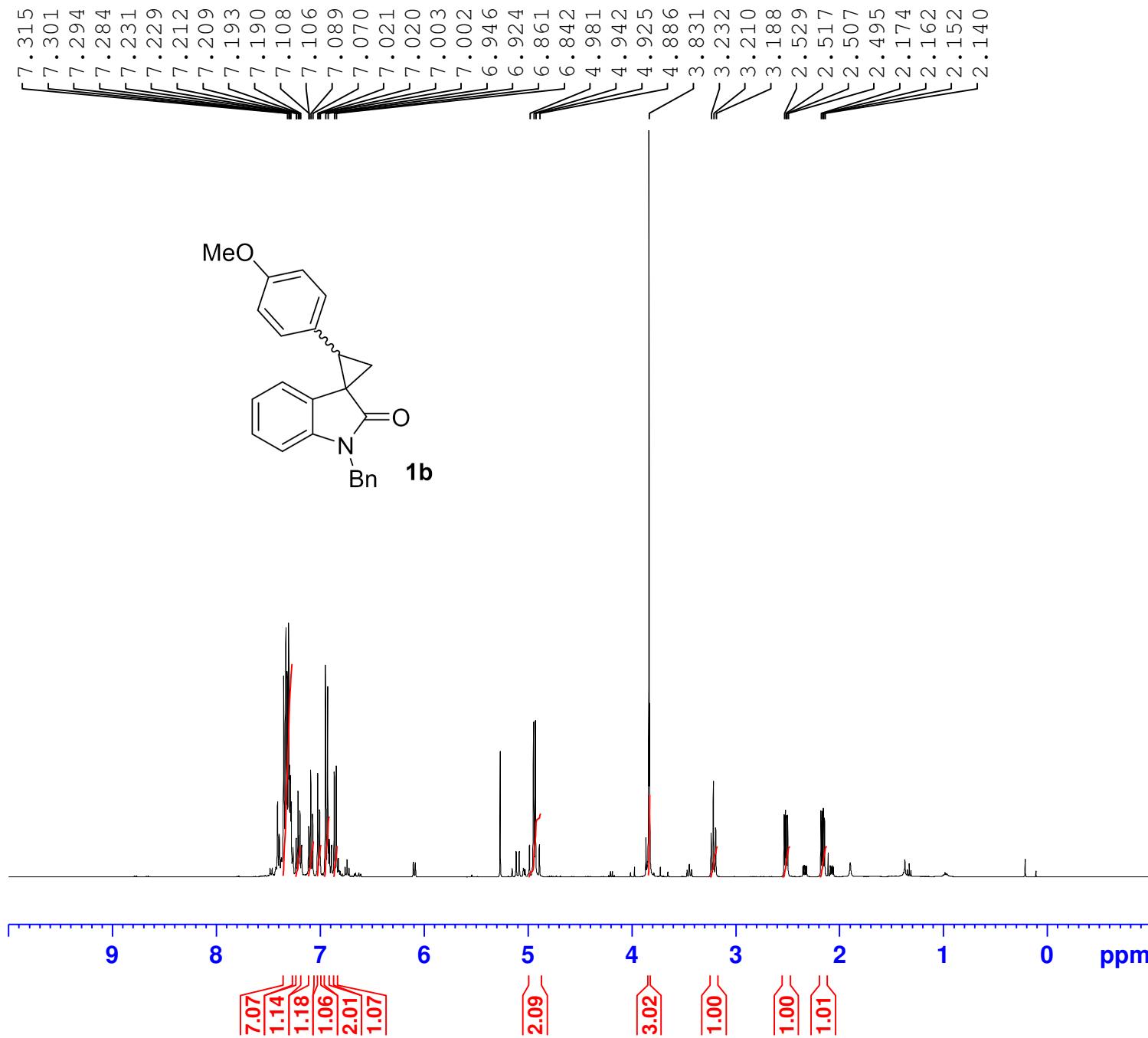


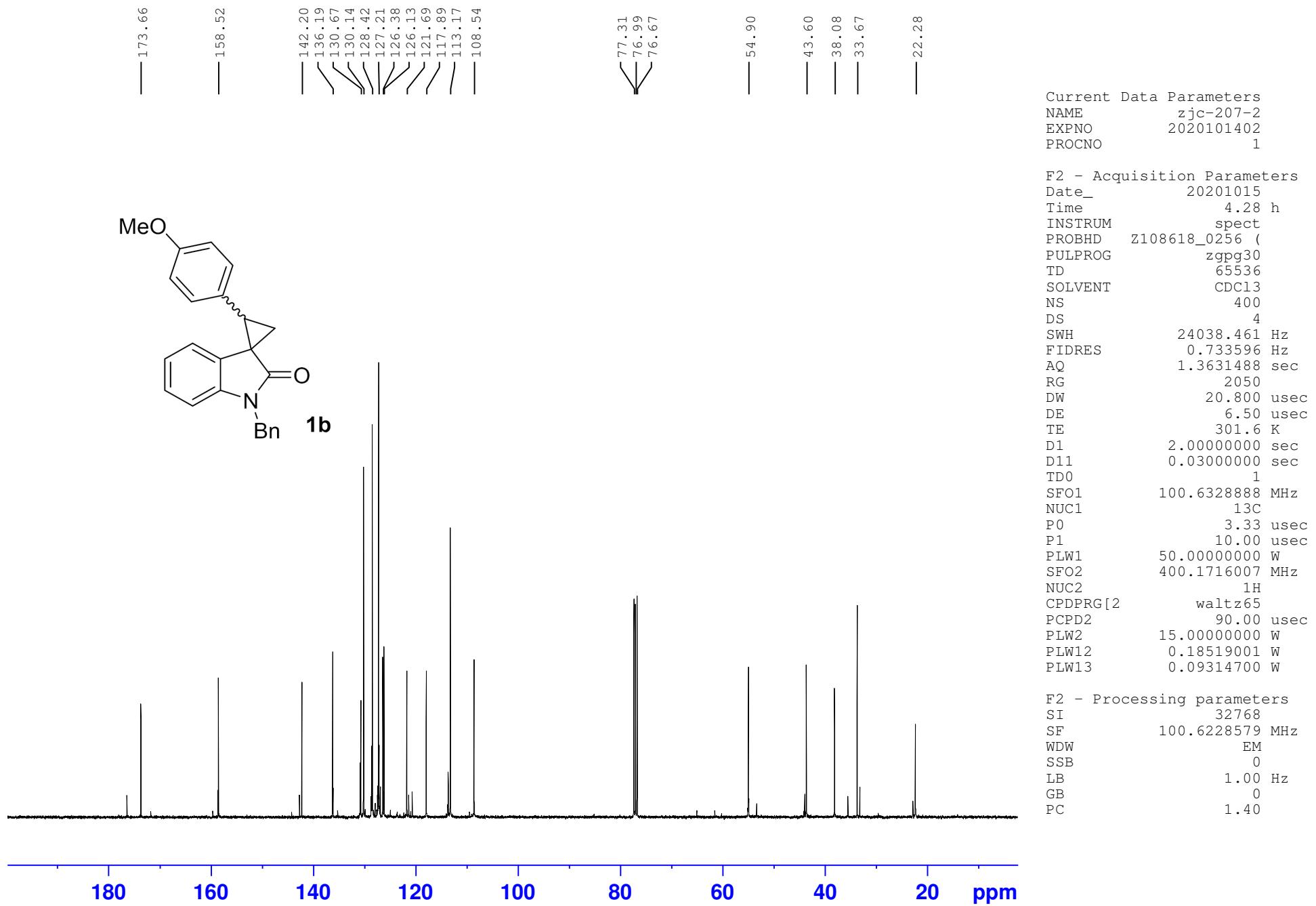
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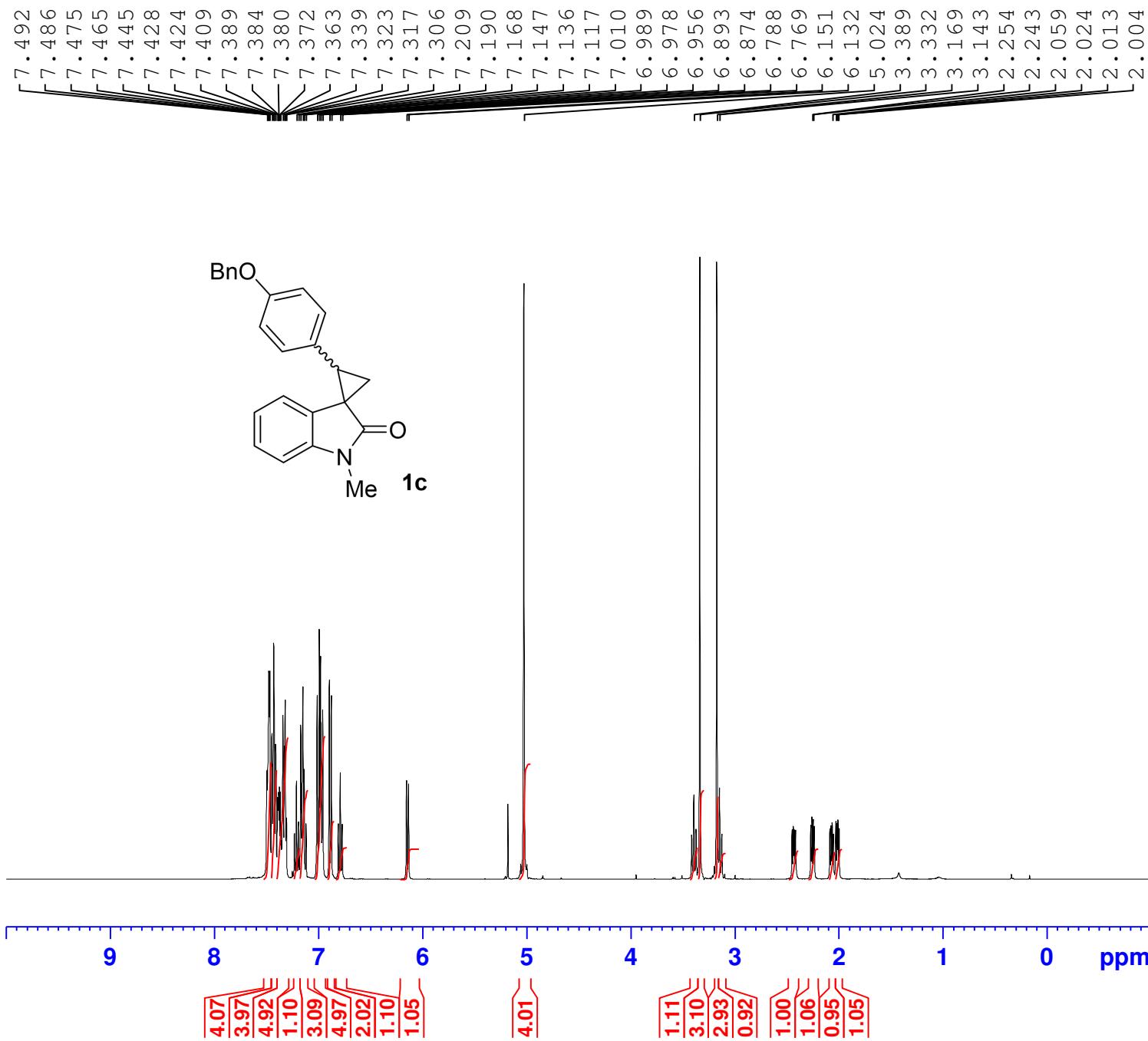
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SWH 0.250967 Hz  
FIDRES 3.9845889 sec  
AQ 25.4  
RG 60.800 usec  
DE 6.50 usec  
TE 301.1 K  
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P0 3.33 usec  
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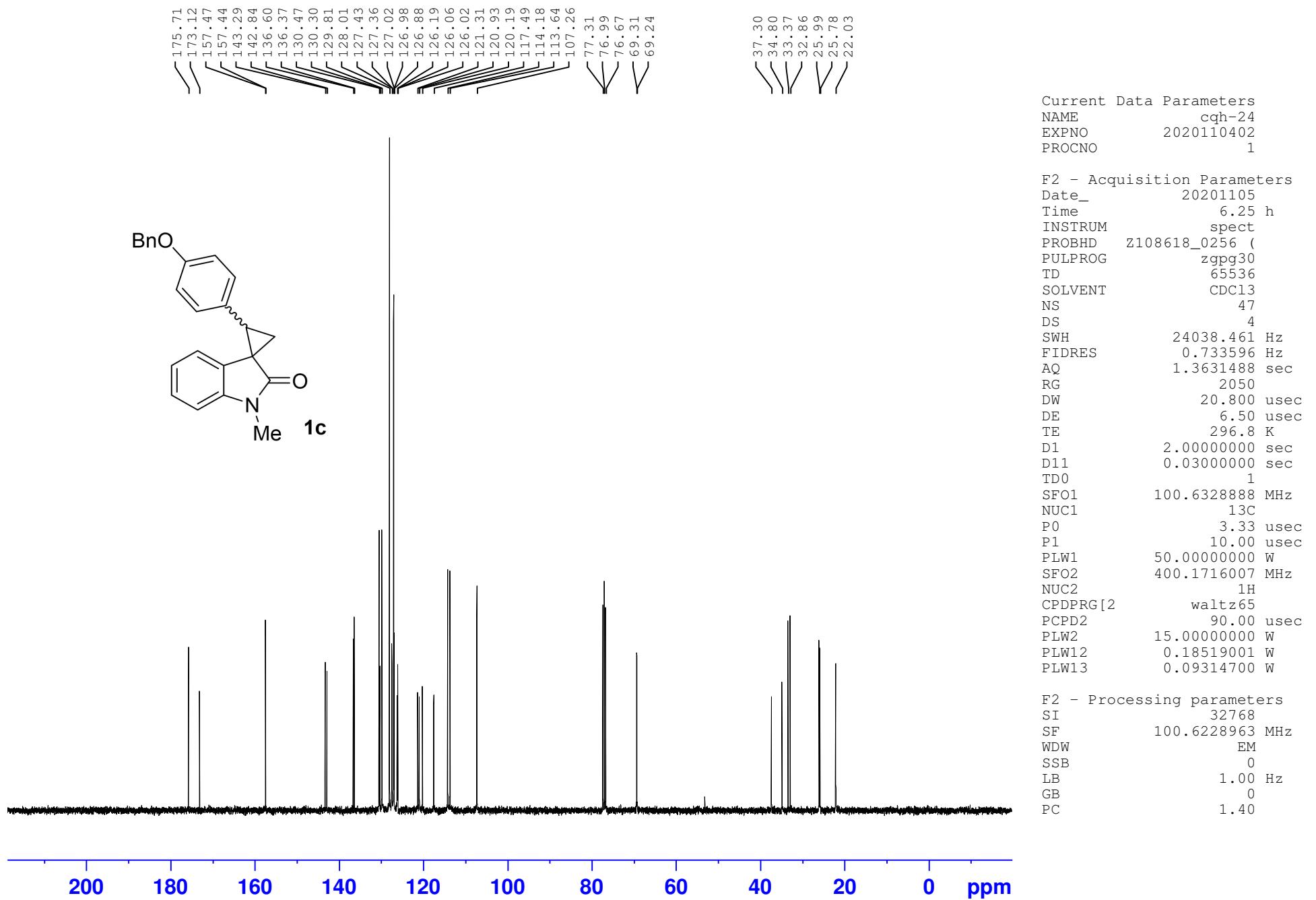


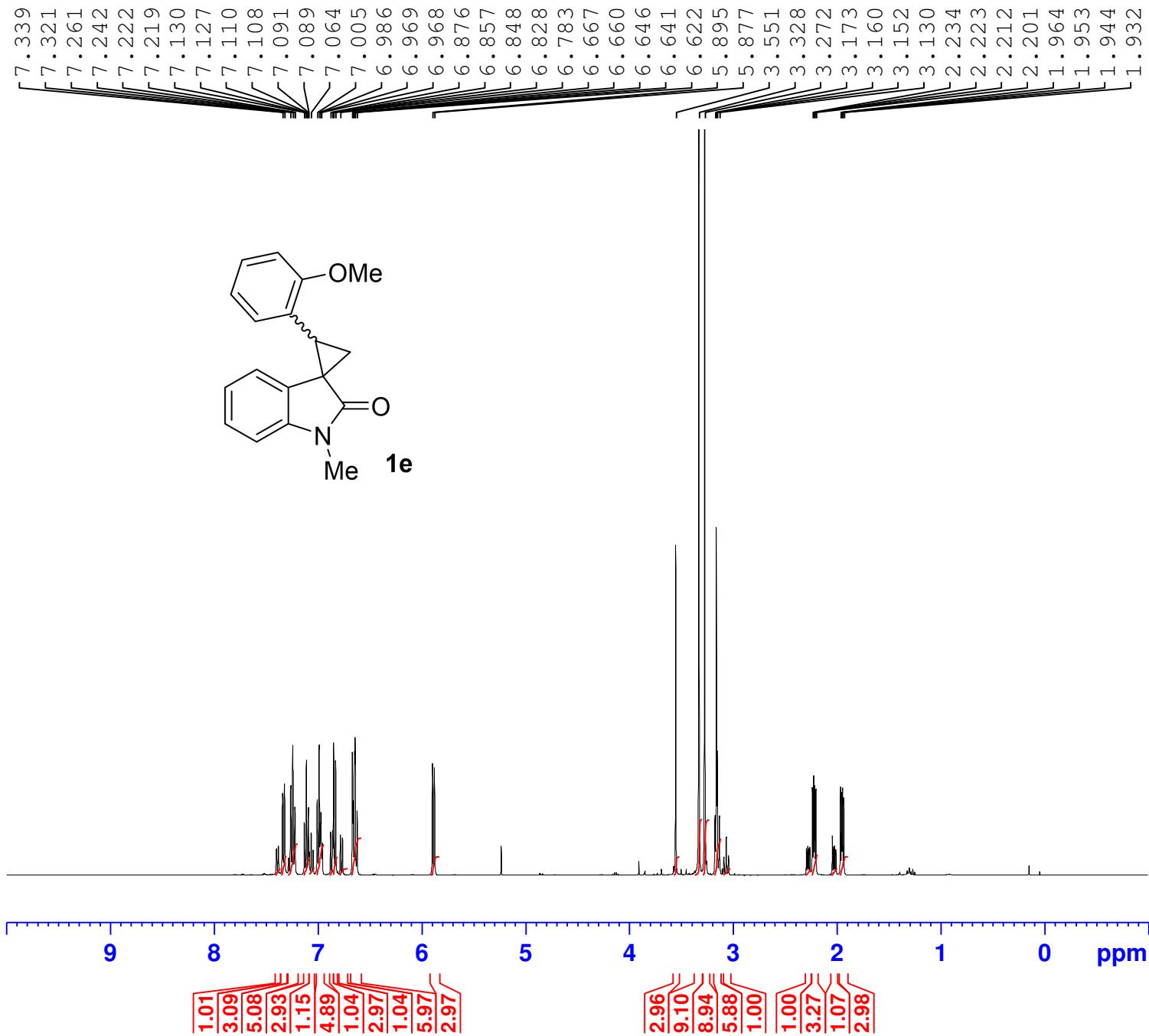


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P0 3.33 usec  
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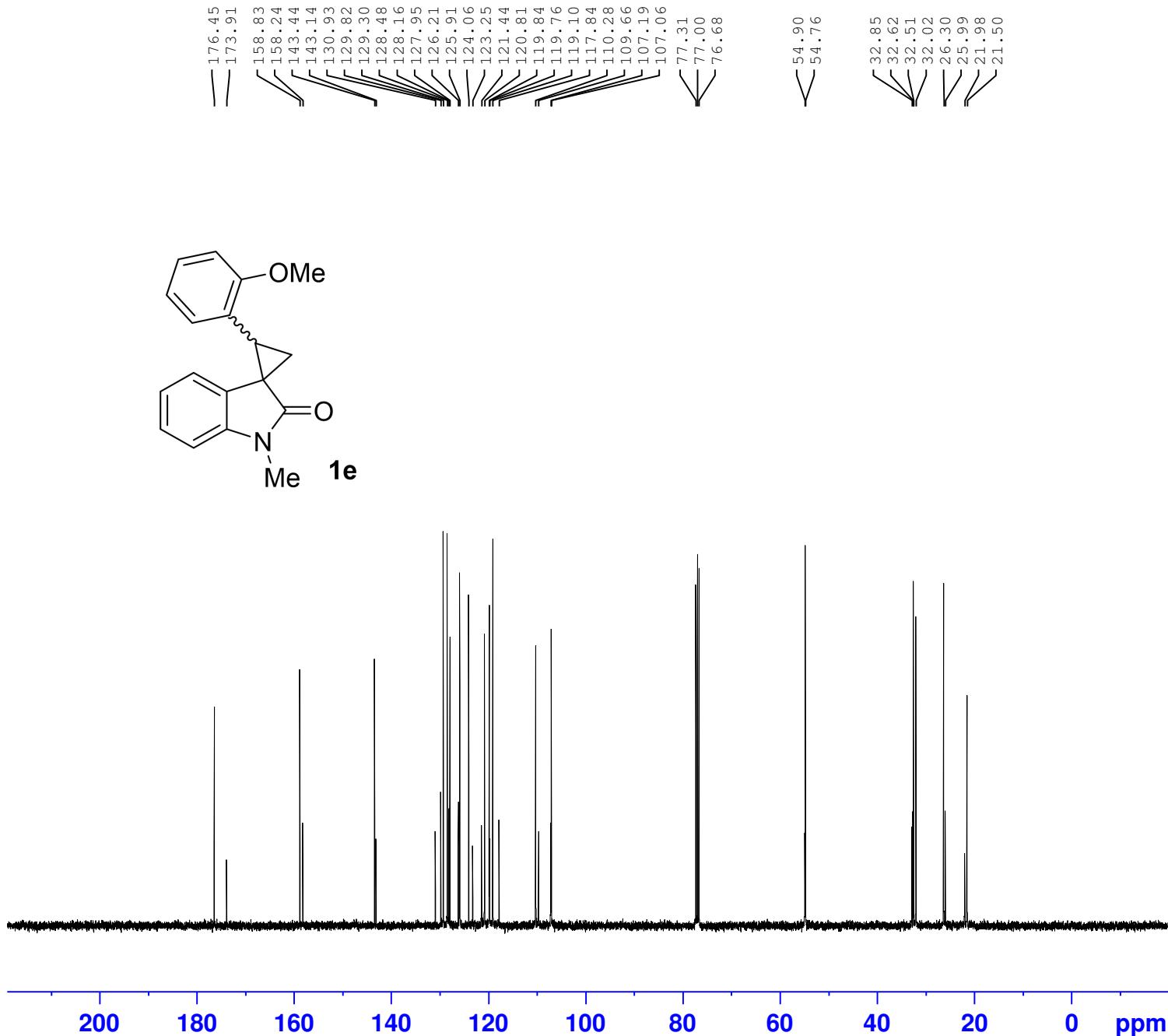


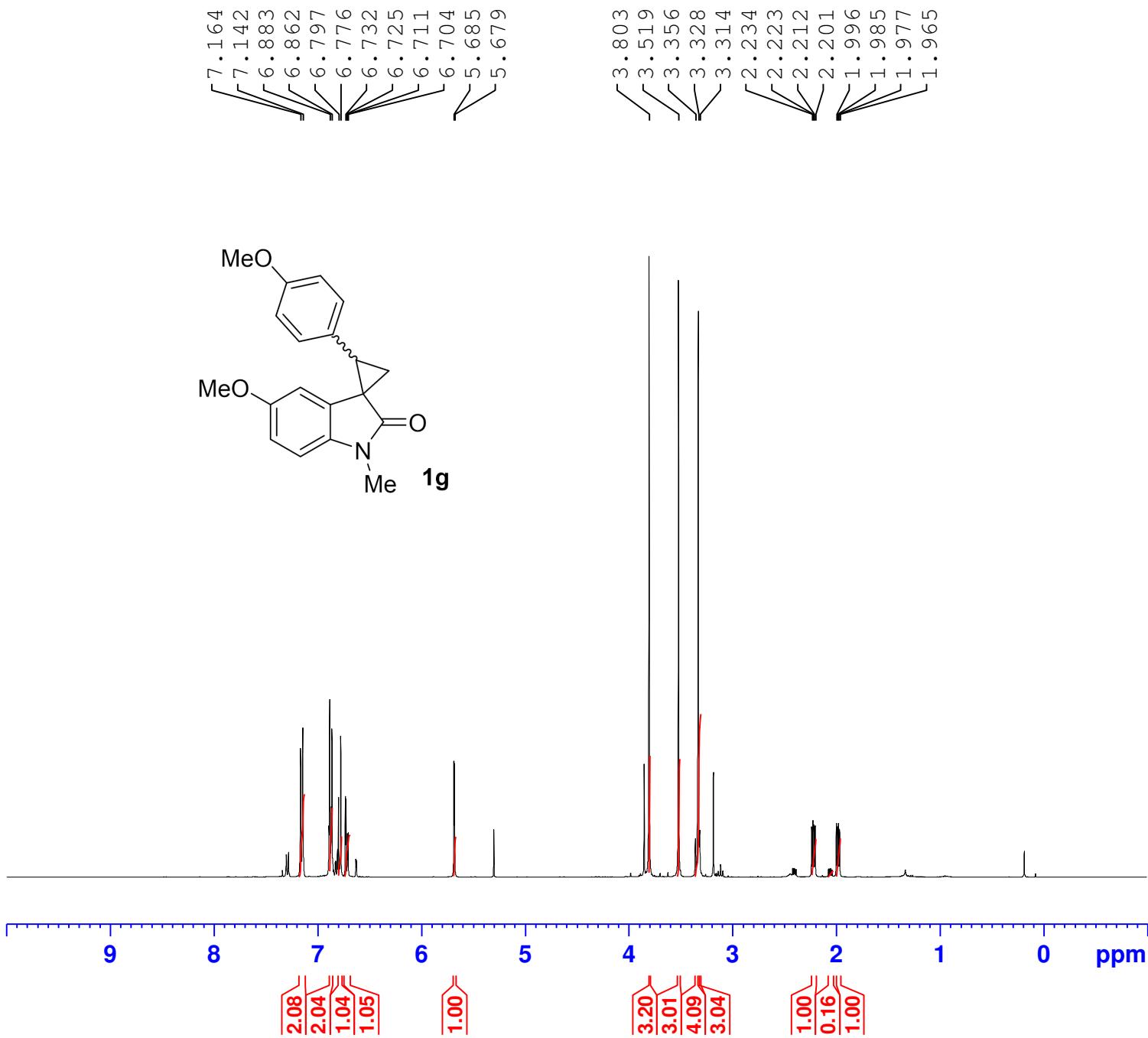


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 SWH 8223.685 Hz  
 FIDRES 0.250967 Hz  
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 DE 6.50 usec  
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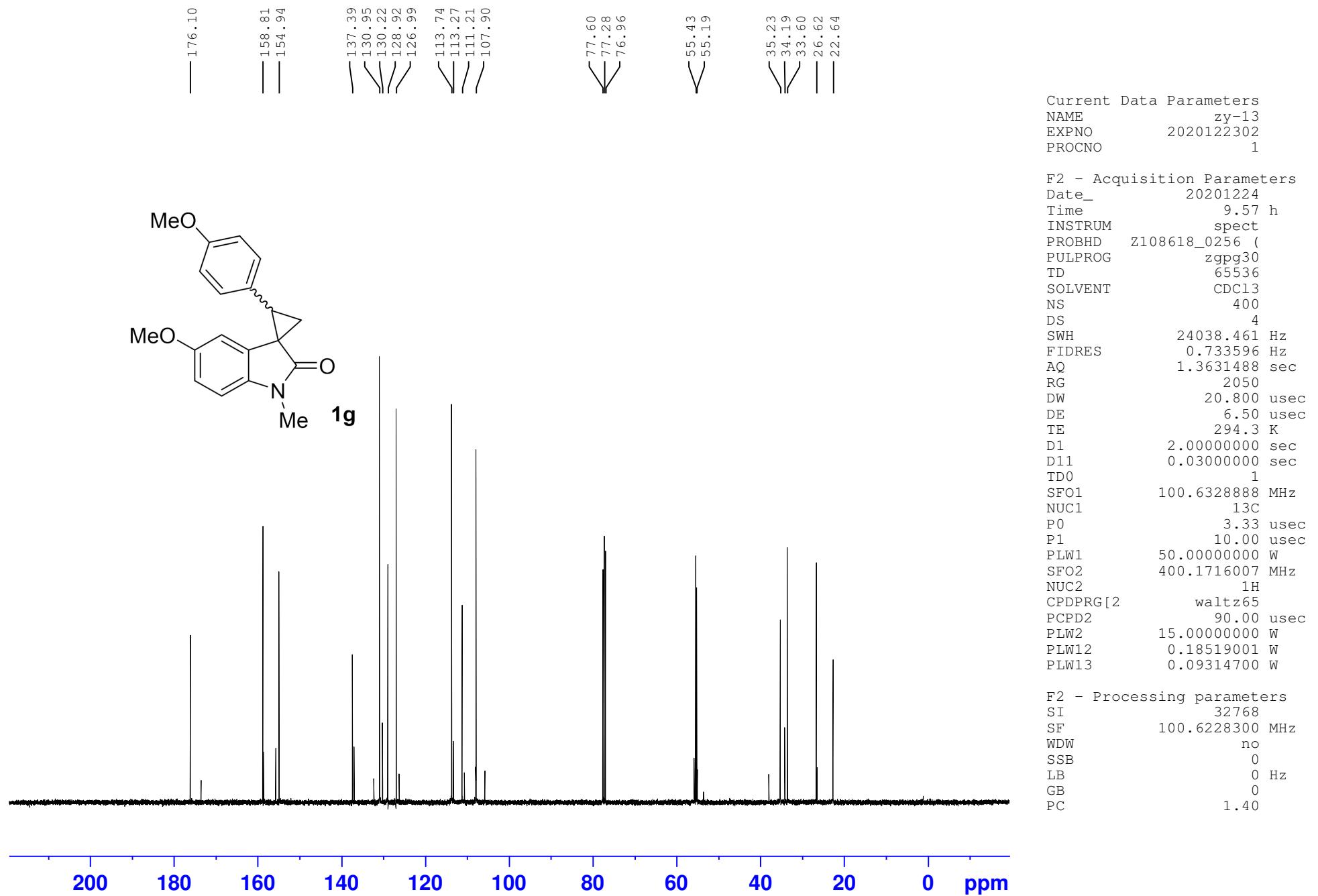


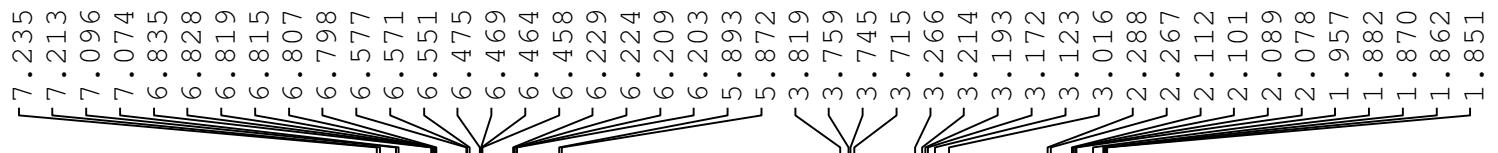


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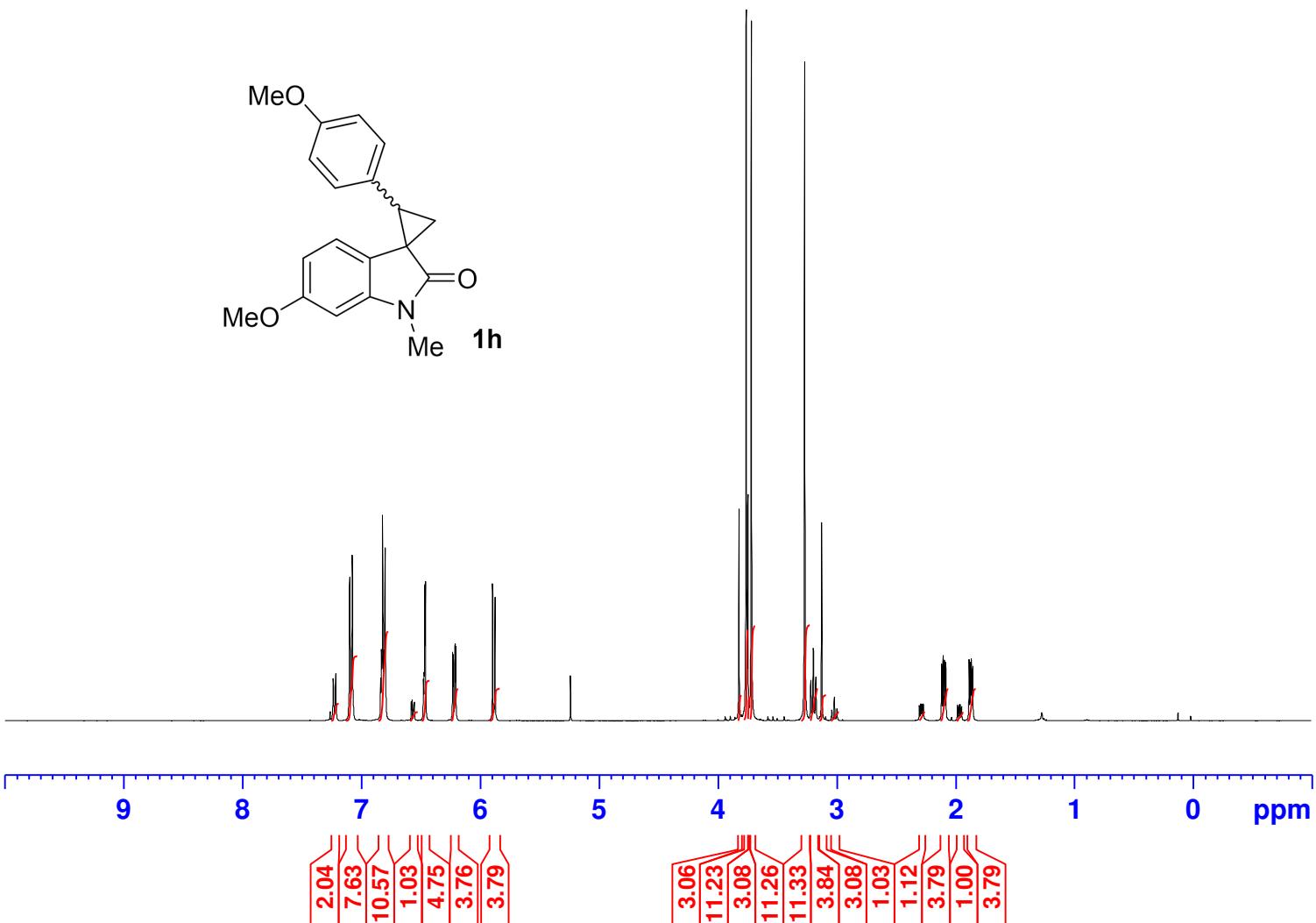


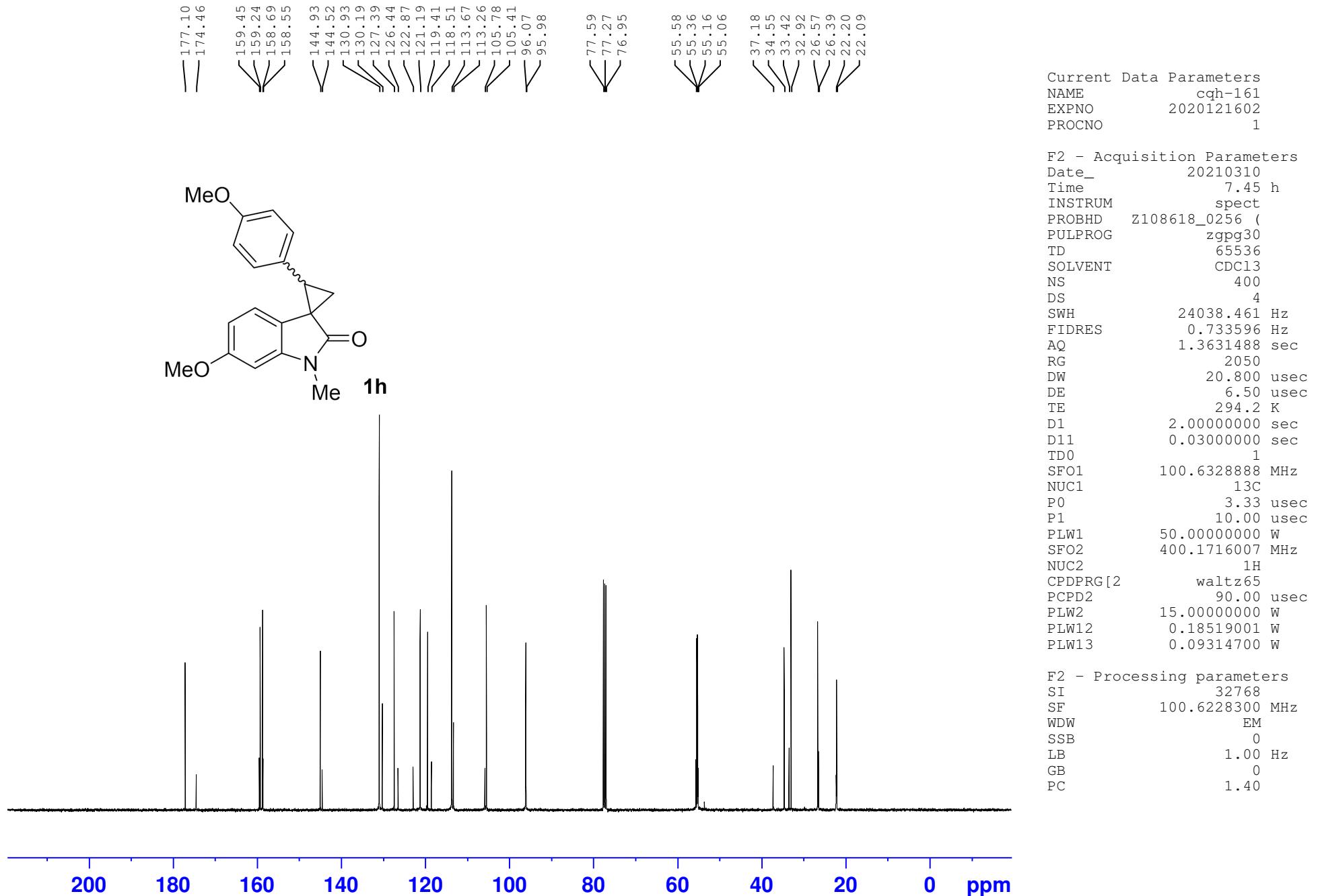


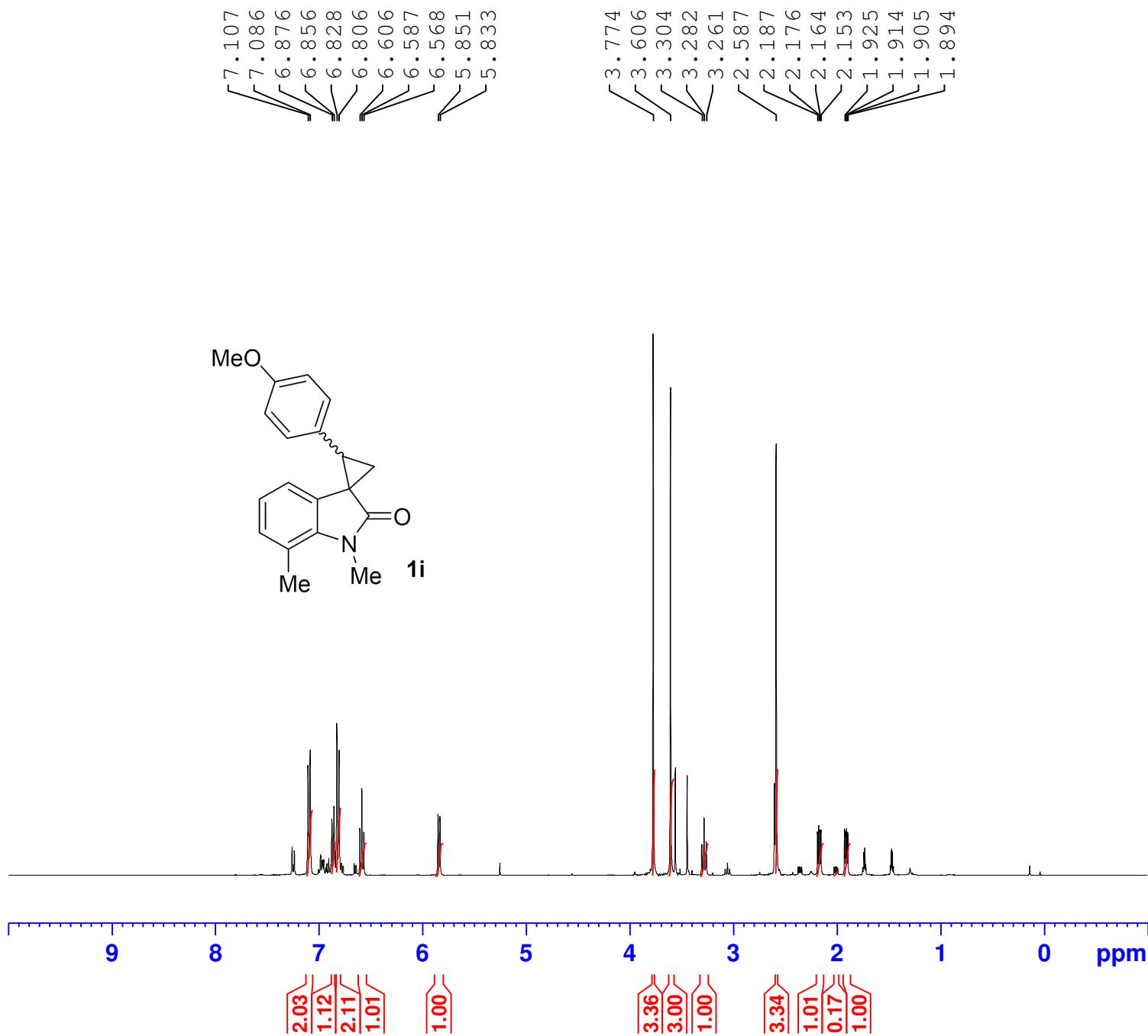
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 NS 16  
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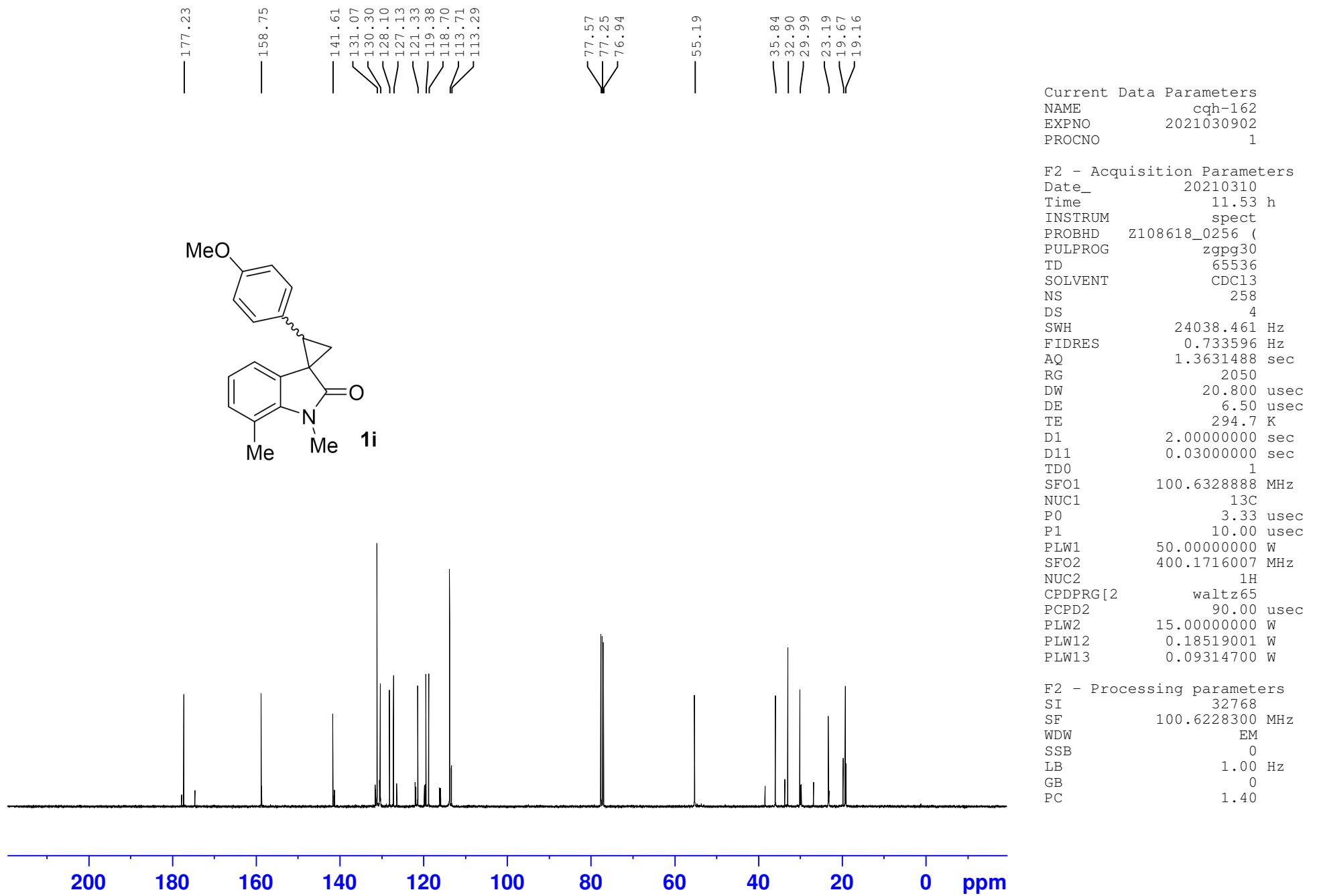


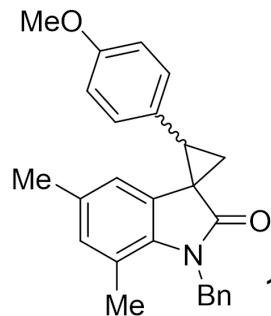
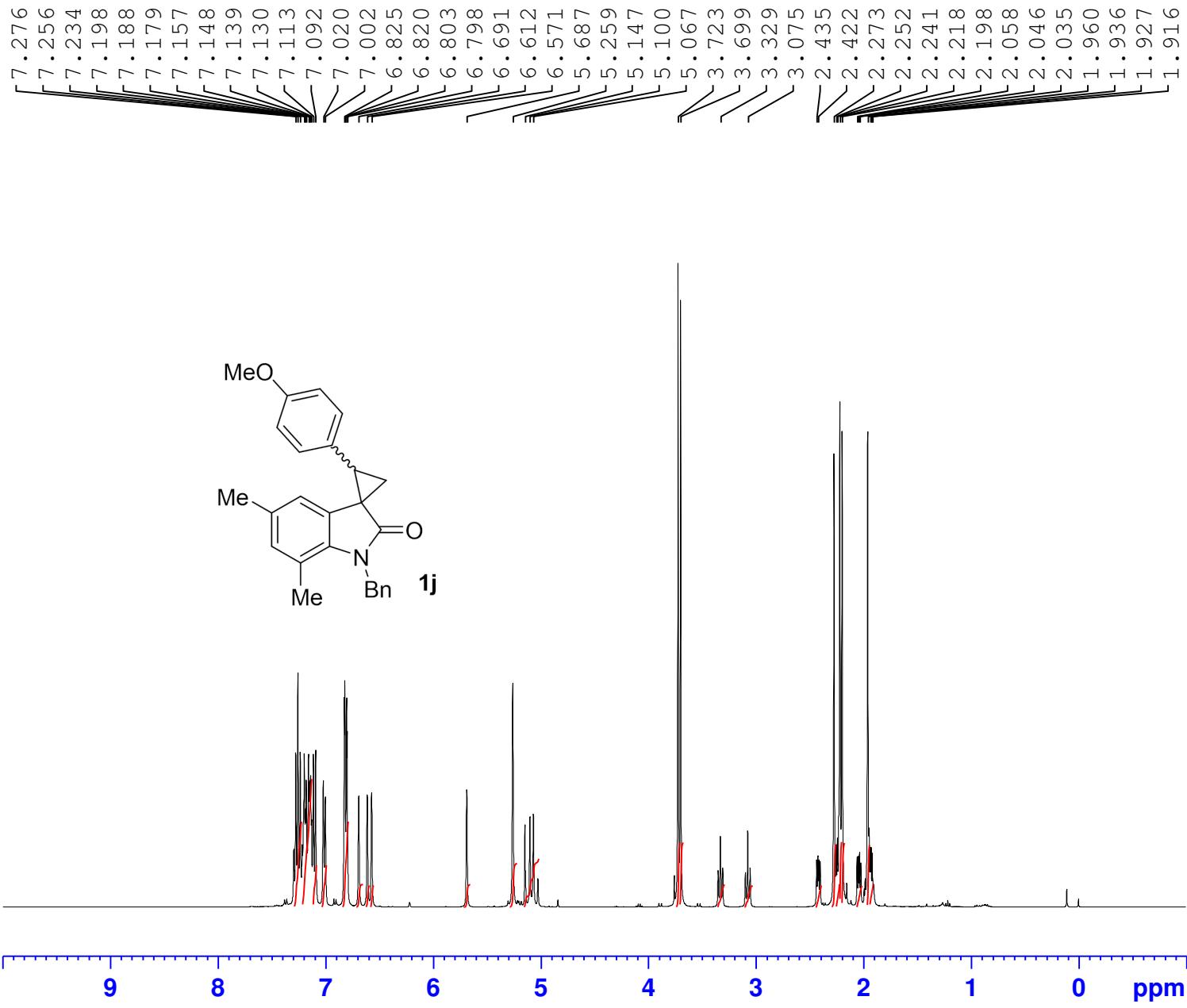


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 AQ 3.9845889 sec  
 RG 32  
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 TE 294.5 K  
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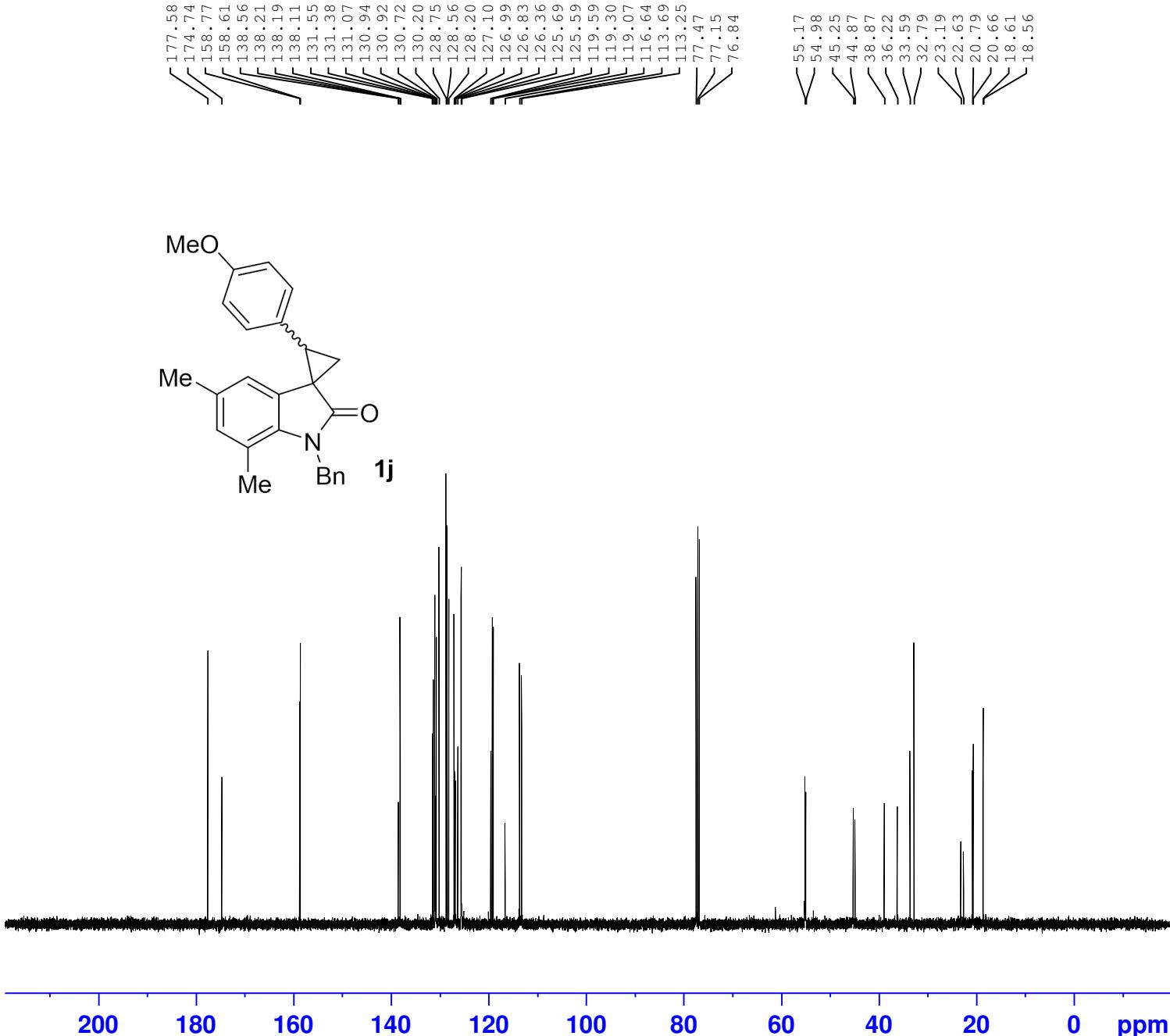
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TE              298.4 K
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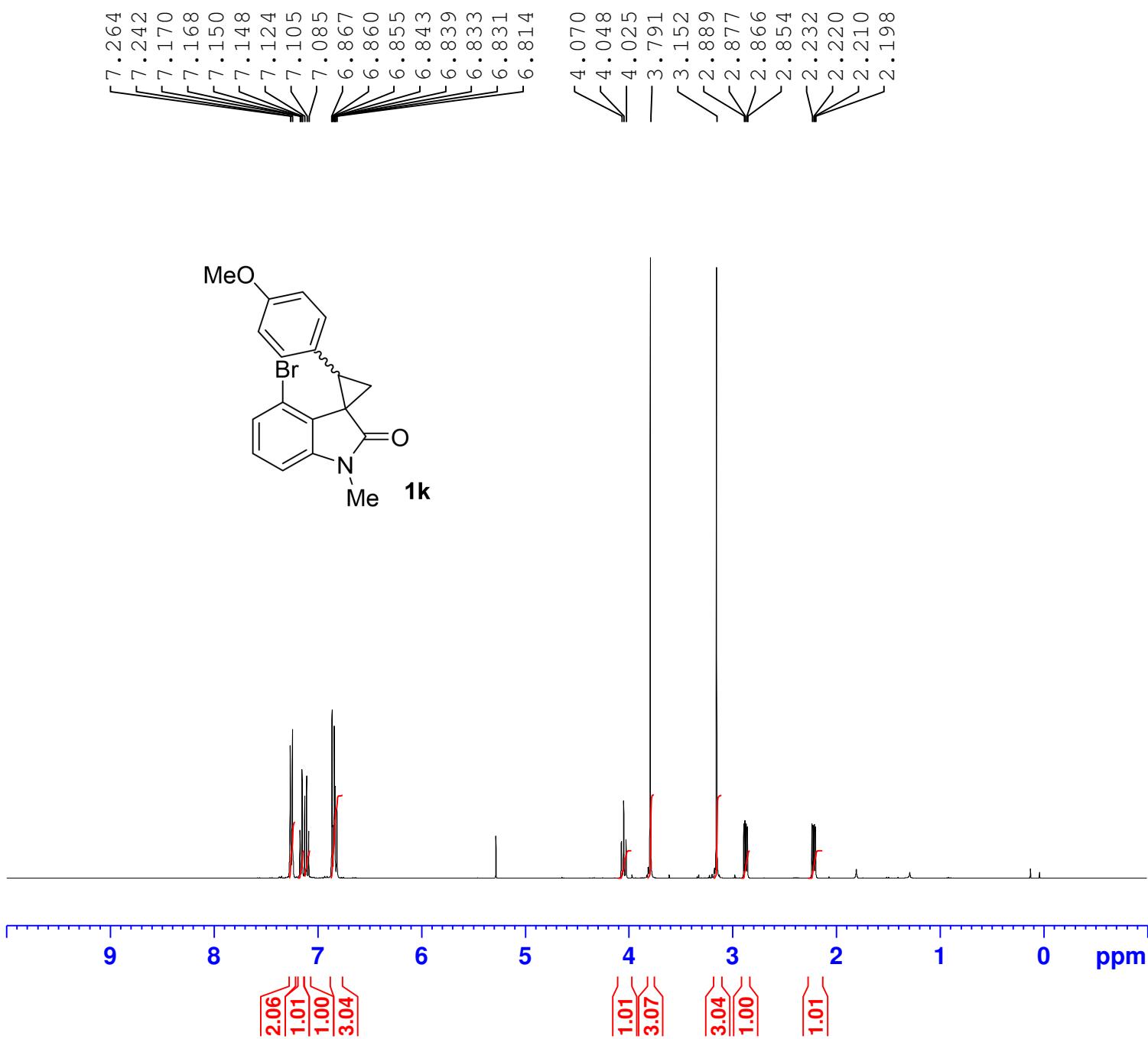
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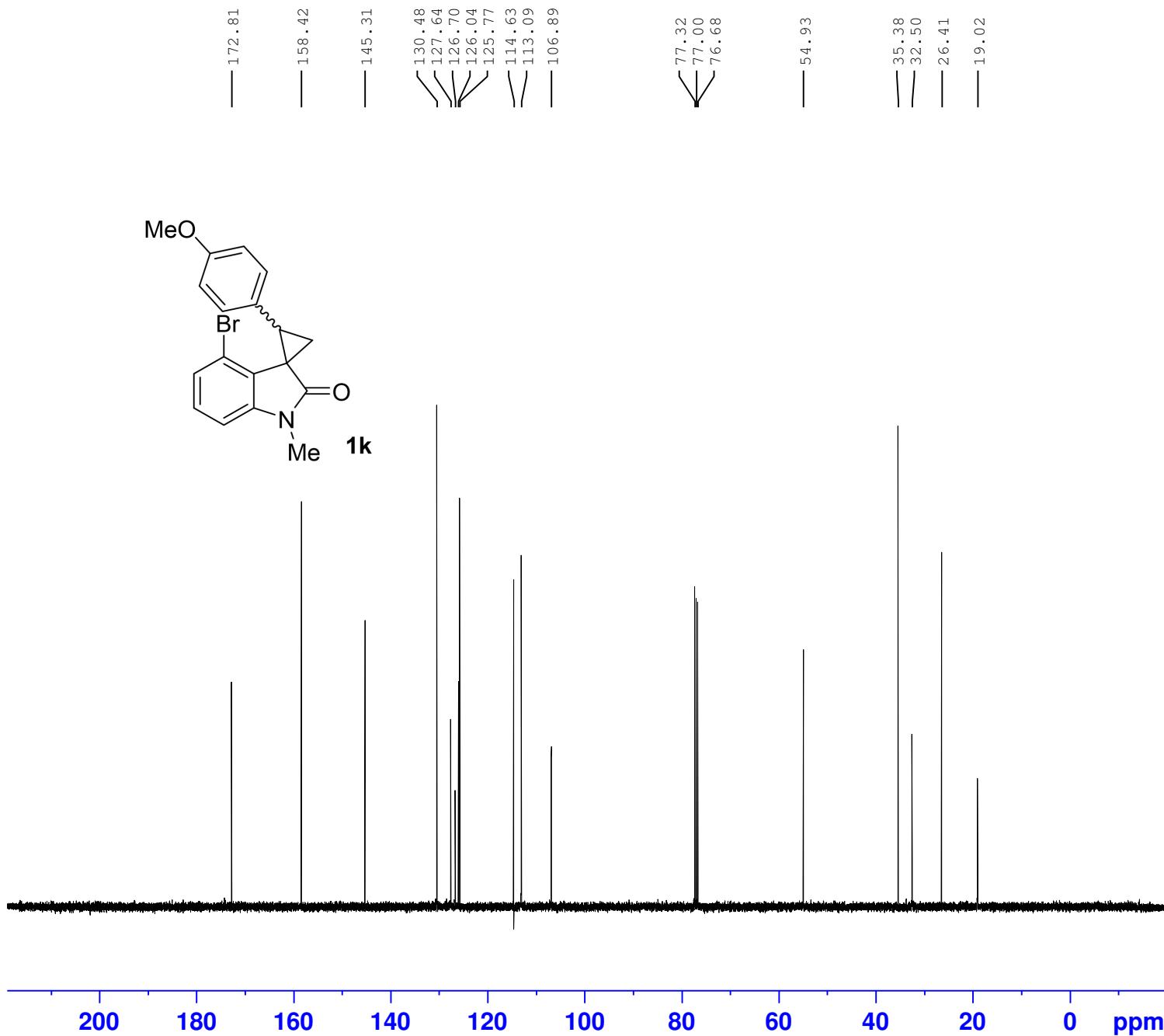
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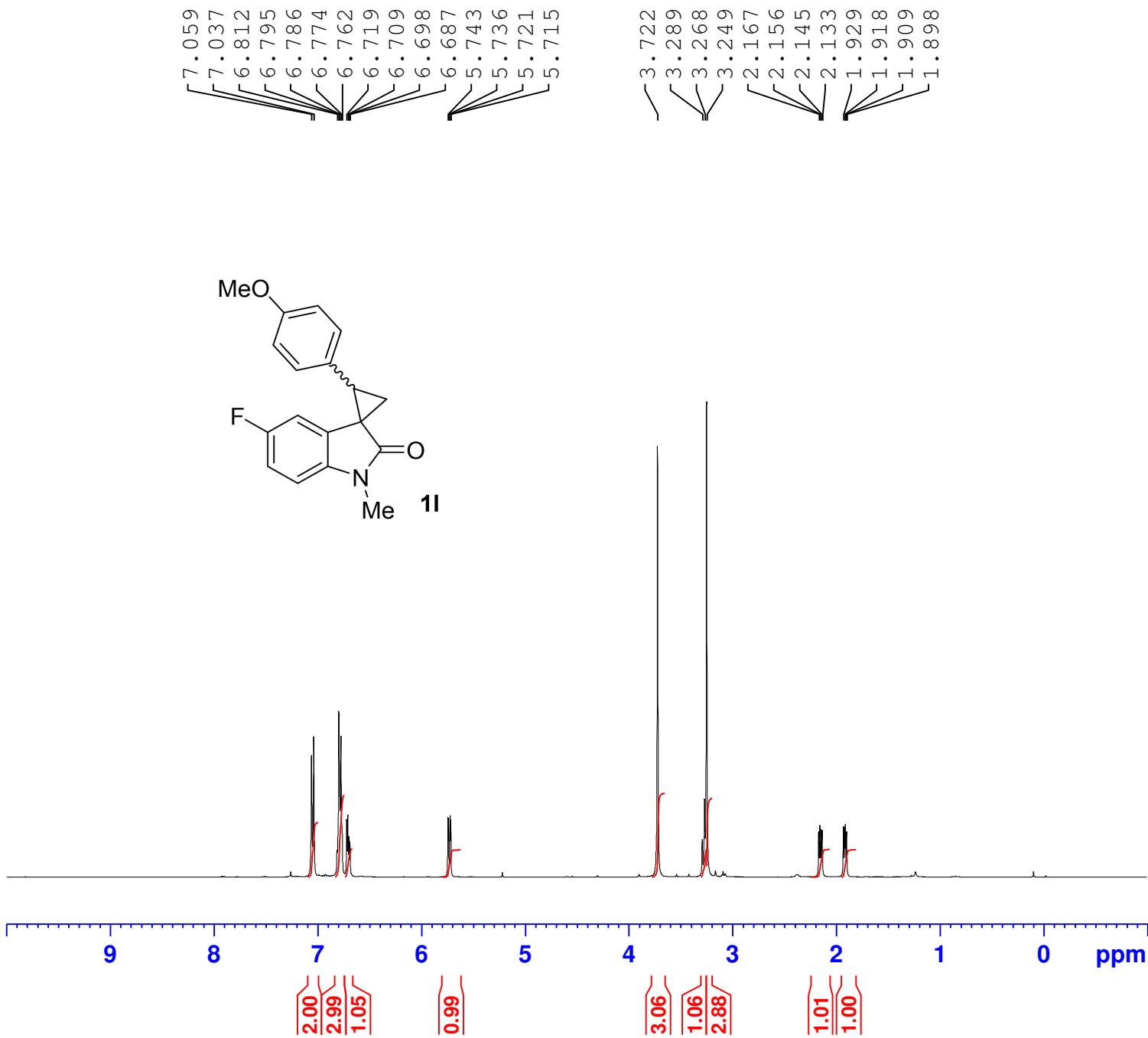


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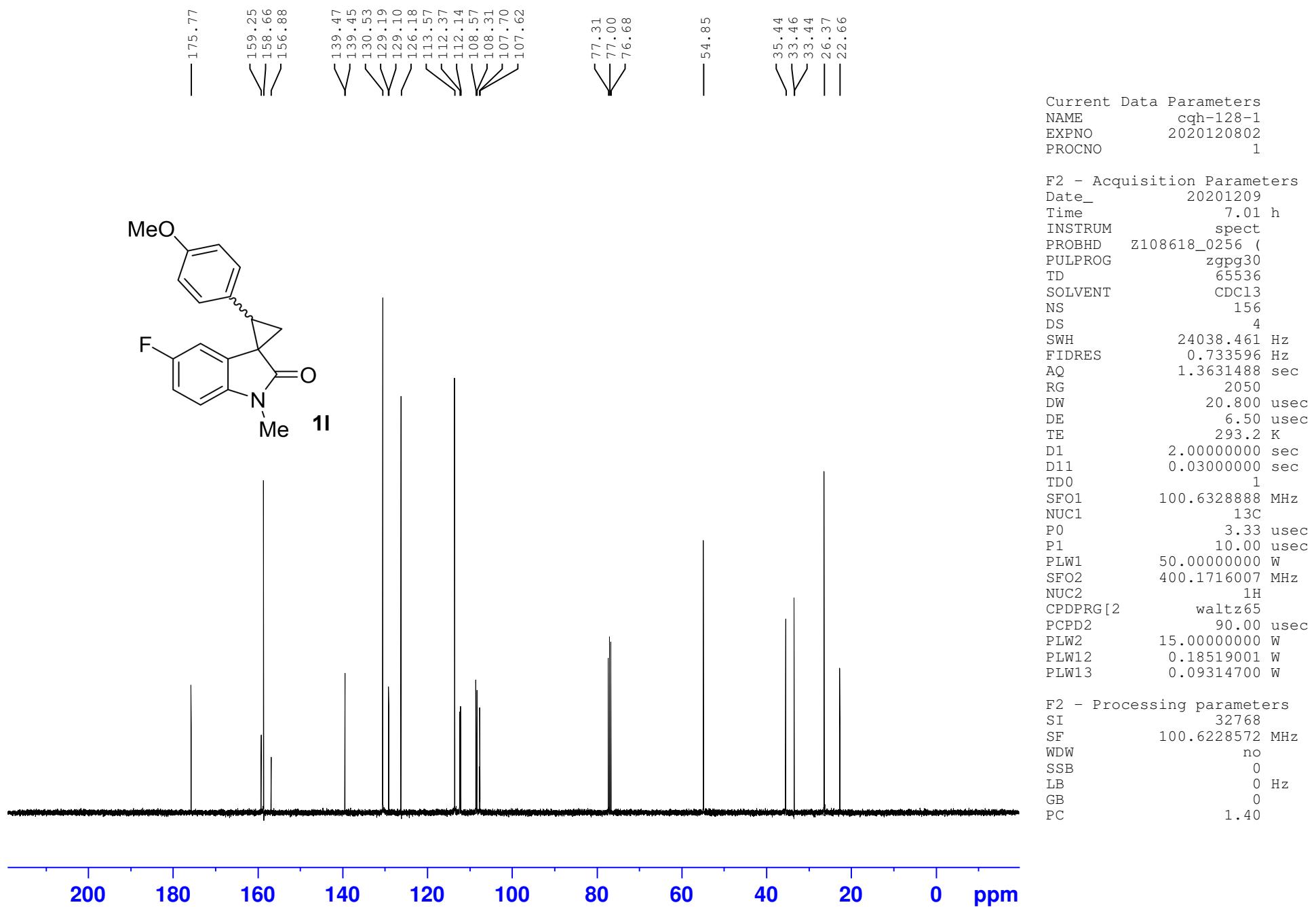


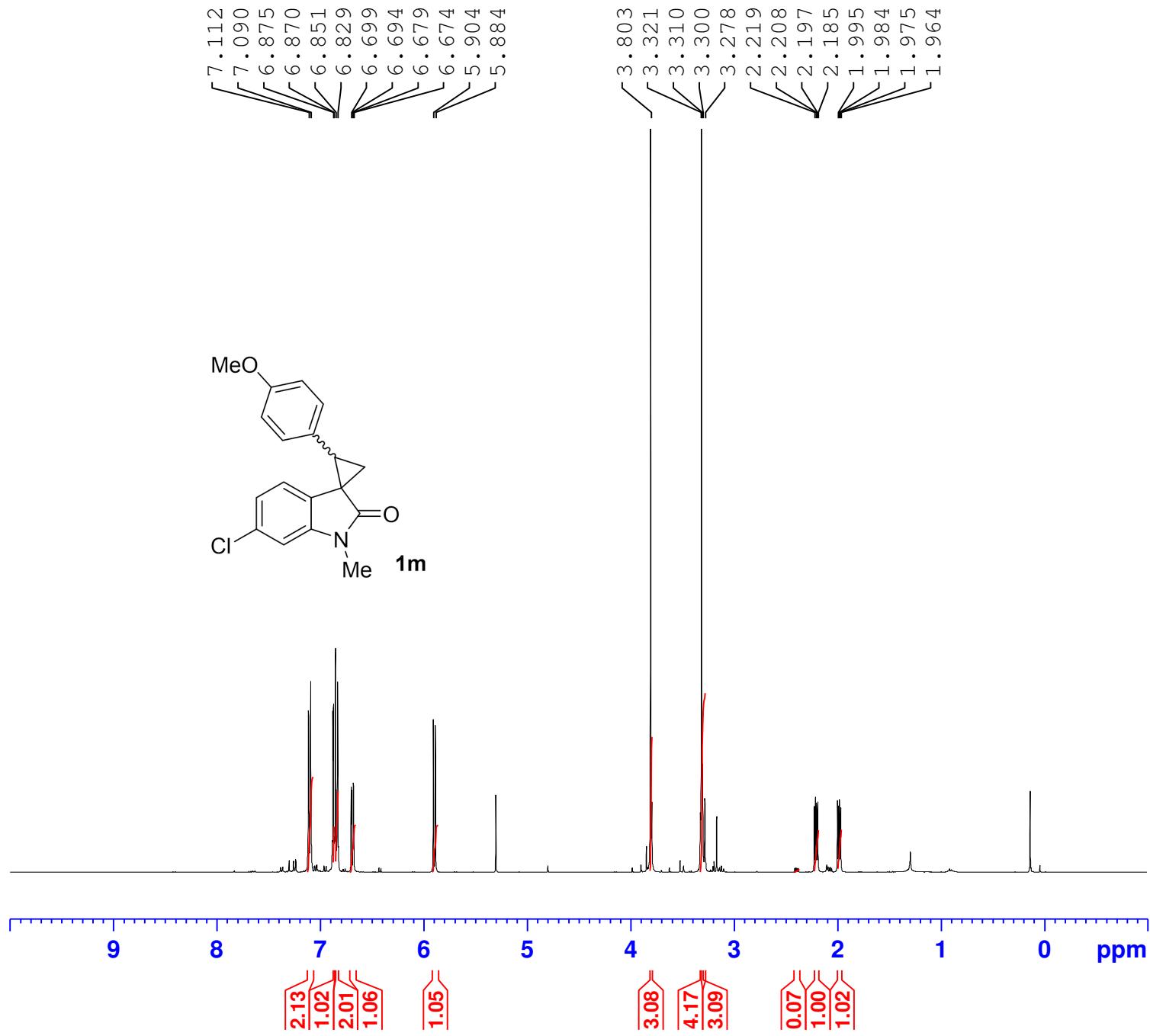


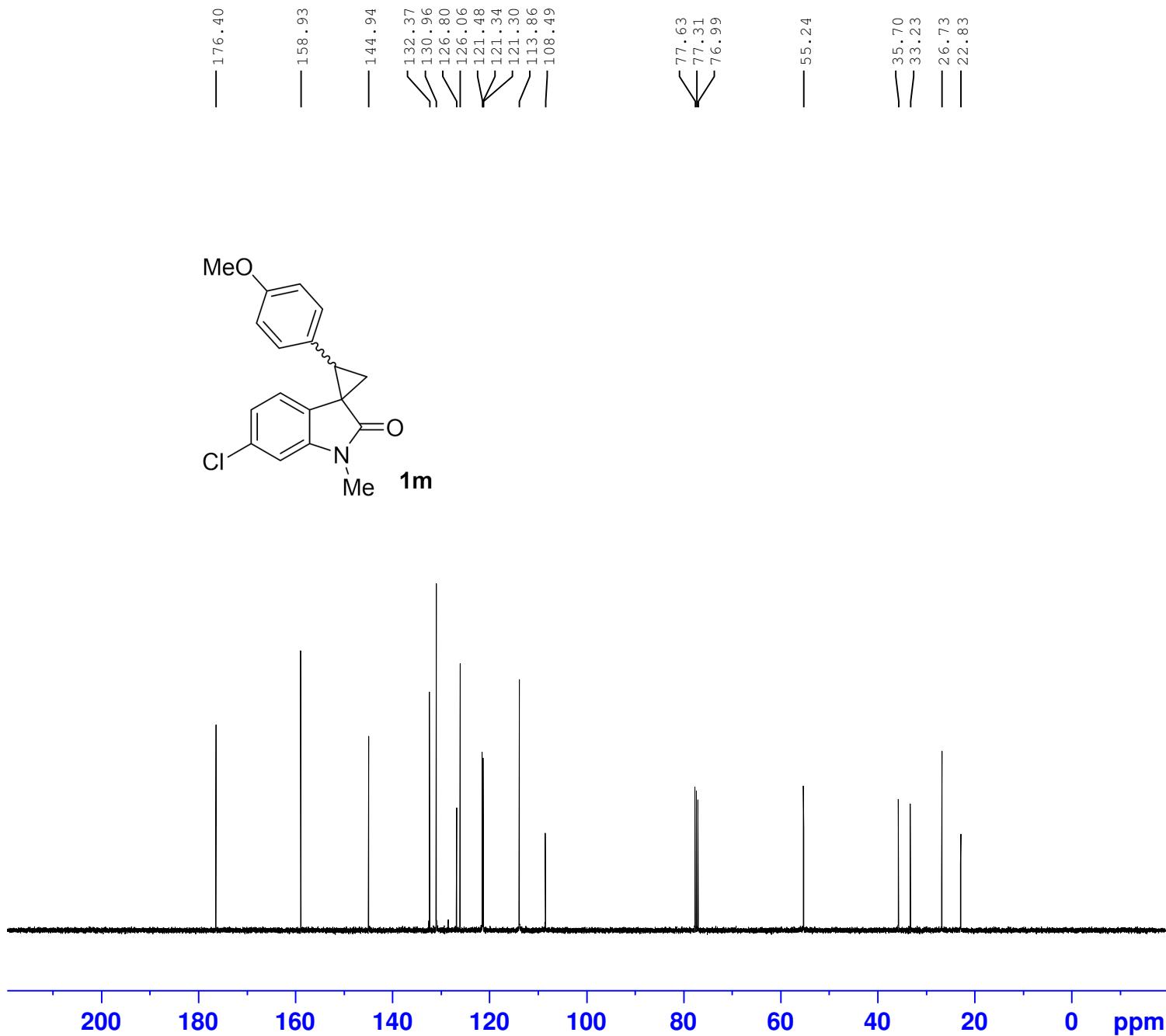
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 PC 1.00



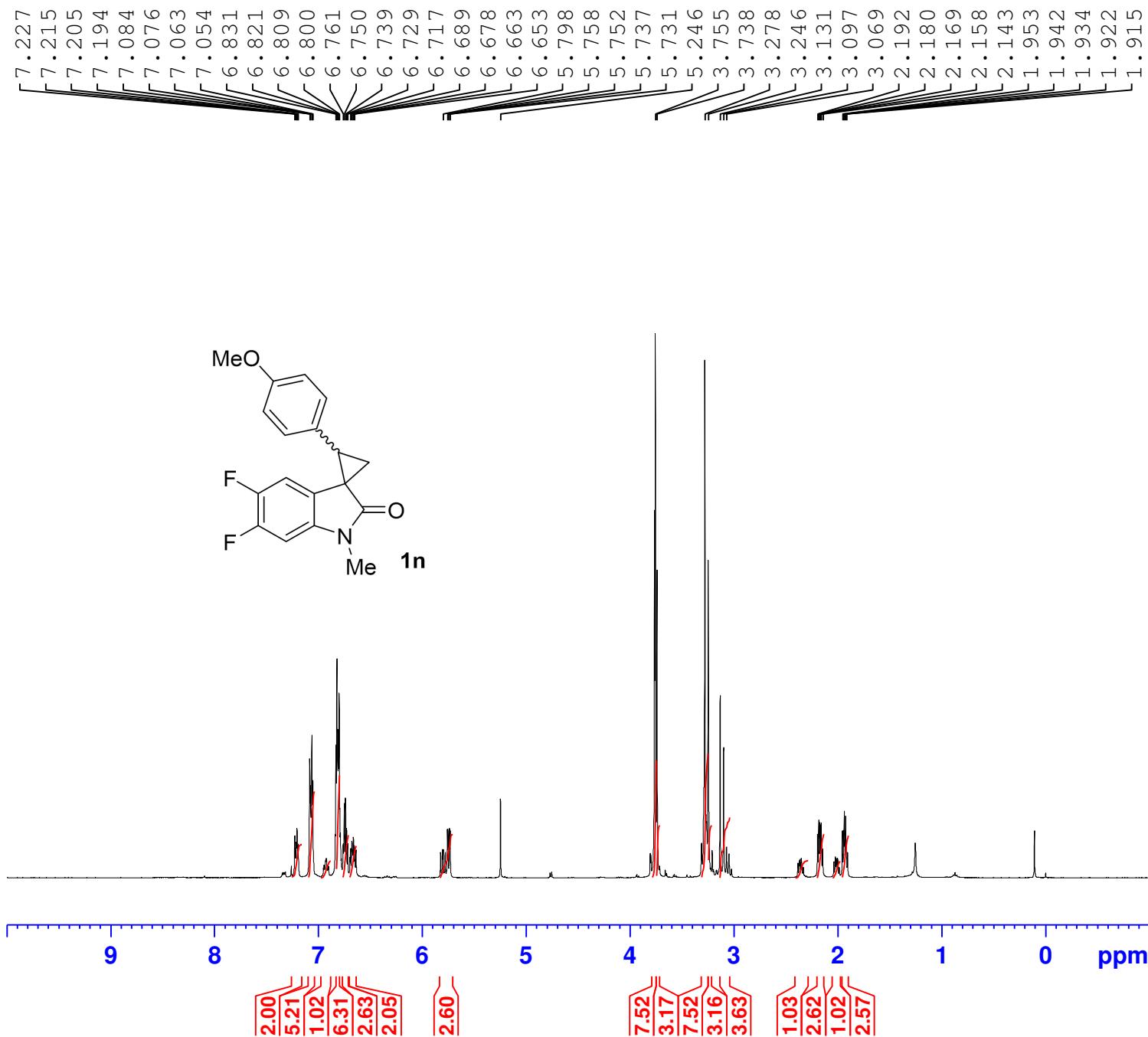


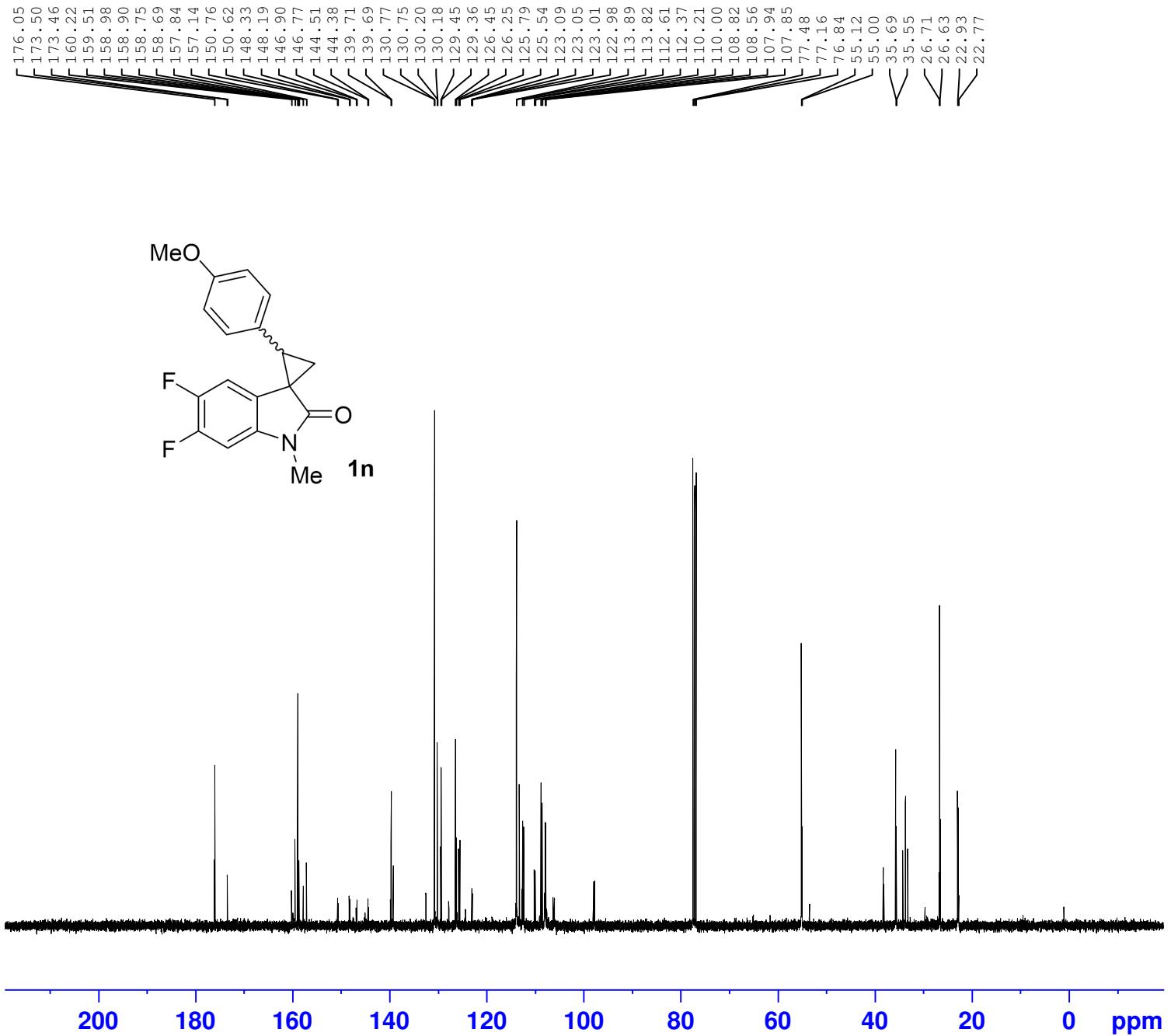


Current Data Parameters  
 NAME cgh-138  
 EXPNO 2020121602  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20201217  
 Time 9.38 h  
 INSTRUM spect  
 PROBHD Z108618\_0256 (zgpg30  
 PULPROG zgpg30  
 TD 65536  
 SOLVENT CDCl3  
 NS 150  
 DS 4  
 SWH 24038.461 Hz  
 FIDRES 0.733596 Hz  
 AQ 1.3631488 sec  
 RG 2050  
 DW 20.800 usec  
 DE 6.50 usec  
 TE 294.1 K  
 D1 2.00000000 sec  
 D11 0.03000000 sec  
 TDO 1  
 SFO1 100.6328888 MHz  
 NUC1 13C  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 50.00000000 W  
 SFO2 400.1716007 MHz  
 NUC2 1H  
 CPDPRG[2] waltz65  
 PCPD2 90.00 usec  
 PLW2 15.00000000 W  
 PLW12 0.18519001 W  
 PLW13 0.09314700 W

F2 - Processing parameters  
 SI 32768  
 SF 100.6228219 MHz  
 WDW no  
 SSB 0  
 LB 0 Hz  
 GB 0  
 PC 1.40





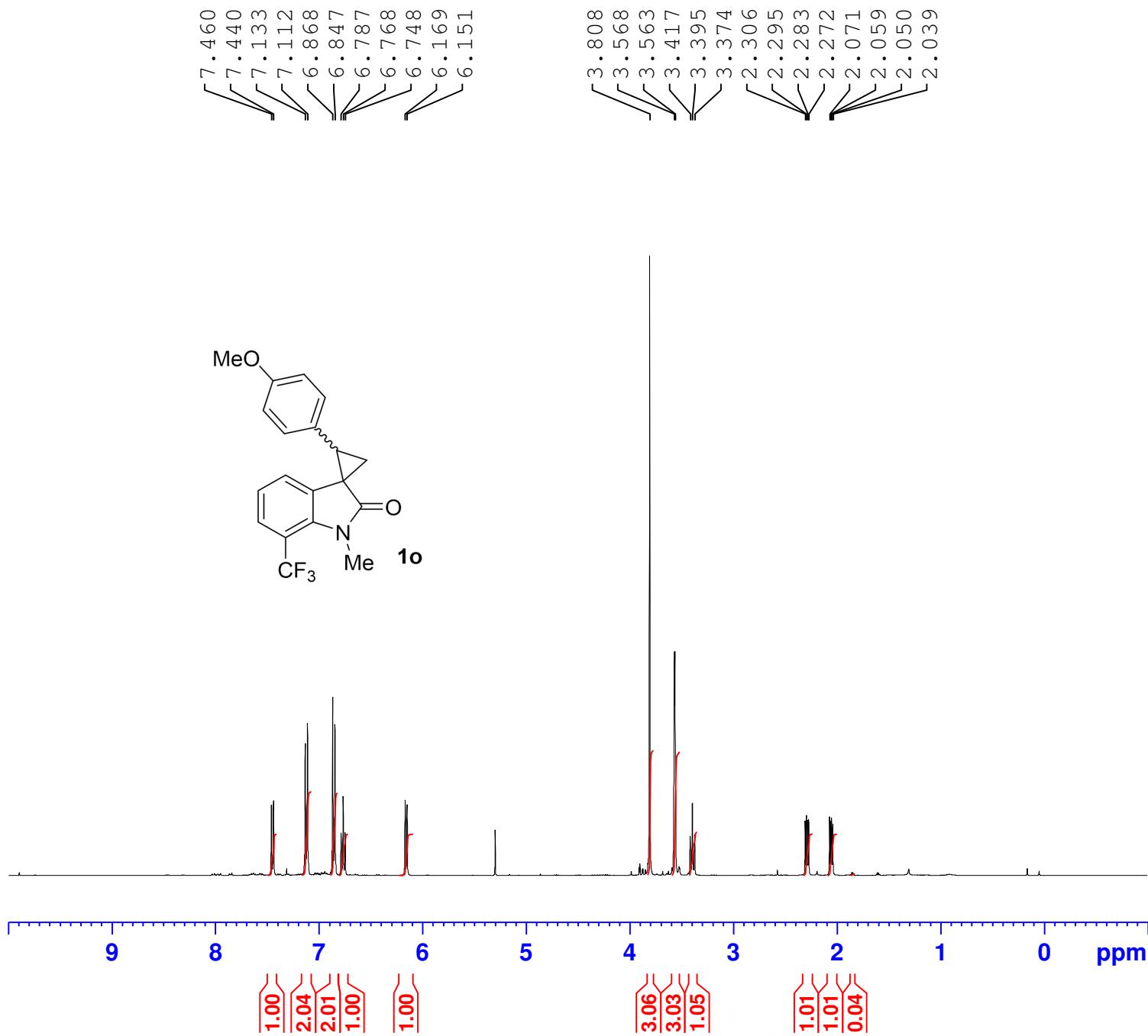
Current	Data	Parameters
NAME		dhq-18
EXPNO		2020122302
PROCNO		1

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F2 - Acquisition Parameters
Date_           20201224
Time            11.10 h
INSTRUM         spect
PROBHD         Z108618_0256 (
PULPROG        zgpg30
TD              65536
SOLVENT         CDC13
NS              377
DS              4
SWH             24038.461 Hz
FIDRES         0.733596 Hz
AQ              1.3631488 sec
RG              2050
DW              20.800 usec
DE              6.50  usec
TE              294.3 K
D1              2.00000000 sec
D11             0.03000000 sec
TD0              1
SFO1            100.6328888 MHz
NUC1            13C
P0              3.33  usec
P1              10.00  usec
PLW1            50.00000000 W
SFO2            400.1716007 MHz
NUC2            1H
CPDPRG[2]       waltz65
PCPD2           90.00  usec
PLW2            15.00000000 W
PLW12           0.18519001 W
PLW13           0.09314700 W

```

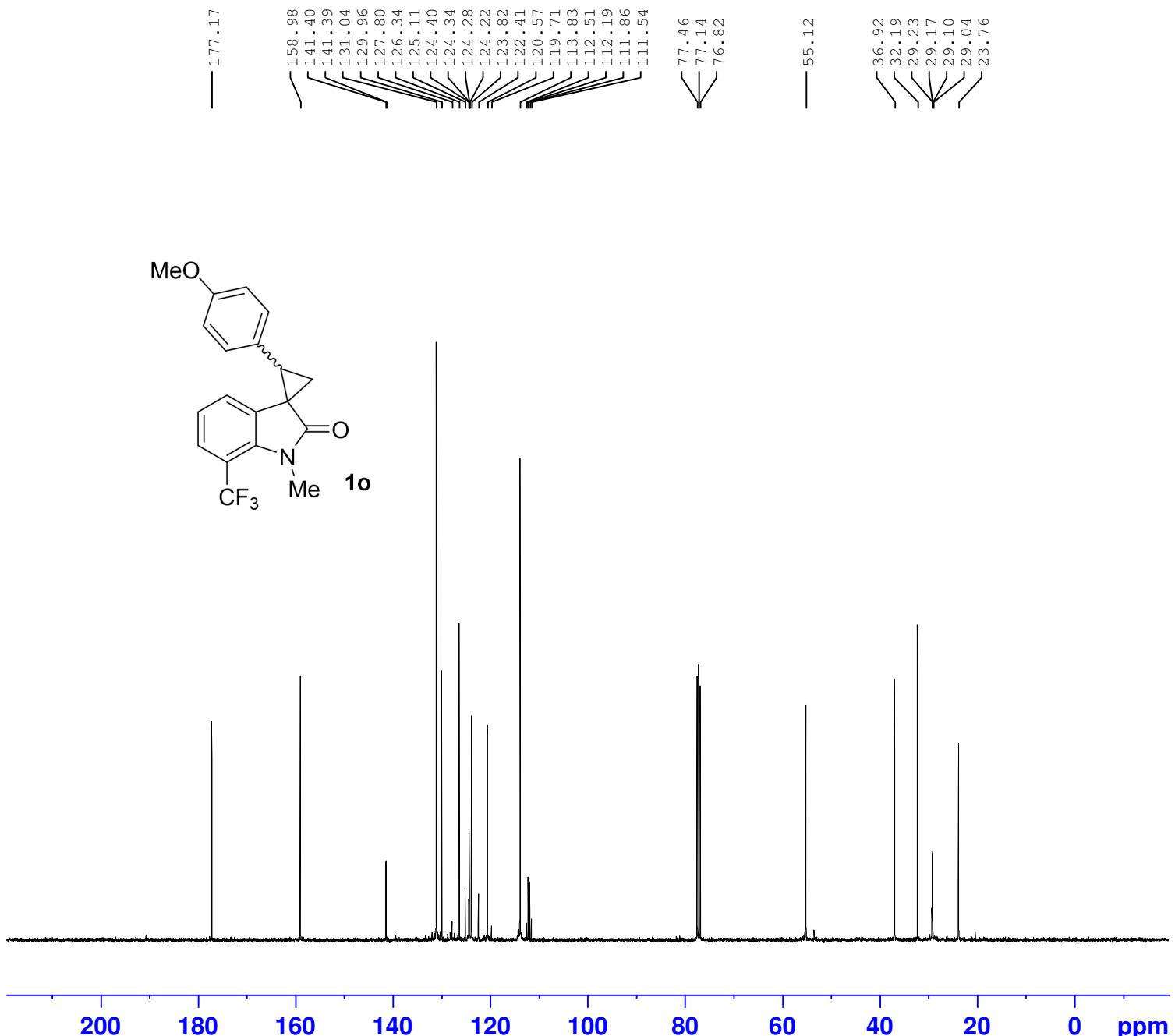
F2 - Processing parameters  
SI 32768  
SF 100.6228345 MHz  
WDW no  
SSB 0  
LB 0 Hz  
GB 0  
PC 1.40



Current Data Parameters  
NAME cgh-139  
EXPNO 2020121601  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20201217  
Time 6.13 h  
INSTRUM spect  
PROBHD z108618\_0256 (zg30  
PULPROG zg30  
TD 65536  
SOLVENT CDCl<sub>3</sub>  
NS 16  
DS 2  
SWH 8223.685 Hz  
FIDRES 0.250967 Hz  
AQ 3.9845889 sec  
RG 50.8  
DW 60.800 usec  
DE 6.50 usec  
TE 293.7 K  
D1 1.00000000 sec  
TD0 1  
SFO1 400.1724712 MHz  
NUC1 1H  
P0 3.33 usec  
P1 10.00 usec  
PLW1 15.00000000 W

F2 - Processing parameters  
SI 32768  
SF 400.1699854 MHz  
WDW no  
SSB 0  
LB 0 Hz  
GB 0  
PC 1.00



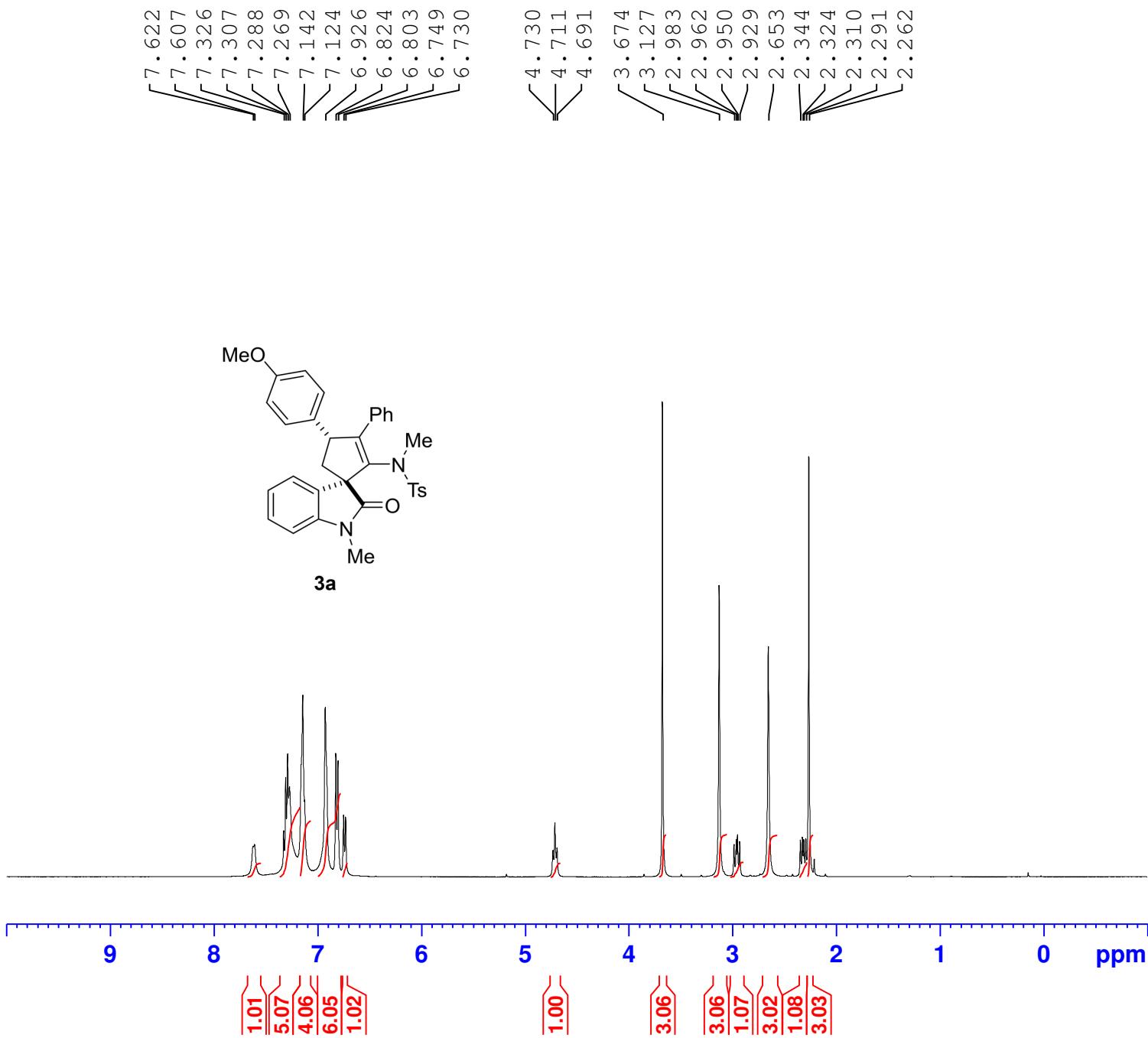
Current	Data	Parameters
NAME	cgh-139	
EXPNO	2020121602	
PROCNO		1

```

F2 - Acquisition Parameters
Date_           20201217
Time            9.56 h
INSTRUM        spect
PROBHD         Z108618_0256 (
PULPROG        zgppg30
TD              65536
SOLVENT         CDC13
NS              274
DS                            4
SWH             24038.461 Hz
FIDRES         0.733596 Hz
AQ              1.3631488 sec
RG              2050
DW              20.800 usec
DE              6.50 usec
TE              294.1 K
D1              2.00000000 sec
D11             0.03000000 sec
TD0                           1
SFO1            100.6328888 MHz
NUC1            13C
P0              3.33 usec
P1              10.00 usec
PLW1            50.00000000 W
SFO2            400.1716007 MHz
NUC2            1H
CPDPRG[2]      waltz65
PCPD2           90.00 usec
PLW2            15.00000000 W
PLW12           0.18519001 W
PLW13           0.09314700 W

```

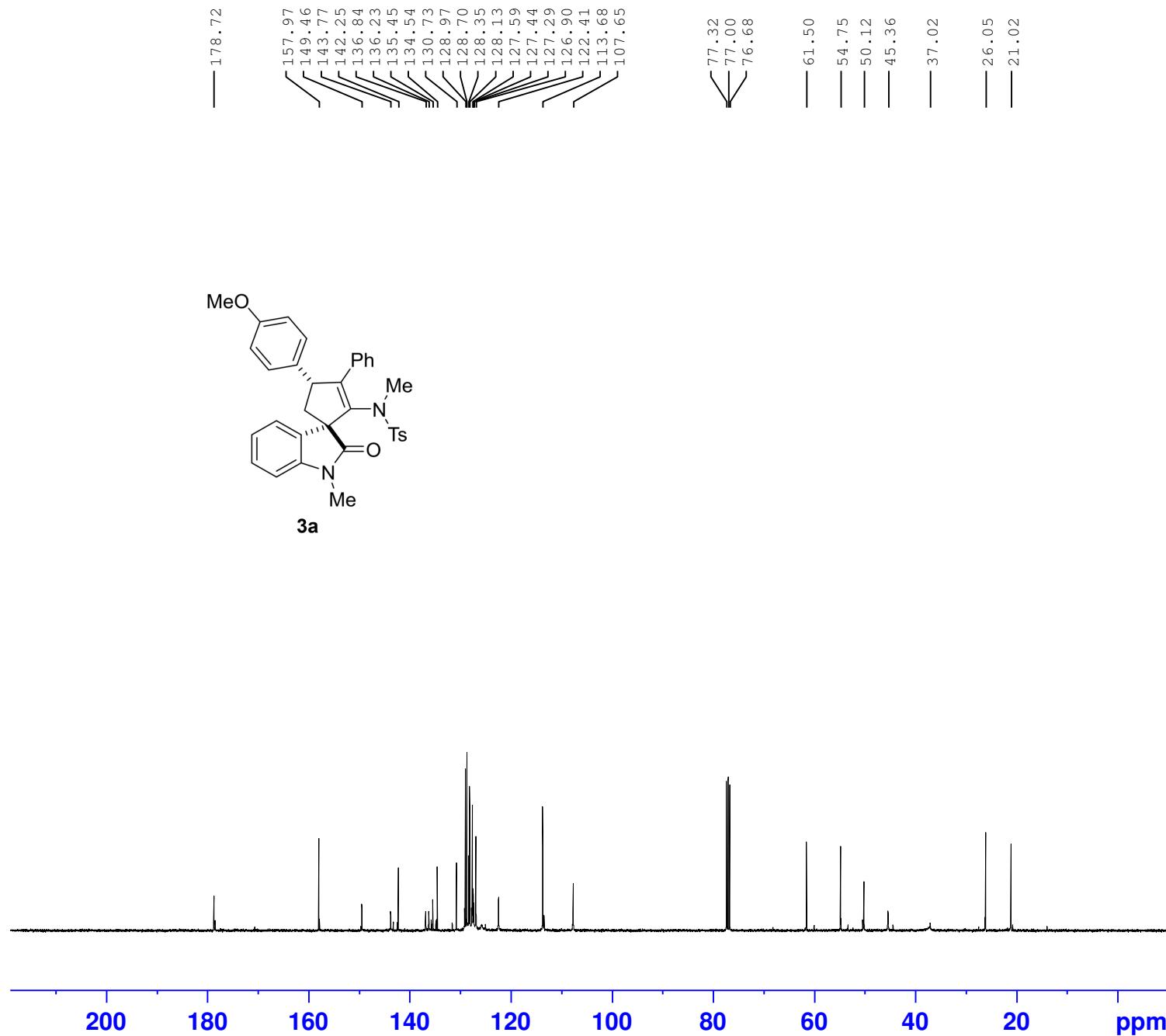
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F2 - Processing parameters
SI           32768
SF          100.6228300 MHz
WDW          EM
SSB            0
LB           1.00 Hz
GB            0
PC          1.40
```



Current Data Parameters  
 NAME cgh-313  
 EXPNO 2021091601  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20210916  
 Time 3.34 h  
 INSTRUM spect  
 PROBHD z108618\_0256 (zg30  
 PULPROG 65536  
 TD CDC13  
 NS 16  
 DS 2  
 SWH 8223.685 Hz  
 FIDRES 0.250967 Hz  
 AQ 3.9845889 sec  
 RG 18  
 DW 60.800 usec  
 DE 6.50 usec  
 TE 296.4 K  
 D1 1.00000000 sec  
 TD0 1  
 SFO1 400.1724712 MHz  
 NUC1 1H  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 15.00000000 W

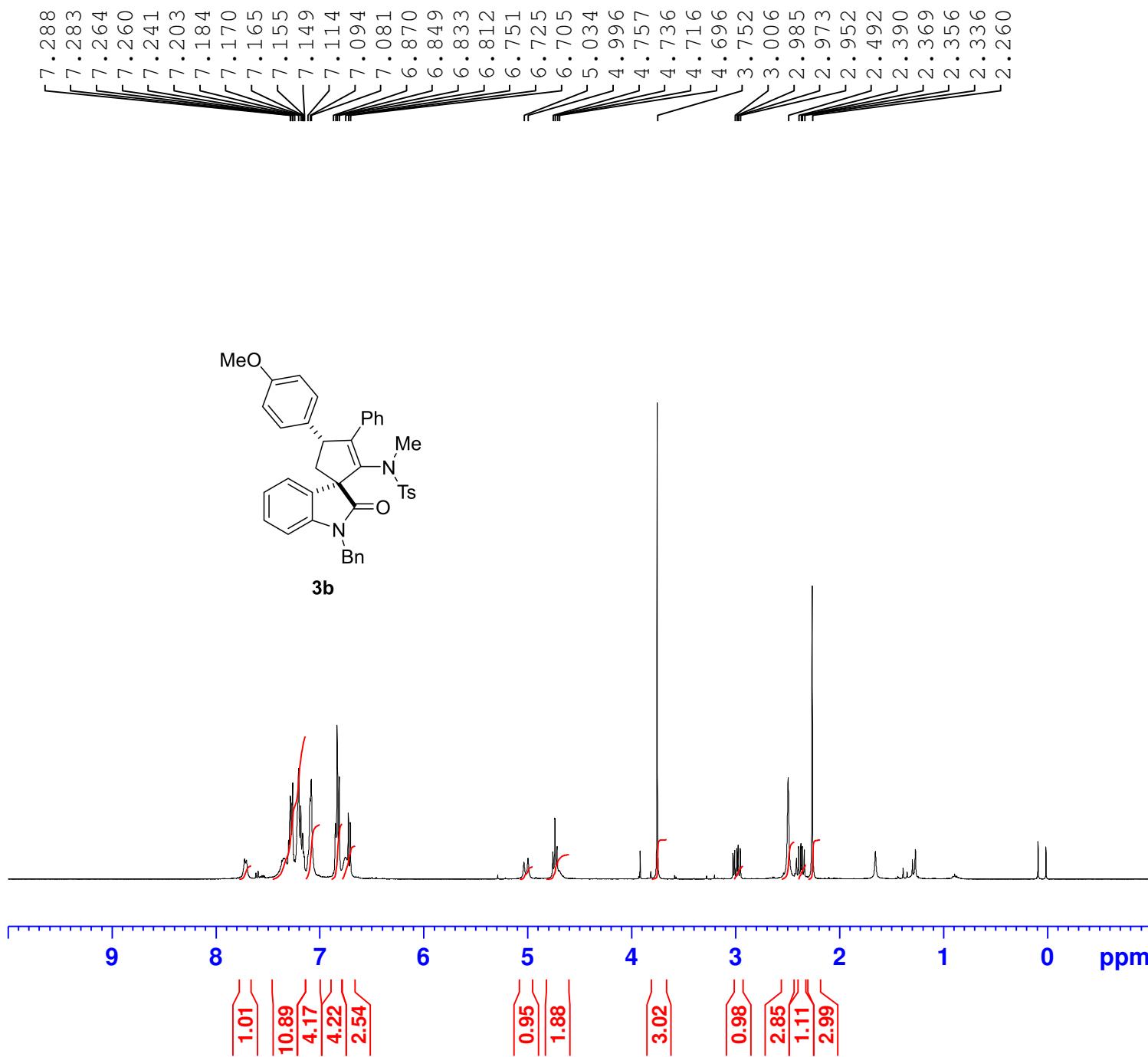
F2 - Processing parameters  
 SI 32768  
 SF 400.1700072 MHz  
 WDW no  
 SSB 0  
 LB 0 Hz  
 GB 0  
 PC 1.00



Current Data Parameters  
NAME zy-2-2  
EXPNO 2020102202  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20201023  
Time 12.39 h  
INSTRUM spect  
PROBHD Z108618\_0256 (   
PULPROG zgpg30  
TD 65536  
SOLVENT CDC13  
NS 240  
DS 4  
SWH 24038.461 Hz  
FIDRES 0.733596 Hz  
AQ 1.3631488 sec  
RG 2050  
DW 20.800 usec  
DE 6.50 usec  
TE 298.5 K  
D1 2.00000000 sec  
D11 0.03000000 sec  
TD0 1  
SFO1 100.6328888 MHz  
NUC1 <sup>13</sup>C  
P0 3.33 usec  
P1 10.00 usec  
PLW1 50.00000000 W  
SFO2 400.1716007 MHz  
NUC2 <sup>1</sup>H  
CPDPRG[2] waltz65  
PCPD2 90.00 usec  
PLW2 15.00000000 W  
PLW12 0.18519001 W  
PLW13 0.09314700 W

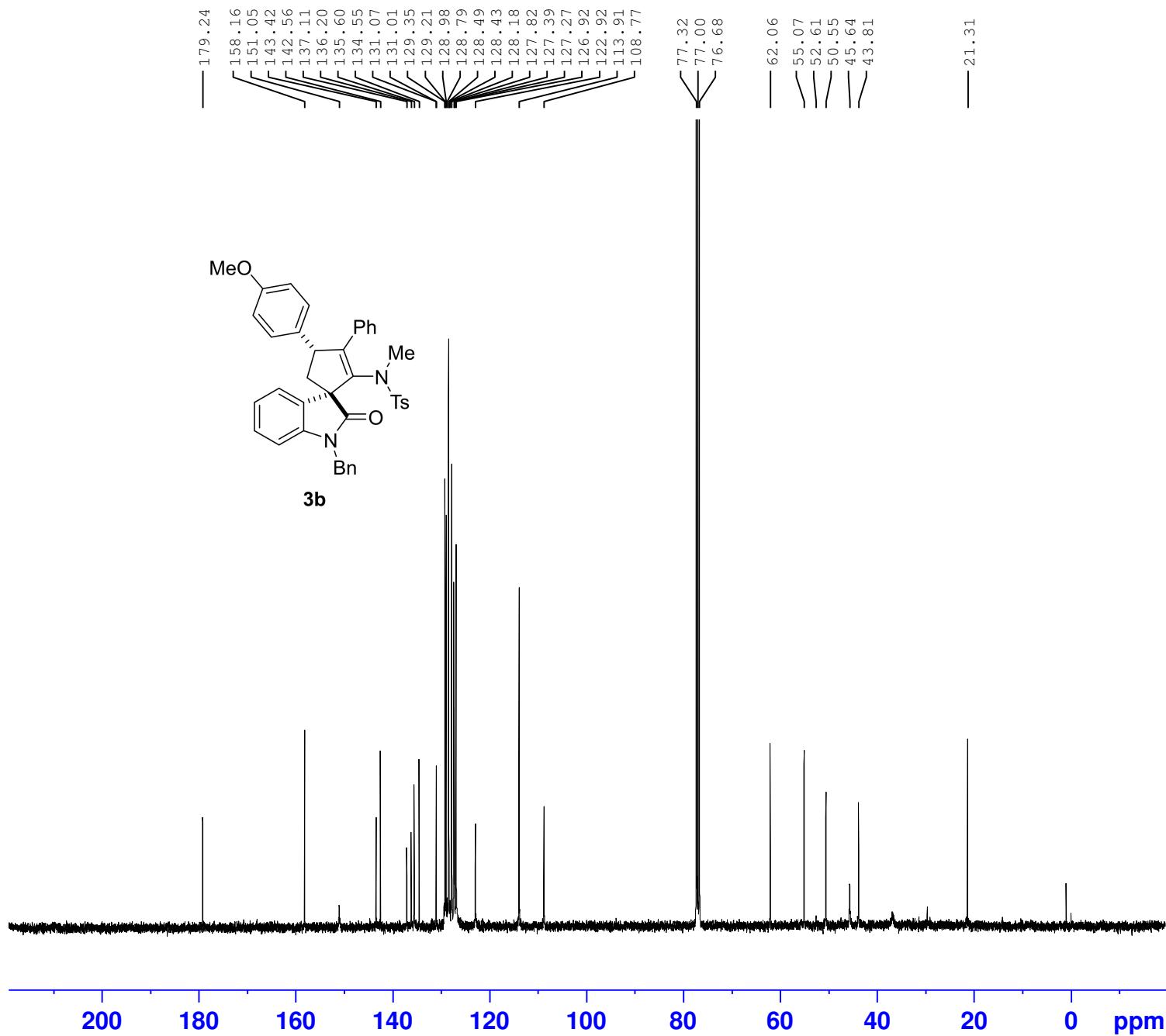
F2 - Processing parameters  
SI 32768  
SF 100.6228660 MHz  
WDW EM  
SSB 0  
LB 1.00 Hz  
GB 0  
PC 1.40



Current Data Parameters  
NAME cgh-103  
EXPNO 2020111701  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20201118  
Time 5.44 h  
INSTRUM spect  
PROBHD Z108618\_0256 (zg30)  
PULPROG 65536  
SOLVENT CDCl<sub>3</sub>  
NS 16  
DS 2  
SWH 8223.685 Hz  
FIDRES 0.250967 Hz  
AQ 3.9845889 sec  
RG 144  
DW 60.800 usec  
DE 6.50 usec  
TE 296.4 K  
D1 1.00000000 sec  
TD0 1  
SFO1 400.1724712 MHz  
NUC1 1H  
P0 3.33 usec  
P1 10.00 usec  
PLW1 15.00000000 W

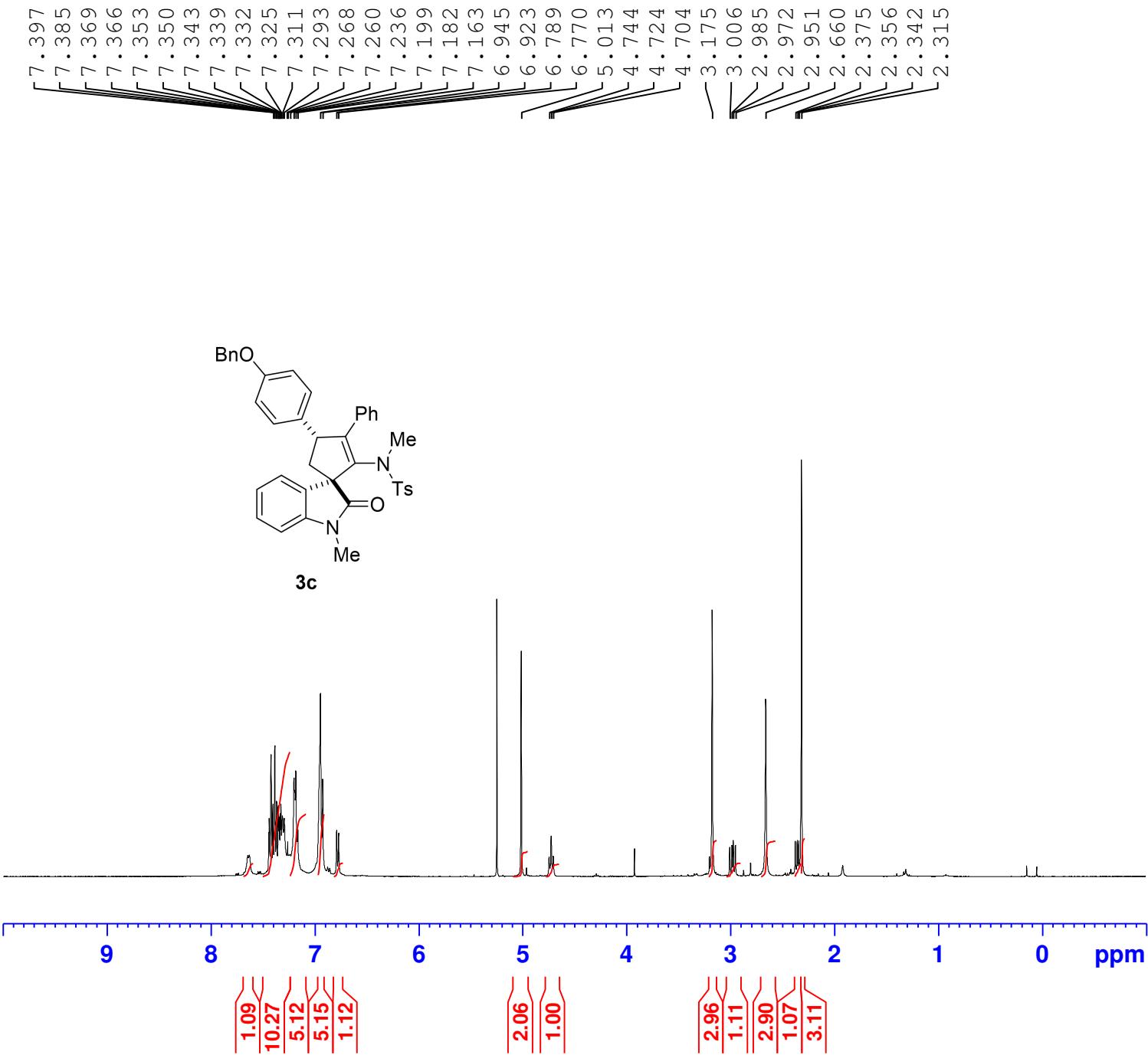
F2 - Processing parameters  
SI 32768  
SF 400.1700088 MHz  
WDW no  
SSB 0  
LB 0 Hz  
GB 0  
PC 1.00



Current Data Parameters  
NAME cgh-103-2  
EXPNO 2021101902  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20211019  
Time 10.29 h  
INSTRUM spect  
PROBHD Z108618\_0256 (zgpg30  
PULPROG 65536  
TD 4096  
SOLVENT CDCl<sub>3</sub>  
NS 800  
DS 4  
SWH 24038.461 Hz  
FIDRES 0.733596 Hz  
AQ 1.3631488 sec  
RG 2050  
DW 20.800 usec  
DE 6.50 usec  
TE 293.2 K  
D1 2.00000000 sec  
D11 0.03000000 sec  
TD0 1  
SFO1 100.6328888 MHz  
NUC1 13C  
P0 3.33 usec  
P1 10.00 usec  
PLW1 50.00000000 W  
SFO2 400.1716007 MHz  
NUC2 1H  
CPDPRG[2] waltz65  
PCPD2 90.00 usec  
PLW2 15.00000000 W  
PLW12 0.18519001 W  
PLW13 0.09314700 W

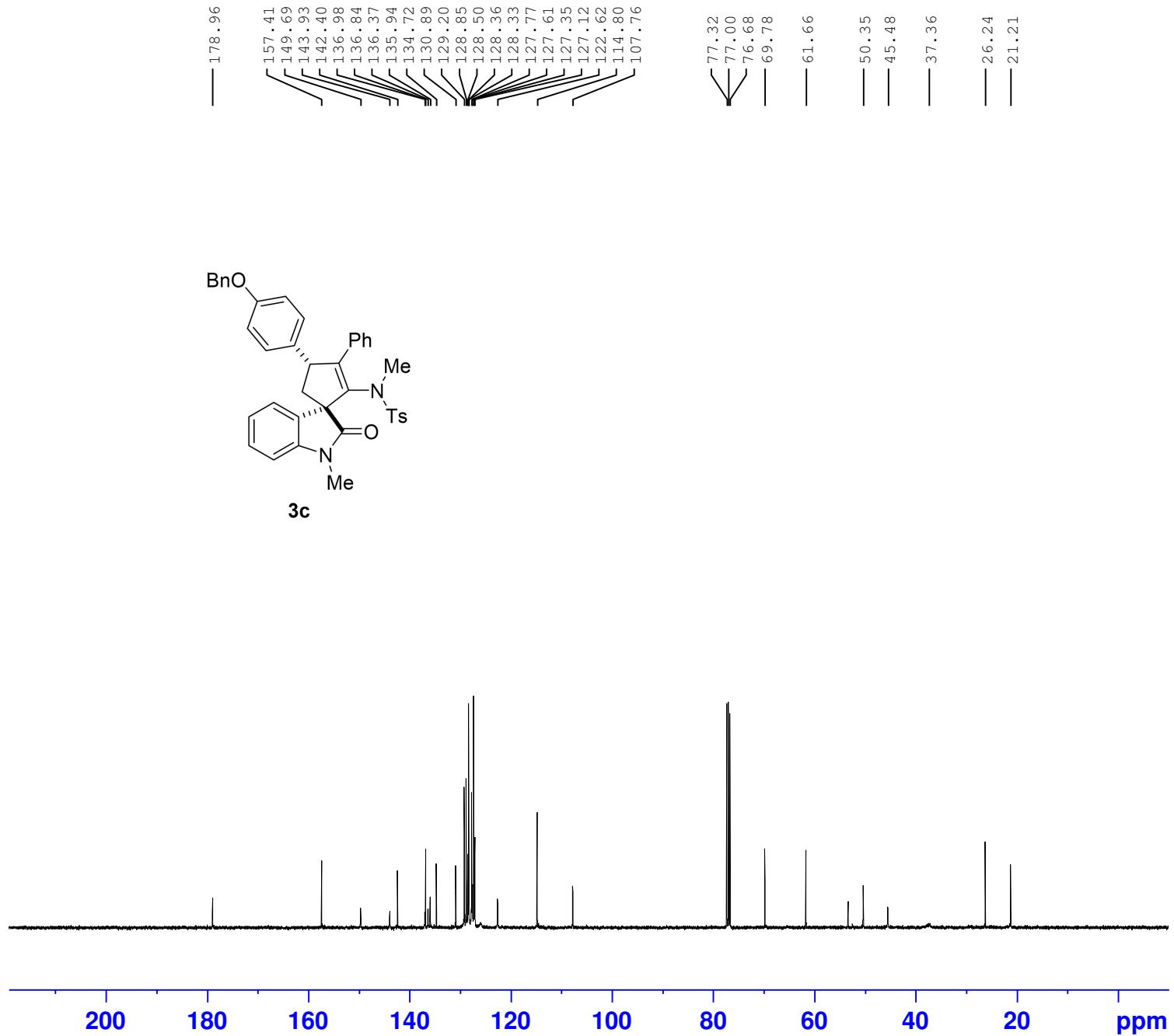
F2 - Processing parameters  
SI 32768  
SF 100.6228388 MHz  
WDW EM  
SSB 0  
LB 1.00 Hz  
GB 0  
PC 1.40



Current Data Parameters  
NAME cgh-82  
EXPNO 2020102901  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20201030  
Time 11.57 h  
INSTRUM spect  
PROBHD Z108618\_0256 (zg30)  
PULPROG zg30  
TD 65536  
SOLVENT CDCl<sub>3</sub>  
NS 16  
DS 2  
SWH 8223.685 Hz  
FIDRES 0.250967 Hz  
AQ 3.9845889 sec  
RG 36  
DW 60.800 usec  
DE 6.50 usec  
TE 297.6 K  
D1 1.00000000 sec  
TD0 1  
SFO1 400.1724712 MHz  
NUC1 1H  
P0 3.33 usec  
P1 10.00 usec  
PLW1 15.00000000 W

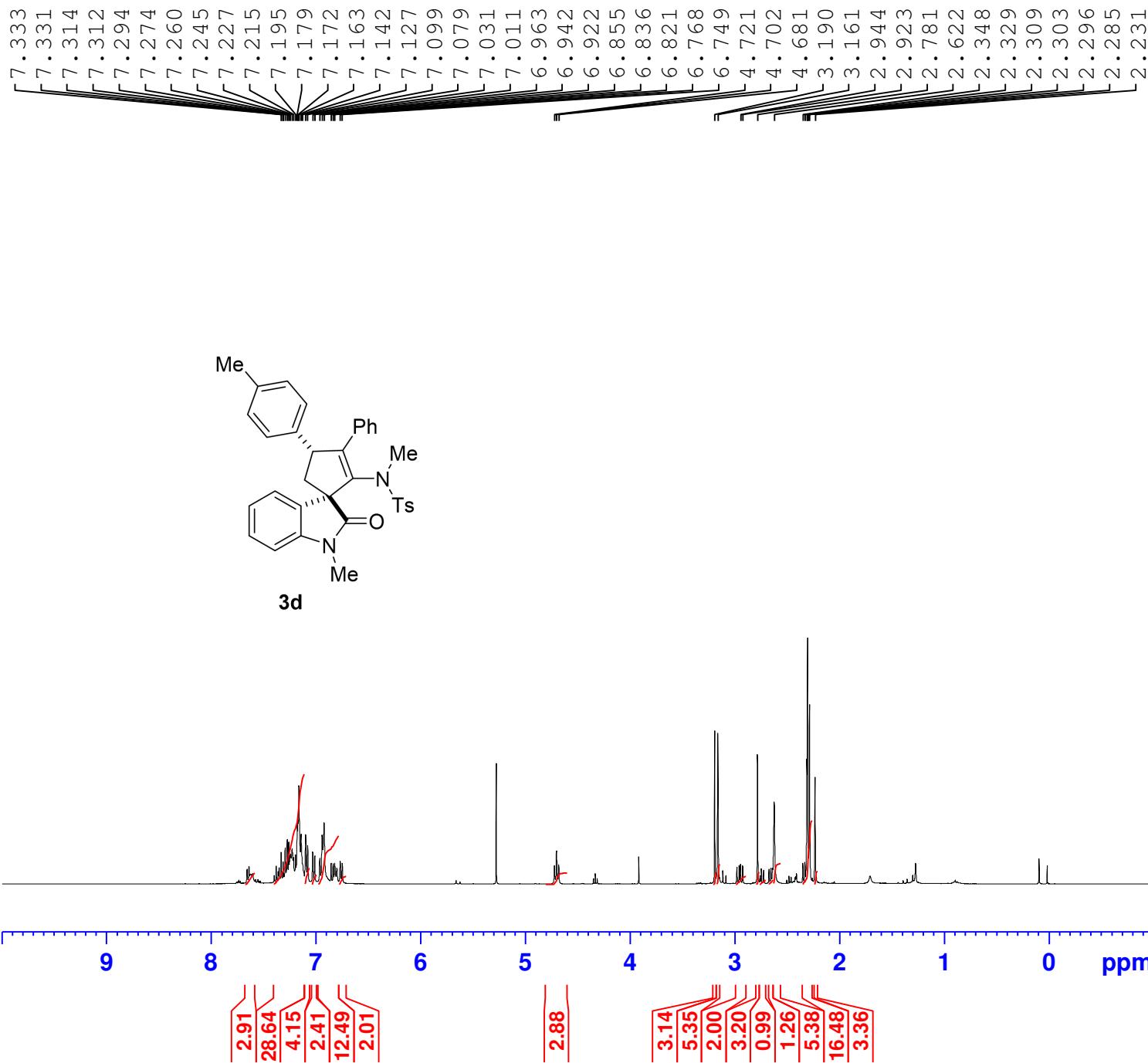
F2 - Processing parameters  
SI 32768  
SF 400.1700072 MHz  
WDW no  
SSB 0  
LB 0 Hz  
GB 0  
PC 1.00



Current Data Parameters  
 NAME cqh-82  
 EXPNO 2020102902  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20201030  
 Time 12.44 h  
 INSTRUM spect  
 PROBHD Z108618\_0256 (   
 PULPROG zgpg30  
 TD 65536  
 SOLVENT CDC13  
 NS 250  
 DS 4  
 SWH 24038.461 Hz  
 FIDRES 0.733596 Hz  
 AQ 1.3631488 sec  
 RG 2050  
 DW 20.800 usec  
 DE 6.50 usec  
 TE 298.1 K  
 D1 2.00000000 sec  
 D11 0.03000000 sec  
 TD0 1  
 SFO1 100.6328888 MHz  
 NUC1 13C  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 50.00000000 W  
 SFO2 400.1716007 MHz  
 NUC2 1H  
 CPDPRG[2] waltz65  
 PCPD2 90.00 usec  
 PLW2 15.00000000 W  
 PLW12 0.18519001 W  
 PLW13 0.09314700 W

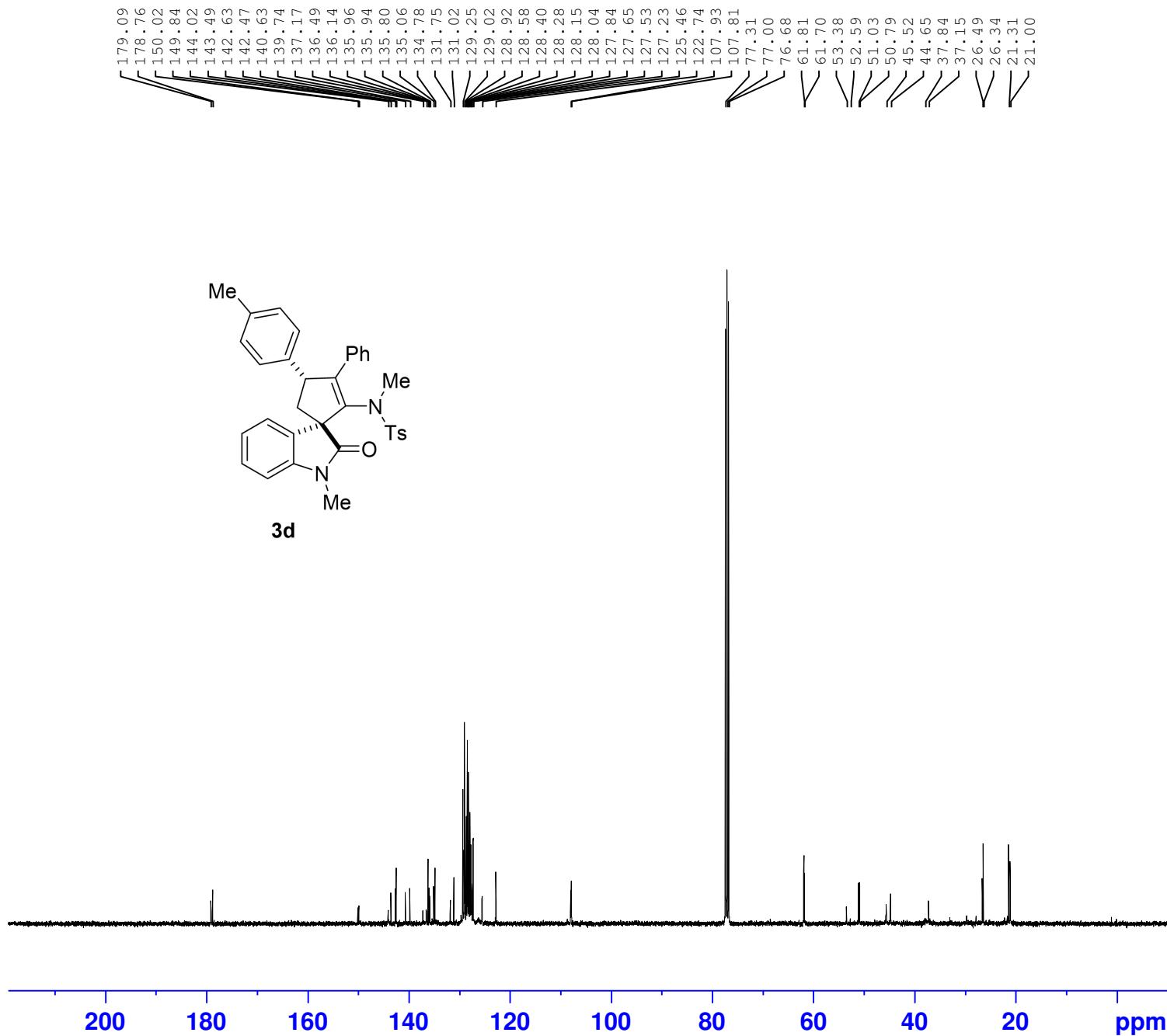
F2 - Processing parameters  
 SI 32768  
 SF 100.6228506 MHz  
 WDW EM  
 SSB 0  
 LB 1.00 Hz  
 GB 0  
 PC 1.40



Current Data Parameters  
 NAME cqh-93  
 EXPNO 2020111001  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20201111  
 Time 5.15 h  
 INSTRUM spect  
 PROBHD Z108618\_0256 (zg30)  
 PULPROG zg30  
 TD 65536  
 SOLVENT CDCl<sub>3</sub>  
 NS 16  
 DS 2  
 SWH 8223.685 Hz  
 FIDRES 0.250967 Hz  
 AQ 3.9845889 sec  
 RG 114  
 DW 60.800 usec  
 DE 6.50 usec  
 TE 297.0 K  
 D1 1.00000000 sec  
 TD0 1  
 SFO1 400.1724712 MHz  
 NUC1 1H  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 15.00000000 W

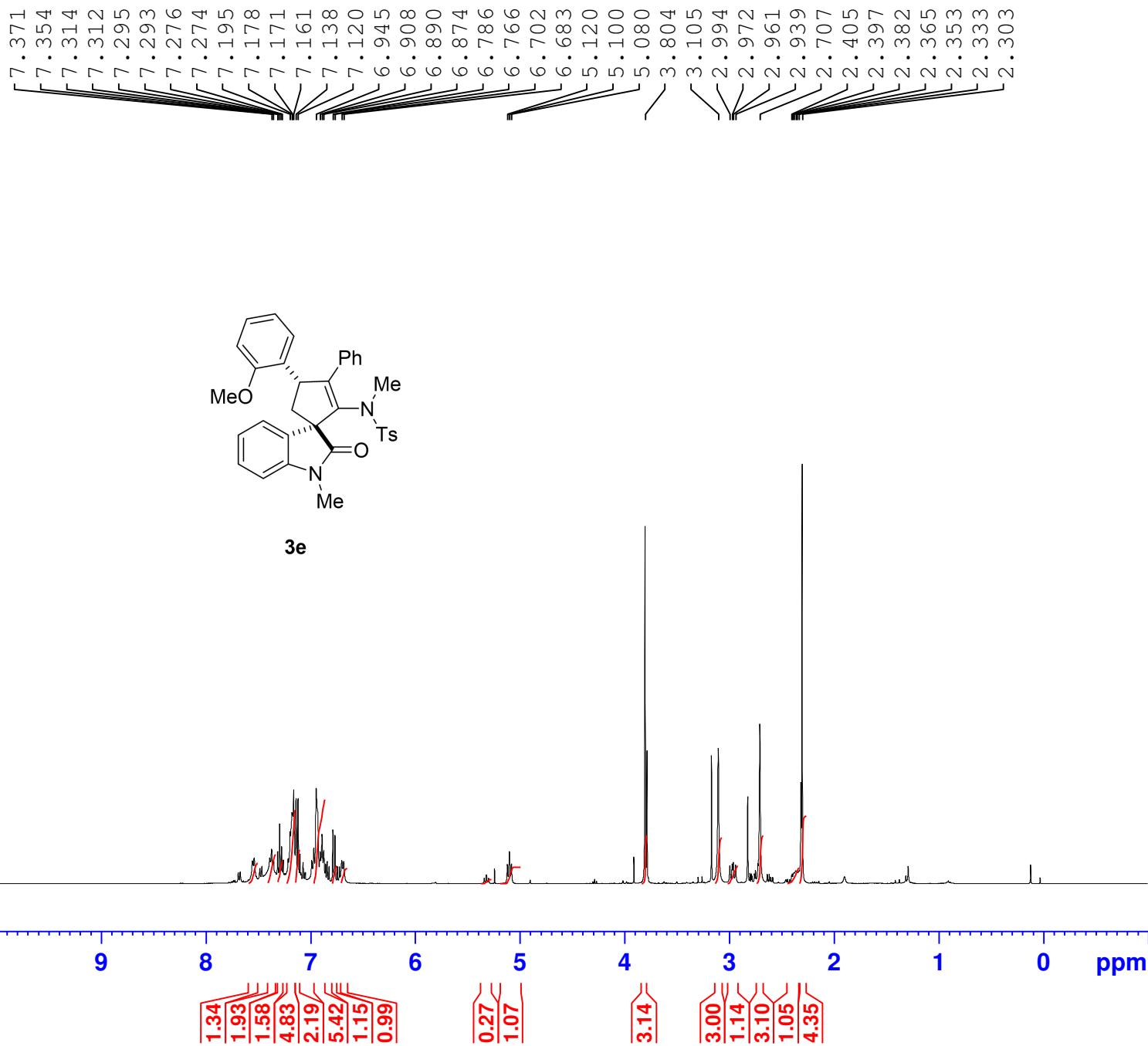
F2 - Processing parameters  
 SI 32768  
 SF 400.1700081 MHz  
 WDW no  
 SSB 0  
 LB 0 Hz  
 GB 0  
 PC 1.00



Current Data Parameters  
 NAME cgh-93  
 EXPNO 2020111002  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20201111  
 Time 8.19 h  
 INSTRUM spect  
 PROBHD Z108618\_0256 (zgpg30)  
 PULPROG 65536  
 TD 800  
 SOLVENT CDC13  
 DS 4  
 SWH 24038.461 Hz  
 FIDRES 0.733596 Hz  
 AQ 1.3631488 sec  
 RG 2050  
 DW 20.800 usec  
 DE 6.50 usec  
 TE 297.0 K  
 D1 2.00000000 sec  
 D11 0.03000000 sec  
 TD0 1  
 SFO1 100.6328888 MHz  
 NUC1 <sup>13</sup>C  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 50.00000000 W  
 SFO2 400.1716007 MHz  
 NUC2 1H  
 CPDPRG[2] waltz65  
 PCPD2 90.00 usec  
 PLW2 15.00000000 W  
 PLW12 0.18519001 W  
 PLW13 0.09314700 W

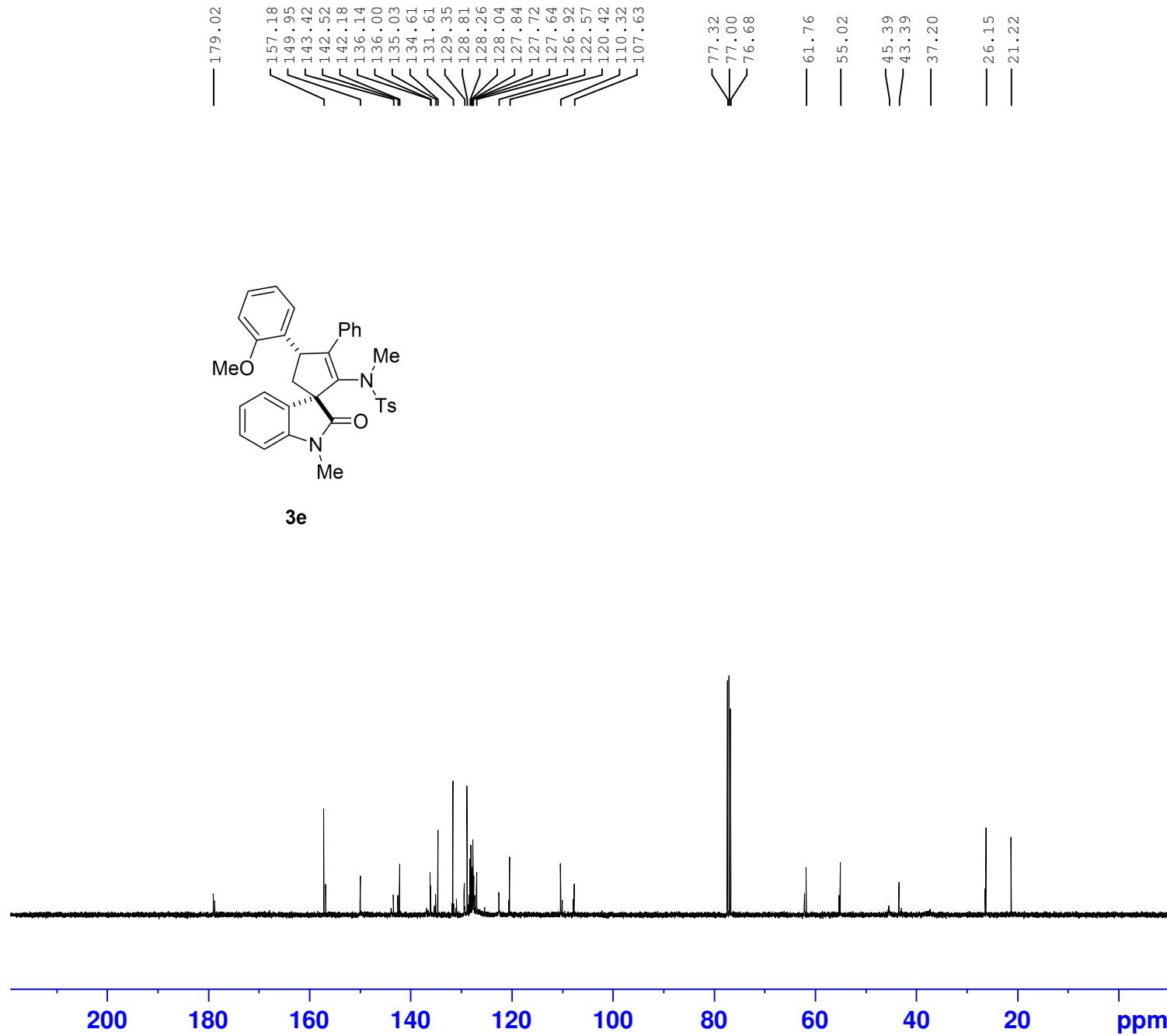
F2 - Processing parameters  
 SI 32768  
 SF 100.6228366 MHz  
 WDW EM  
 SSB 0  
 LB 1.00 Hz  
 GB 0  
 PC 1.40



Current Data Parameters  
NAME cgh-95  
EXPNO 2020111002  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20201111  
Time 5.20 h  
INSTRUM spect  
PROBHD Z108618\_0256 (zg30)  
PULPROG zg30  
TD 65536  
SOLVENT CDCl<sub>3</sub>  
NS 16  
DS 2  
SWH 8223.685 Hz  
FIDRES 0.250967 Hz  
AQ 3.9845889 sec  
RG 40.3  
DW 60.800 usec  
DE 6.50 usec  
TE 296.9 K  
D1 1.00000000 sec  
TD0 1  
SFO1 400.1724712 MHz  
NUC1 1H  
P0 3.33 usec  
P1 10.00 usec  
PLW1 15.00000000 W

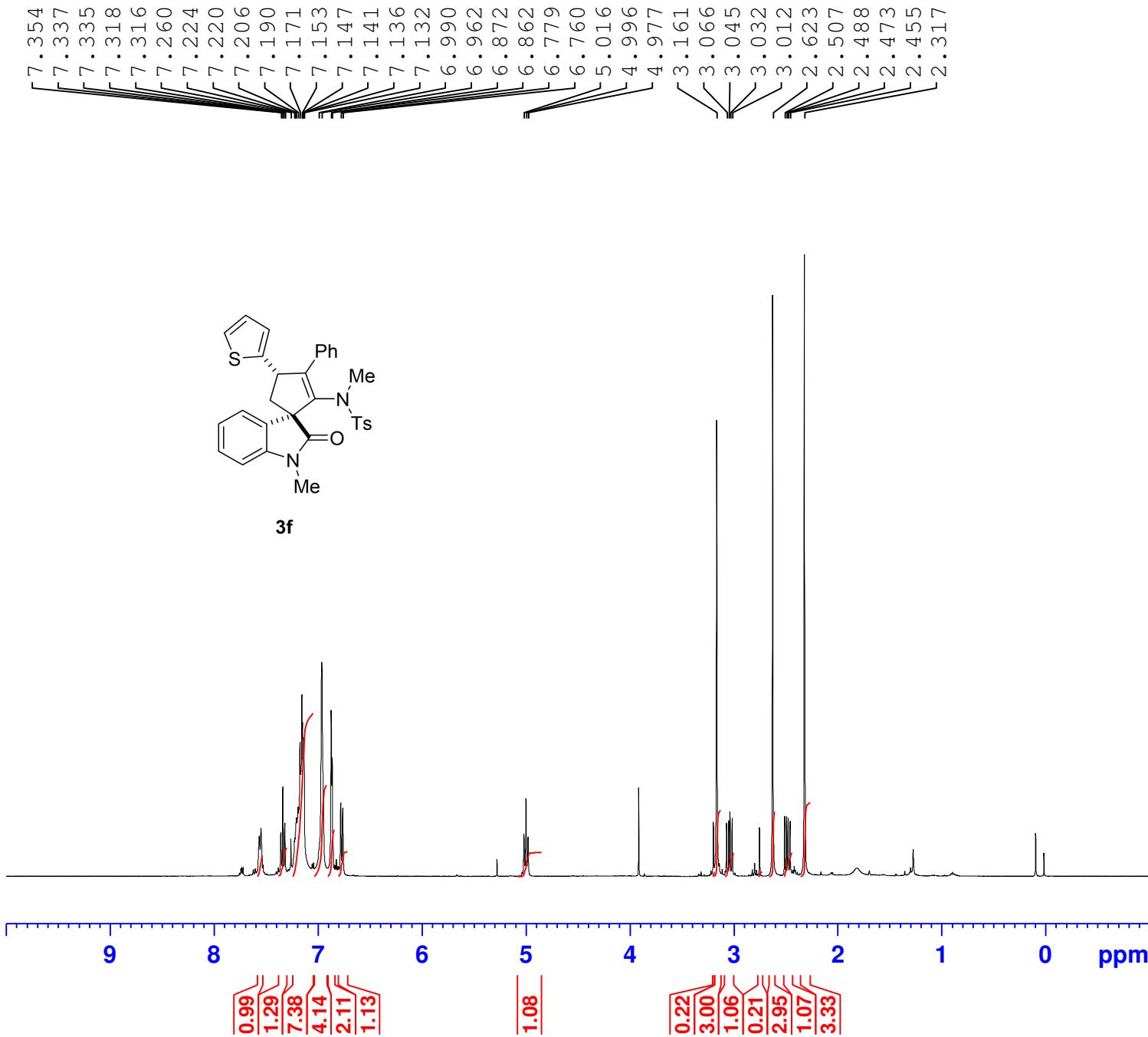
F2 - Processing parameters  
SI 32768  
SF 400.1700072 MHz  
WDW no  
SSB 0  
LB 0 Hz  
GB 0  
PC 1.00



Current Data Parameters  
NAME cqh-95  
EXPNO 2020111003  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20201111  
Time 8.44 h  
INSTRUM spect  
PROBHD Z108618\_0256 (zgpg30  
PULPROG zgpg30  
TD 65536  
SOLVENT CDCl<sub>3</sub>  
NS 351  
DS 4  
SWH 24038.461 Hz  
FIDRES 0.733596 Hz  
AQ 1.3631488 sec  
RG 2050  
DW 20.800 usec  
DE 6.50 usec  
TE 296.9 K  
D1 2.00000000 sec  
D11 0.03000000 sec  
TD0 1  
SFO1 100.6328888 MHz  
NUC1 13C  
P0 3.33 usec  
P1 10.00 usec  
PLW1 50.00000000 W  
SFO2 400.1716007 MHz  
NUC2 1H  
CPDPRG[2] waltz65  
PCPD2 90.00 usec  
PLW2 15.00000000 W  
PLW12 0.18519001 W  
PLW13 0.09314700 W

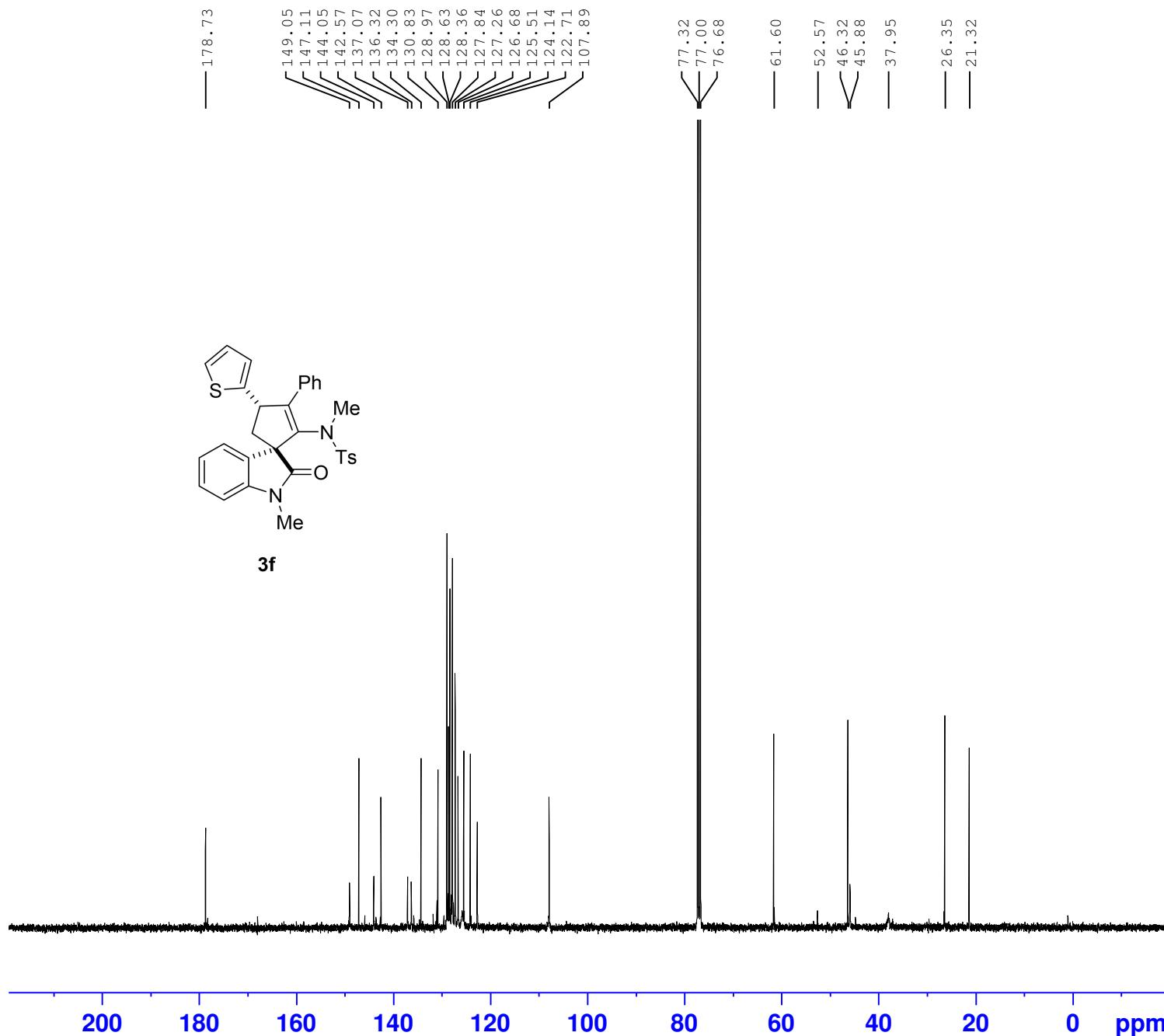
F2 - Processing parameters  
SI 32768  
SF 100.6228476 MHz  
WDW no  
SSB 0  
LB 0 Hz  
GB 0  
PC 1.40



Current Data Parameters  
 NAME cqh-71-1  
 EXPNO 2020092301  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20200924  
 Time 23.39 h  
 INSTRUM spect  
 PROBHD Z108618\_0256 (br)  
 PULPROG zg30  
 TD 65536  
 SOLVENT CDCl<sub>3</sub>  
 NS 16  
 DS 2  
 SWH 8223.685 Hz  
 FIDRES 0.250967 Hz  
 AQ 3.9845889 sec  
 RG 144  
 DW 60.800 usec  
 DE 6.50 usec  
 TE 297.5 K  
 D1 1.00000000 sec  
 TDO 1  
 SFO1 400.1724712 MHz  
 NUC1 1H  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 15.00000000 W

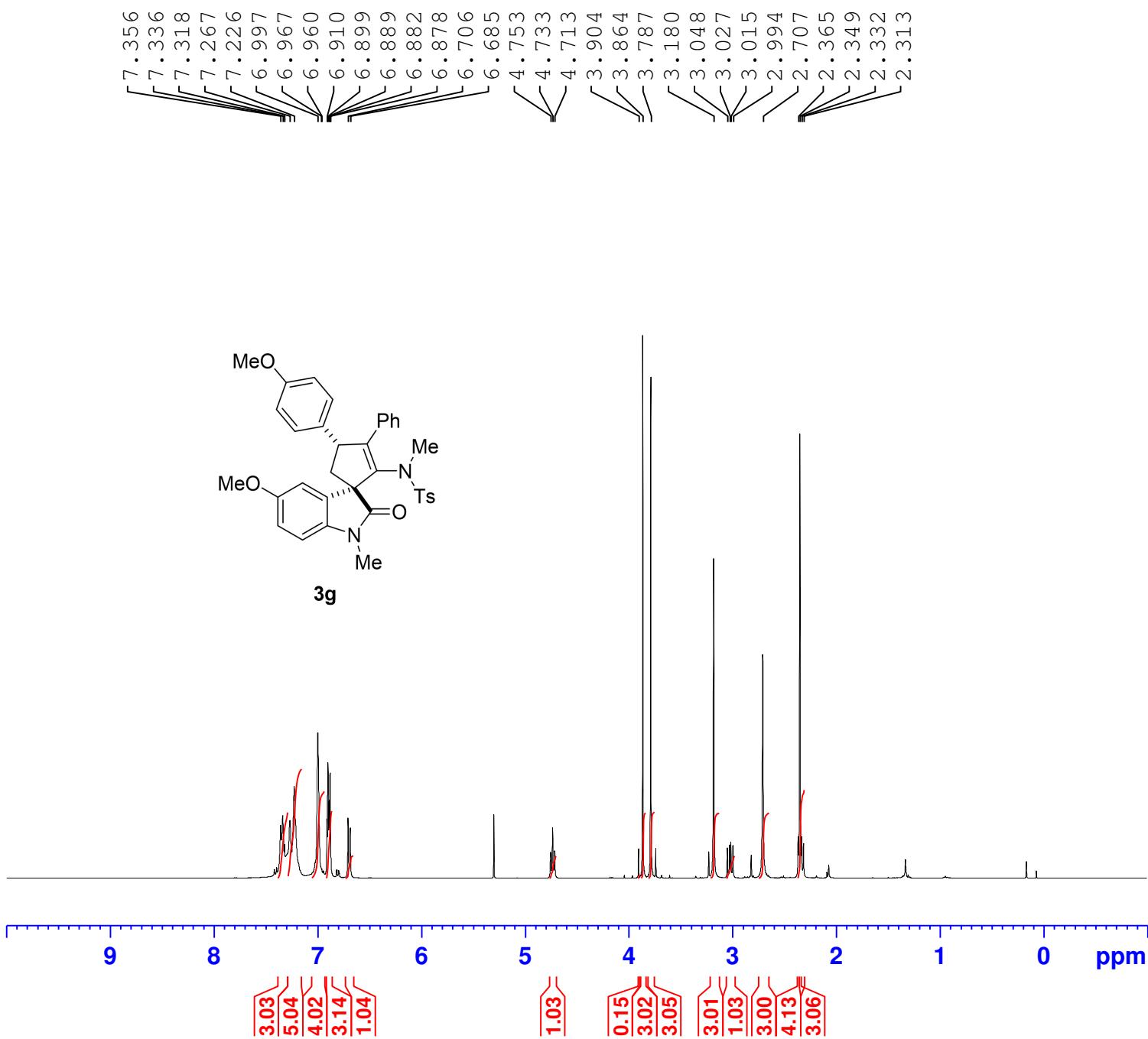
F2 - Processing parameters  
 SI 32768  
 SF 400.1700087 MHz  
 WDW no  
 SSB 0  
 LB 0 Hz  
 GB 0  
 PC 1.00



Current Data Parameters  
 NAME cgh-71-1  
 EXPNO 2020092302  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20200925  
 Time 0.54 h  
 INSTRUM spect  
 PROBHD Z108618\_0256 (zgpg30  
 PULPROG 65536  
 TD 1  
 SOLVENT CDCl3  
 NS 661  
 DS 4  
 SWH 24038.461 Hz  
 FIDRES 0.733596 Hz  
 AQ 1.3631488 sec  
 RG 2050  
 DW 20.800 usec  
 DE 6.50 usec  
 TE 298.0 K  
 D1 2.00000000 sec  
 D11 0.03000000 sec  
 TD0 1  
 SFO1 100.6328888 MHz  
 NUC1 13C  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 50.00000000 W  
 SFO2 400.1716007 MHz  
 NUC2 1H  
 CPDPRG[2] waltz65  
 PCPD2 90.00 usec  
 PLW2 15.00000000 W  
 PLW12 0.18519001 W  
 PLW13 0.09314700 W

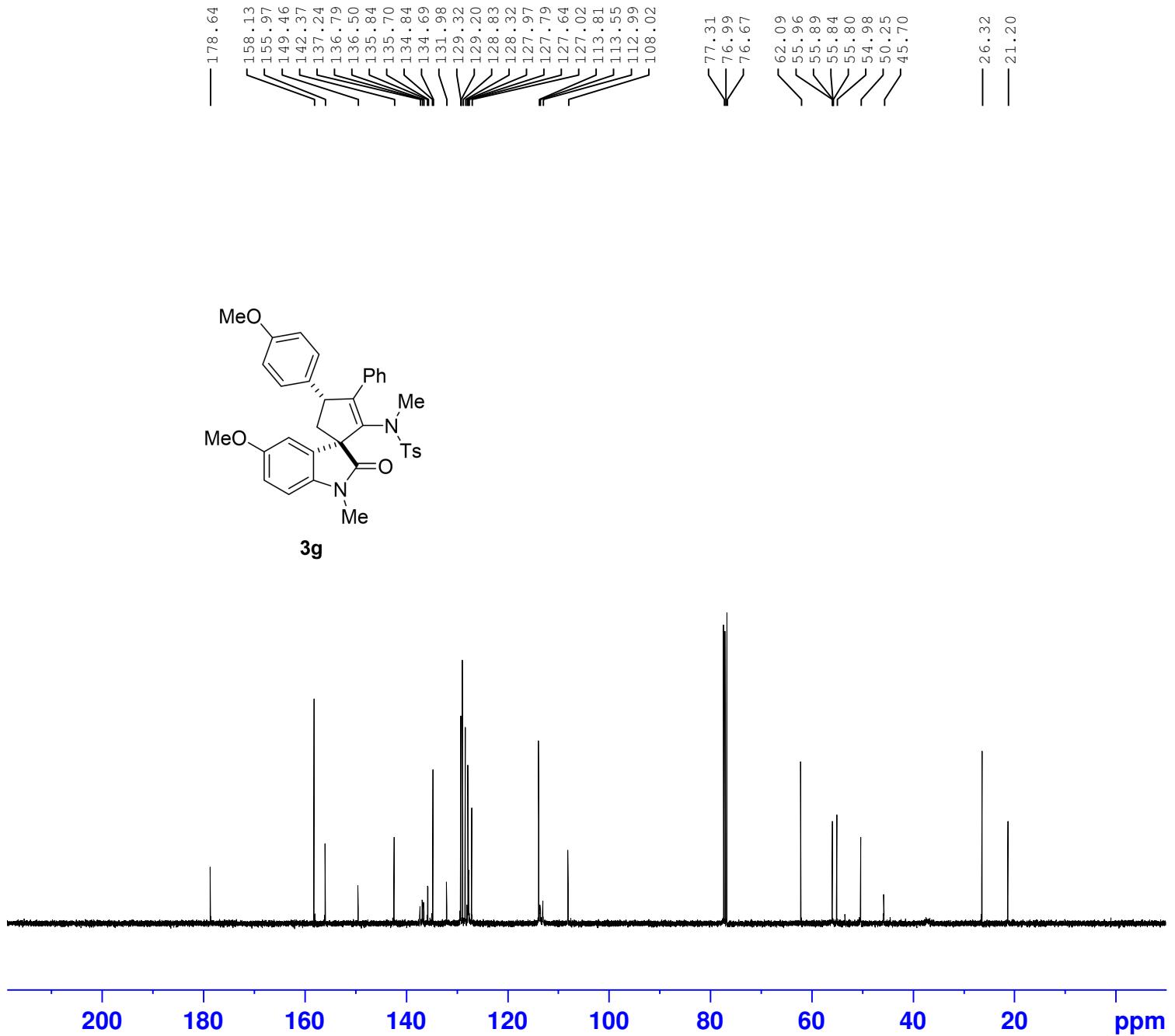
F2 - Processing parameters  
 SI 32768  
 SF 100.6228381 MHz  
 WDW EM  
 SSB 0  
 LB 1.00 Hz  
 GB 0  
 PC 1.40

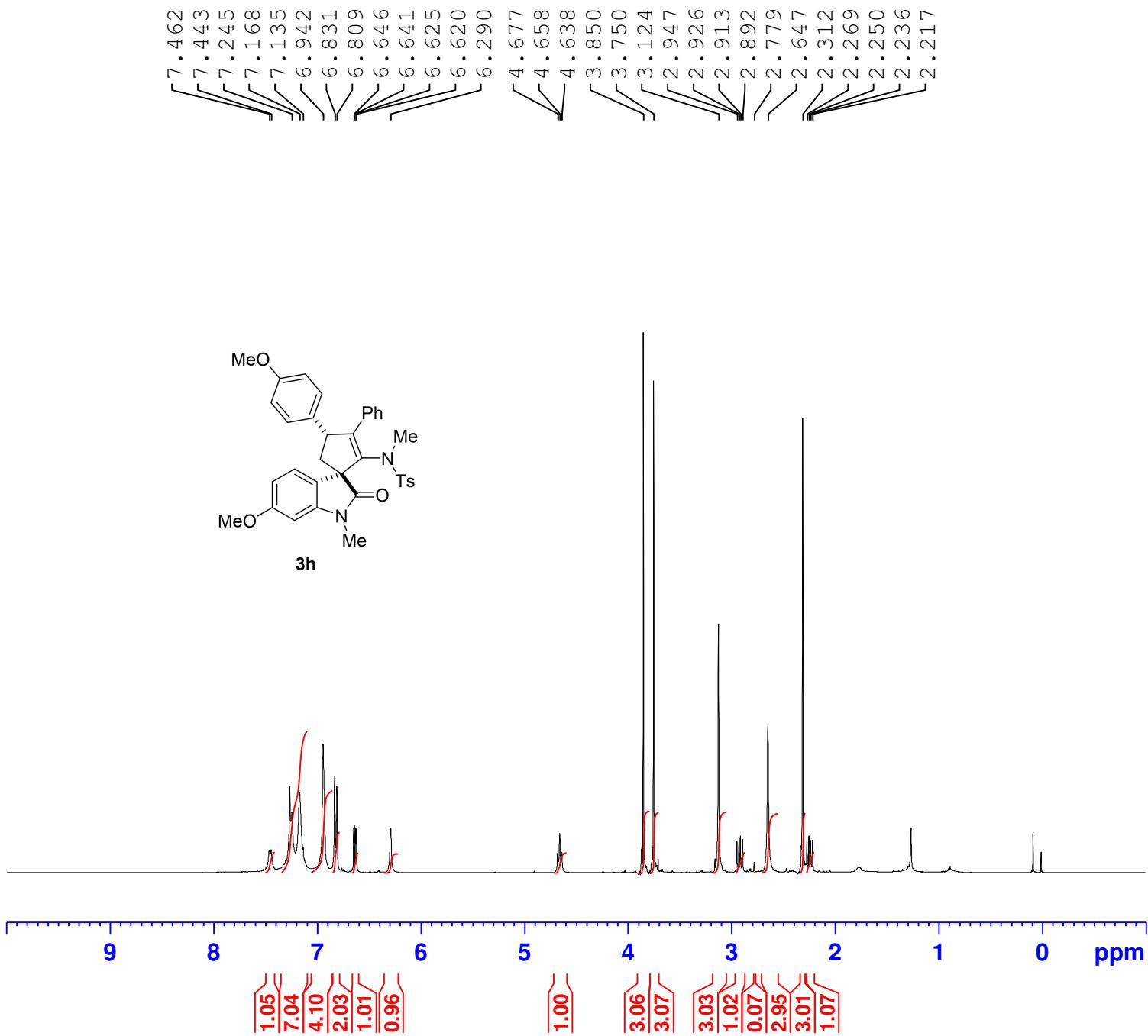


Current Data Parameters  
NAME cgh-144  
EXPNO 2020122301  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20201224  
Time 9.08 h  
INSTRUM spect  
PROBHD z108618\_0256 (   
PULPROG zg30  
TD 65536  
SOLVENT CDCl<sub>3</sub>  
NS 16  
DS 2  
SWH 8223.685 Hz  
FIDRES 0.250967 Hz  
AQ 3.9845889 sec  
RG 32  
DW 60.800 usec  
DE 6.50 usec  
TE 294.0 K  
D1 1.00000000 sec  
TD0 1  
SFO1 400.1724712 MHz  
NUC1 1H  
P0 3.33 usec  
P1 10.00 usec  
PLW1 15.00000000 W

F2 - Processing parameters  
SI 32768  
SF 400.1699841 MHz  
WDW no  
SSB 0  
LB 0 Hz  
GB 0  
PC 1.00

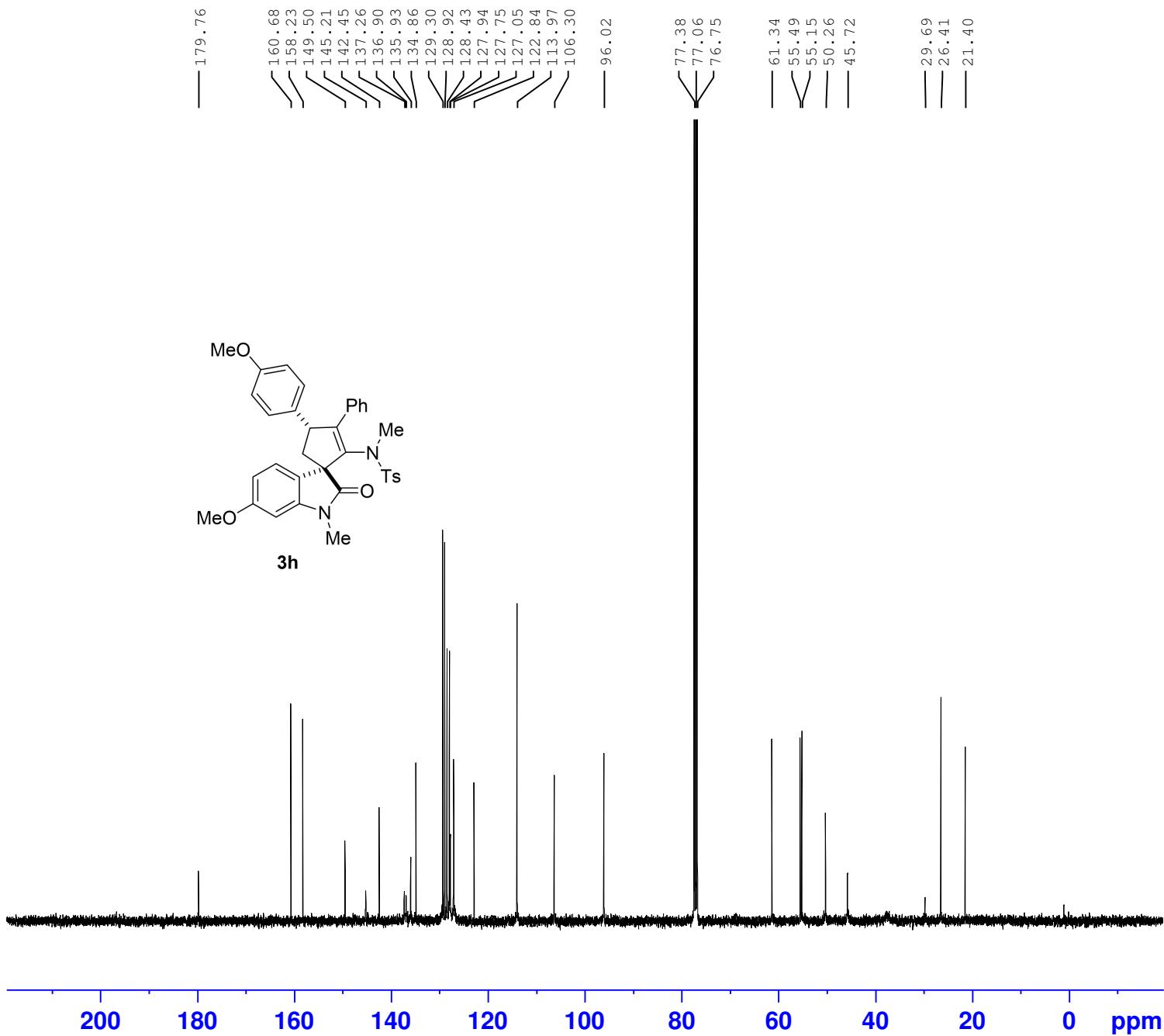




Current Data Parameters  
NAME cqh-163  
EXPNO 2021030901  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20210310  
Time 6.30 h  
INSTRUM spect  
PROBHD Z108618\_0256 (   
PULPROG zg30  
TD 65536  
SOLVENT CDCl<sub>3</sub>  
NS 16  
DS 2  
SWH 8223.685 Hz  
FIDRES 0.250967 Hz  
AQ 3.9845889 sec  
RG 90.5  
DW 60.800 usec  
DE 6.50 usec  
TE 294.6 K  
D1 1.0000000 sec  
TD0 1  
SFO1 400.1724712 MHz  
NUC1 1H  
P0 3.33 usec  
P1 10.00 usec  
PLW1 15.00000000 W

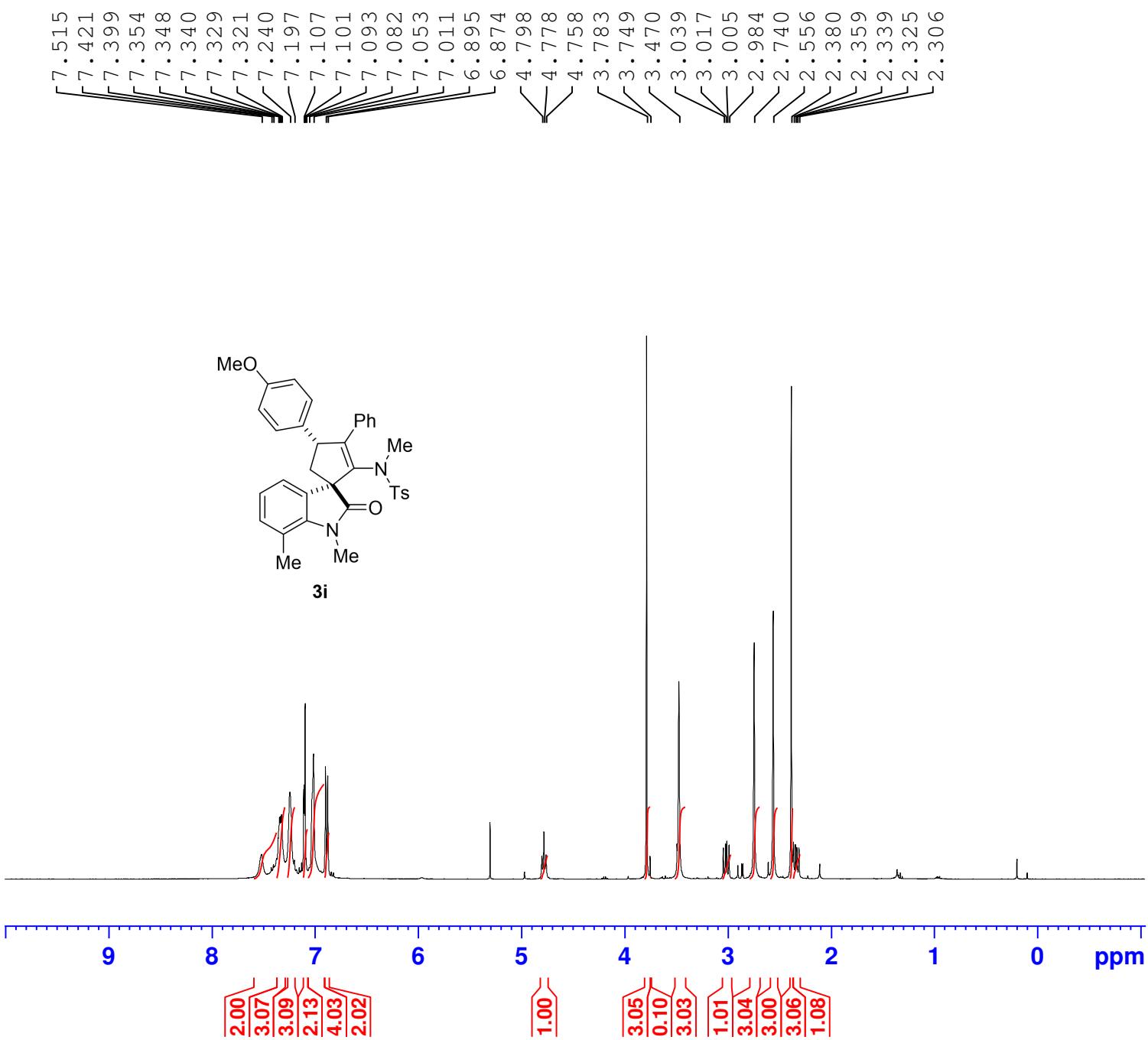
F2 - Processing parameters  
SI 32768  
SF 400.1700072 MHz  
WDW EM  
SSB 0  
LB 0.30 Hz  
GB 0  
PC 1.00

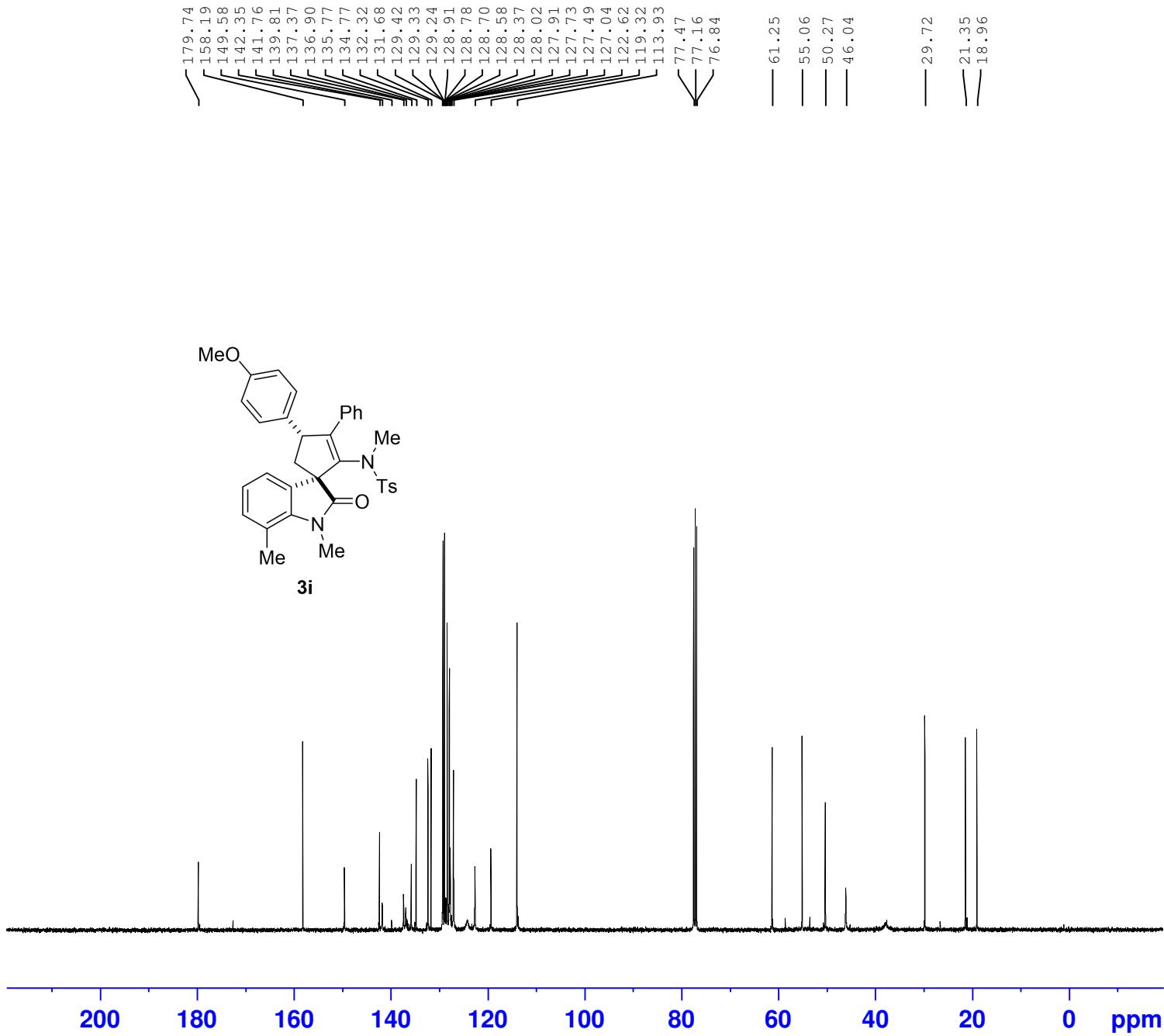


Current Data Parameters  
 NAME cgh-163  
 EXPNO 2021030902  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20210310  
 Time 11.35 h  
 INSTRUM spect  
 PROBHD z108618\_0256 (   
 PULPROG zgpg30  
 TD 65536  
 SOLVENT CDCl3  
 NS 506  
 DS 4  
 SWH 24038.461 Hz  
 FIDRES 0.733596 Hz  
 AQ 1.3631488 sec  
 RG 2050  
 DW 20.800 usec  
 DE 6.50 usec  
 TE 294.8 K  
 D1 2.00000000 sec  
 D11 0.03000000 sec  
 TDO 1  
 SFO1 100.6328888 MHz  
 NUC1 13C  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 50.00000000 W  
 SFO2 400.1716007 MHz  
 NUC2 1H  
 CPDPRG[2] waltz65  
 PCPD2 90.00 usec  
 PLW2 15.00000000 W  
 PLW12 0.18519001 W  
 PLW13 0.09314700 W

F2 - Processing parameters  
 SI 32768  
 SF 100.6228300 MHz  
 WDW EM  
 SSB 0  
 LB 1.00 Hz  
 GB 0  
 PC 1.40

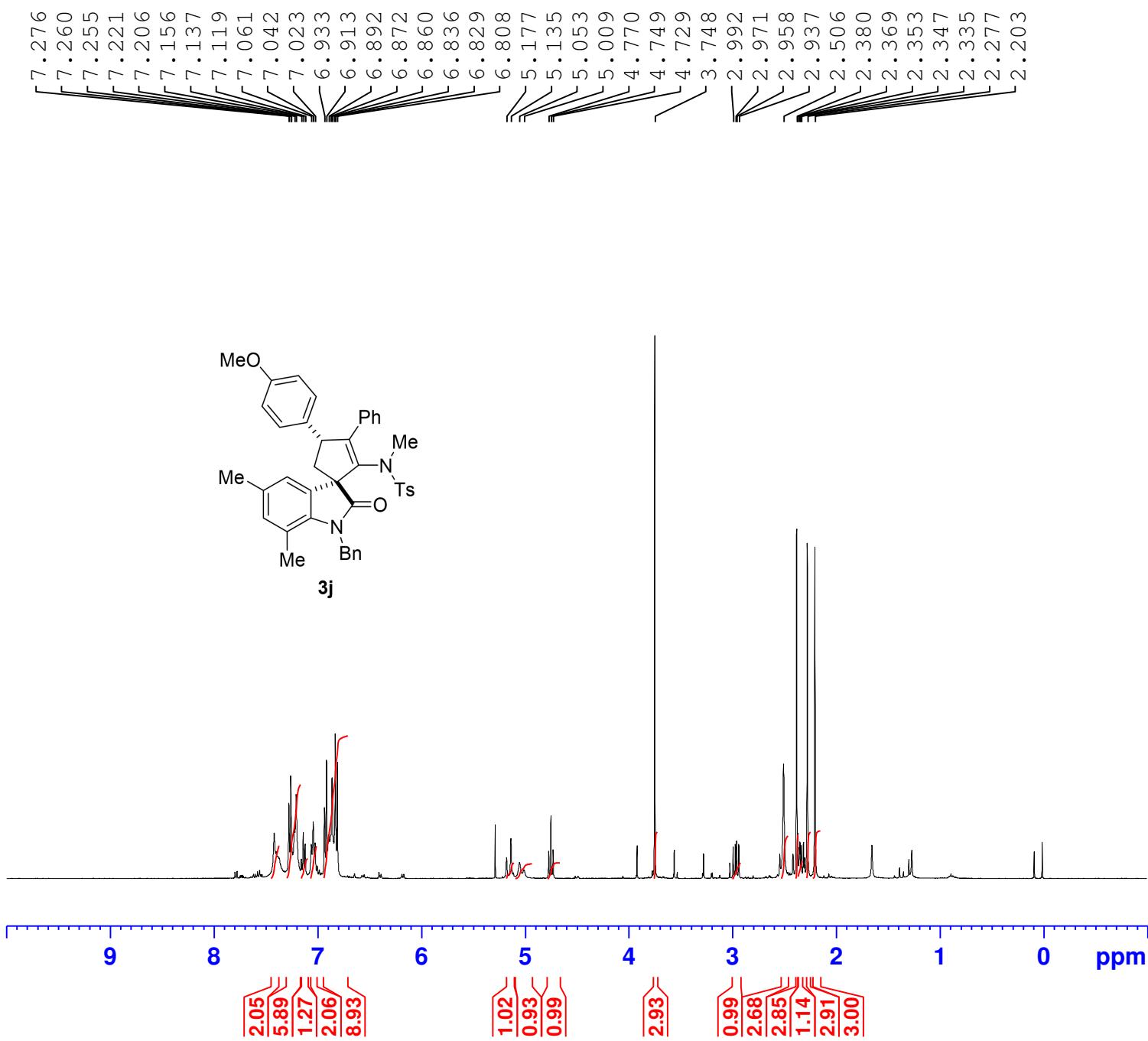


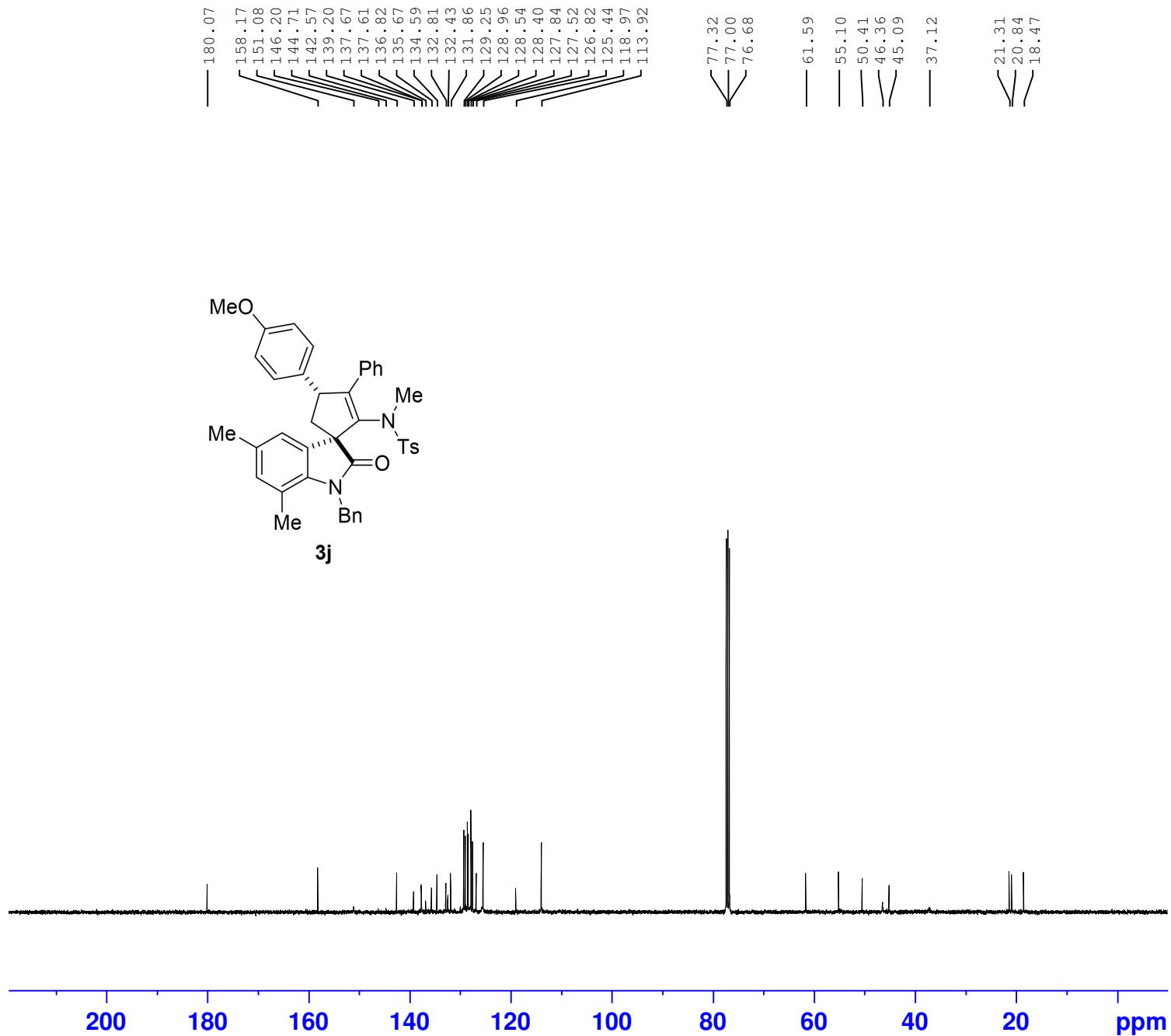


Current Data Parameters  
 NAME cgh-164  
 EXPNO 2021010802  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20210109  
 Time 6.47 h  
 INSTRUM spect  
 PROBHD Z108618\_0256 (   
 PULPROG zgpg30  
 TD 65536  
 SOLVENT CDCl<sub>3</sub>  
 NS 519  
 DS 4  
 SWH 24038.461 Hz  
 FIDRES 0.733596 Hz  
 AQ 1.3631488 sec  
 RG 2050  
 DW 20.800 usec  
 DE 6.50 usec  
 TE 294.2 K  
 D1 2.00000000 sec  
 D11 0.03000000 sec  
 TDO 1  
 SFO1 100.6328888 MHz  
 NUC1 <sup>13</sup>C  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 50.00000000 W  
 SFO2 400.1716007 MHz  
 NUC2 <sup>1</sup>H  
 CPDPRG[2] waltz65  
 PCPD2 90.00 usec  
 PLW2 15.00000000 W  
 PLW12 0.18519001 W  
 PLW13 0.09314700 W

F2 - Processing parameters  
 SI 32768  
 SF 100.6228367 MHz  
 WDW EM  
 SSB 0  
 LB 1.00 Hz  
 GB 0  
 PC 1.40

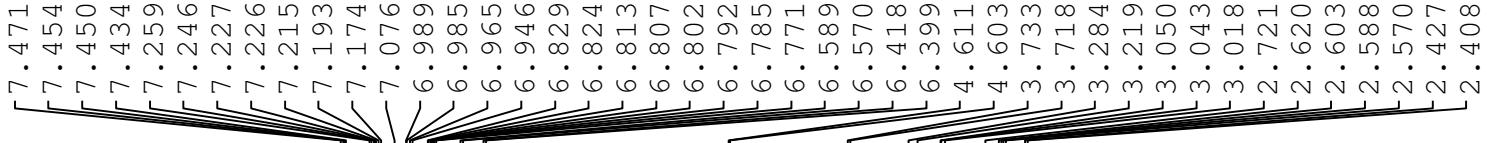




Current Data Parameters  
 NAME zy-8  
 EXPNO 2020111702  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20201118  
 Time 7.34 h  
 INSTRUM spect  
 PROBHD Z108618\_0256 (zgpg30)  
 PULPROG 65536  
 TD 500  
 SOLVENT CDCl<sub>3</sub>  
 NS 4  
 DS 24038.461 Hz  
 SWH 0.733596 Hz  
 FIDRES 1.3631488 sec  
 AQ 2050  
 RG 20.800 usec  
 DE 6.50 usec  
 TE 296.8 K  
 D1 2.000000000 sec  
 D11 0.030000000 sec  
 TD0 1  
 SFO1 100.6328888 MHz  
 NUC1 <sup>13</sup>C  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 50.00000000 W  
 SFO2 400.1716007 MHz  
 NUC2 <sup>1</sup>H  
 CPDPRG[2] waltz65  
 PCPD2 90.00 usec  
 PLW2 15.00000000 W  
 PLW12 0.18519001 W  
 PLW13 0.09314700 W

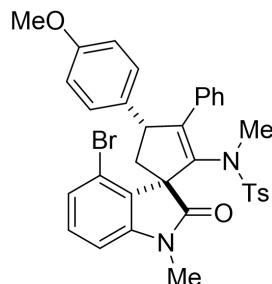
F2 - Processing parameters  
 SI 32768  
 SF 100.6228337 MHz  
 WDW EM  
 SSB 0  
 LB 1.00 Hz  
 GB 0  
 PC 1.40



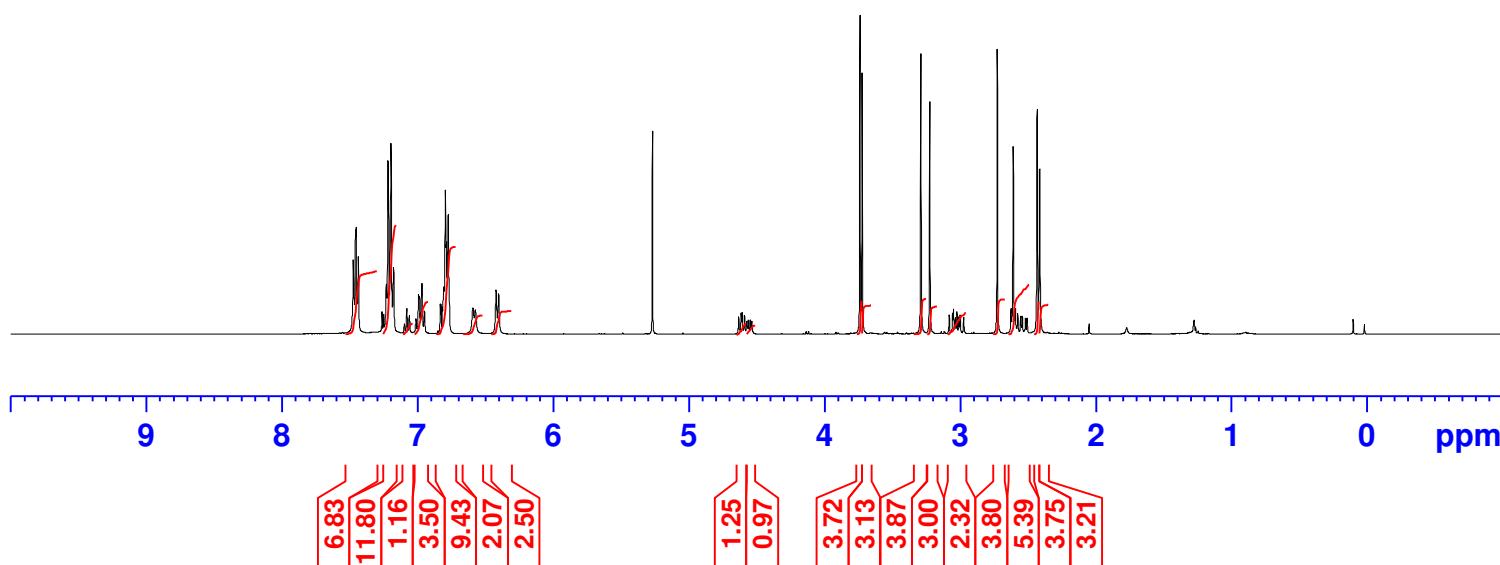
Current Data Parameters  
 NAME cqh-131  
 EXPNO 2020121601  
 PROCNO 1

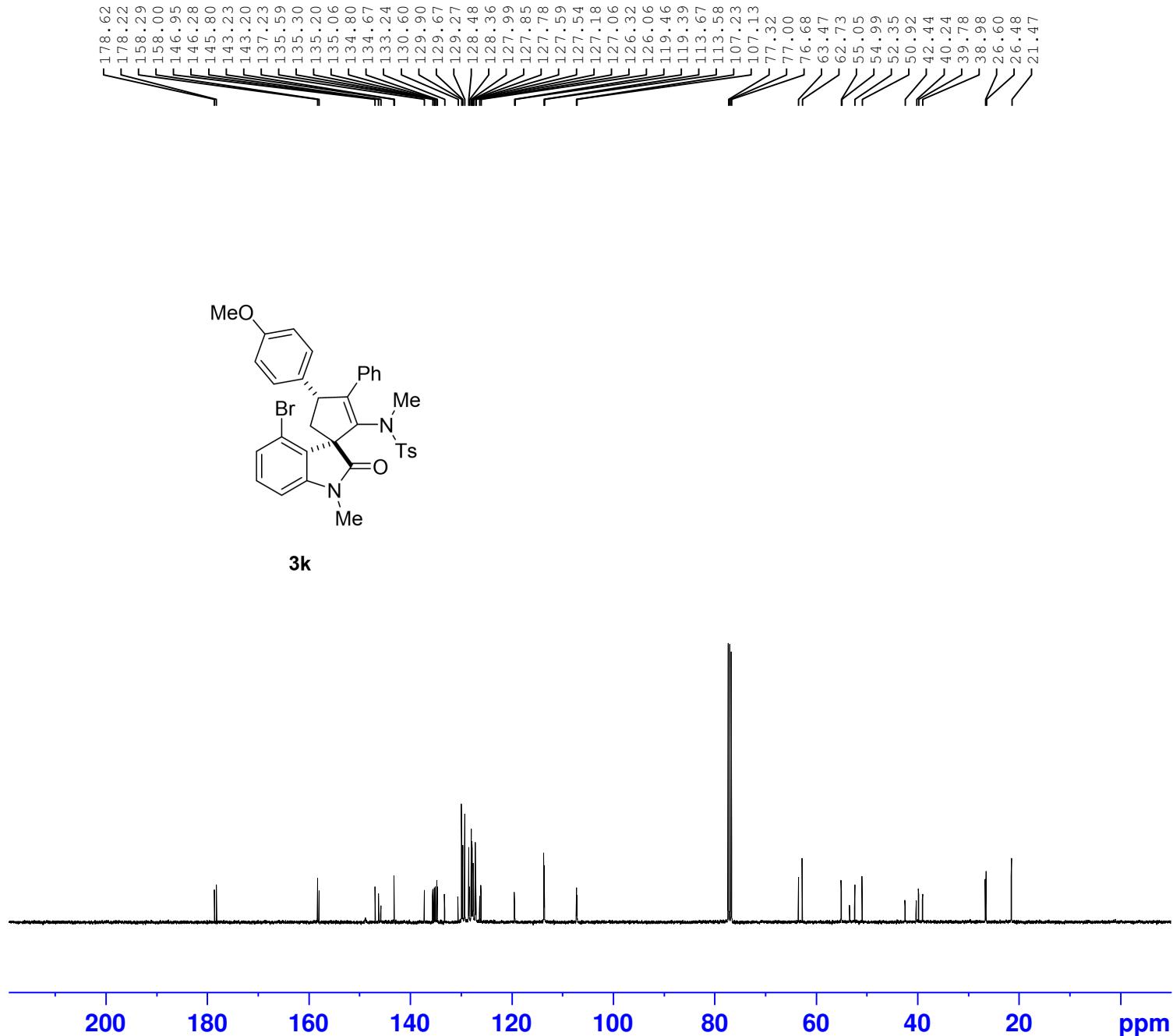
F2 - Acquisition Parameters  
 Date\_ 20201217  
 Time 6.05 h  
 INSTRUM spect  
 PROBHD Z108618\_0256 (   
 PULPROG zg30  
 TD 65536  
 SOLVENT CDCl3  
 NS 16  
 DS 2  
 SWH 8223.685 Hz  
 FIDRES 0.250967 Hz  
 AQ 3.9845889 sec  
 RG 50.8  
 DW 60.800 usec  
 DE 6.50 usec  
 TE 293.8 K  
 D1 1.00000000 sec  
 TD0 1  
 SFO1 400.1724712 MHz  
 NUC1 1H  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 15.00000000 W

F2 - Processing parameters  
 SI 32768  
 SF 400.1700072 MHz  
 WDW no  
 SSB 0  
 LB 0 Hz  
 GB 0  
 PC 1.00



**3k**

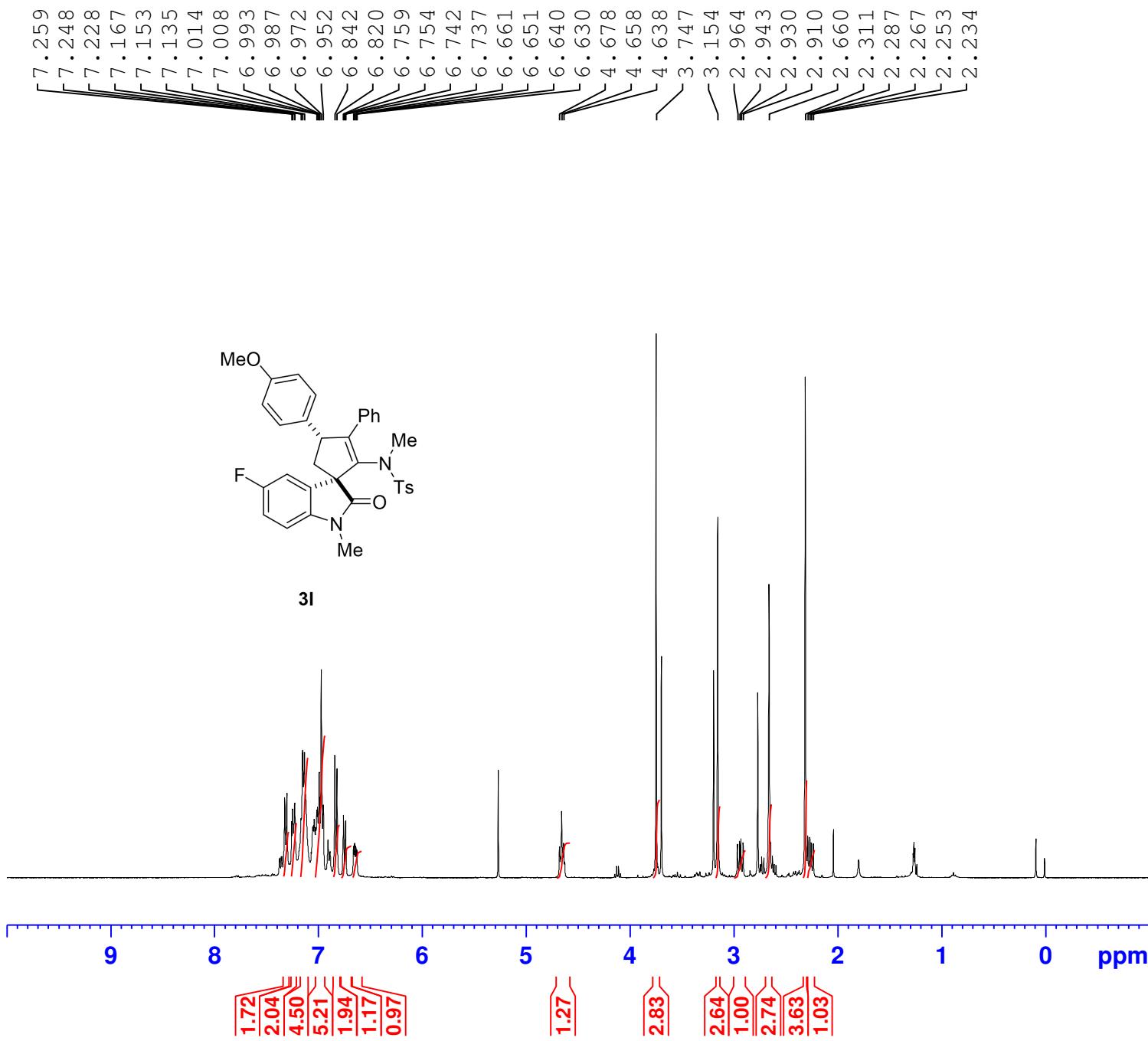


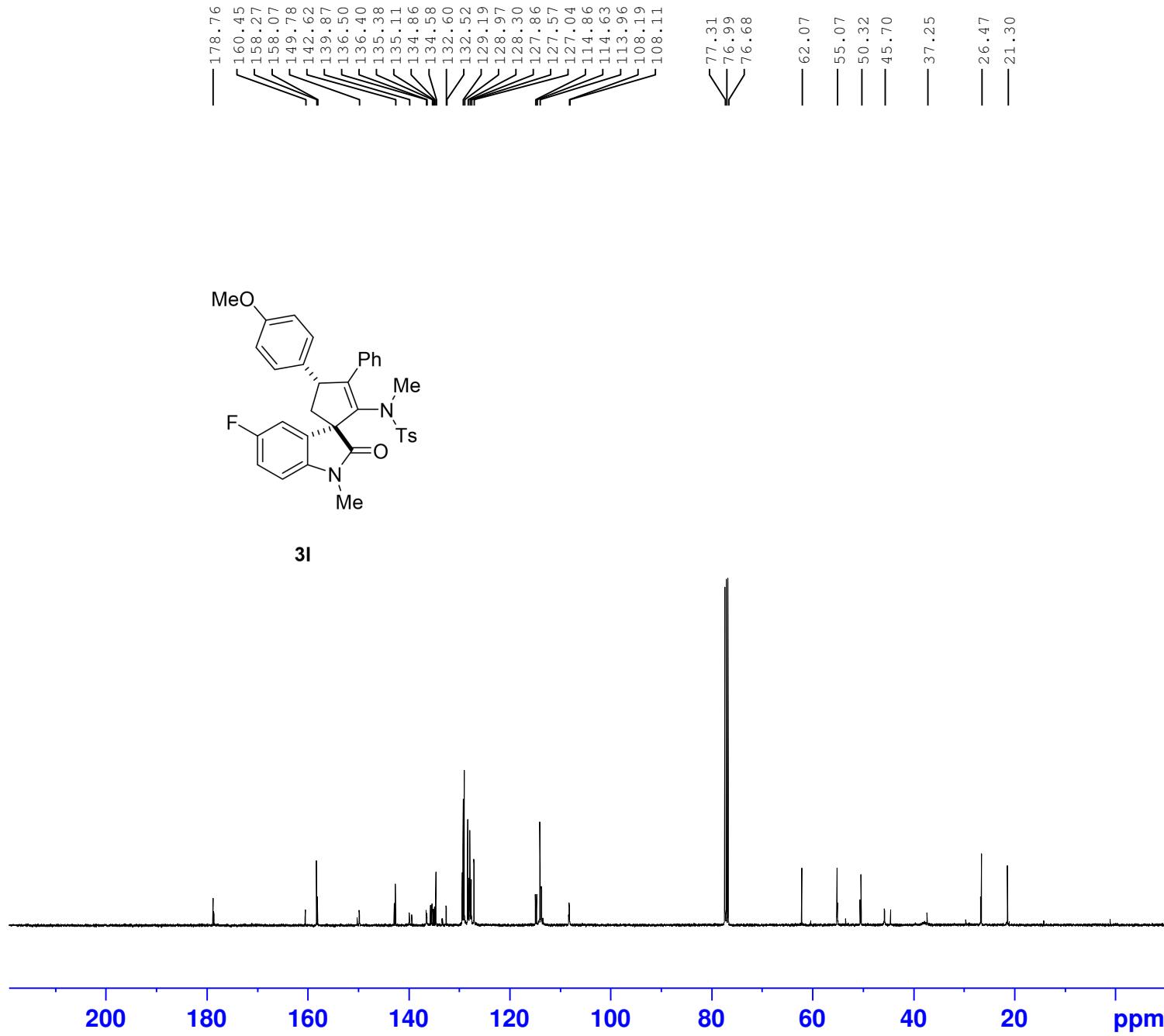


Current Data Parameters  
NAME cgh-131  
EXPNO 2020121602  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20201217  
Time 9.24 h  
INSTRUM spect  
PROBHD Z108618\_0256 (zgpg30)  
TD 65536  
SOLVENT CDCl3  
NS 400  
DS 4  
SWH 24038.461 Hz  
FIDRES 0.733596 Hz  
AQ 1.3631488 sec  
RG 2050  
DW 20.800 usec  
DE 6.50 usec  
TE 293.8 K  
D1 2.00000000 sec  
D11 0.03000000 sec  
TD0 1  
SFO1 100.6328888 MHz  
NUC1 13C  
P0 3.33 usec  
P1 10.00 usec  
PLW1 50.00000000 W  
SFO2 400.1716007 MHz  
NUC2 1H  
CPDPG[2] waltz65  
PCPD2 90.00 usec  
PLW2 15.00000000 W  
PLW12 0.18519001 W  
PLW13 0.09314700 W

F2 - Processing parameters  
SI 32768  
SF 100.6228410 MHz  
WDW EM  
SSB 0  
LB 1.00 Hz  
GB 0  
PC 1.40

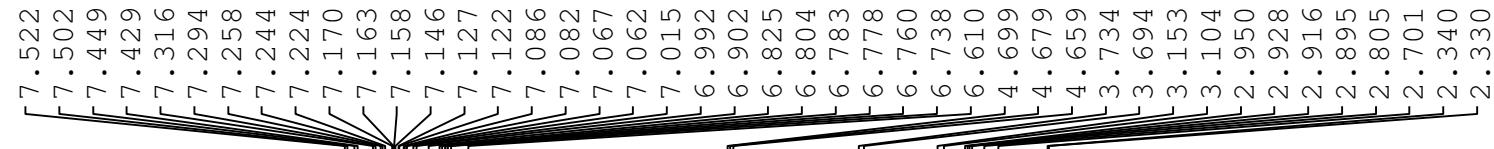




Current Data Parameters  
NAME cgh-130  
EXPNO 2020121602  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20201217  
Time 8.38 h  
INSTRUM spect  
PROBHD Z108618\_0256 (zpg30)  
PULPROG 65536  
TD CDC13  
NS 1000  
DS 4  
SWH 24038.461 Hz  
FIDRES 0.733596 Hz  
AQ 1.3631488 sec  
RG 2050  
DW 20.800 usec  
DE 6.50 usec  
TE 293.7 K  
D1 2.00000000 sec  
D11 0.03000000 sec  
TD0 1  
SFO1 100.6328888 MHz  
NUC1 13C  
P0 3.33 usec  
P1 10.00 usec  
PLW1 50.00000000 W  
SFO2 400.1716007 MHz  
NUC2 1H  
CPDPGRG[2] waltz65  
PCPD2 90.00 usec  
PLW2 15.00000000 W  
PLW12 0.18519001 W  
PLW13 0.09314700 W

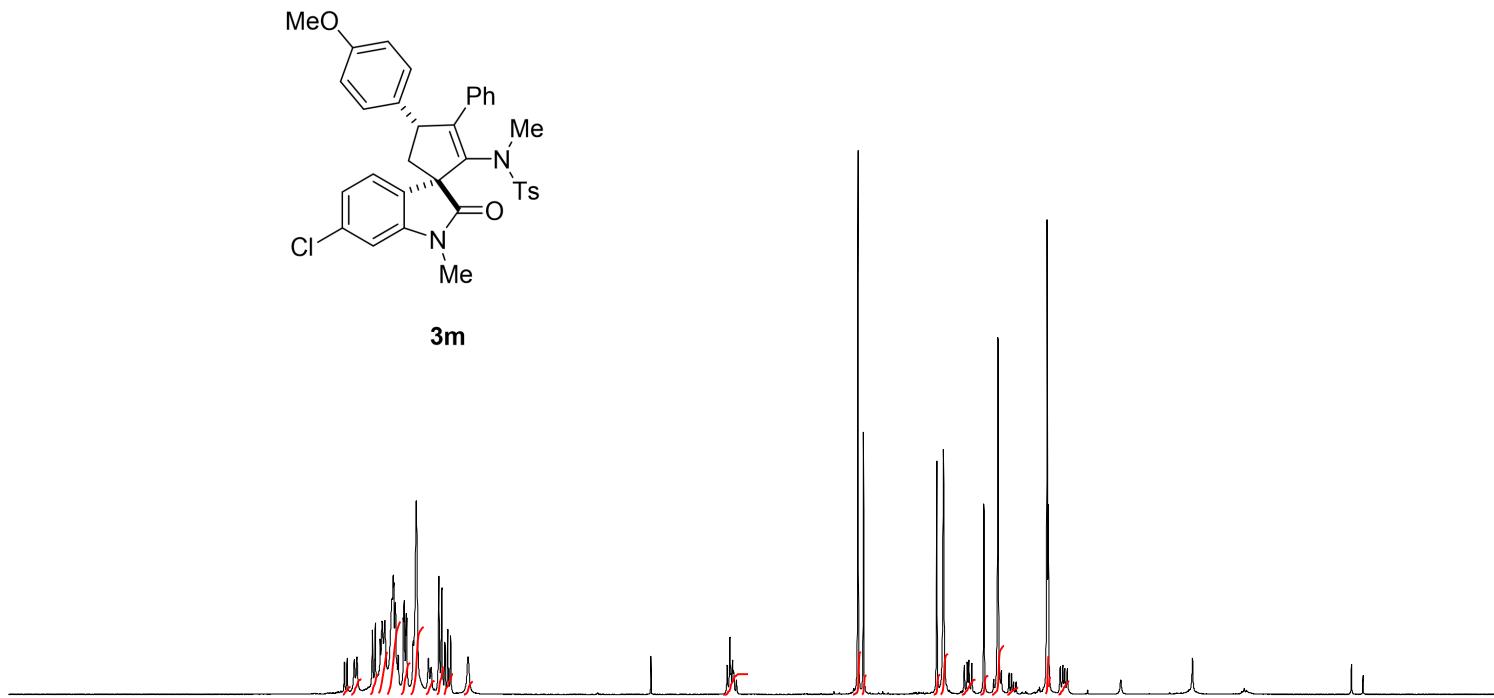
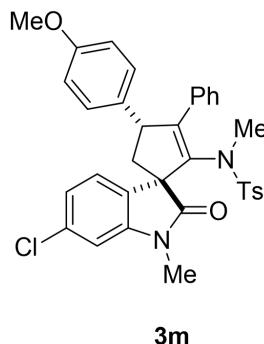
F2 - Processing parameters  
SI 32768  
SF 100.6228403 MHz  
WDW EM  
SSB 0  
LB 1.00 Hz  
GB 0  
PC 1.40

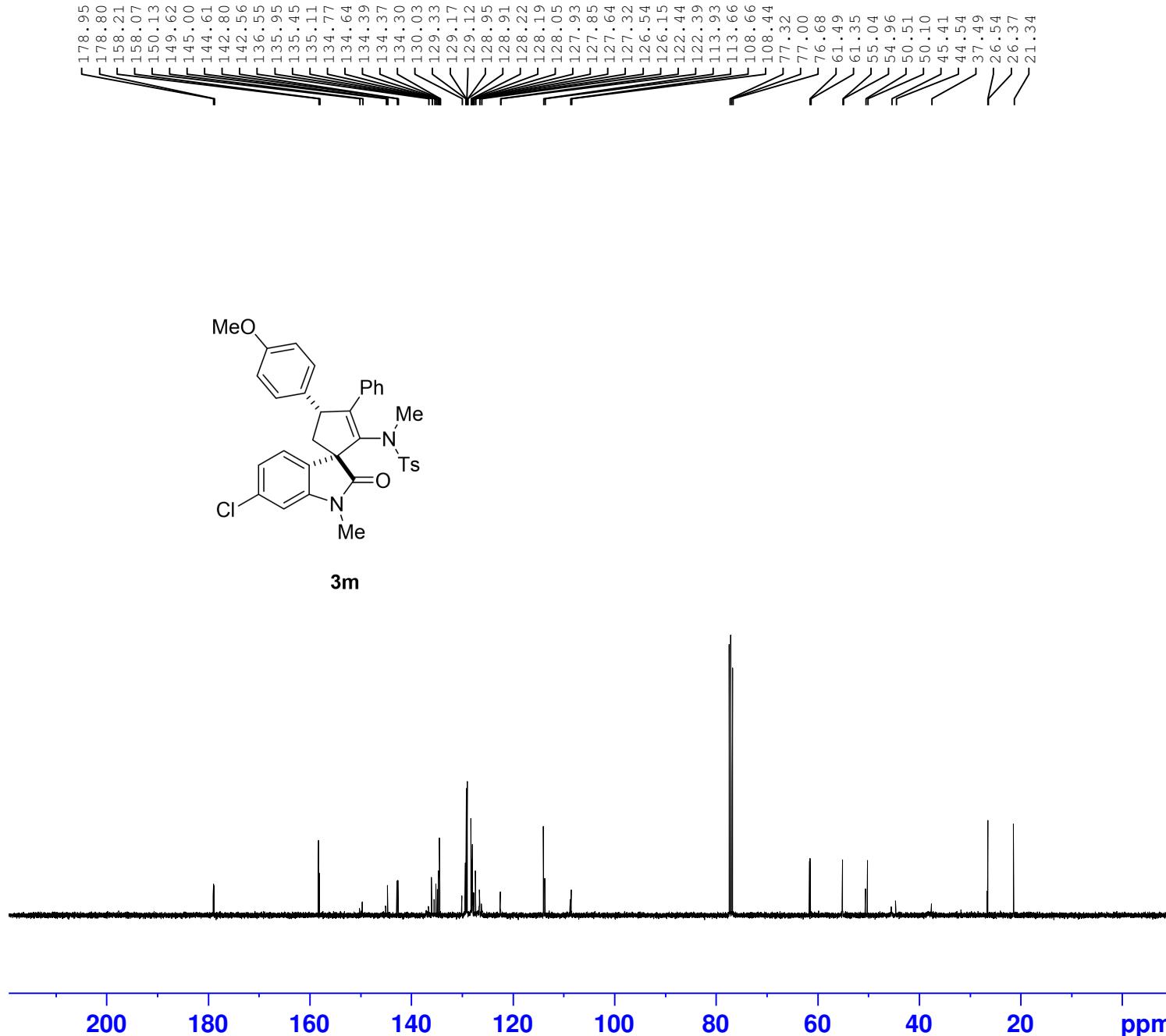


Current Data Parameters  
NAME zy-10  
EXPNO 2020121601  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20201217  
Time 6.26 h  
INSTRUM spect  
PROBHD Z108618\_0256 (zg30  
PULPROG 65536  
TD CDC13  
NS 16  
DS 2  
SWH 8223.685 Hz  
FIDRES 0.250967 Hz  
AQ 3.9845889 sec  
RG 45.2  
DW 60.800 usec  
DE 6.50 usec  
TE 293.4 K  
D1 1.00000000 sec  
TD0 1  
SFO1 400.1724712 MHz  
NUC1 1H  
P0 3.33 usec  
P1 10.00 usec  
PLW1 15.00000000 W

F2 - Processing parameters  
SI 32768  
SF 400.1700072 MHz  
WDW no  
SSB 0  
LB 0 Hz  
GB 0  
PC 1.00

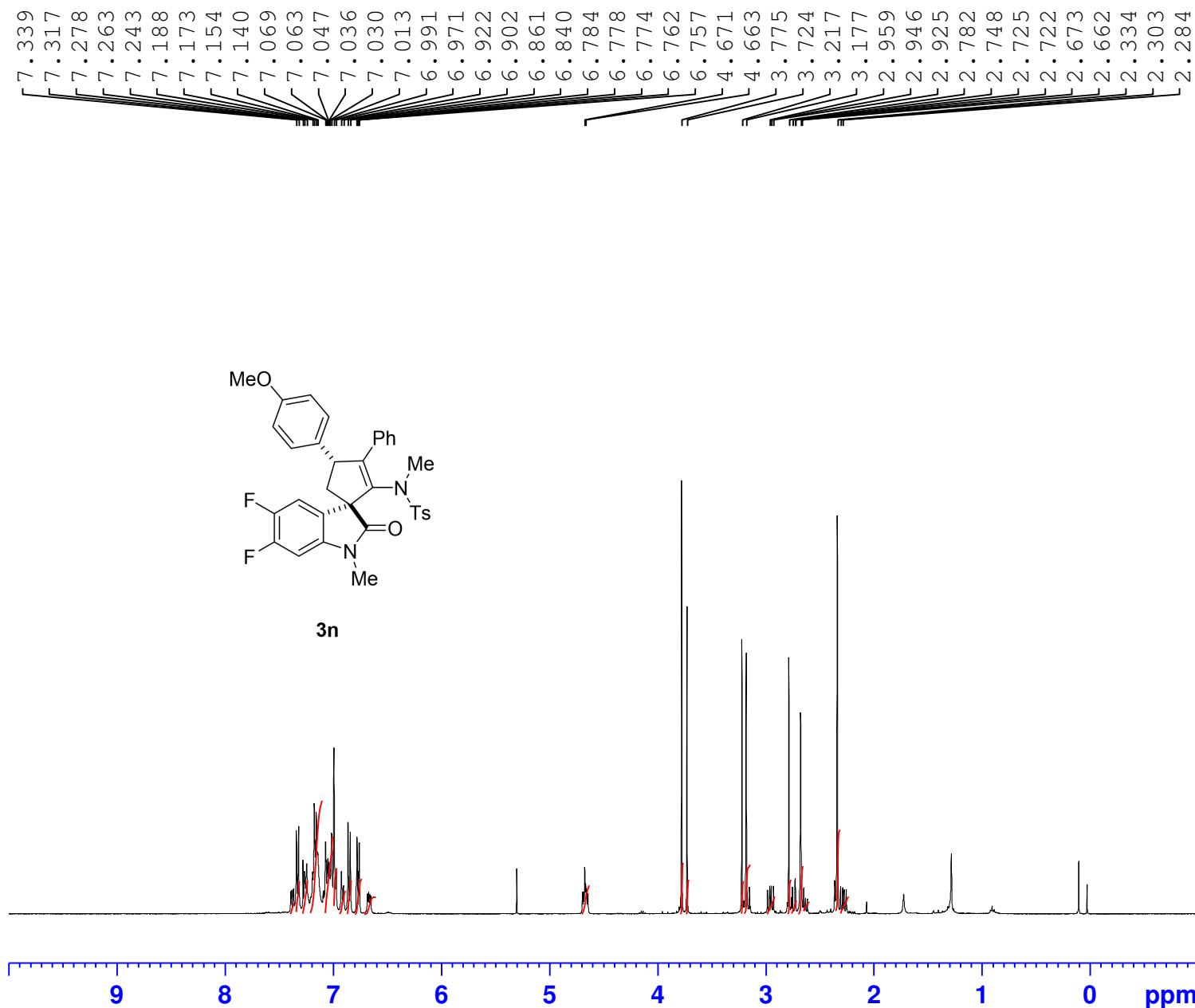




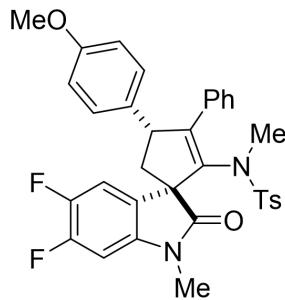
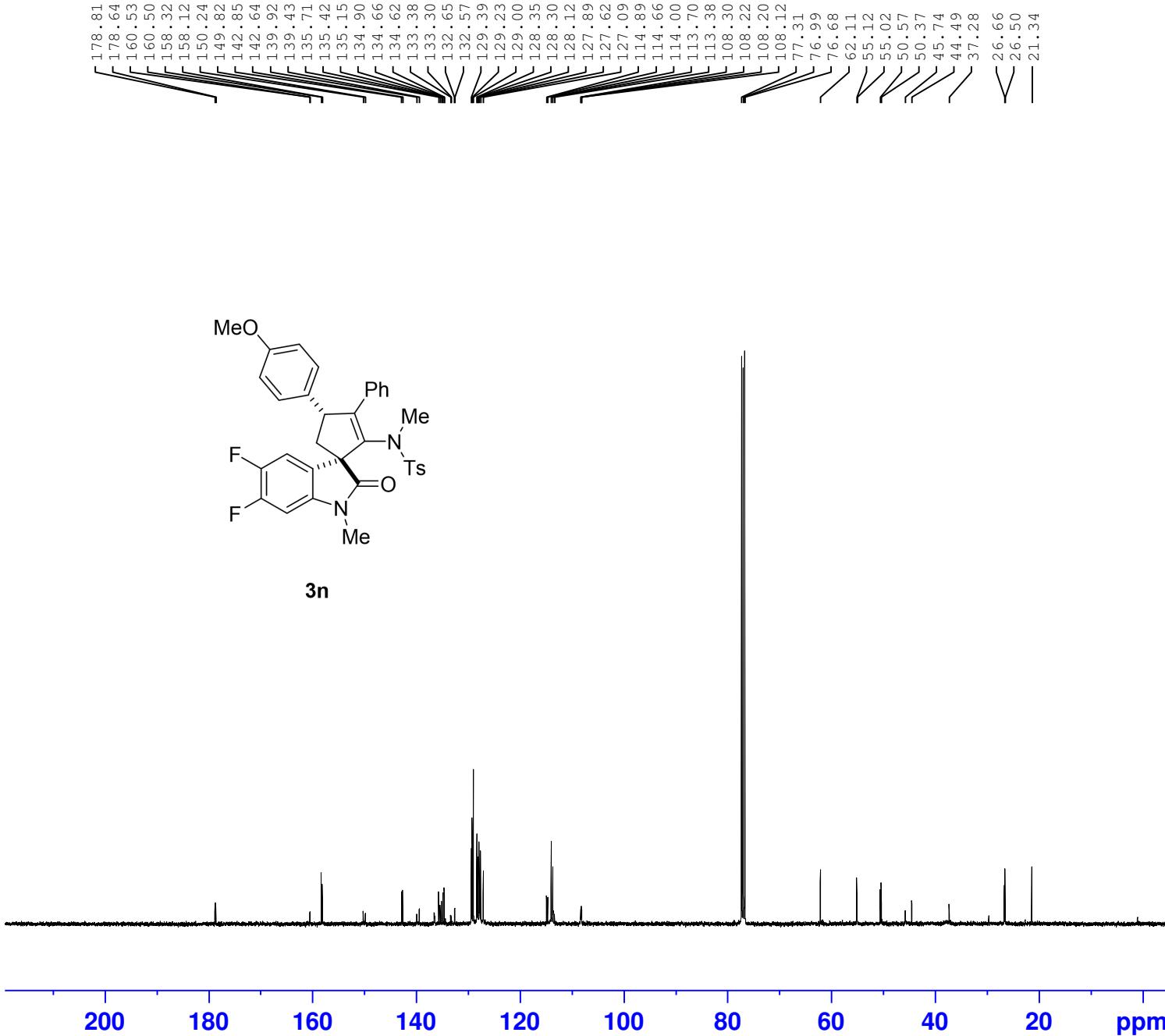
Current Data Parameters  
 NAME zy-10  
 EXPNO 2020121602  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20201217  
 Time 11.01 h  
 INSTRUM spect  
 PROBHD Z108618\_0256 (zgpg30)  
 PULPROG 65536  
 SOLVENT CDC13  
 NS 400  
 DS 4  
 SWH 24038.461 Hz  
 FIDRES 0.733596 Hz  
 AQ 1.3631488 sec  
 RG 2050  
 DW 20.800 usec  
 DE 6.50 usec  
 TE 293.7 K  
 D1 2.00000000 sec  
 D11 0.03000000 sec  
 TDO 1  
 SFO1 100.6328888 MHz  
 NUC1 13C  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 50.00000000 W  
 SFO2 400.1716007 MHz  
 NUC2 1H  
 CPDPRG[2] waltz65  
 PCPD2 90.00 usec  
 PLW2 15.00000000 W  
 PLW12 0.18519001 W  
 PLW13 0.09314700 W

F2 - Processing parameters  
 SI 32768  
 SF 100.6228418 MHz  
 WDW no  
 SSB 0  
 LB 0 Hz  
 GB 0  
 PC 1.40



1.14  
3.13  
3.17  
10.06  
7.14  
4.23  
2.05  
3.09  
3.07  
1.50  
2.51  
4.54  
3.00  
2.94  
1.55  
1.24  
4.24  
1.08  
7.43  
1.50



3n

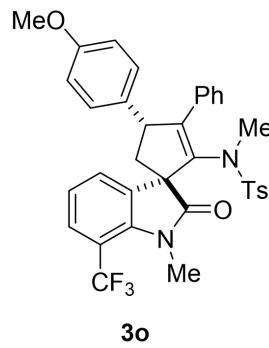
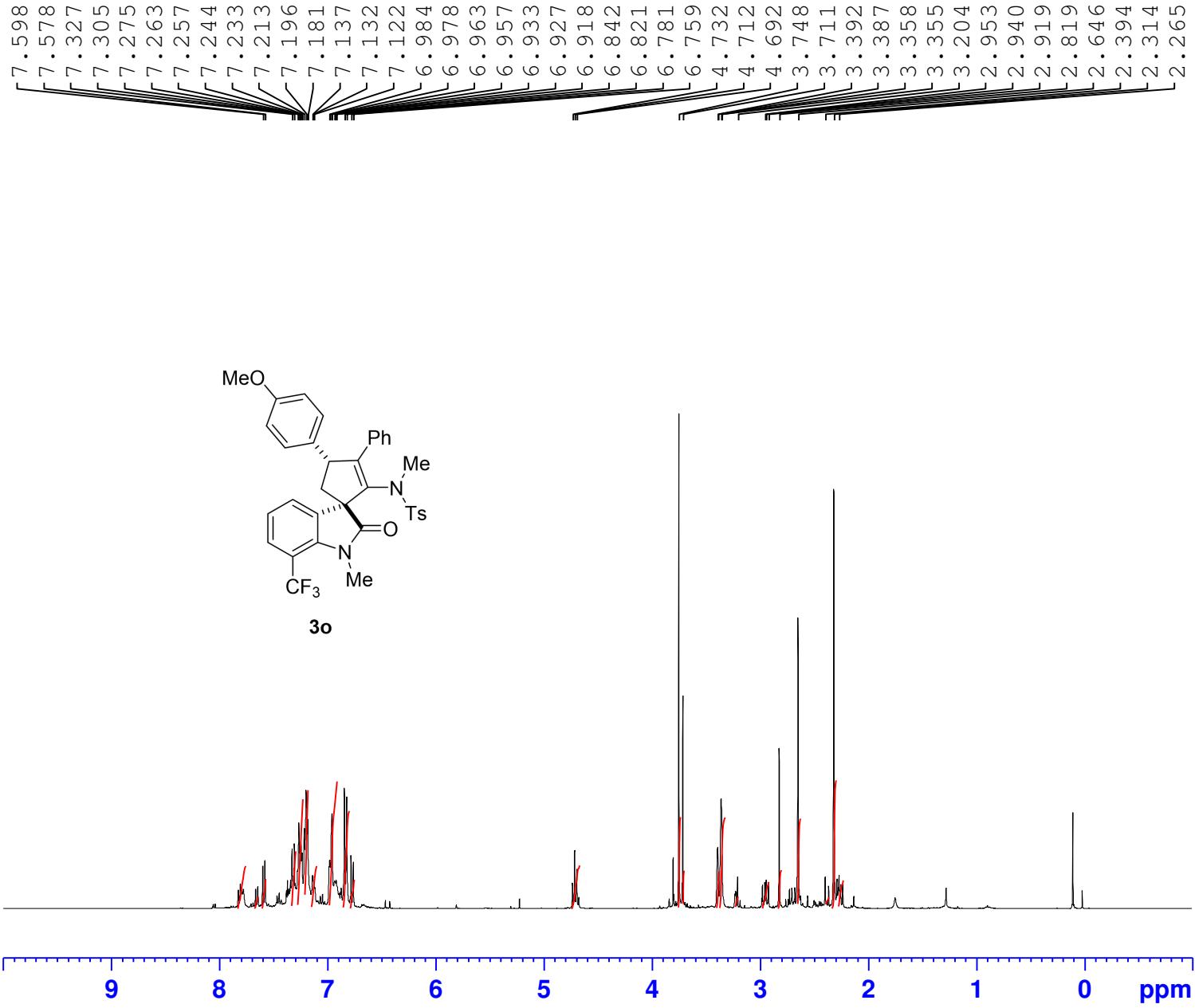
Current	Data	Parameters
NAME	cqh-145	
EXPNO	2020122302	
PROCNO		1

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F2 - Acquisition Parameters
Date_           20201224
Time            13.27 h
INSTRUM        spect
PROBHD         Z108618_0256 (
PULPROG        zgppg30
TD              65536
SOLVENT         CDC13
NS              800
DS              4
SWH             24038.461 Hz
FIDRES         0.733596 Hz
AQ              1.3631488 sec
RG              2050
DW              20.800 usec
DE              6.50  usec
TE              294.6 K
D1              2.00000000 sec
D11             0.03000000 sec
TDO              1
SFO1            100.6328888 MHz
NUC1            13C
P0              3.33  usec
P1              10.00 usec
PLW1            50.00000000 W
SFO2            400.1716007 MHz
NUC2              1H
CPDPRG[2       waltz65
PCPD2           90.00 usec
PLW2            15.00000000 W
PLW12           0.18519001 W
PLW13           0.09314700 W

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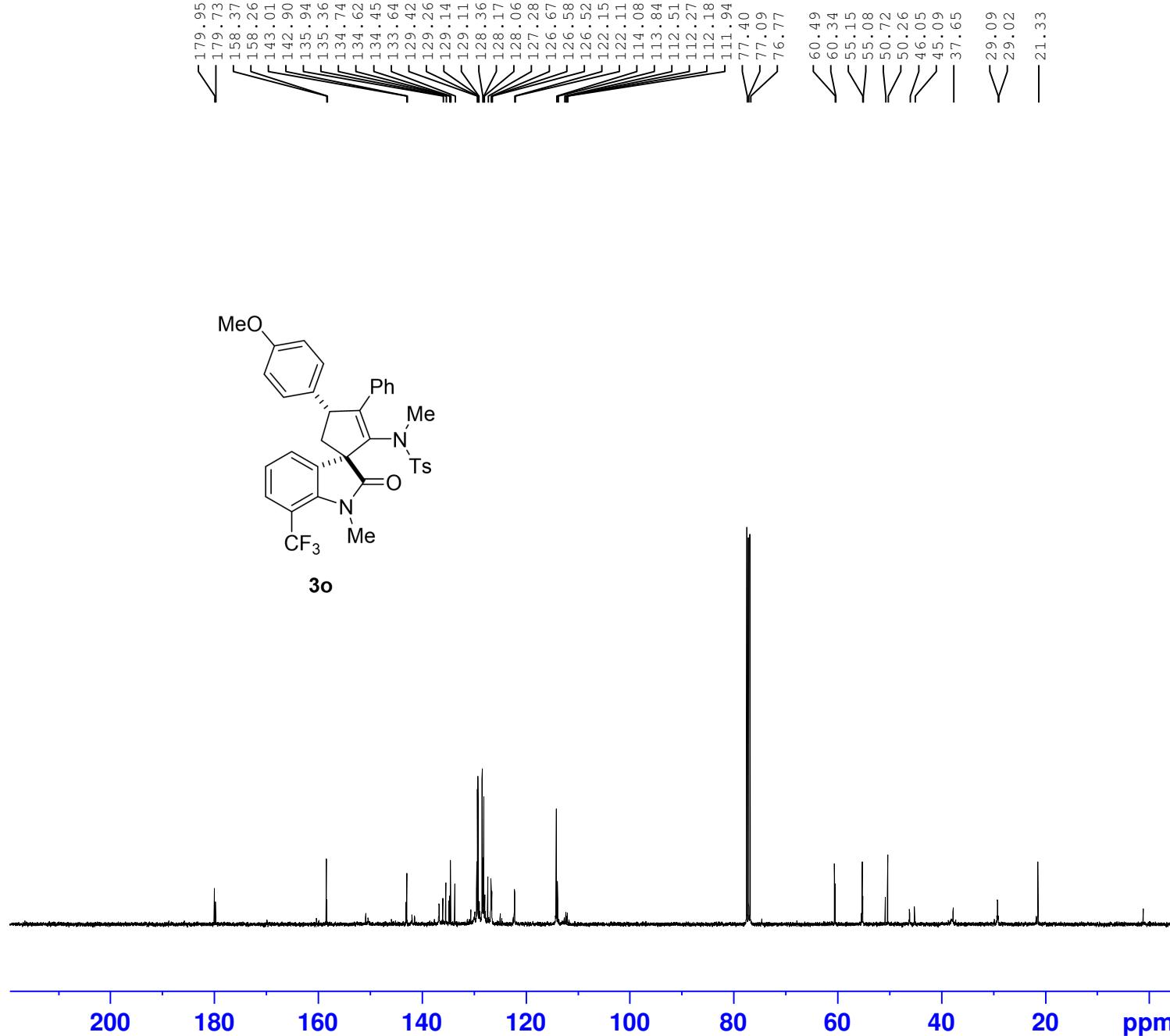
F2 - Processing parameters  
SI 32768  
SF 100.6228359 MHz  
WDW EM  
SSB 0  
LB 1.00 Hz  
GB 0  
PC 1.40



Current Data Parameters  
NAME cgh-167  
EXPNO 2021030901  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20210310  
Time 6.42 h  
INSTRUM spect  
PROBHD Z108618\_0256 (zg30  
PULPROG zg30  
TD 65536  
SOLVENT CDCl3  
NS 16  
DS 2  
SWH 8223.685 Hz  
FIDRES 0.250967 Hz  
AQ 3.9845889 sec  
RG 71.8  
DW 60.800 usec  
DE 6.50 usec  
TE 294.5 K  
D1 1.00000000 sec  
TD0 1  
SFO1 400.1724712 MHz  
NUC1 1H  
P0 3.33 usec  
P1 10.00 usec  
PLW1 15.00000000 W

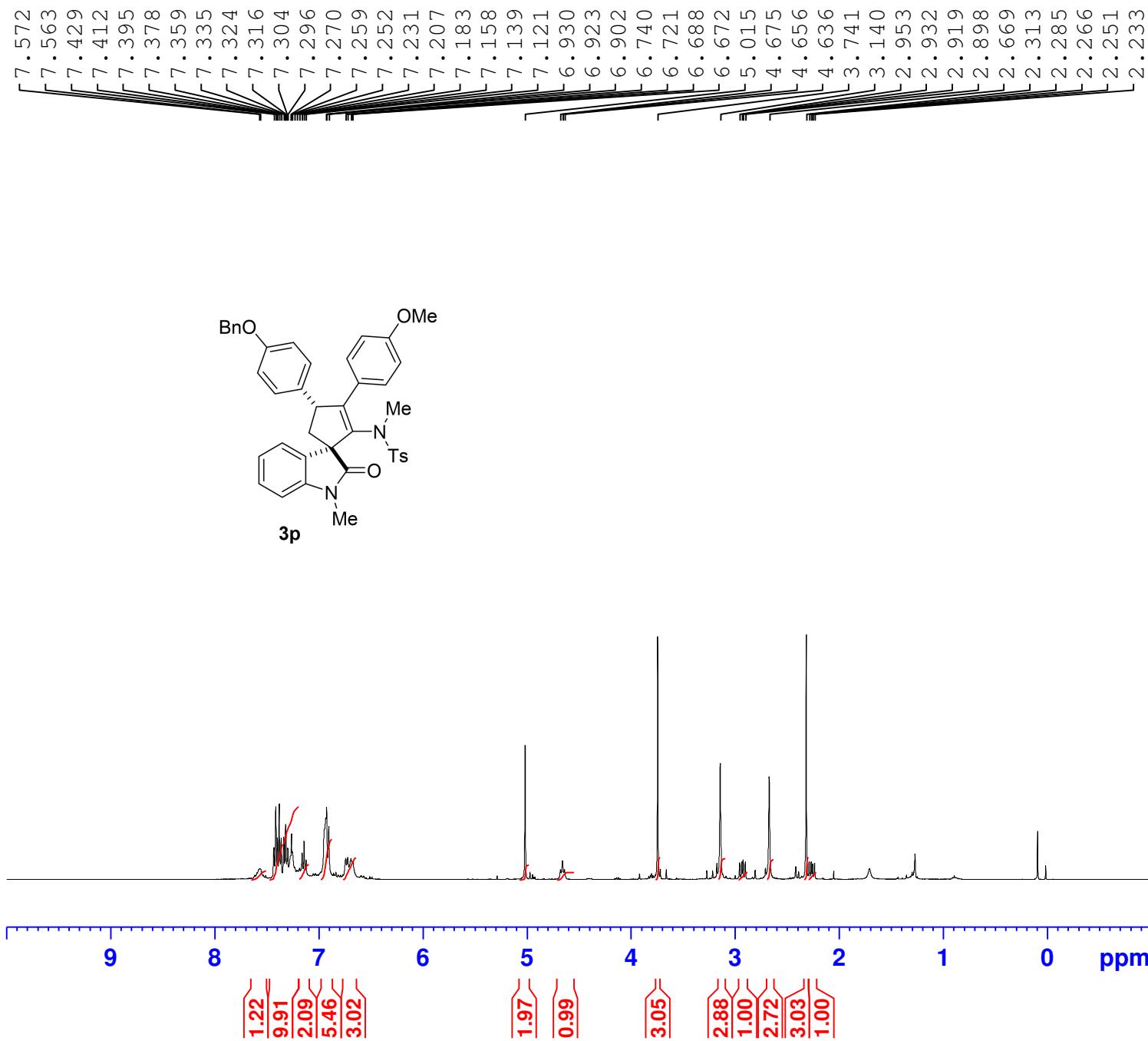
F2 - Processing parameters  
SI 32768  
SF 400.1700072 MHz  
WDW no  
SSB 0  
LB 0 Hz  
GB 0  
PC 1.00

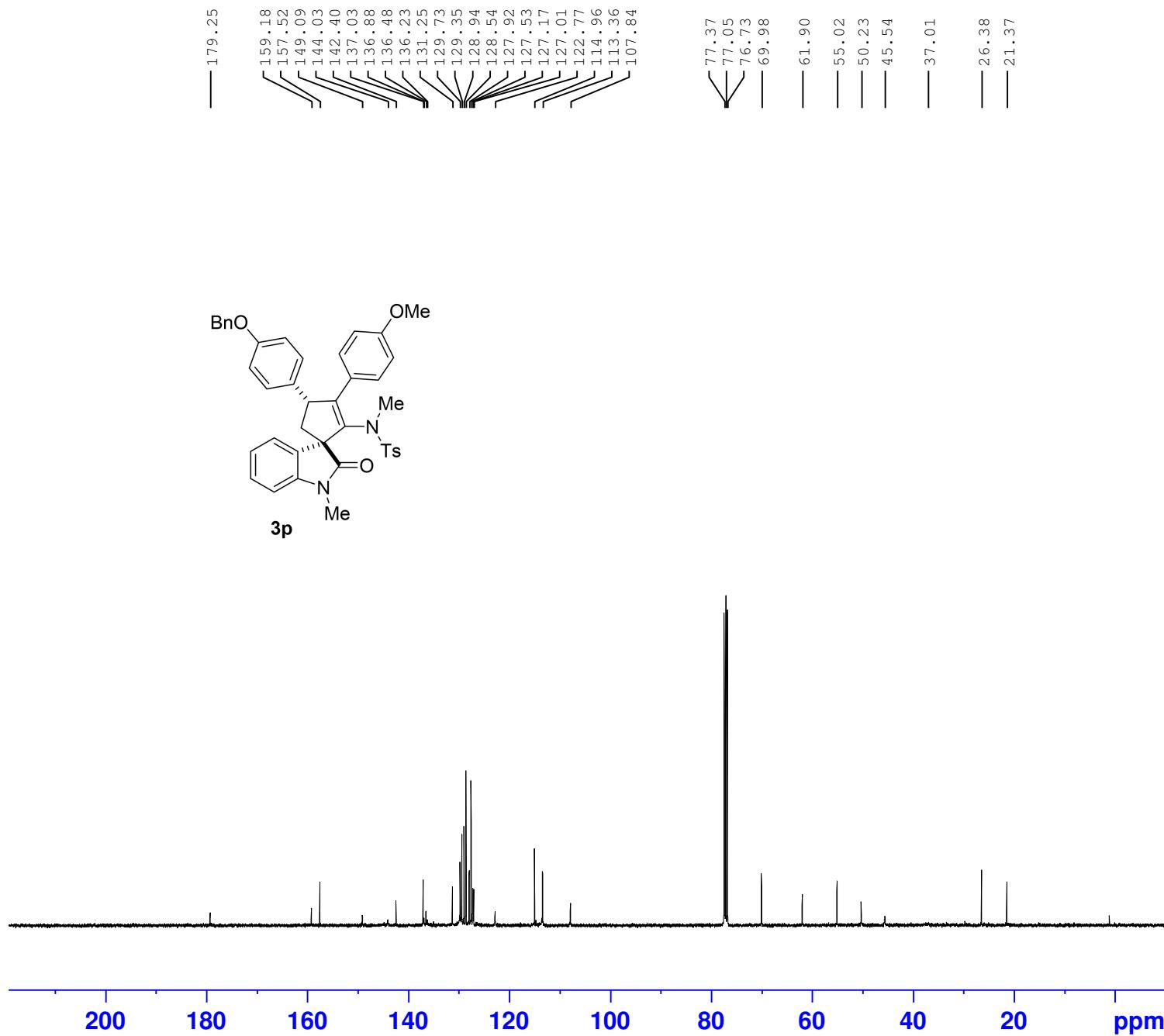


Current Data Parameters  
 NAME cqh-167  
 EXPNO 2021030902  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20210310  
 Time 13.16 h  
 INSTRUM spect  
 PROBHD Z108618\_0256 (zgpg30  
 PULPROG 65536  
 TD 533  
 SOLVENT CDC13  
 NS 4  
 DS 24038.461 Hz  
 SWH 0.733596 Hz  
 FIDRES 1.3631488 sec  
 AQ 2050  
 RG 20.800 usec  
 DE 6.50 usec  
 TE 294.5 K  
 D1 2.00000000 sec  
 D11 0.03000000 sec  
 TDO 1  
 SFO1 100.6328888 MHz  
 NUC1 13C  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 50.00000000 W  
 SFO2 400.1716007 MHz  
 NUC2 1H  
 CPDPRG[2] waltz65  
 PCPD2 90.00 usec  
 PLW2 15.00000000 W  
 PLW12 0.18519001 W  
 PLW13 0.09314700 W

F2 - Processing parameters  
 SI 32768  
 SF 100.6228300 MHz  
 WDW EM  
 SSB 0  
 LB 1.00 Hz  
 GB 0  
 PC 1.40

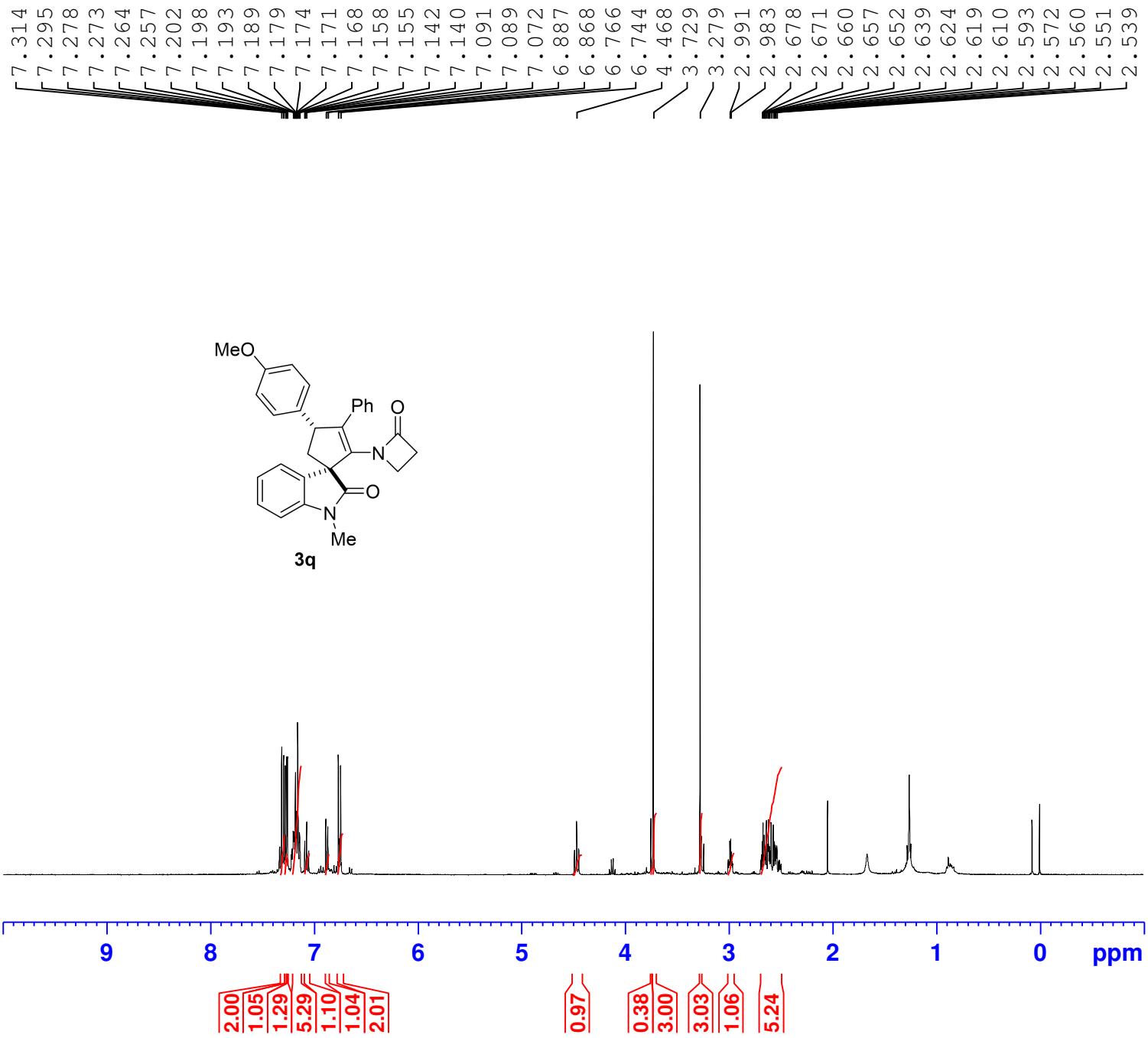




Current Data Parameters  
 NAME zy-1  
 EXPNO 2020092102  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20200925  
 Time 1.56 h  
 INSTRUM spect  
 PROBHD Z108618\_0256 (zgpg30)  
 PULPROG zgpg30  
 TD 65536  
 SOLVENT CDCl<sub>3</sub>  
 NS 866  
 DS 4  
 SWH 24038.461 Hz  
 FIDRES 0.733596 Hz  
 AQ 1.3631488 sec  
 RG 2050  
 DW 20.800 usec  
 DE 6.500 usec  
 TE 298.1 K  
 D1 2.000000000 sec  
 D11 0.030000000 sec  
 TD0 1  
 SFO1 100.6328888 MHz  
 NUC1 13C  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 50.000000000 W  
 SFO2 400.1716007 MHz  
 NUC2 1H  
 CPDPRG[2] waltz65  
 PCPD2 90.00 usec  
 PLW2 15.000000000 W  
 PLW12 0.18519001 W  
 PLW13 0.09314700 W

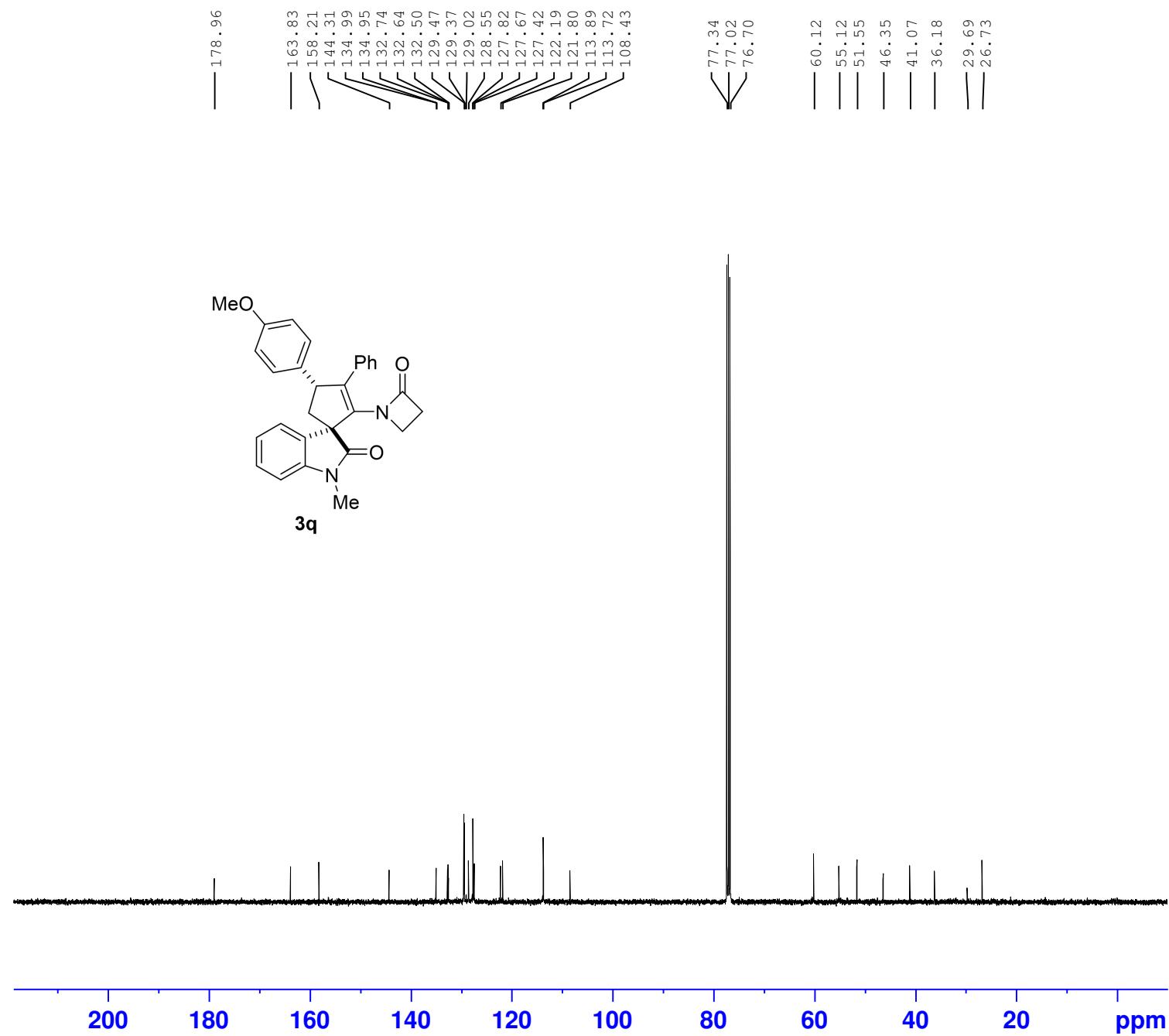
F2 - Processing parameters  
 SI 32768  
 SF 100.6228300 MHz  
 WDW EM  
 SSB 0  
 LB 1.00 Hz  
 GB 0  
 PC 1.40



Current Data Parameters  
 NAME cqh-146  
 EXPNO 2021030901  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20210310  
 Time 5.58 h  
 INSTRUM spect  
 PROBHD Z108618\_0256 (zg30  
 PULPROG zg30  
 TD 65536  
 SOLVENT CDCl3  
 NS 16  
 DS 2  
 SWH 8223.685 Hz  
 FIDRES 0.250967 Hz  
 AQ 3.9845889 sec  
 RG 362  
 DW 60.800 usec  
 DE 6.50 usec  
 TE 294.4 K  
 D1 1.00000000 sec  
 TDO 1  
 SFO1 400.1724712 MHz  
 NUC1 1H  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 15.00000000 W

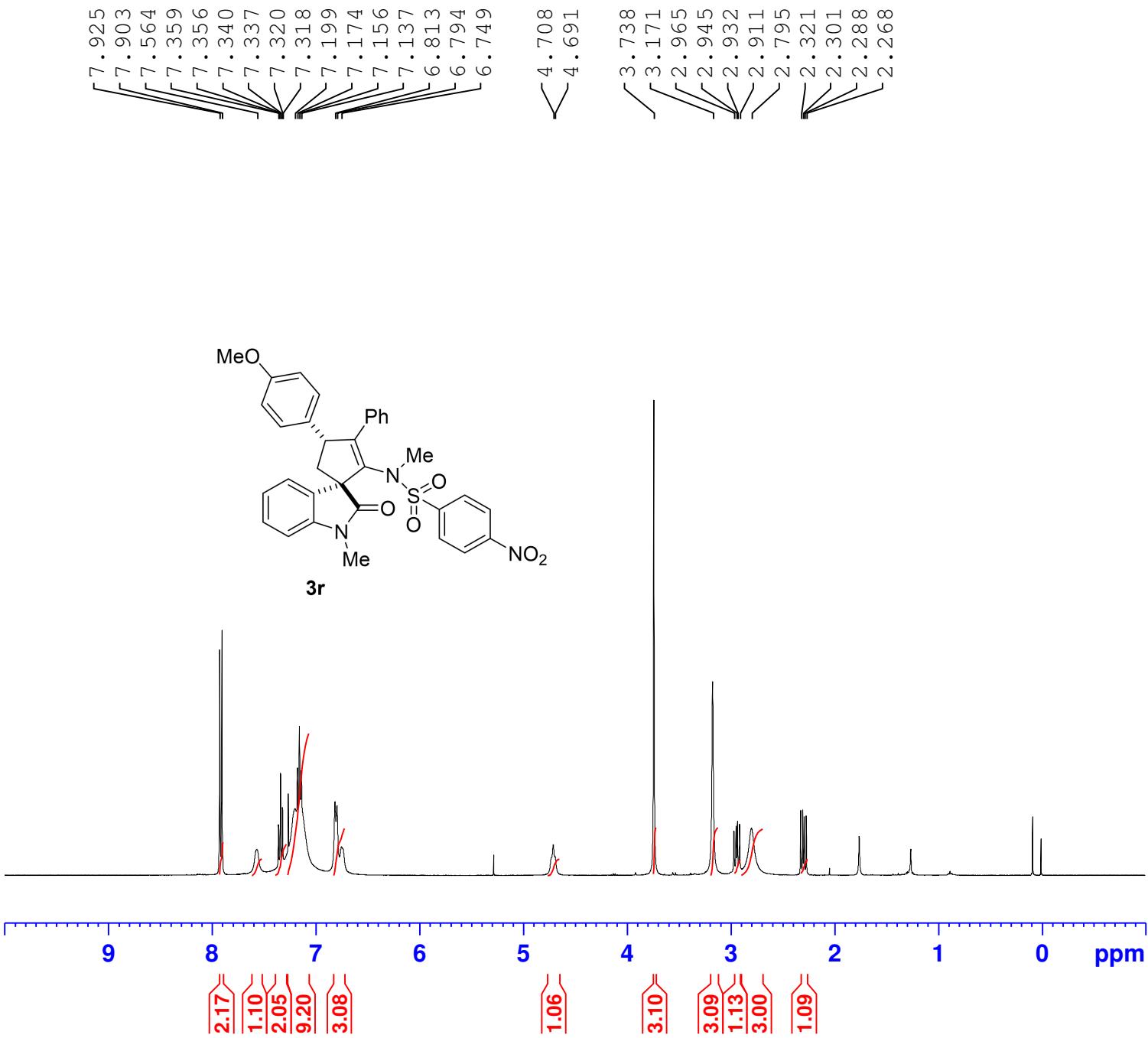
F2 - Processing parameters  
 SI 32768  
 SF 400.1700072 MHz  
 WDW no  
 SSB 0  
 LB 0 Hz  
 GB 0  
 PC 1.00



Current Data Parameters  
 NAME cgh-146  
 EXPNO 2021010802  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20210310  
 Time 9.53 h  
 INSTRUM spect  
 PROBHD Z108618\_0256 (zpg30)  
 PULPROG zpg30  
 TD 65536  
 SOLVENT CDCl3  
 NS 681  
 DS 4  
 SWH 24038.461 Hz  
 FIDRES 0.733596 Hz  
 AQ 1.3631488 sec  
 RG 2050  
 DW 20.800 usec  
 DE 6.50 usec  
 TE 294.7 K  
 D1 2.00000000 sec  
 D11 0.03000000 sec  
 TDO 1  
 SFO1 100.6328888 MHz  
 NUC1 13C  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 50.00000000 W  
 SFO2 400.1716007 MHz  
 NUC2 1H  
 CPDPRG[2] waltz65  
 PCPD2 90.00 usec  
 PLW2 15.00000000 W  
 PLW12 0.18519001 W  
 PLW13 0.09314700 W

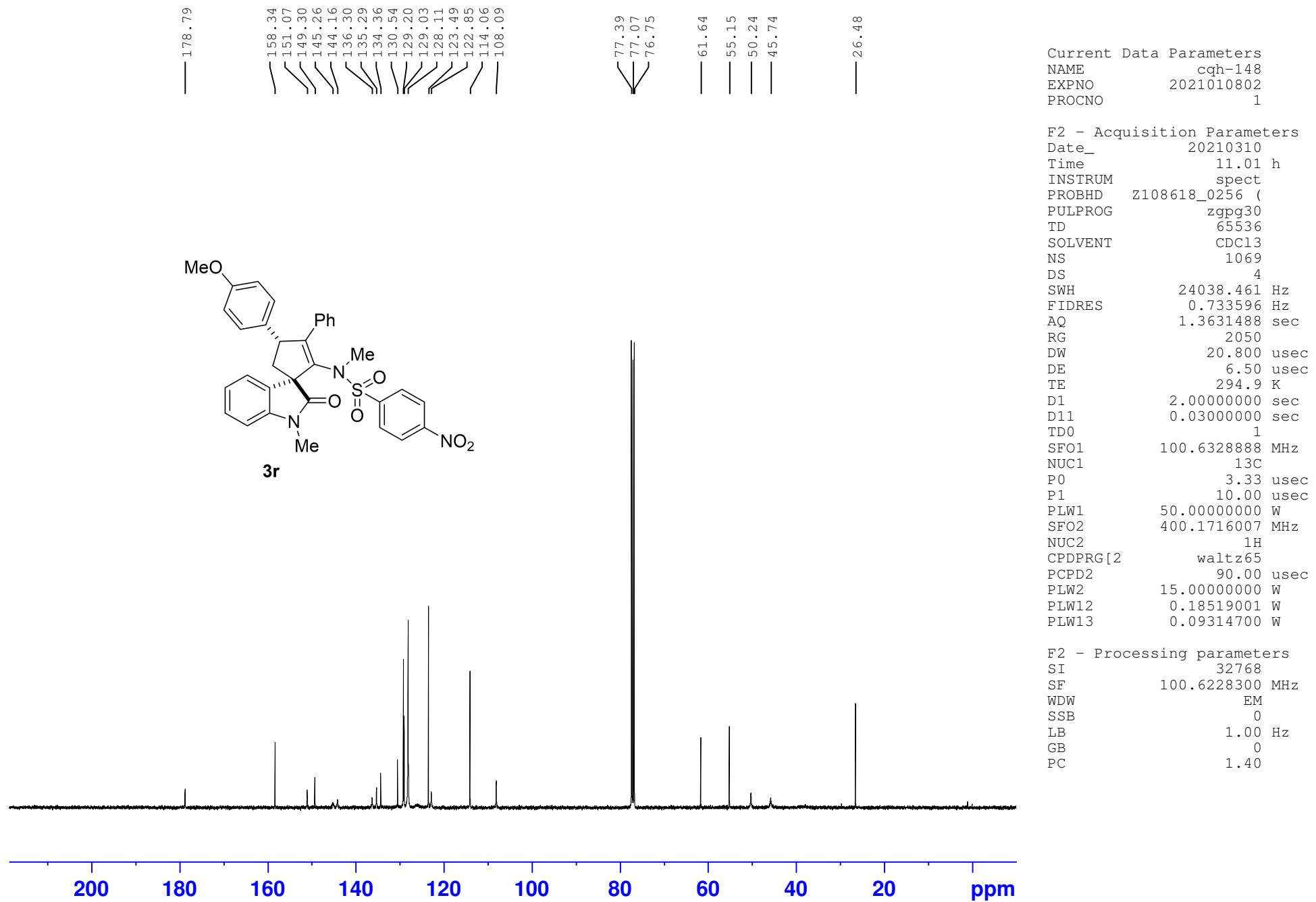
F2 - Processing parameters  
 SI 32768  
 SF 100.6228300 MHz  
 WDW EM  
 SSB 0  
 LB 1.00 Hz  
 GB 0  
 PC 1.40

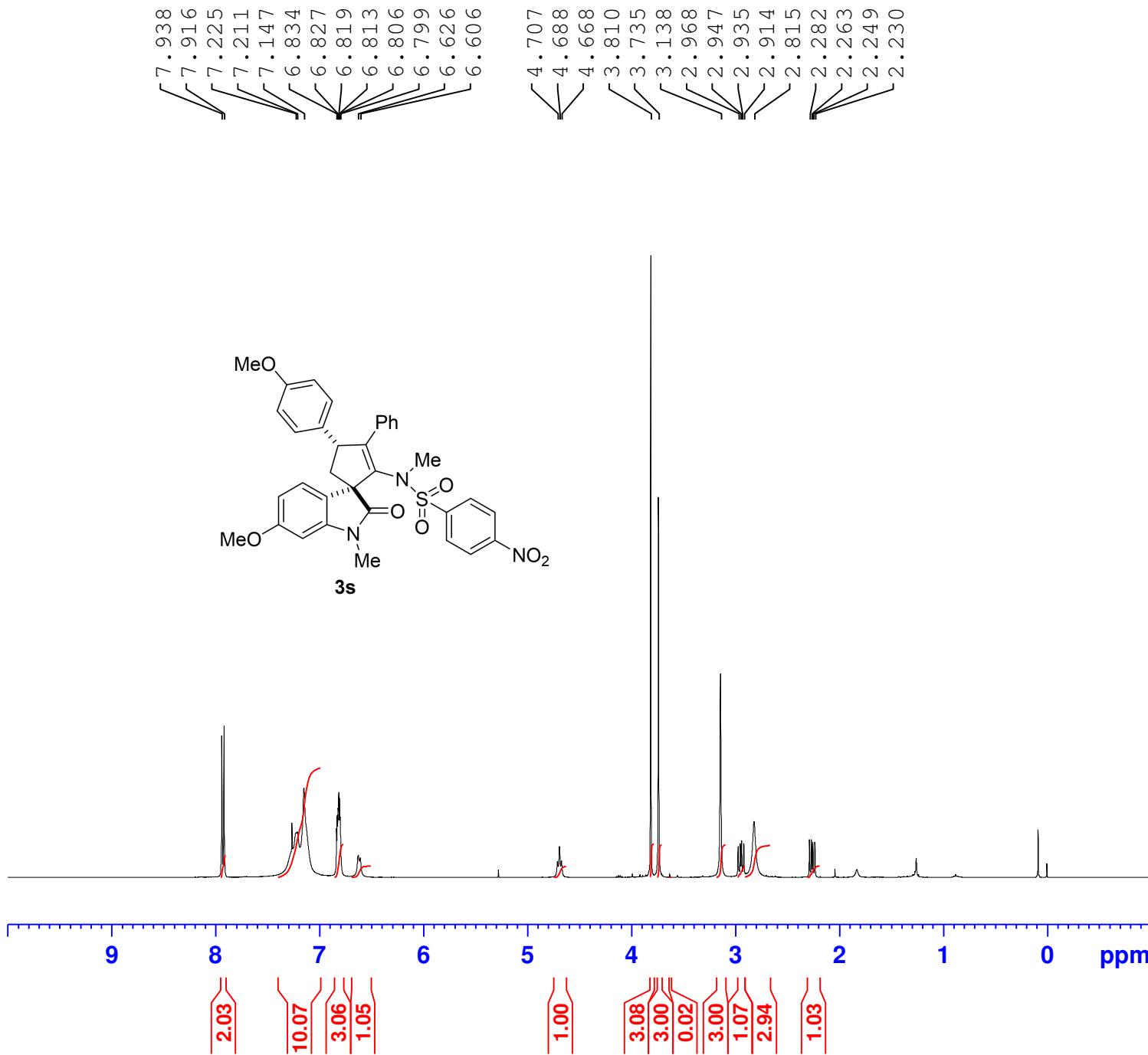


Current Data Parameters  
NAME cgh-148  
EXPNO 2021030901  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20210310  
Time 6.04 h  
INSTRUM spect  
PROBHD Z108618\_0256 (zg30  
PULPROG 65536  
TD CDC13  
NS 16  
DS 2  
SWH 8223.685 Hz  
FIDRES 0.250967 Hz  
AQ 3.9845889 sec  
RG 362  
DW 60.800 usec  
DE 6.50 usec  
TE 294.5 K  
D1 1.000000000 sec  
TD0 1  
SFO1 400.1724712 MHz  
NUC1 1H  
P0 3.33 usec  
P1 10.00 usec  
PLW1 15.000000000 W

F2 - Processing parameters  
SI 32768  
SF 400.1700072 MHz  
WDW no  
SSB 0  
LB 0 Hz  
GB 0  
PC 1.00

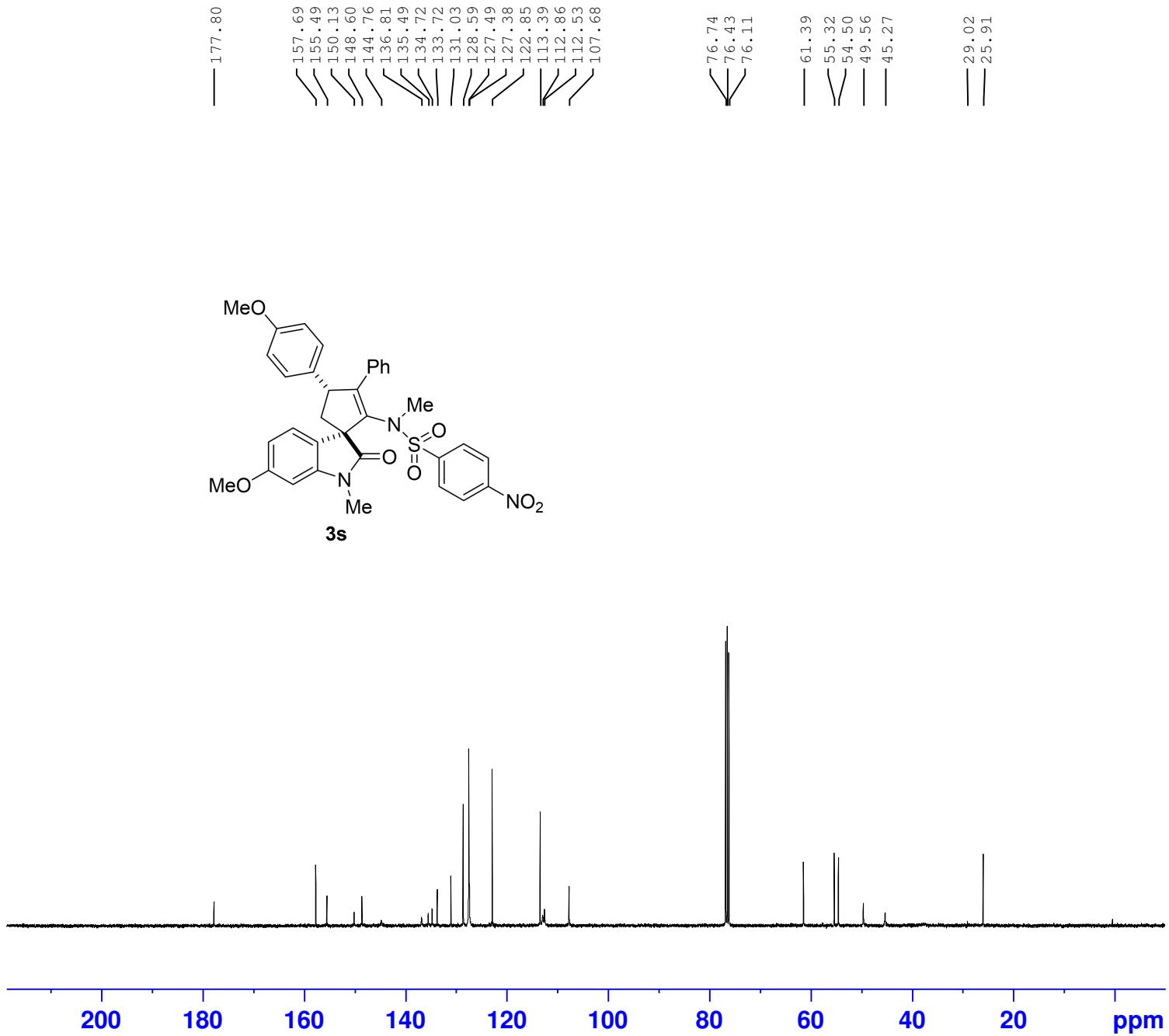




Current Data Parameters  
 NAME cgh-193  
 EXPNO 2021041901  
 PROCNO 1

F2 - Acquisition Parameters  
 Date\_ 20210420  
 Time 2.00 h  
 INSTRUM spect  
 PROBHD Z108618\_0256 (zg30  
 PULPROG zg30  
 TD 65536  
 SOLVENT CDCl<sub>3</sub>  
 NS 16  
 DS 2  
 SWH 8223.685 Hz  
 FIDRES 0.250967 Hz  
 AQ 3.9845889 sec  
 RG 101  
 DW 60.800 usec  
 DE 6.50 usec  
 TE 293.2 K  
 D1 1.00000000 sec  
 TDO 1  
 SFO1 400.1724712 MHz  
 NUC1 1H  
 P0 3.33 usec  
 P1 10.00 usec  
 PLW1 15.00000000 W

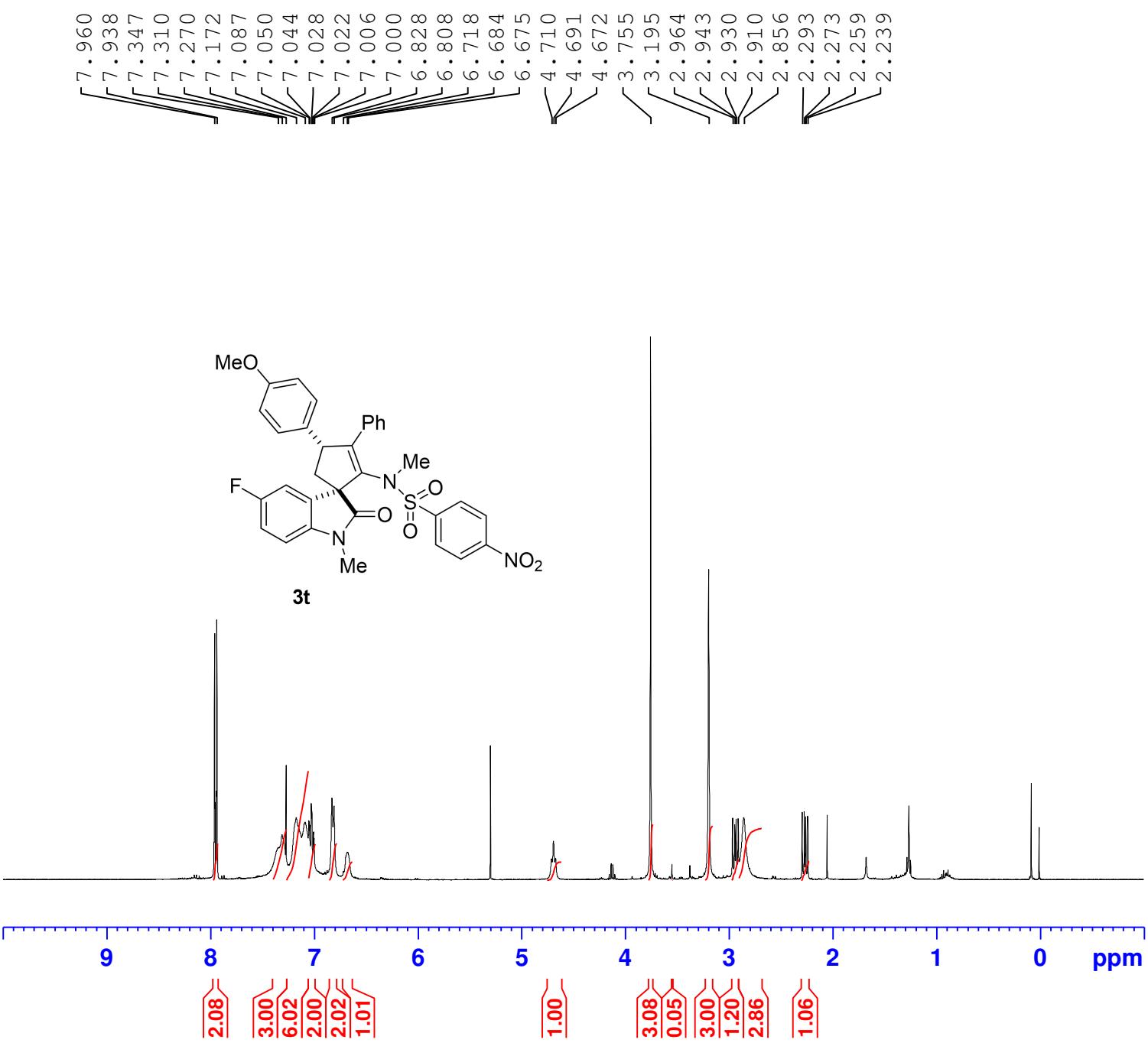
F2 - Processing parameters  
 SI 32768  
 SF 400.1700072 MHz  
 WDW no  
 SSB 0  
 LB 0 Hz  
 GB 0  
 PC 1.00



Current Data Parameters  
NAME cqh-193  
EXPNO 2021041902  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20210420  
Time 3.38 h  
INSTRUM spect  
PROBHD Z108618\_0256 (   
PULPROG zgpg30  
TD 65536  
SOLVENT CDC13  
NS 501  
DS 4  
SWH 24038.461 Hz  
FIDRES 0.733596 Hz  
AQ 1.3631488 sec  
RG 2050  
DW 20.800 usec  
DE 6.50 usec  
TE 293.7 K  
D1 2.00000000 sec  
D11 0.03000000 sec  
TD0 1  
SFO1 100.6328888 MHz  
NUC1 13C  
P0 3.33 usec  
P1 10.00 usec  
PLW1 50.00000000 W  
SFO2 400.1716007 MHz  
NUC2 1H  
CPDPRG[2] waltz65  
PCPD2 90.00 usec  
PLW2 15.00000000 W  
PLW12 0.18519001 W  
PLW13 0.09314700 W

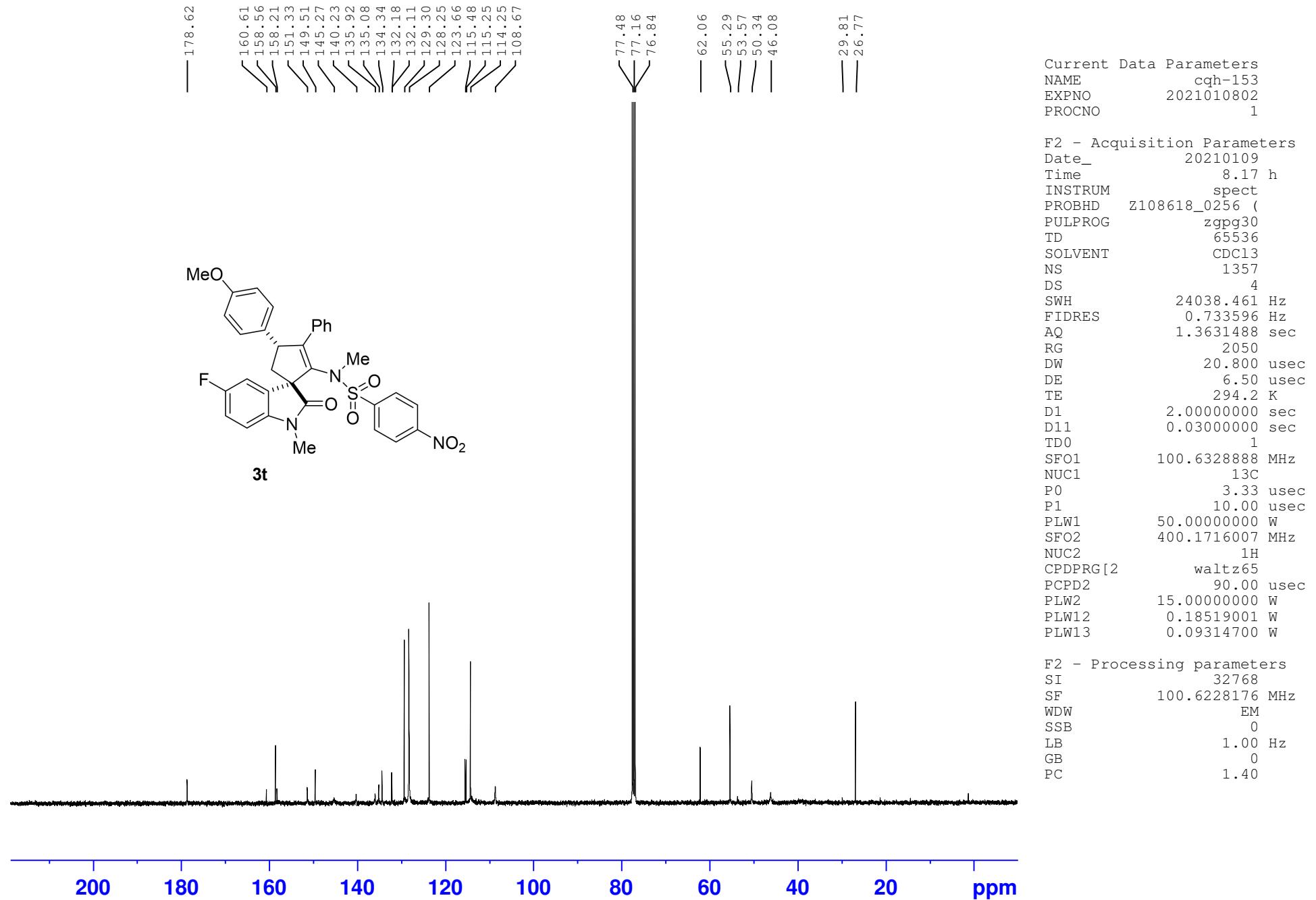
F2 - Processing parameters  
SI 32768  
SF 100.6228963 MHz  
WDW EM  
SSB 0  
LB 1.00 Hz  
GB 0  
PC 1.40

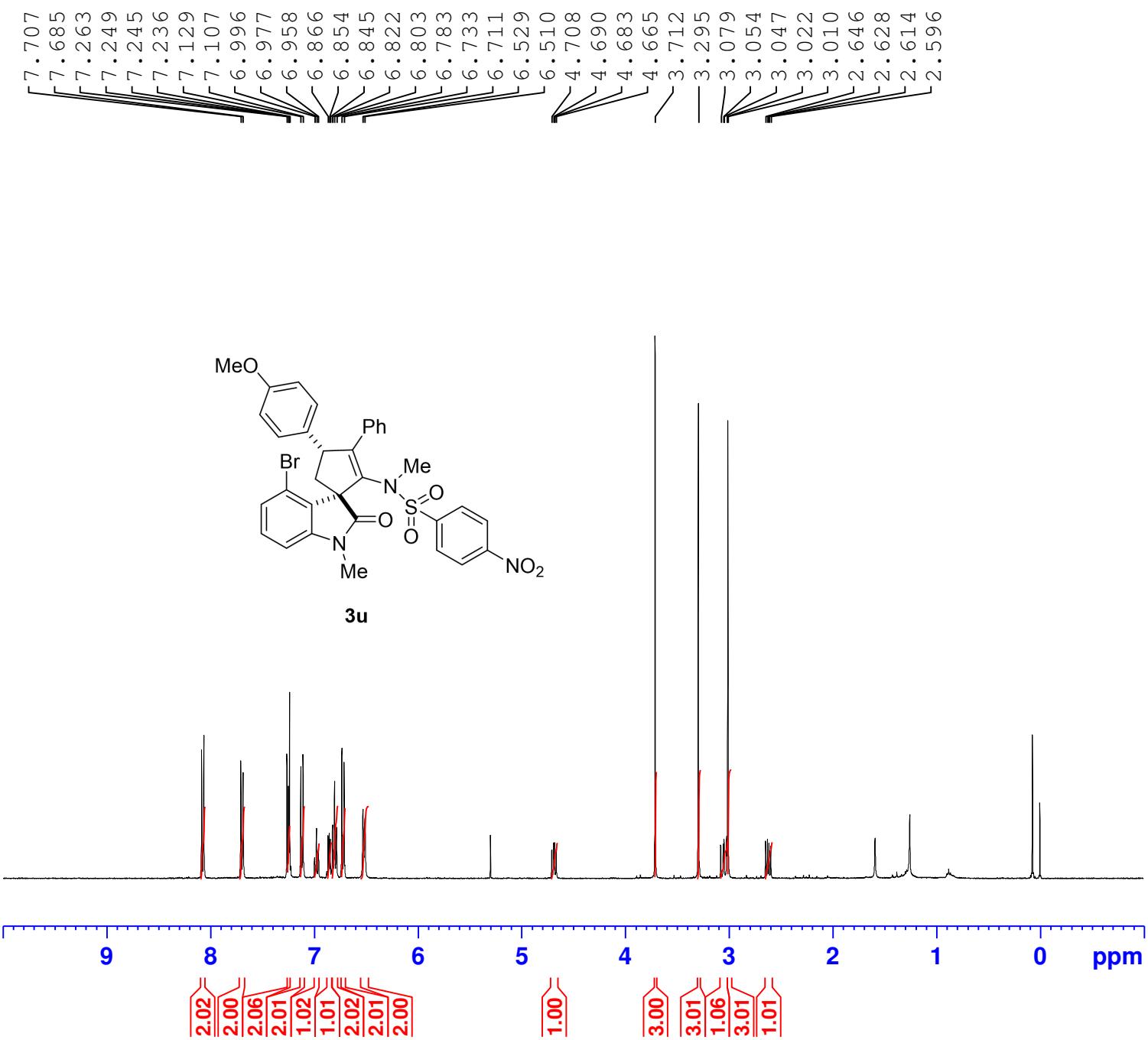


Current Data Parameters  
NAME cgh-153  
EXPNO 2021010801  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20210109  
Time 5.14 h  
INSTRUM spect  
PROBHD Z108618\_0256 (PULPROG zg30  
TD 65536  
SOLVENT CDCl<sub>3</sub>  
NS 16  
DS 2  
SWH 8223.685 Hz  
FIDRES 0.250967 Hz  
AQ 3.9845889 sec  
RG 128  
DW 60.800 usec  
DE 6.50 usec  
TE 293.7 K  
D1 1.0000000 sec  
TD0 1  
SFO1 400.1724712 MHz  
NUC1 1H  
P0 3.33 usec  
P1 10.00 usec  
PLW1 15.00000000 W

F2 - Processing parameters  
SI 32768  
SF 400.1700042 MHz  
WDW no  
SSB 0  
LB 0 Hz  
GB 0  
PC 1.00





Current Data Parameters  
NAME cgh-154  
EXPNO 2021010801  
PROCNO 1

F2 - Acquisition Parameters  
Date\_ 20210109  
Time 5.18 h  
INSTRUM spect  
PROBHD Z108618\_0256 (PULPROG zg30  
TD 65536  
SOLVENT CDCl3  
NS 16  
DS 2  
SWH 8223.685 Hz  
FIDRES 0.250967 Hz  
AQ 3.9845889 sec  
RG 287  
DW 60.800 usec  
DE 6.50 usec  
TE 293.7 K  
D1 1.00000000 sec  
TD0 1  
SFO1 400.1724712 MHz  
NUC1 1H  
P0 3.33 usec  
P1 10.00 usec  
PLW1 15.00000000 W

F2 - Processing parameters  
SI 32768  
SF 400.1700075 MHz  
WDW no  
SSB 0  
LB 0 Hz  
GB 0  
PC 1.00

