

Supporting Information

**K₂CO₃- Accelerated Amidation of Carboxylic Acids with α -
Oxo Ketene-*N*, *S*-Acetals as Amine Surrogates**

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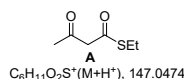
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I. Detection of S-ethyl 3-oxobutanethioate A from LC-MS

1. General Information. The chromatographic separation was performed on an Agilent 1260 Infinity HPLC apparatus equipped with a reverse phase SepaChrom Srl C18 (50 mm × 4.6 mm, 5.0 µm particle size) analytical column. LC flow rate: 3.5 mL min⁻¹, The mass spectrometry shunt flow rate is 0.3 mL min⁻¹. The ultra-pure water was prepared by a Millipore purification device (New York, USA). The column temperature was set at 35 °C. The injection volume was set at 2 µL. HPLC grade Acetonitrile (0.1% HPLC grade methanoic acid) and ultrapure water (0.1% HPLC grade methanoic acid) were used as mobile phase A and mobile phase B, respectively. The gradient was set as follows: 0→5 min, 30%→80% A; 5→6 min, 80% A. A Thermofisher TSQ Quantum Access MAX mass spectrometer was used for the identification and quantification of analytes with a Full Scan mode (m/z 100 – 1000). The parameters of the ESI source (positive ionization mode) were carefully optimized.

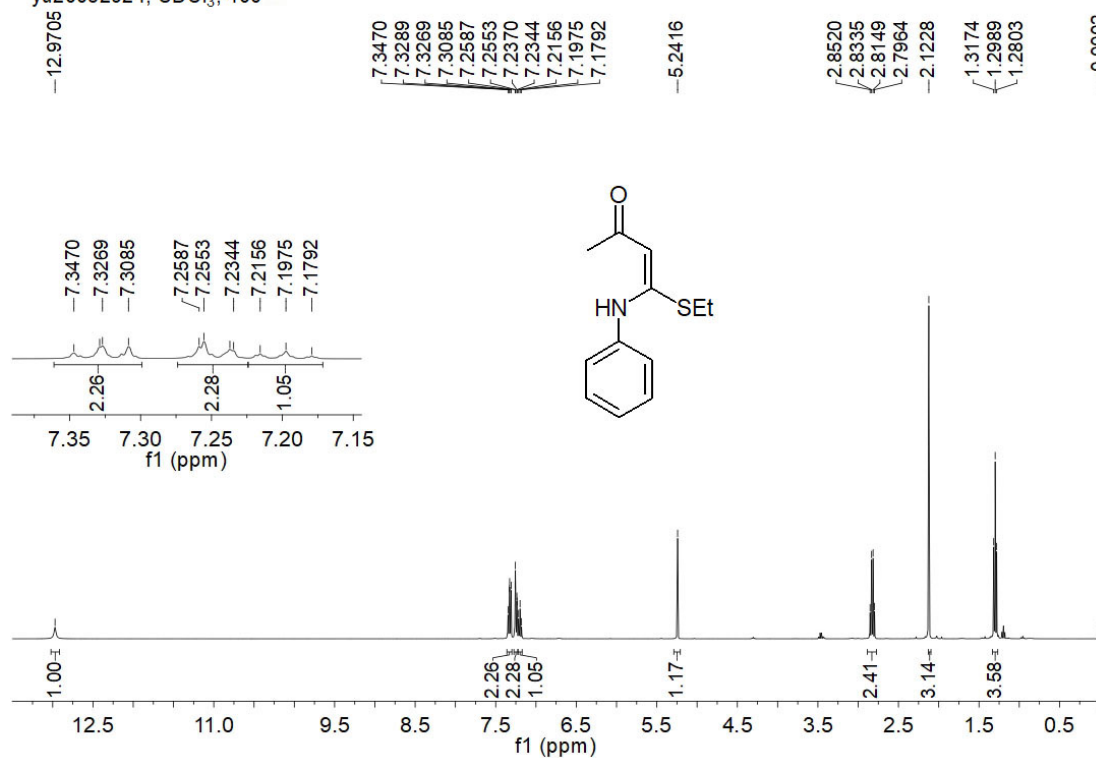
2. Total ion chromatogram

3. The MS of components based on retention time 1 minute

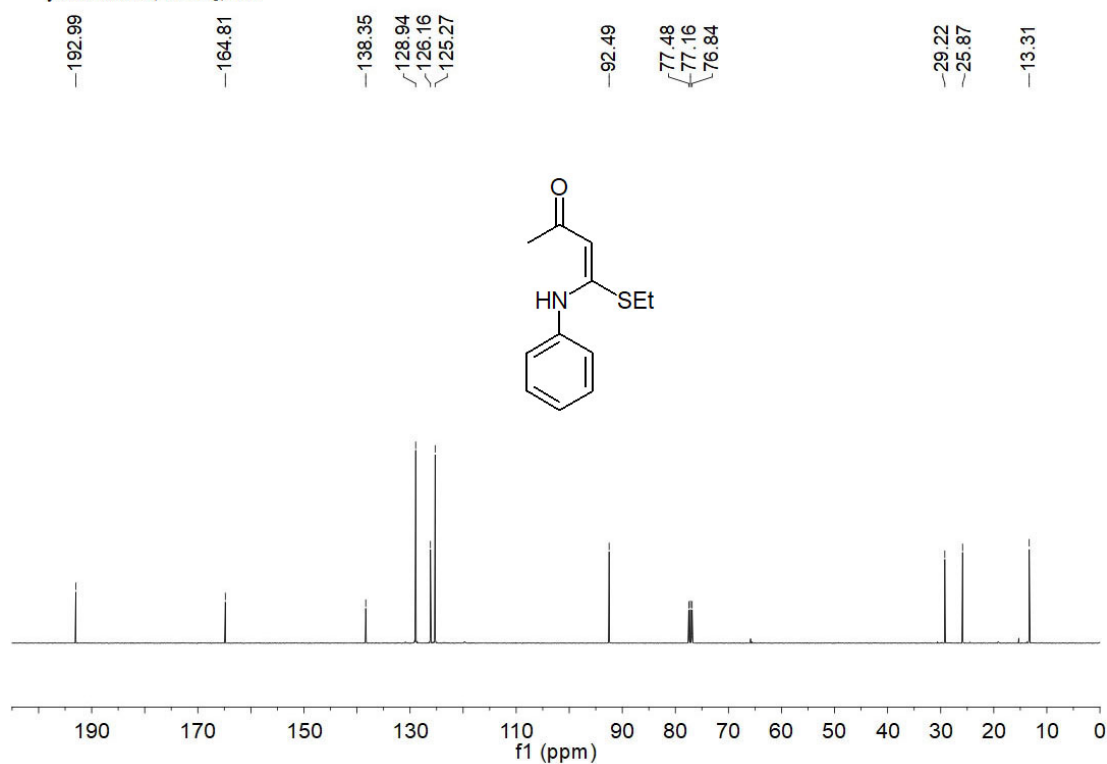


II. Copies of NMR spectra for 1

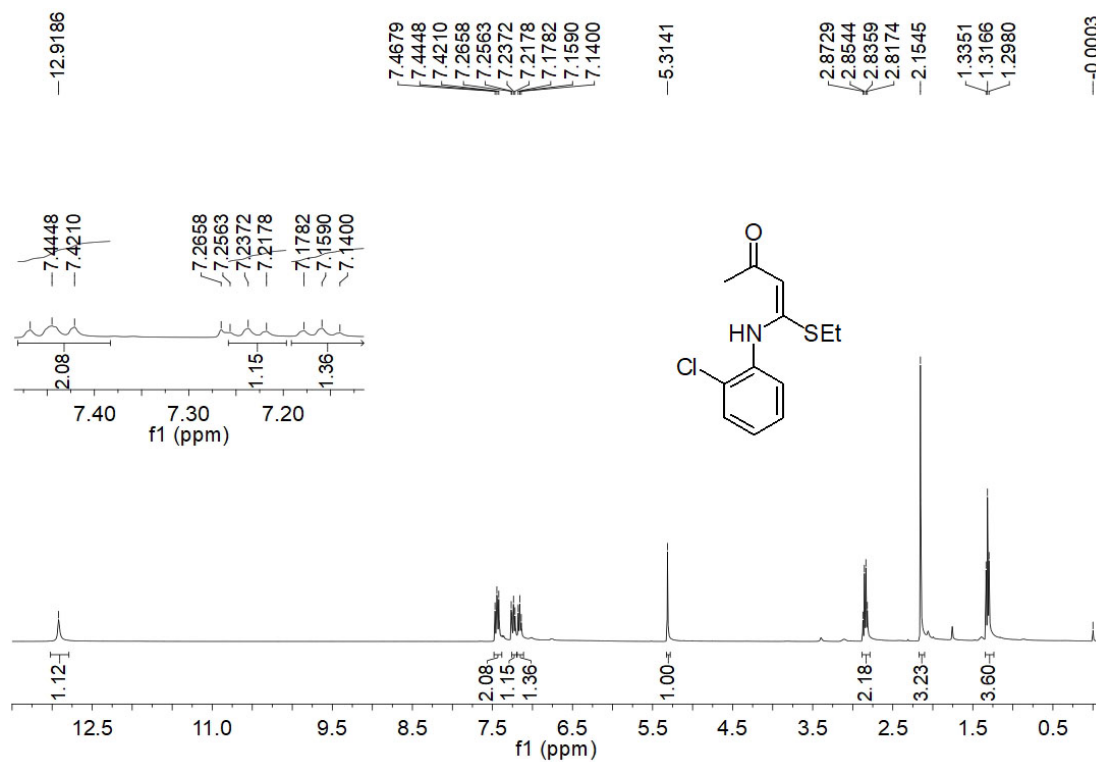
yu20052924, CDCl₃, 400



yu20052924, CDCl₃, 400

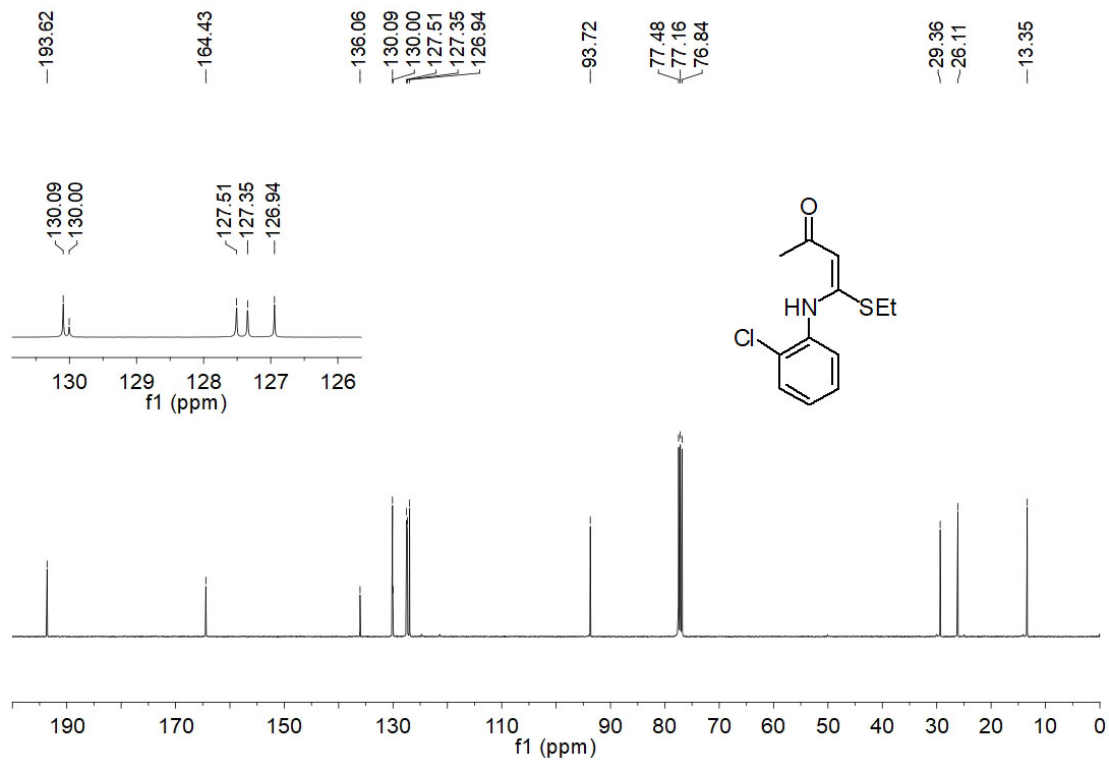


yu22071513, CDCl₃, 400



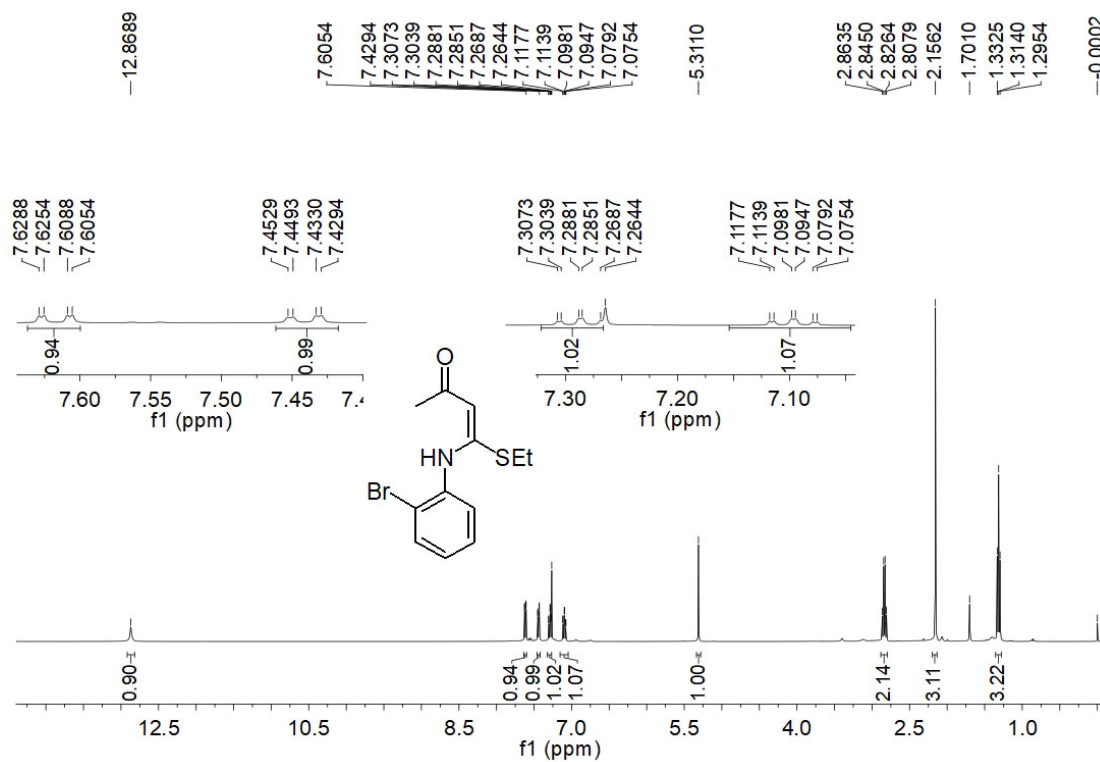
¹H NMR of **1b**

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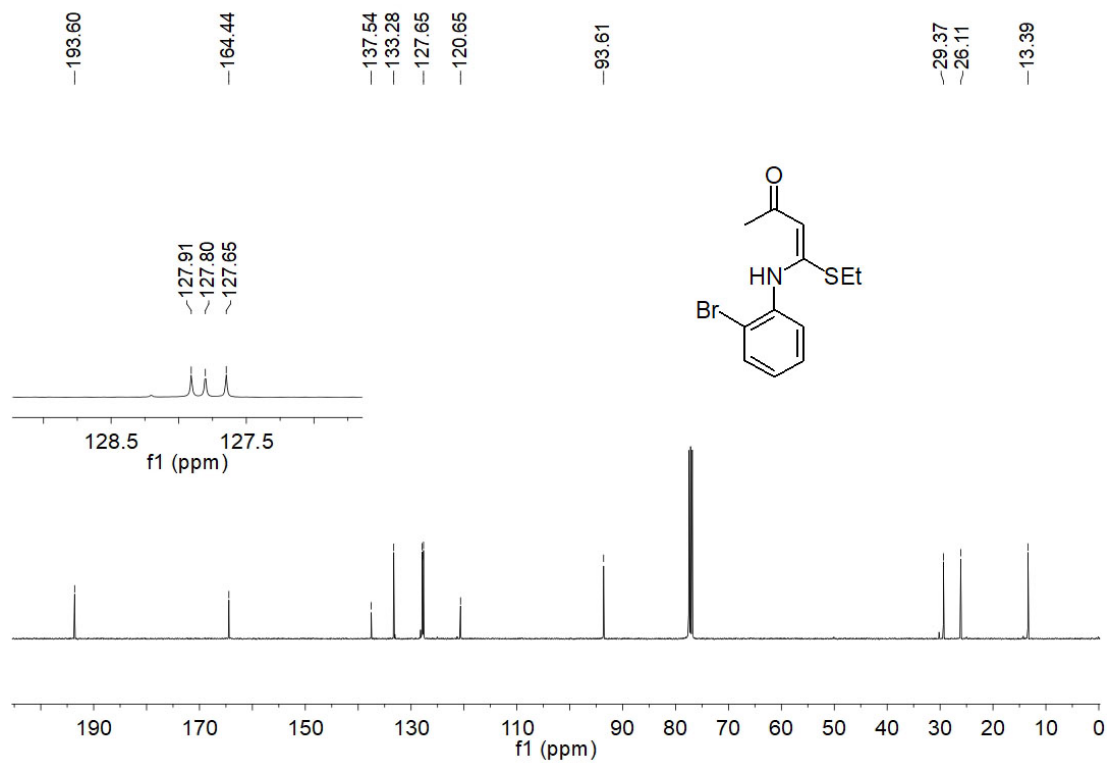
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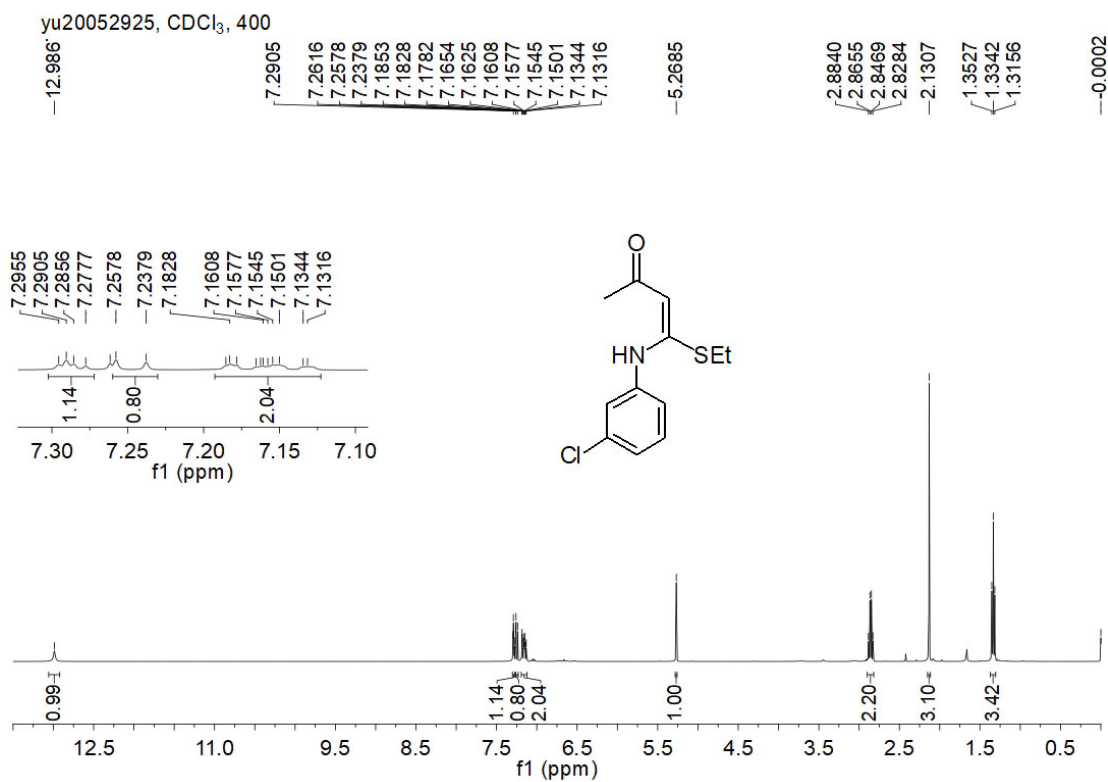


¹HNMR of **1c**

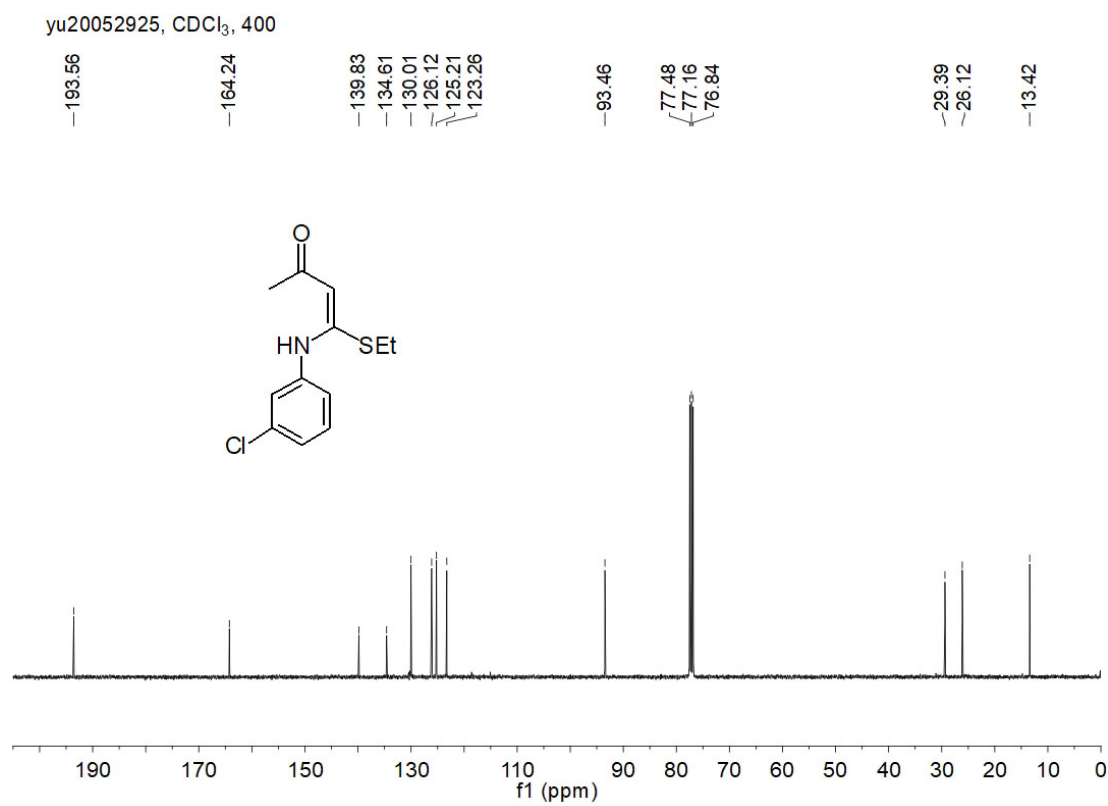
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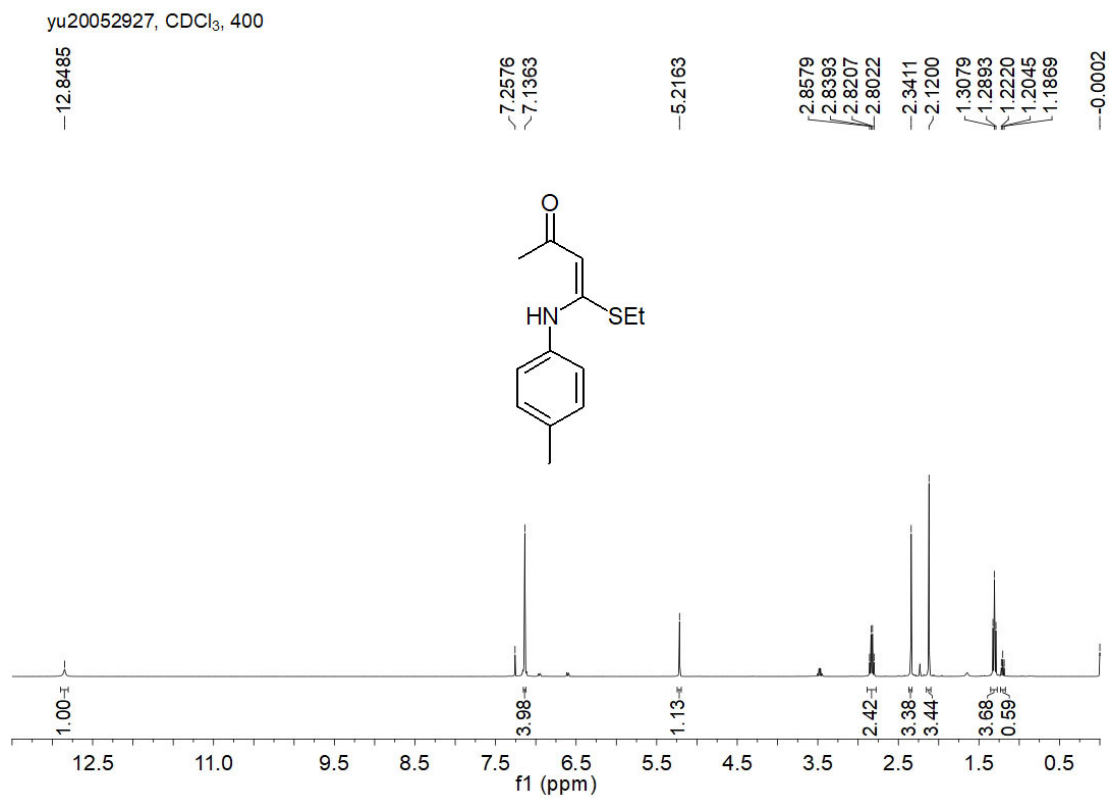
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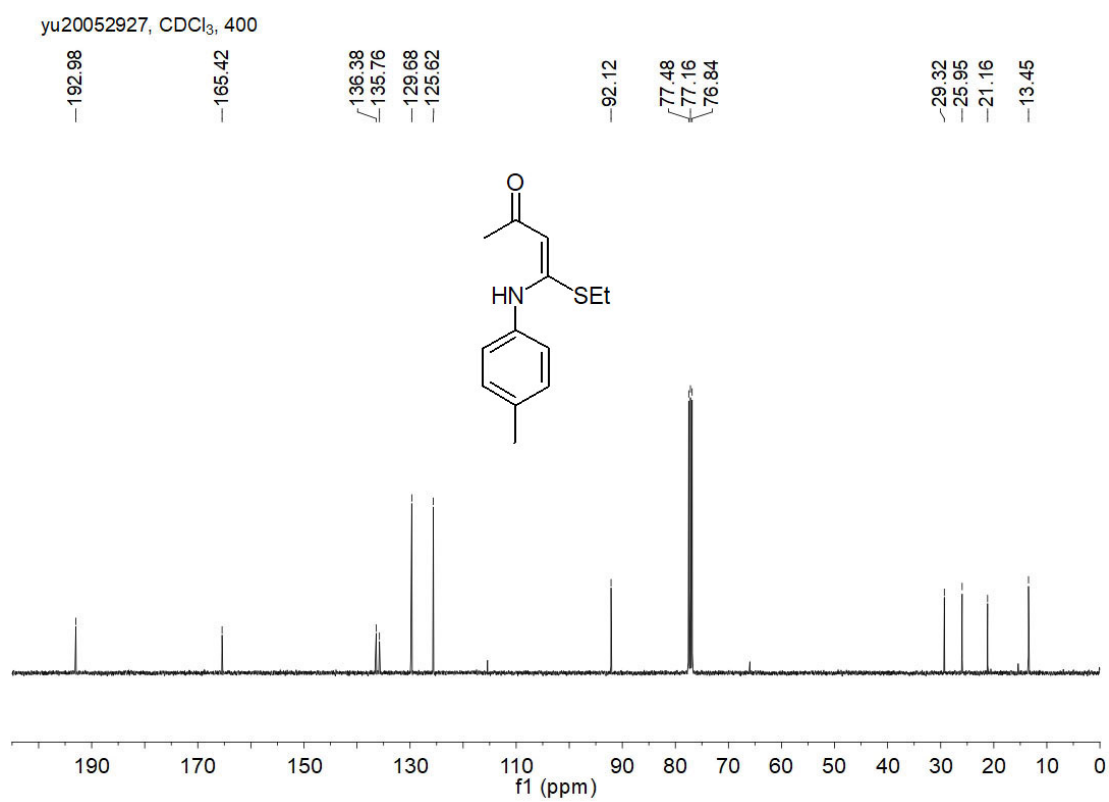
¹HNMR of **1d**



¹³CNMR of **1d**

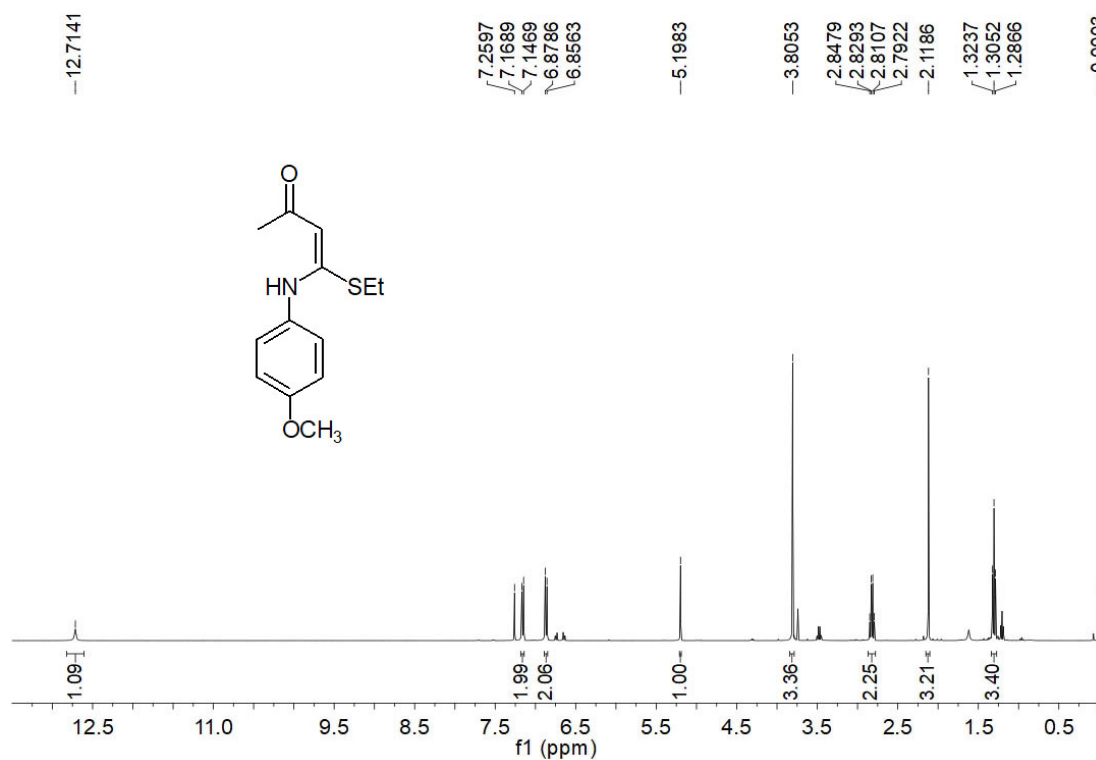


¹HNMR of **1e**



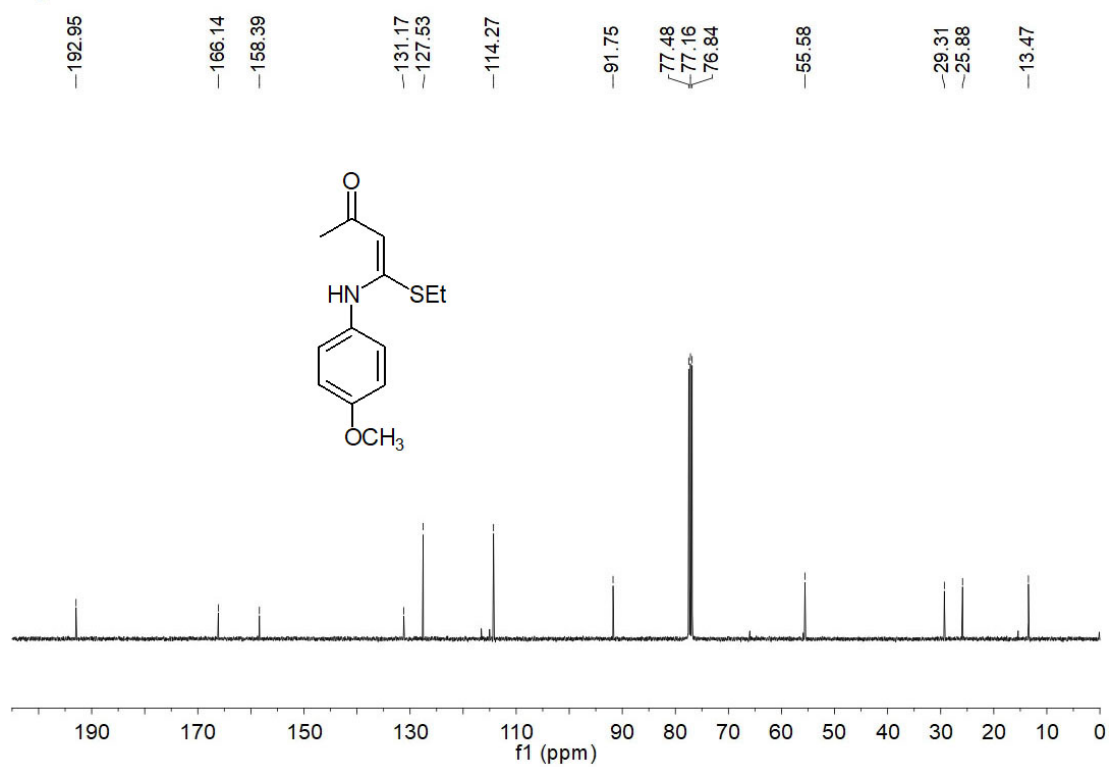
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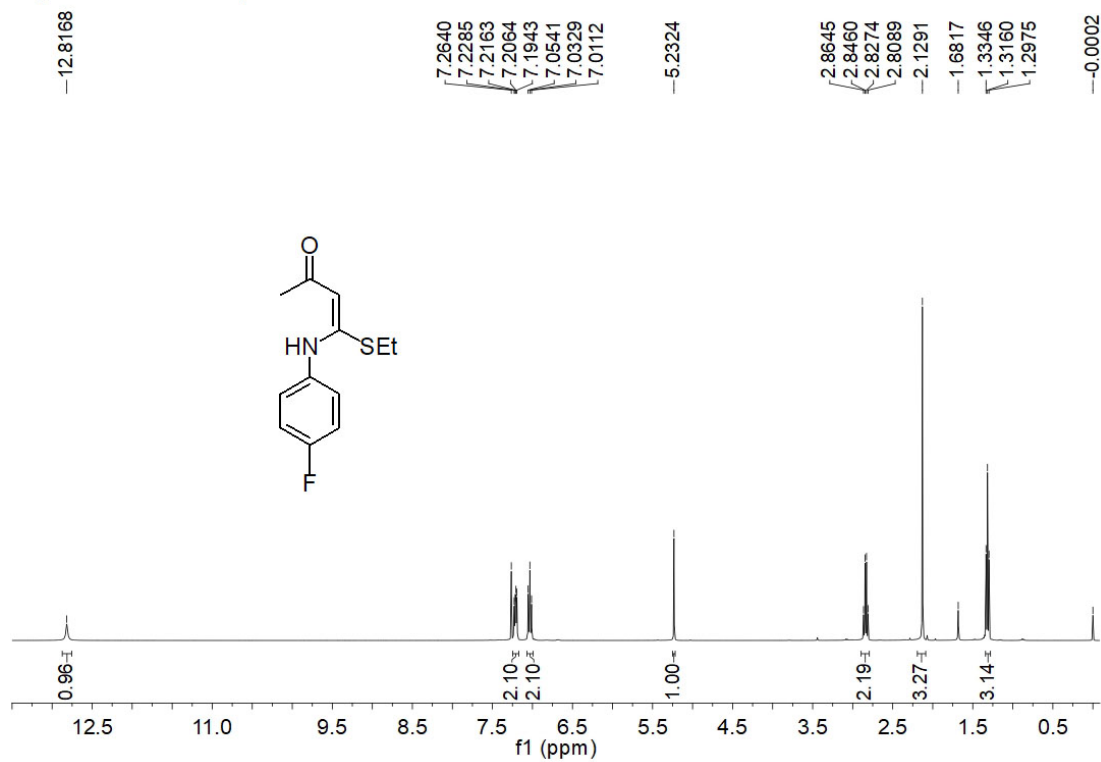
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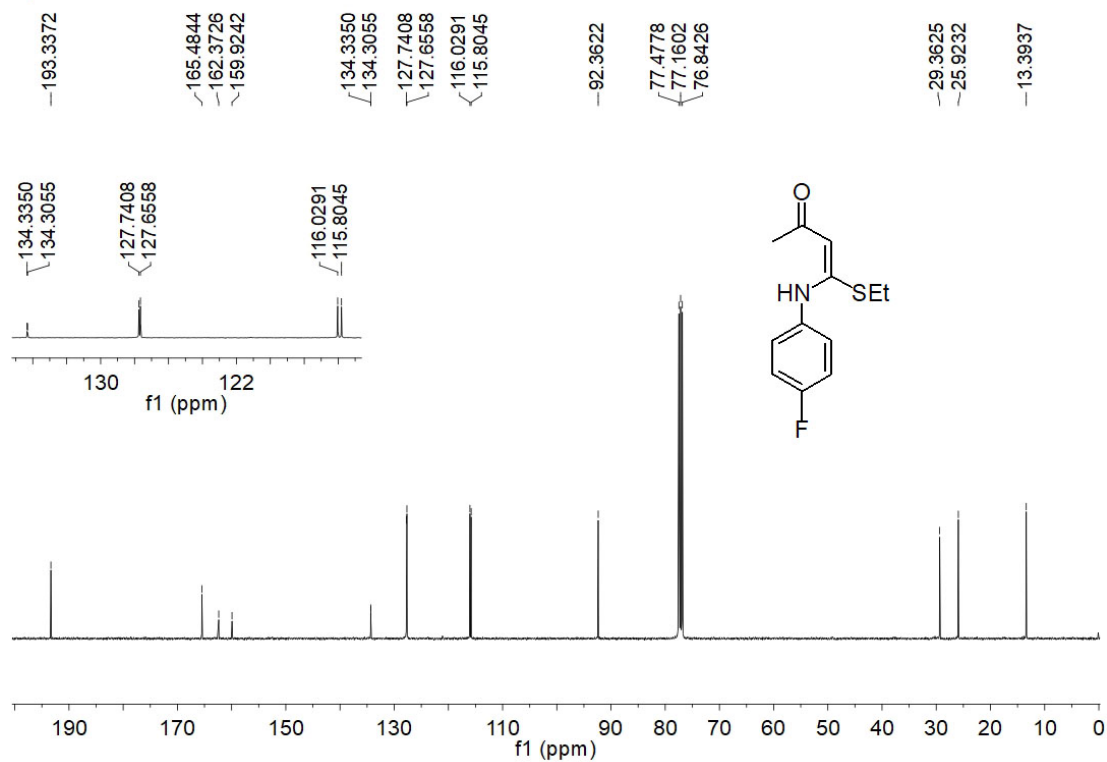
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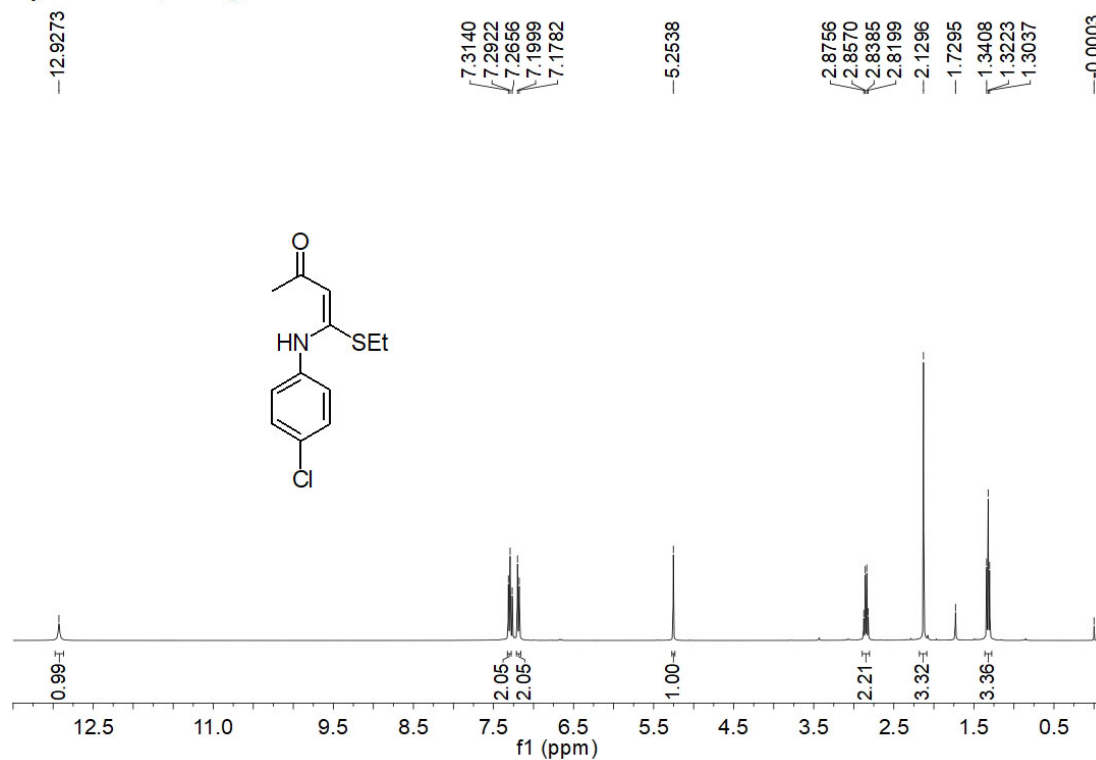
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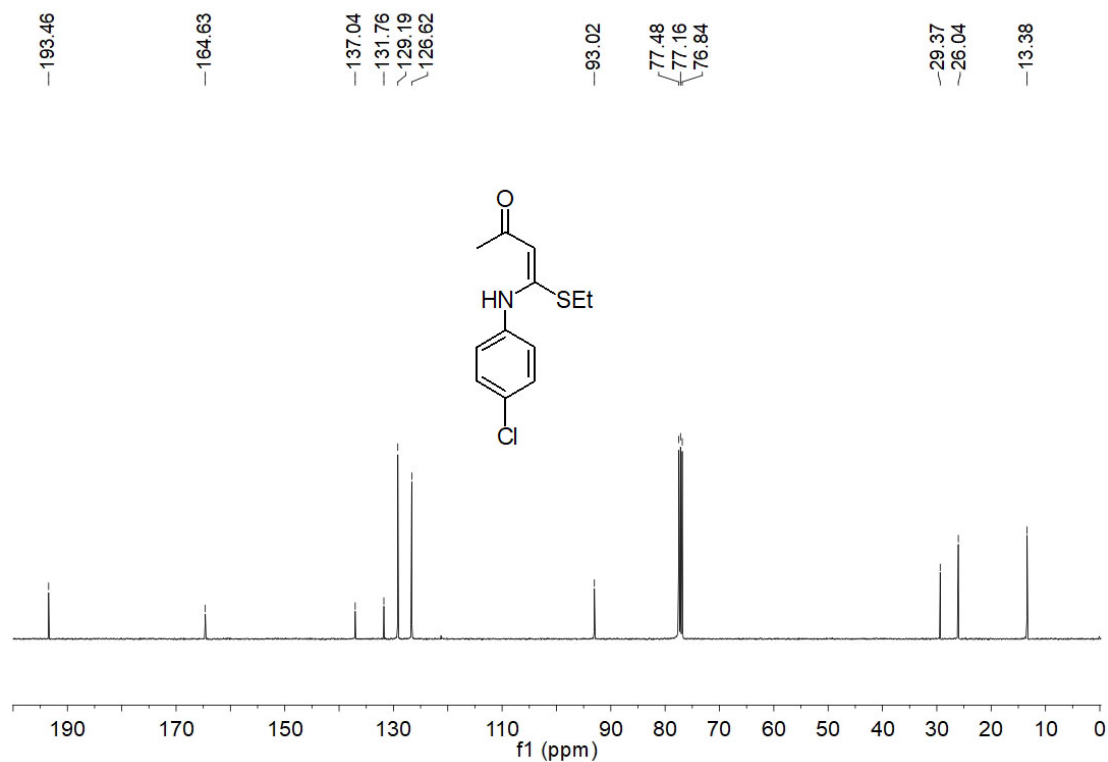
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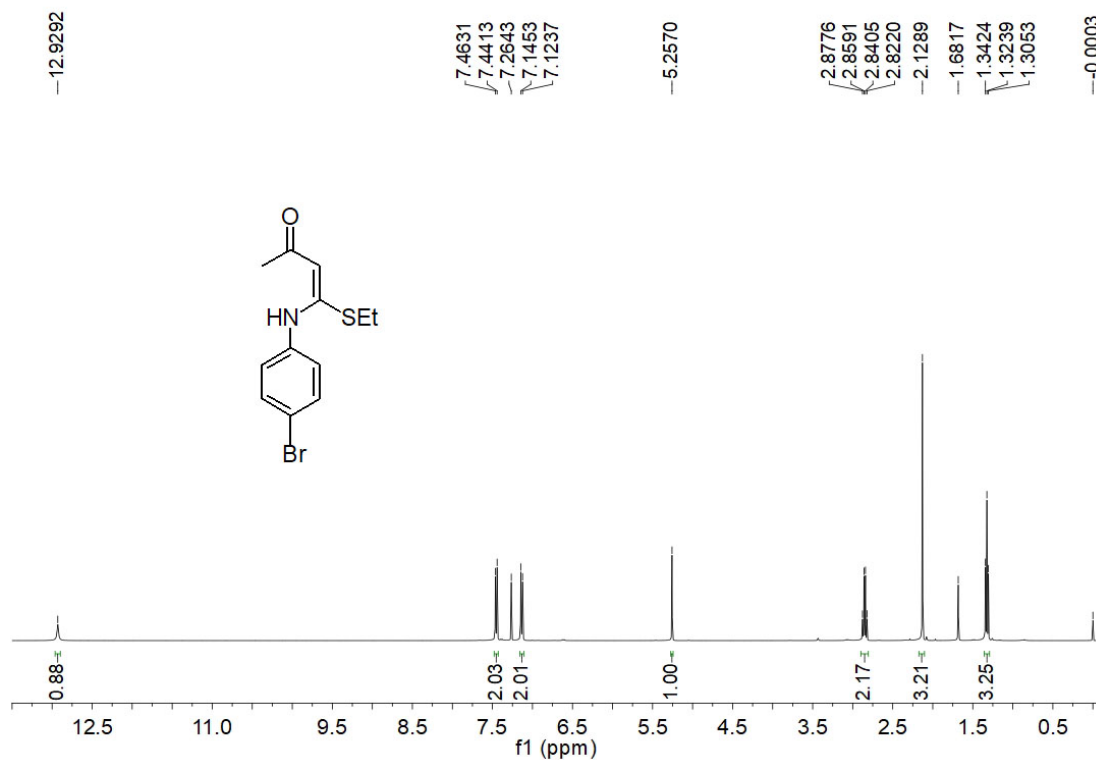
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yu22071514, CDCl₃, 400



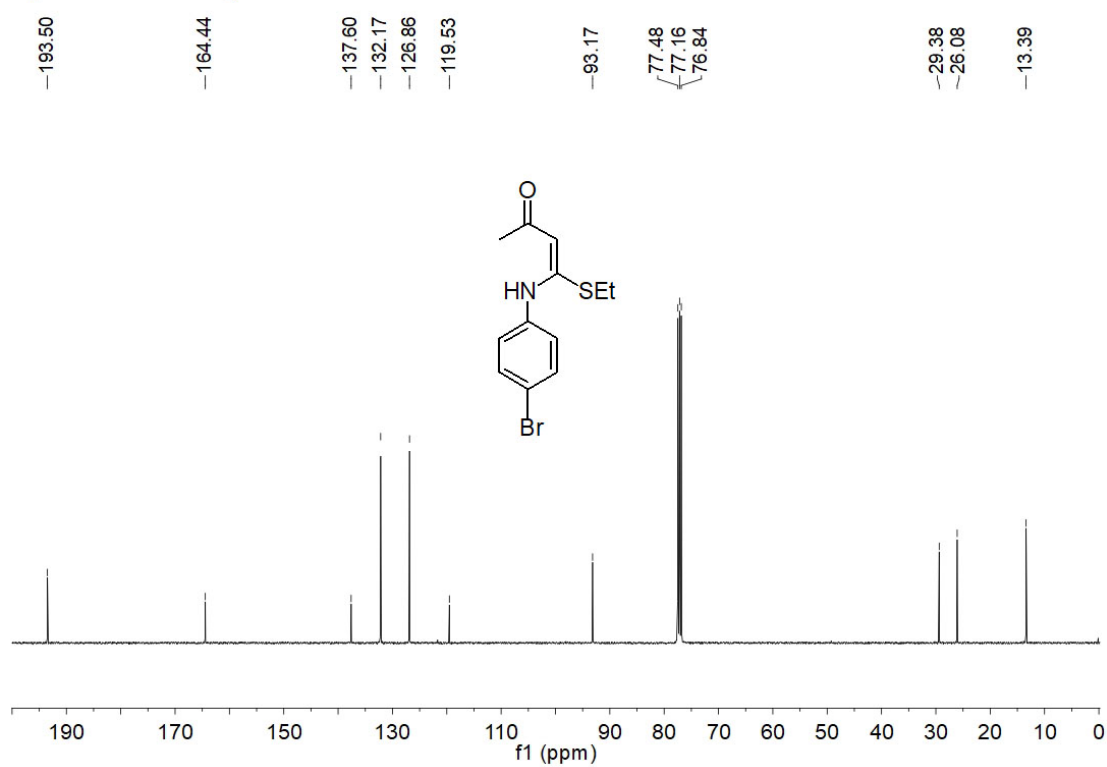
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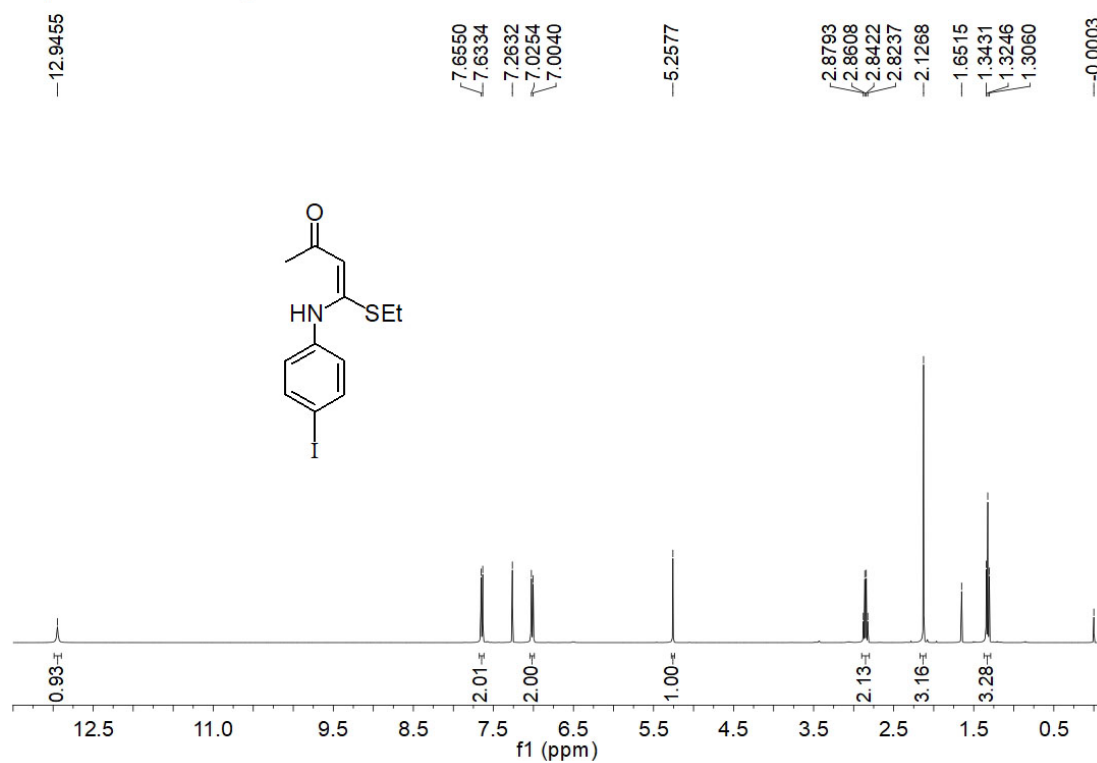
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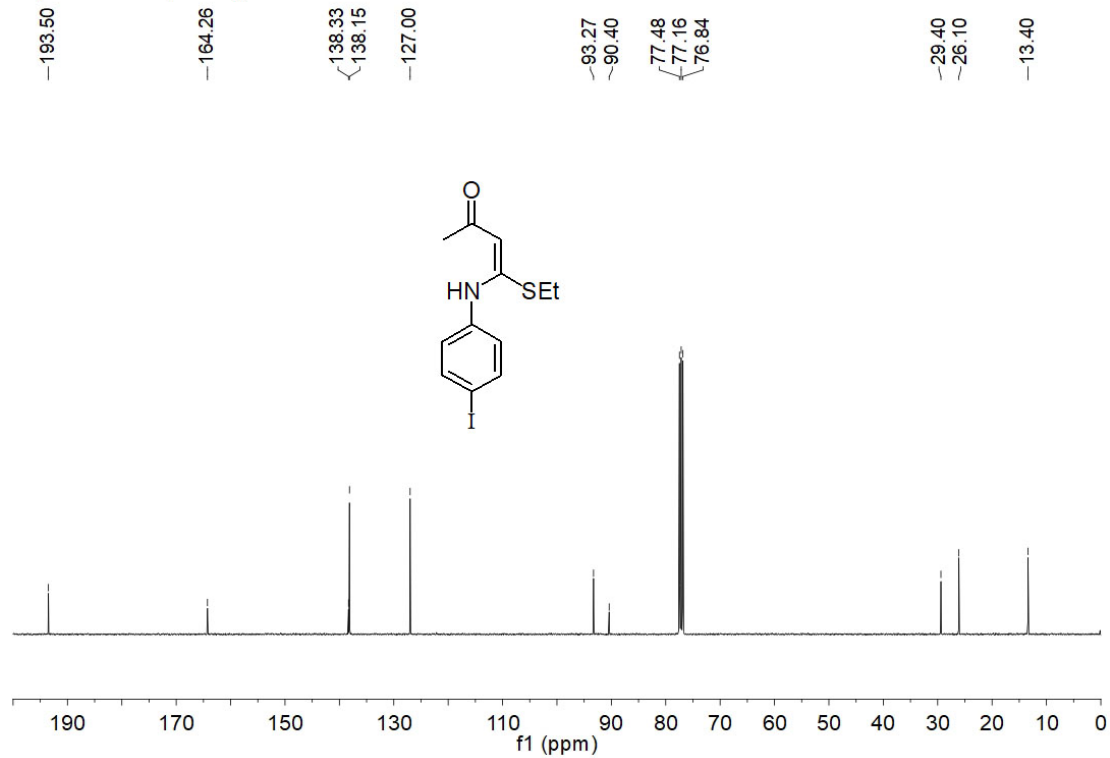
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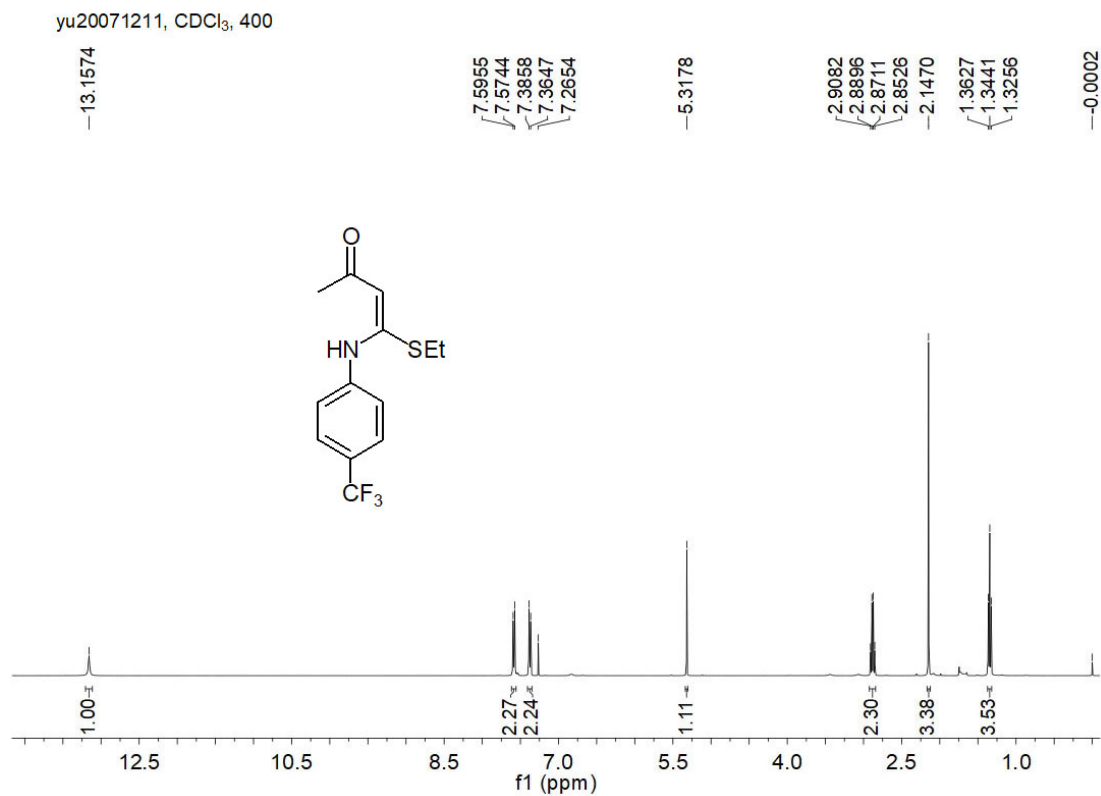


¹H NMR of **1j**

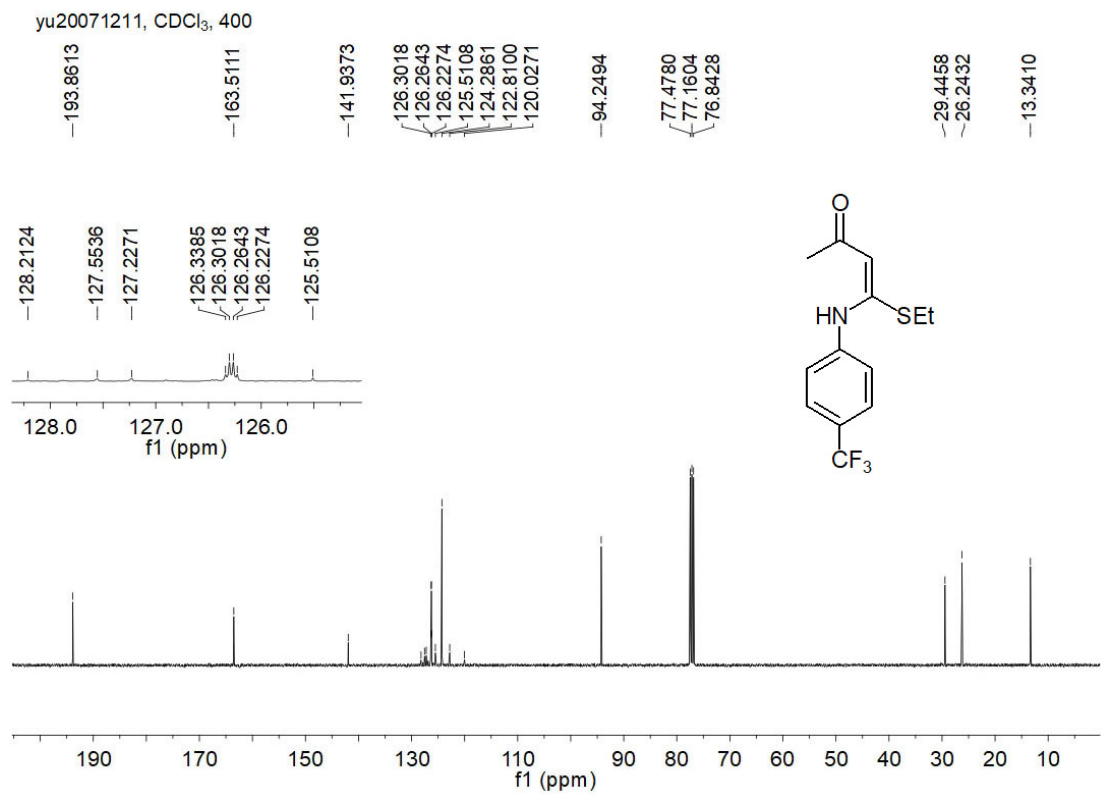
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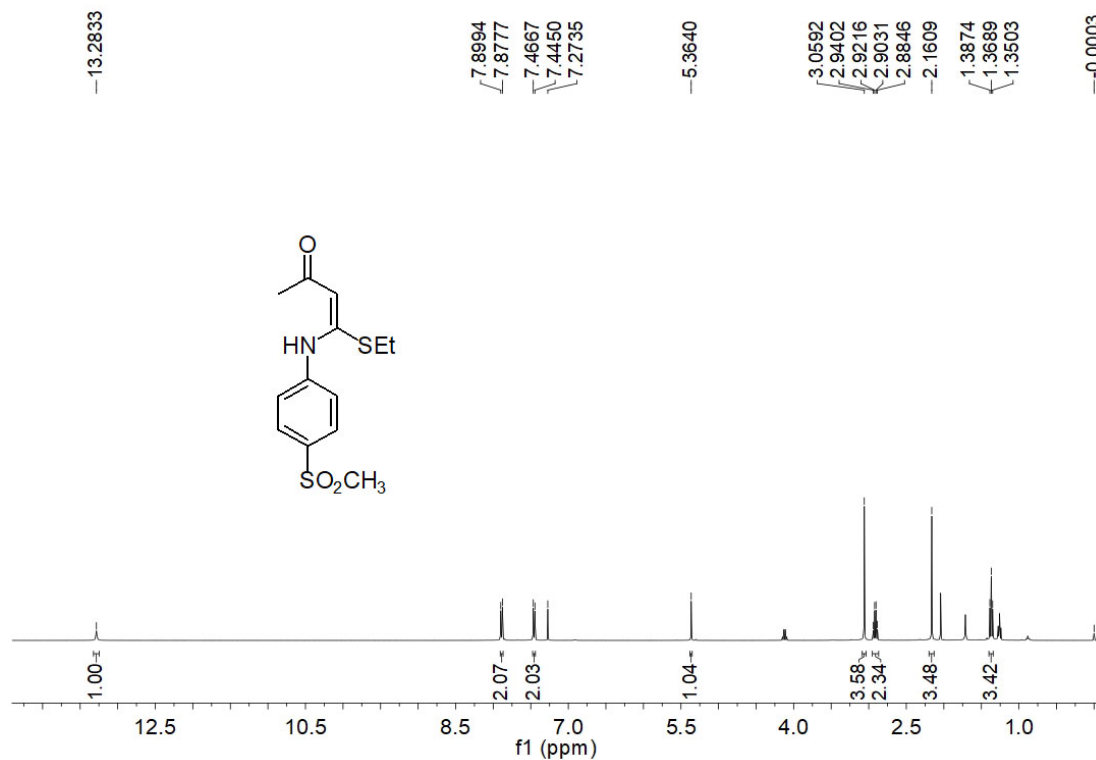


¹HNMR of **1k**



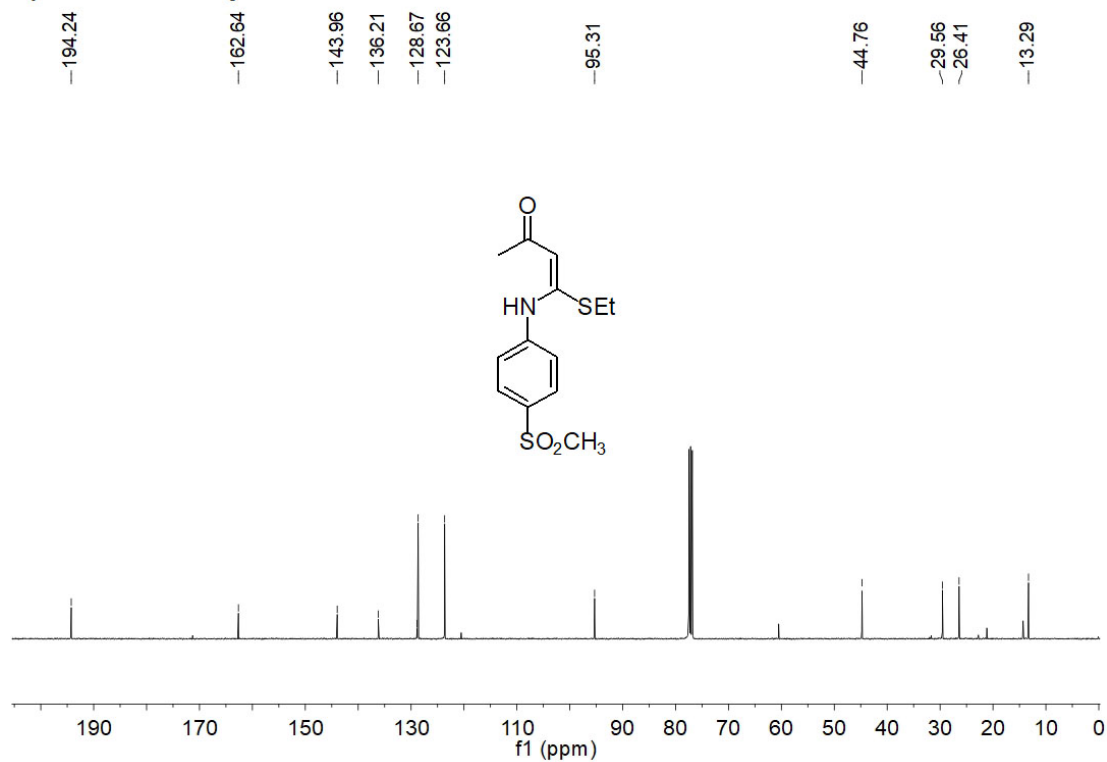
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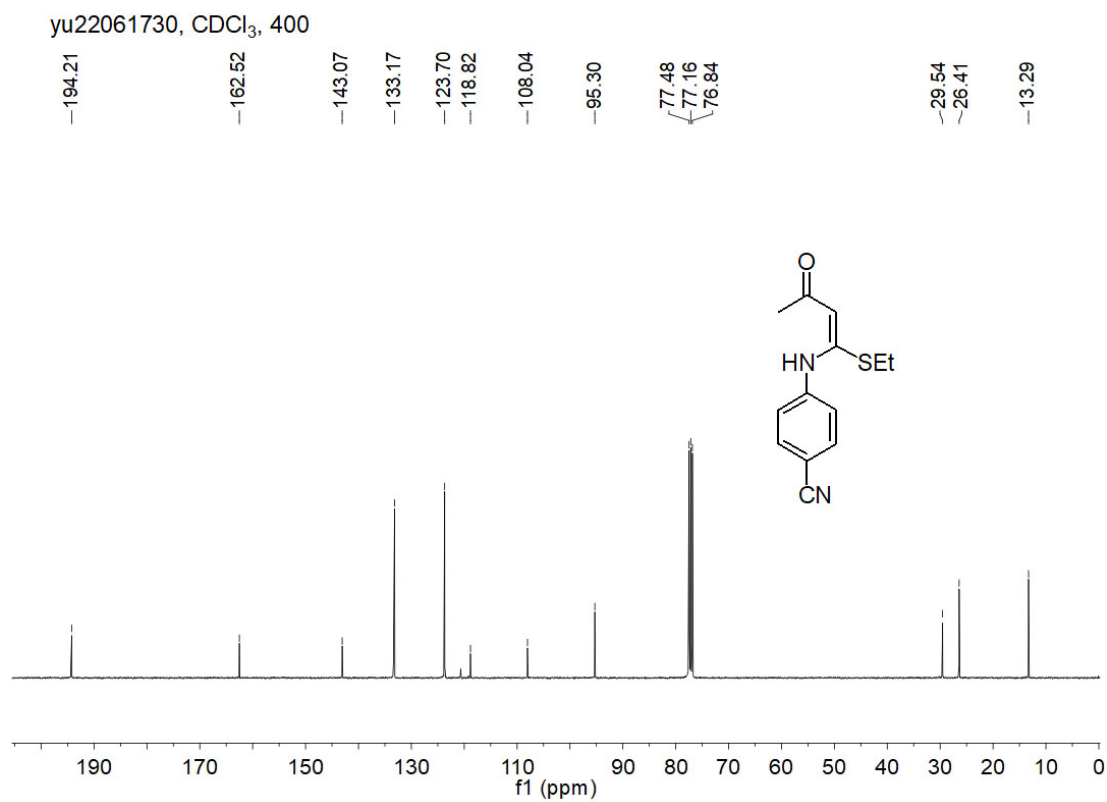
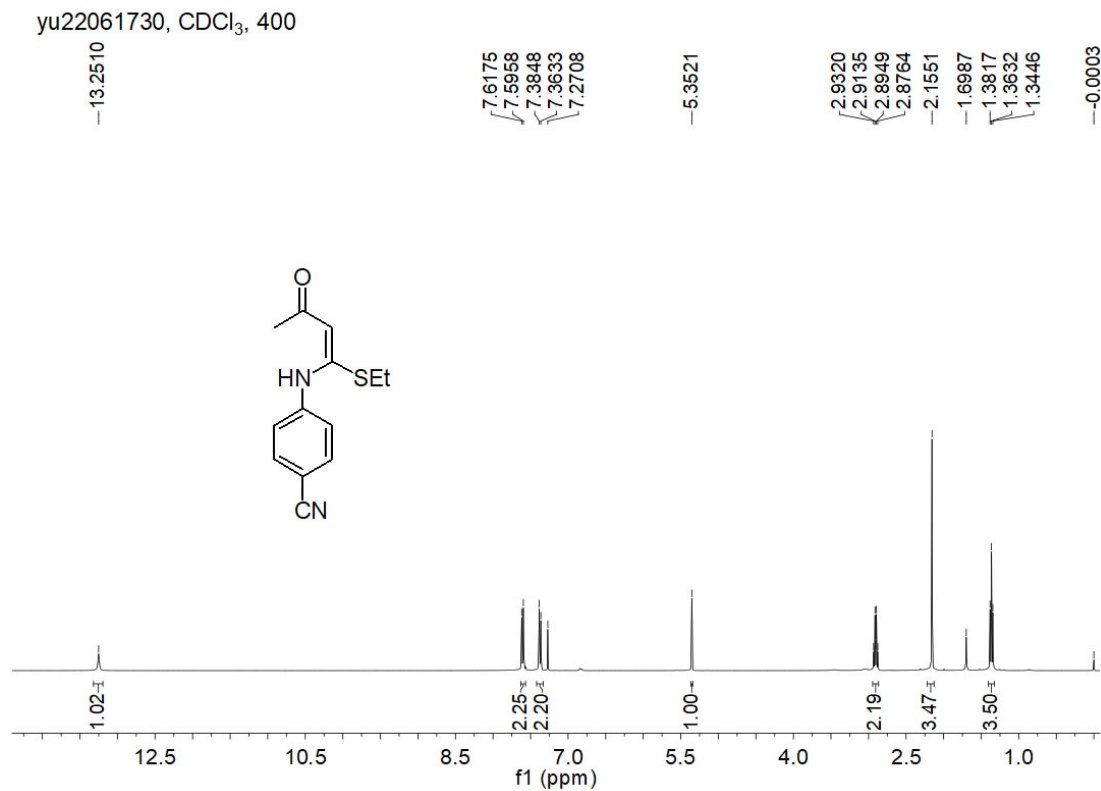


¹HNMR of **11**

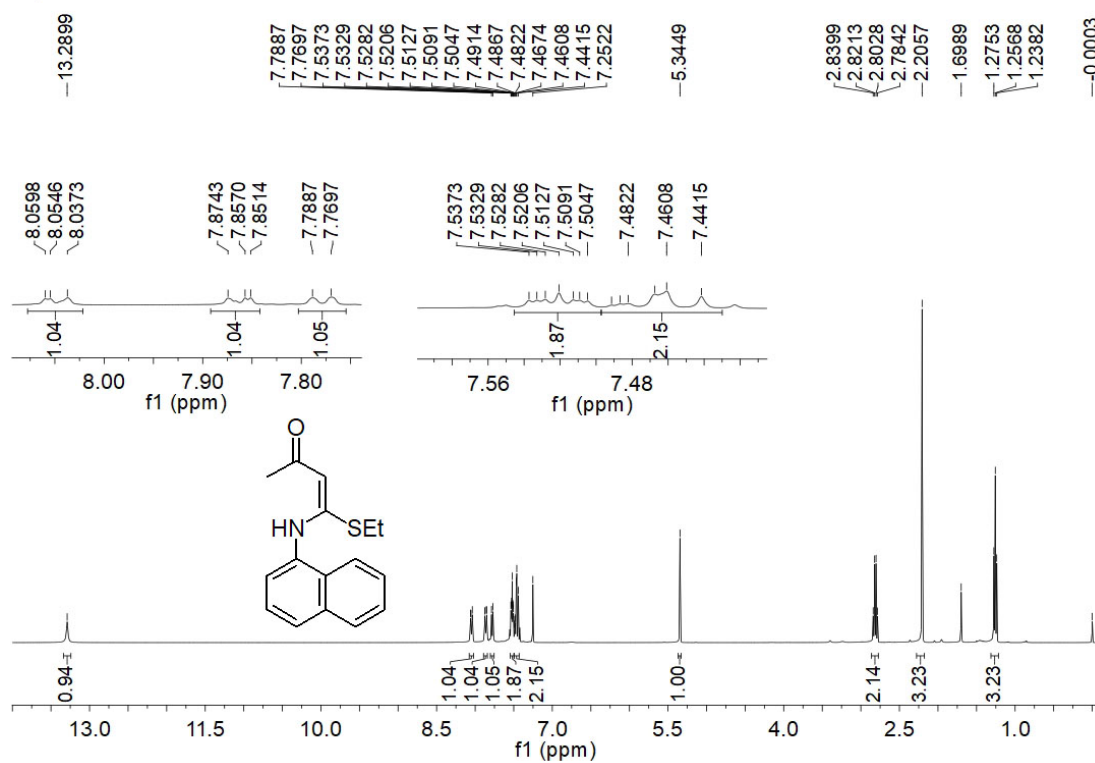
yu22061732, CDCl₃, 400



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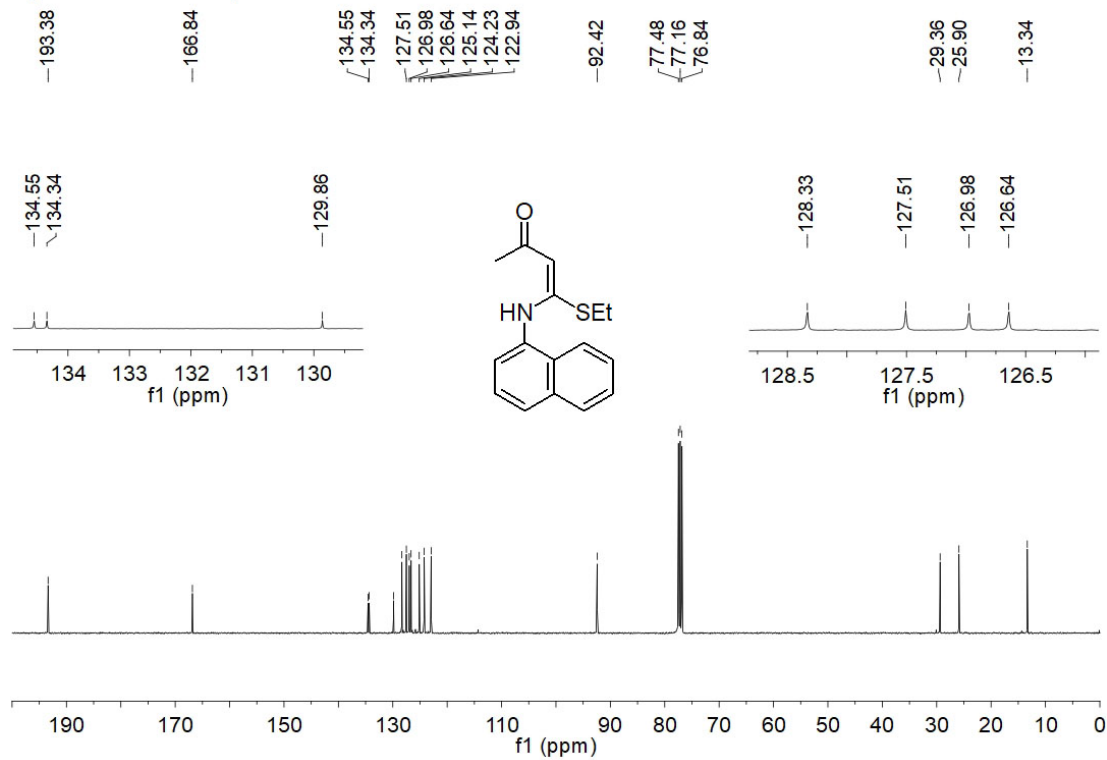


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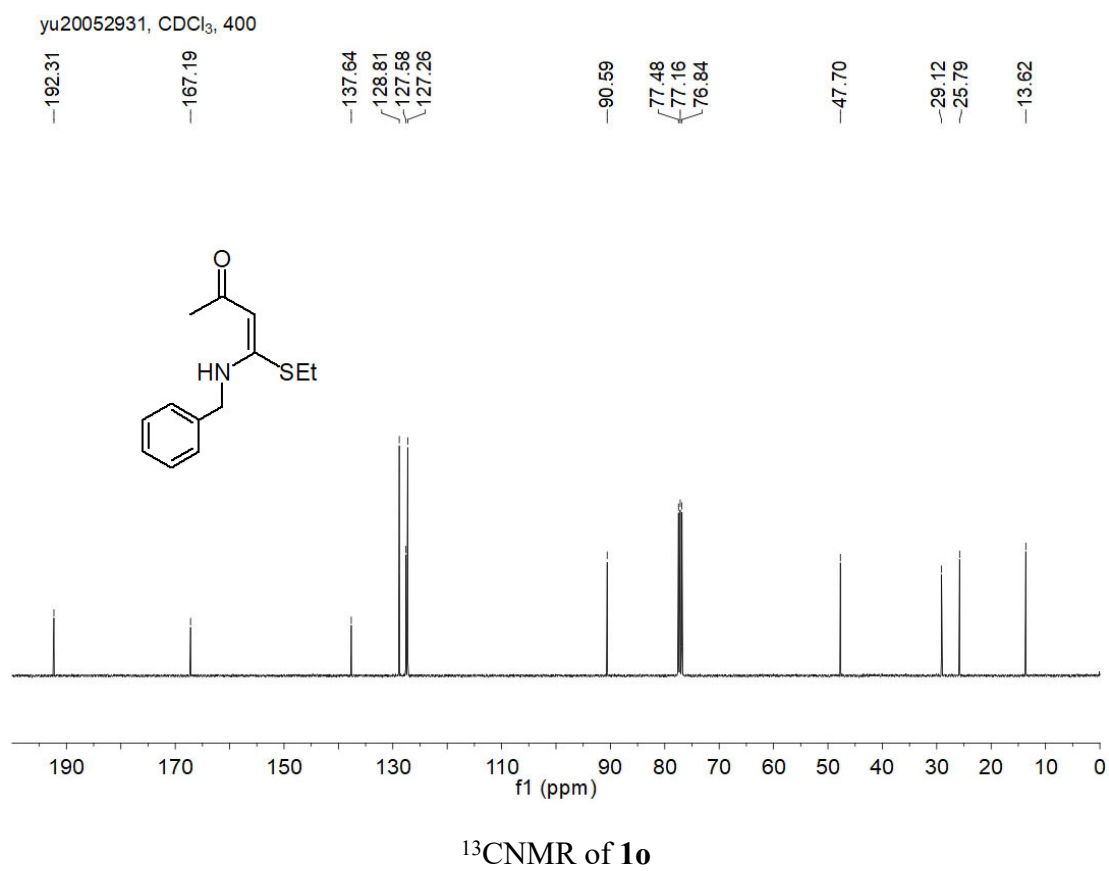
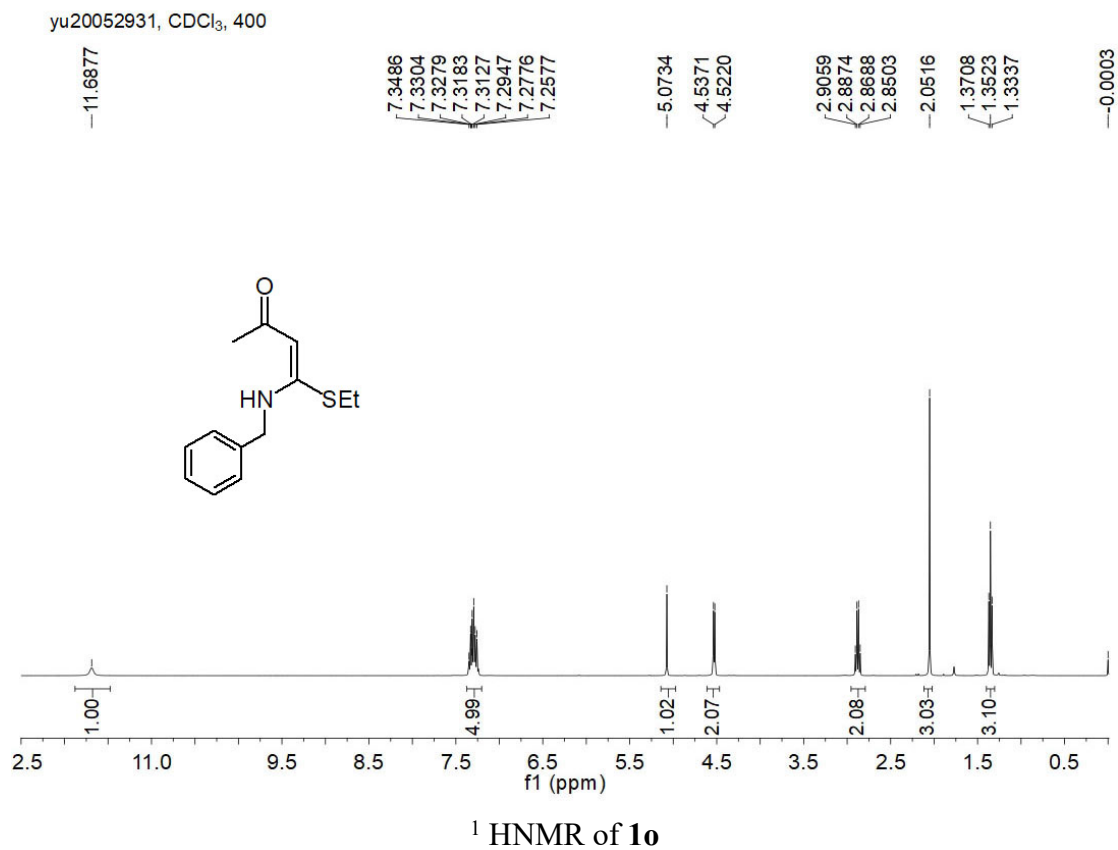


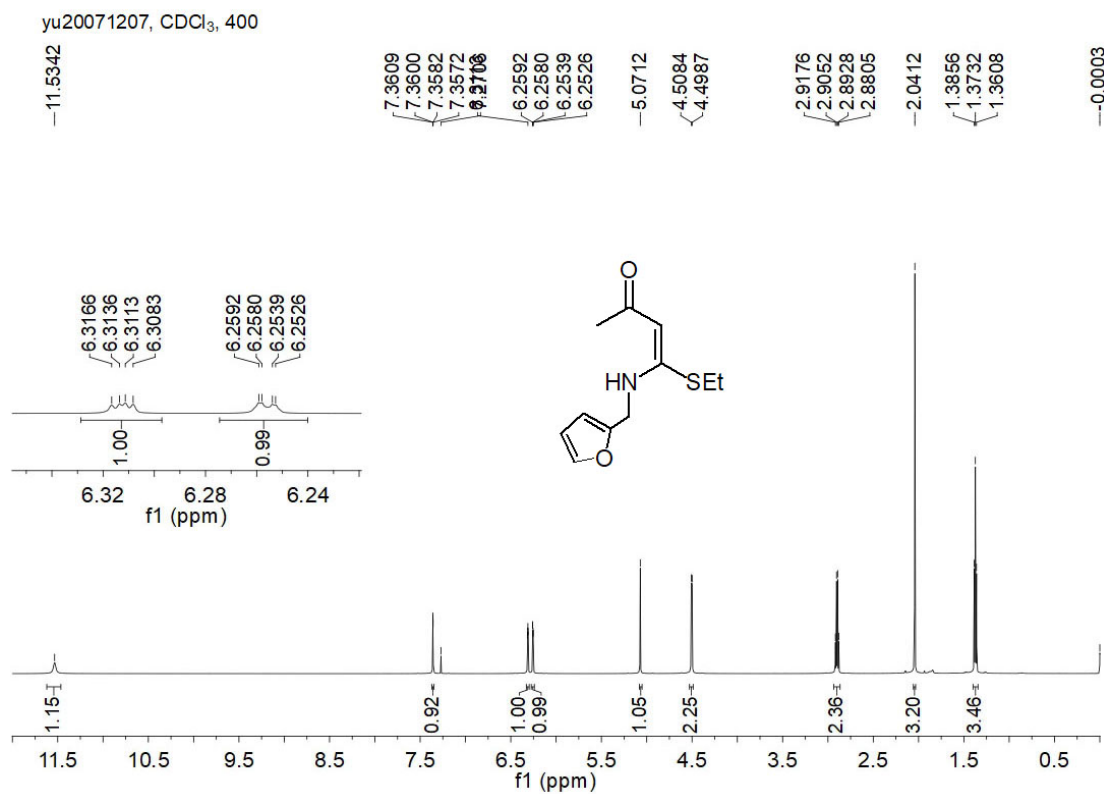
¹HNMR of **1n**

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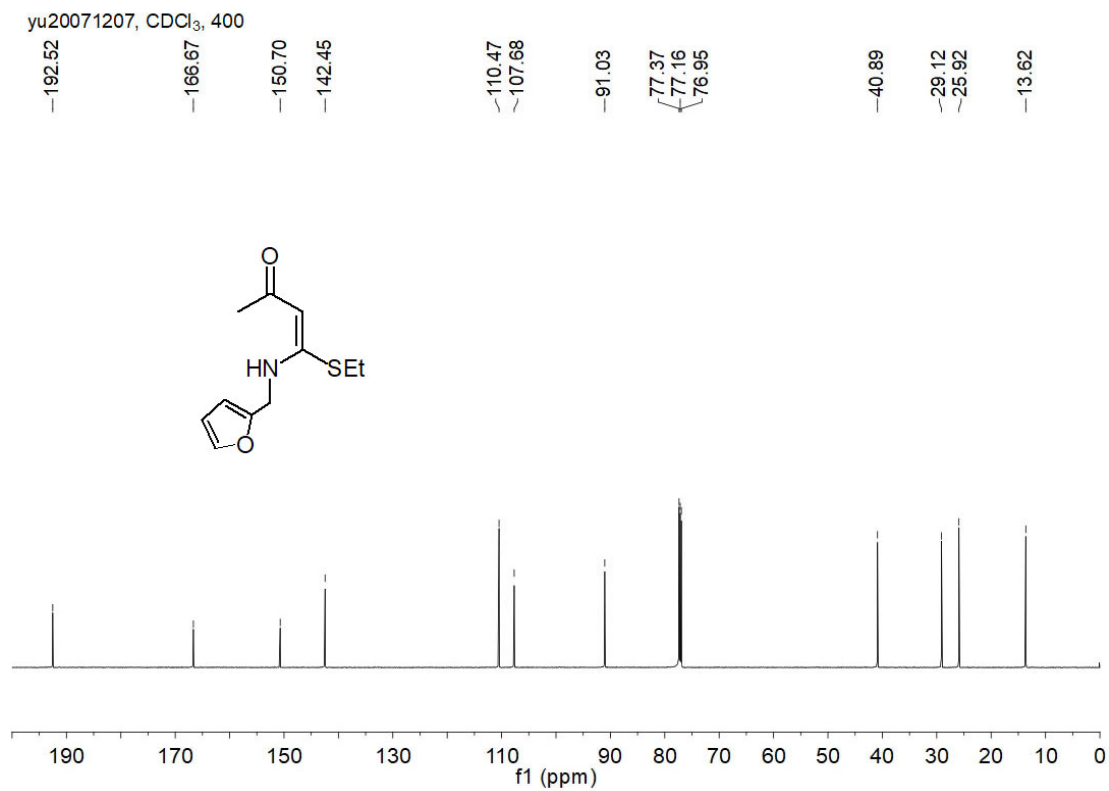


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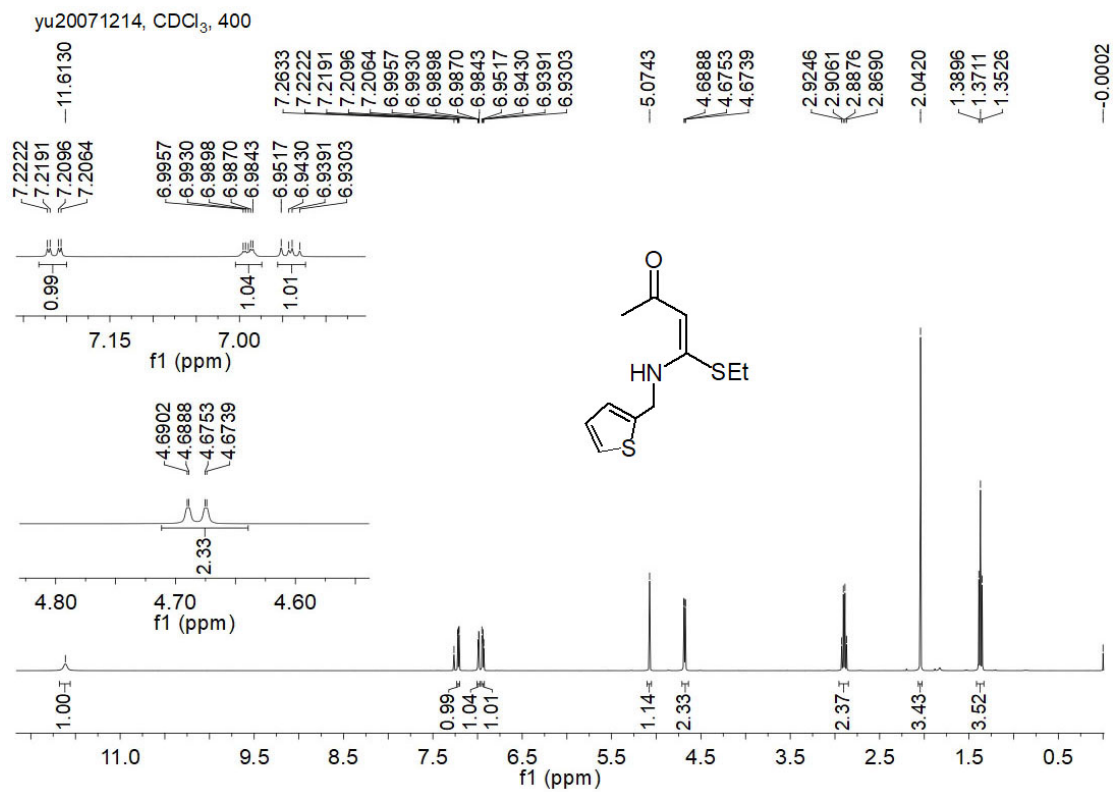




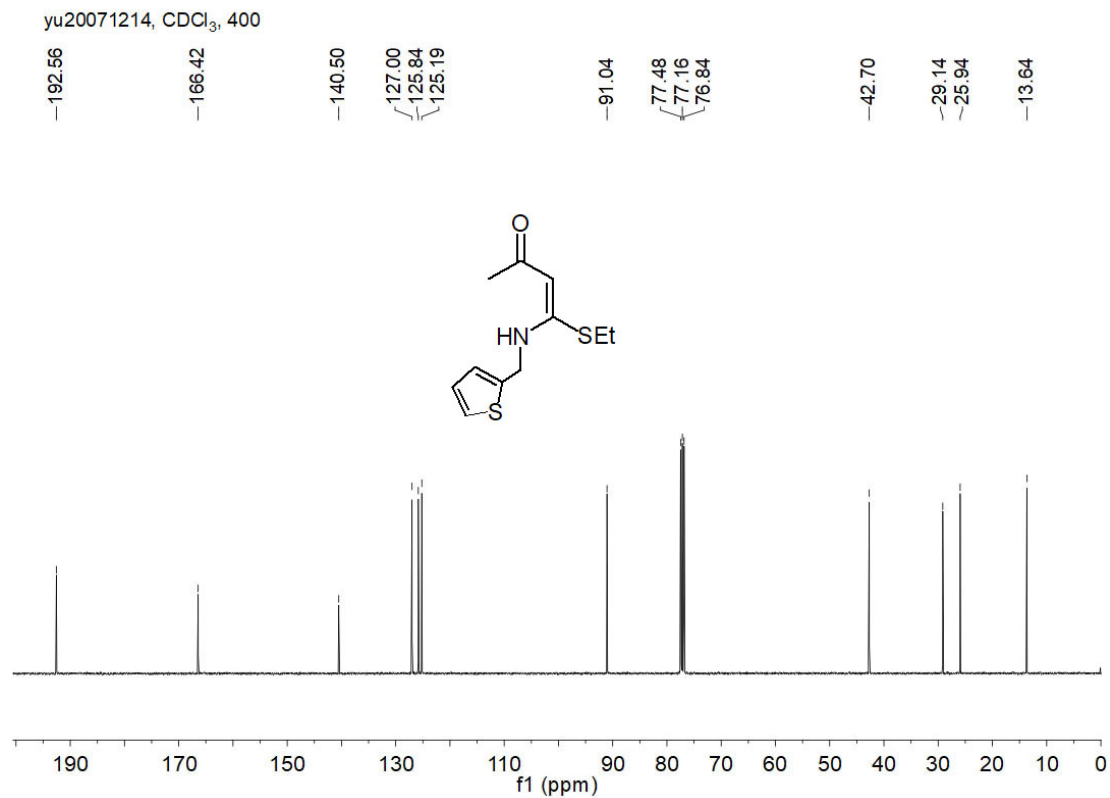
¹H NMR of **1p**



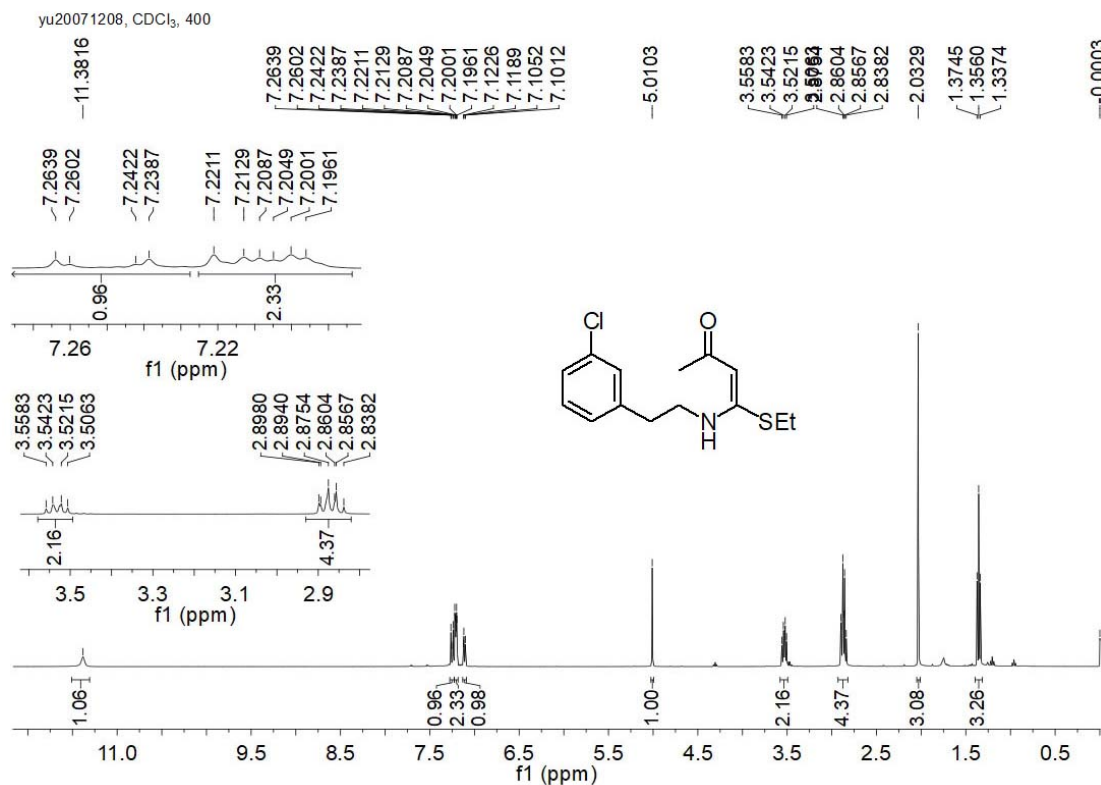
¹³C NMR of **1p**



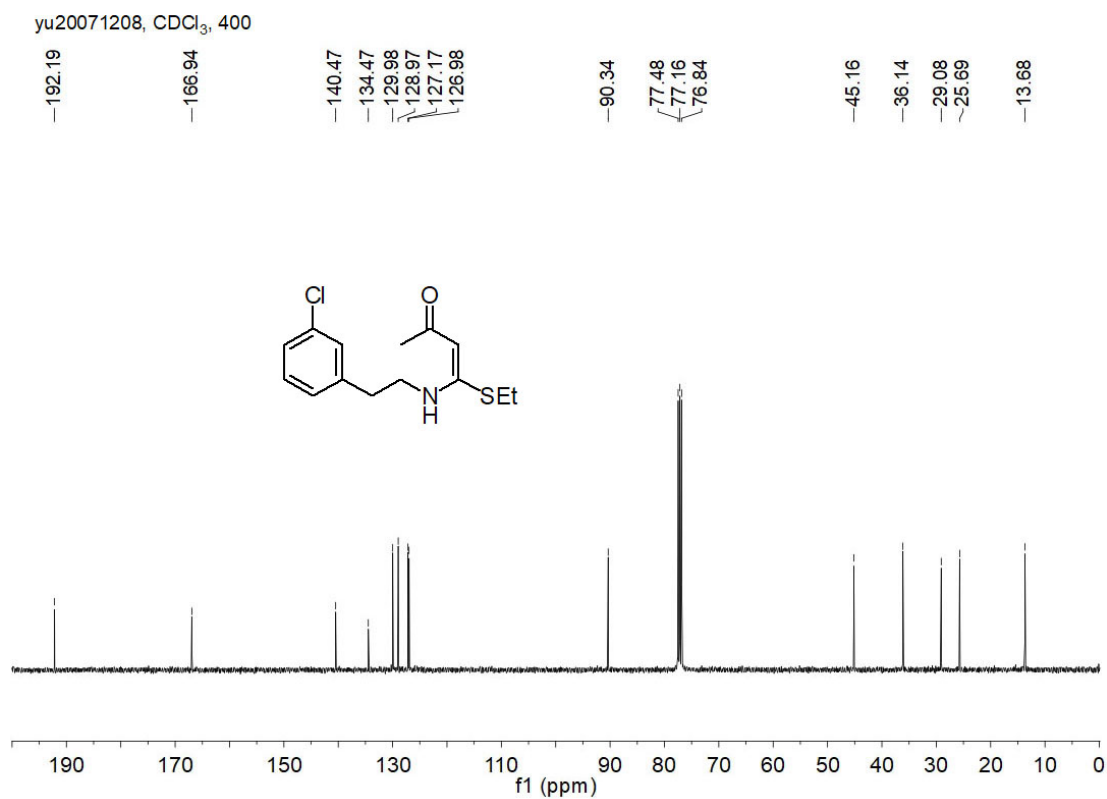
¹H NMR of **1q**



¹³C NMR of **1q**

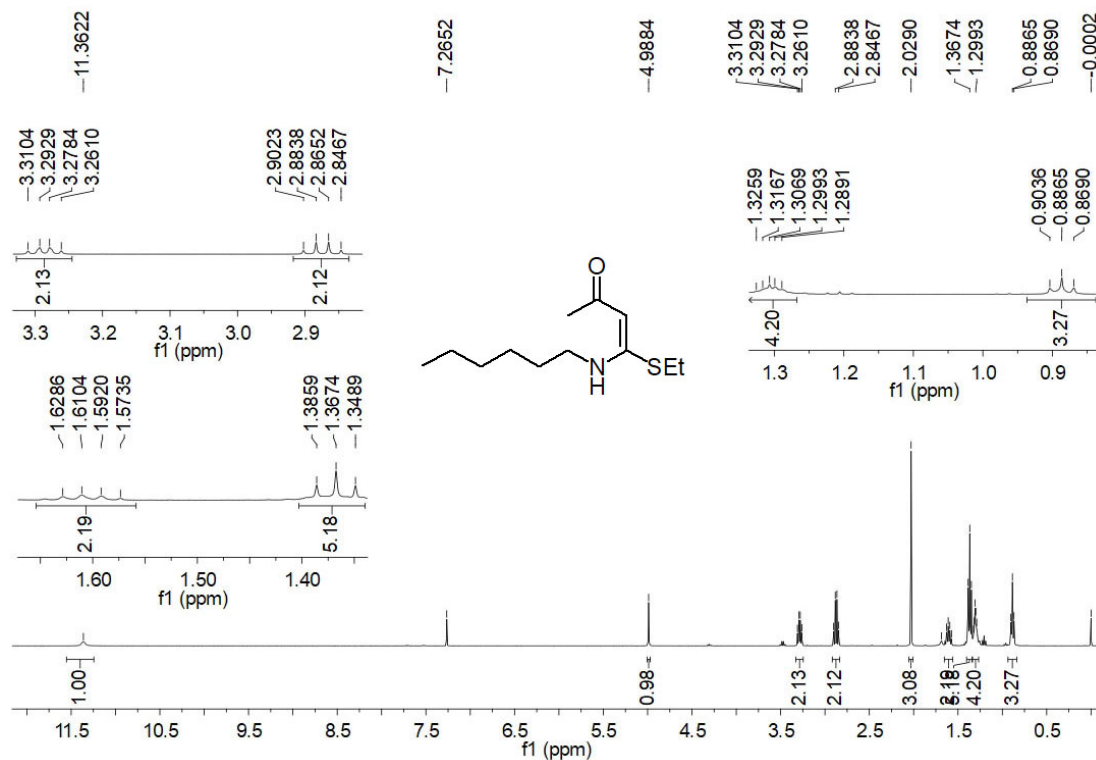


¹H NMR of **1r**



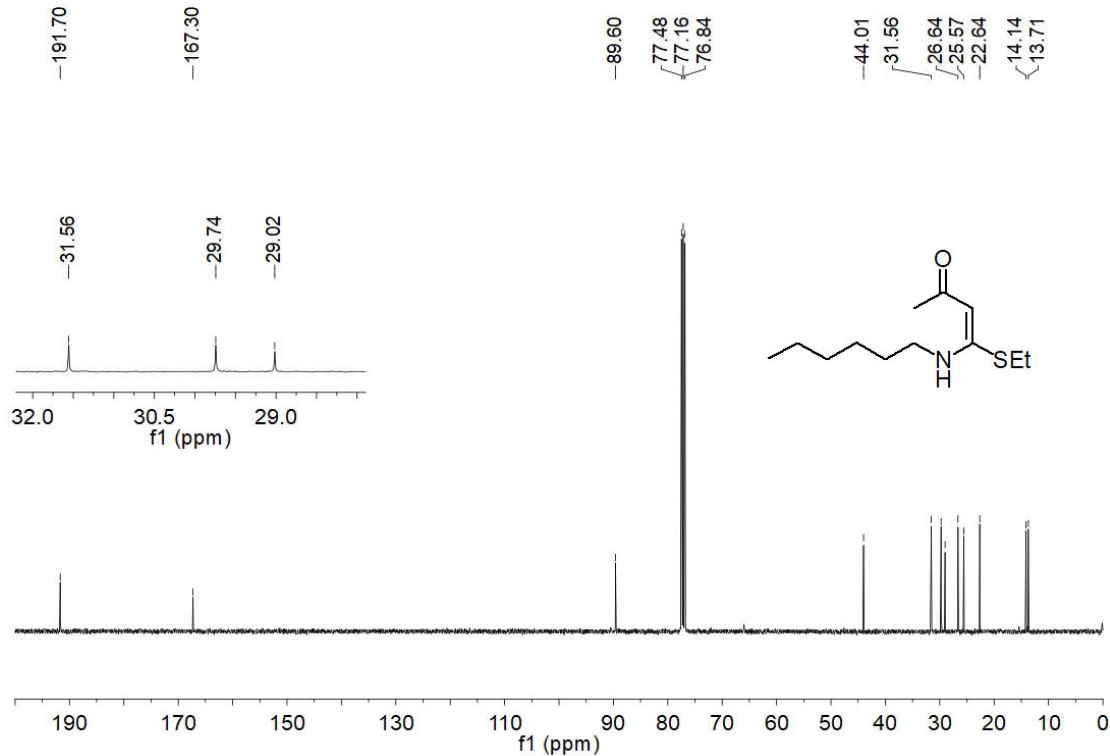
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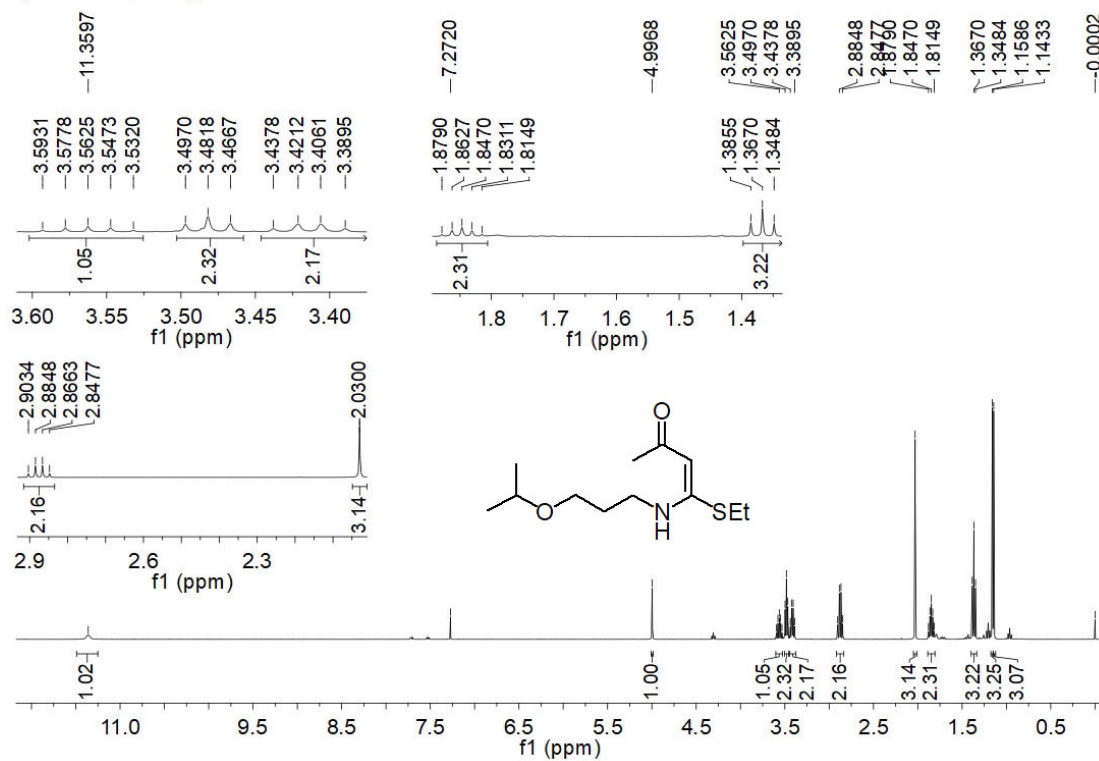
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yu20071213, CDCl₃, 400



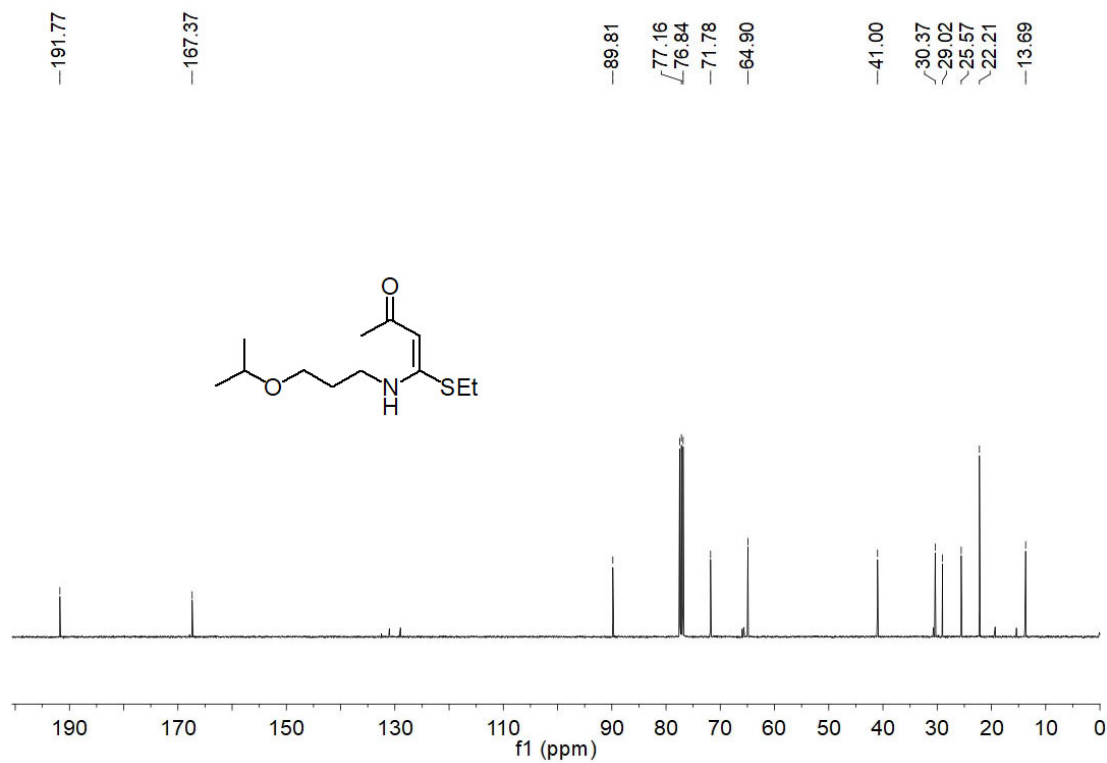
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yu20071212, CDCl₃, 400



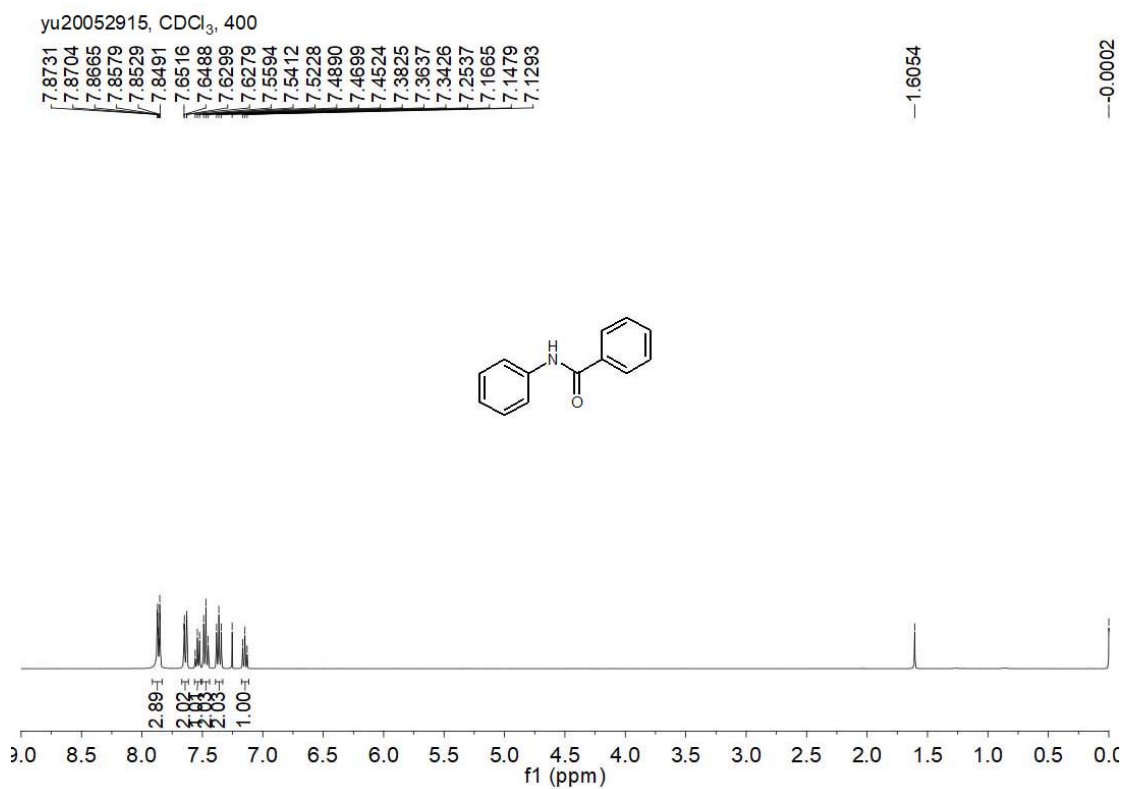
¹H NMR of 1t

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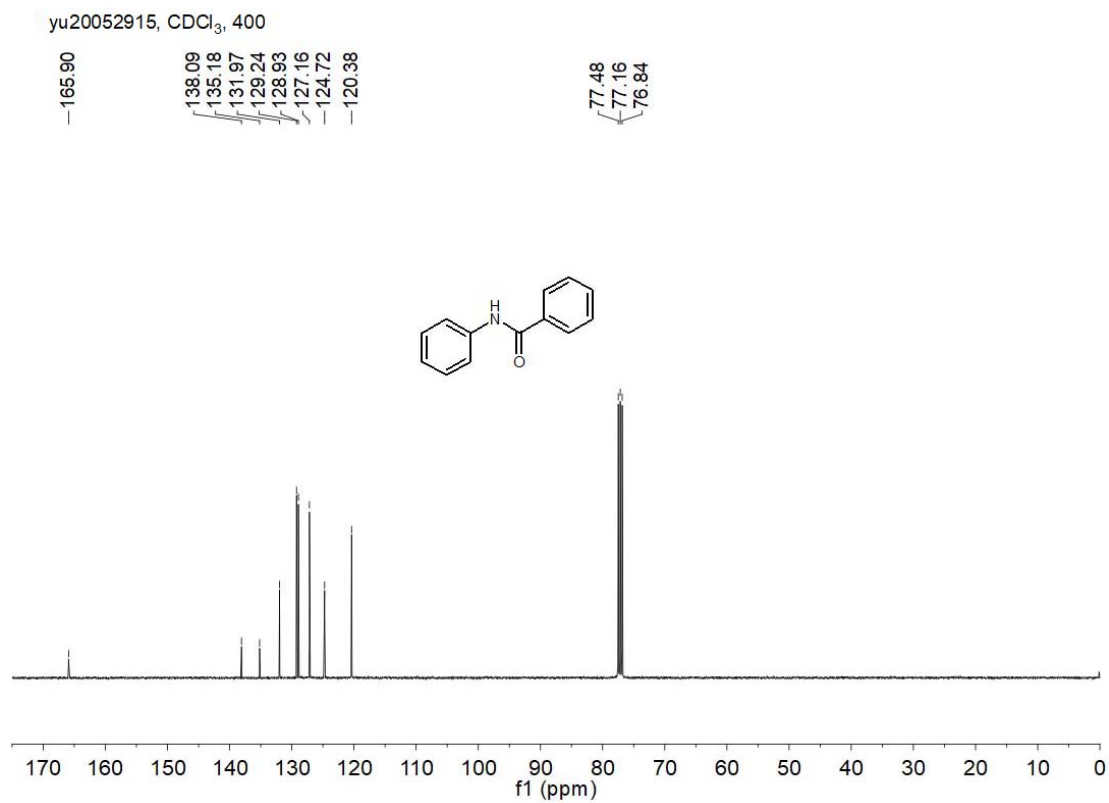


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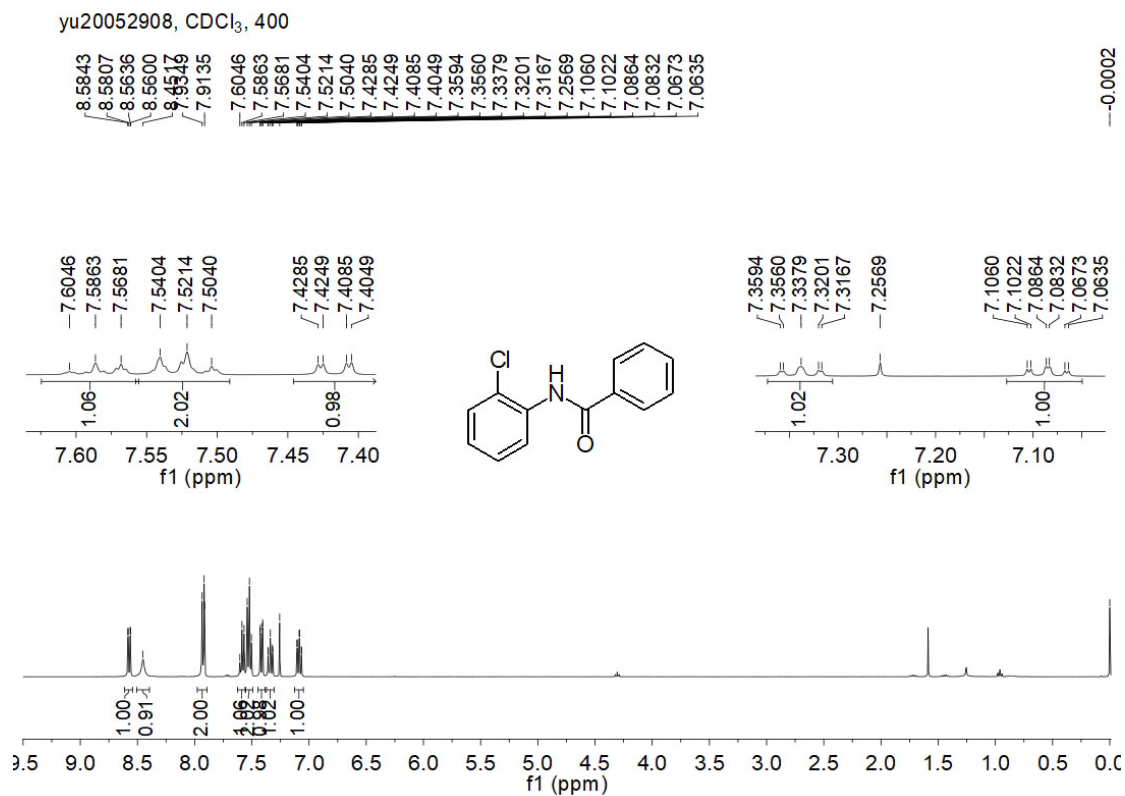
III. Copies of NMR spectra for 3



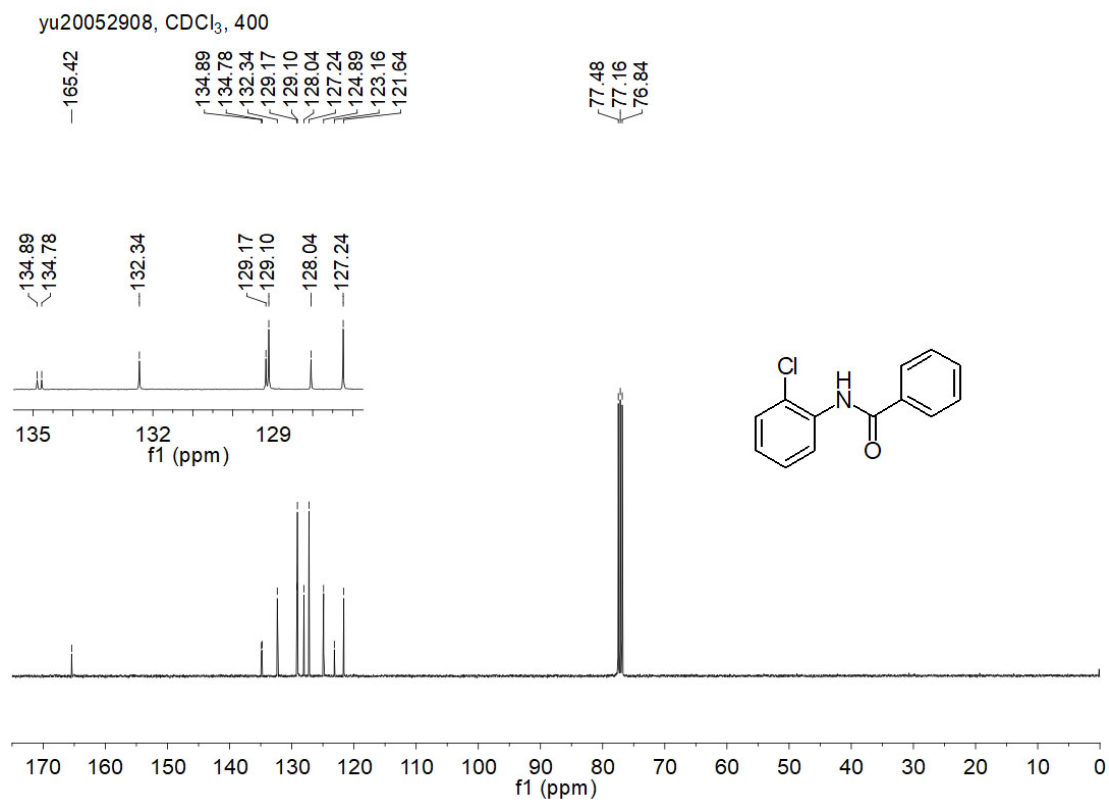
¹H NMR of 3a



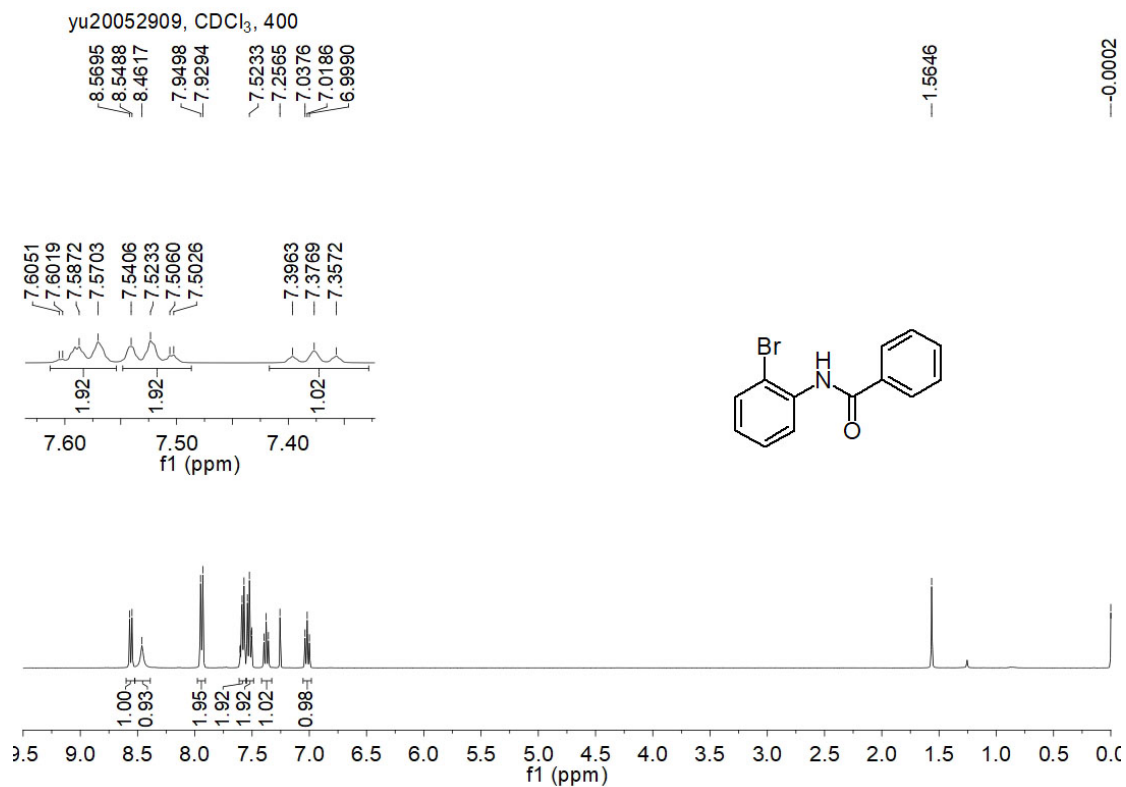
¹³C NMR of 3a



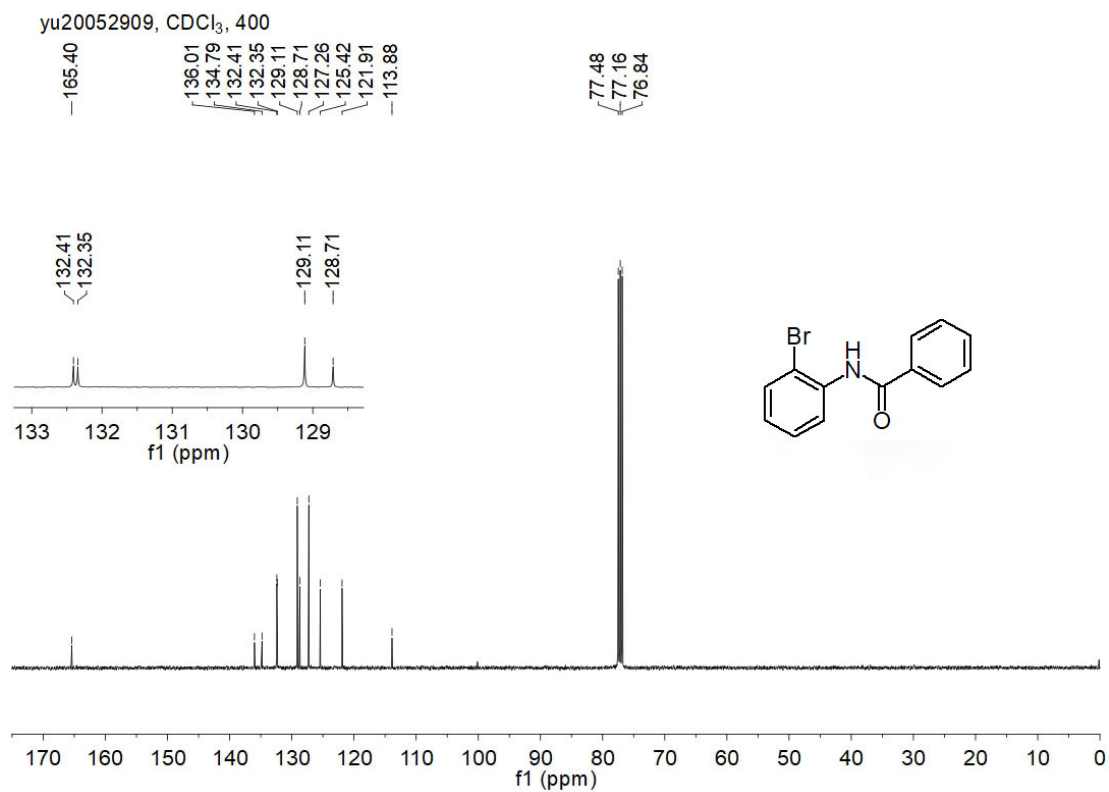
¹HNMR of **3b**



¹³CNMR of **3b**

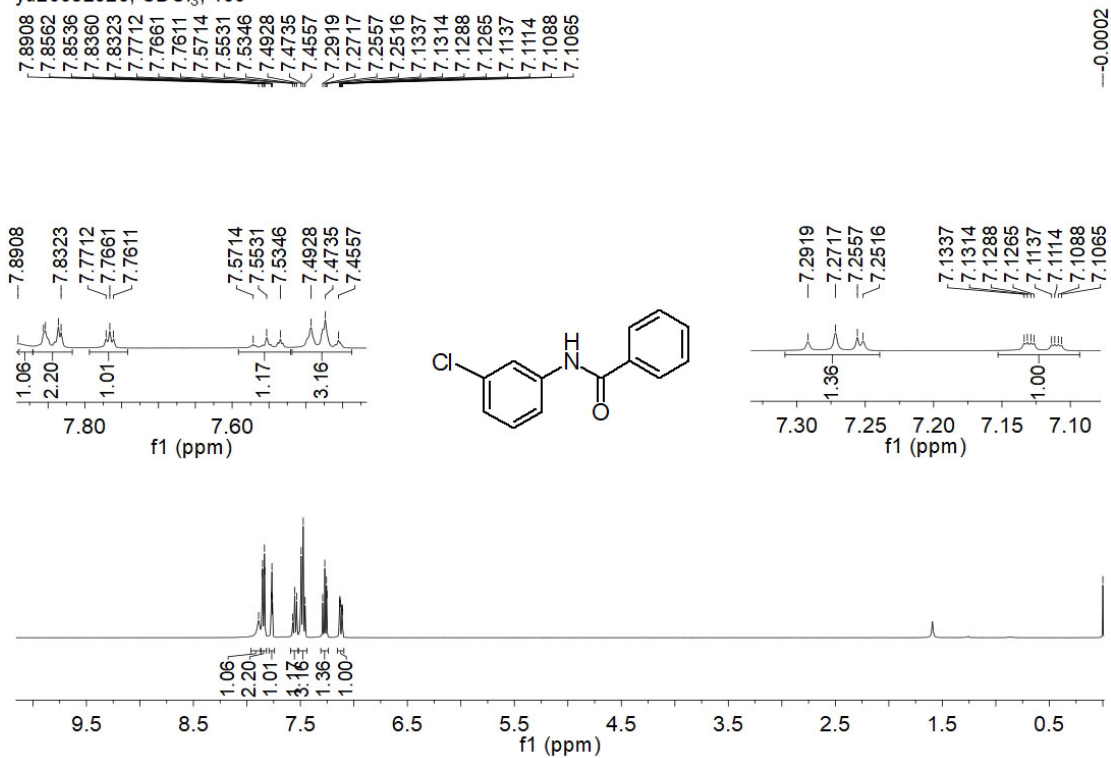


¹H NMR of **3c**



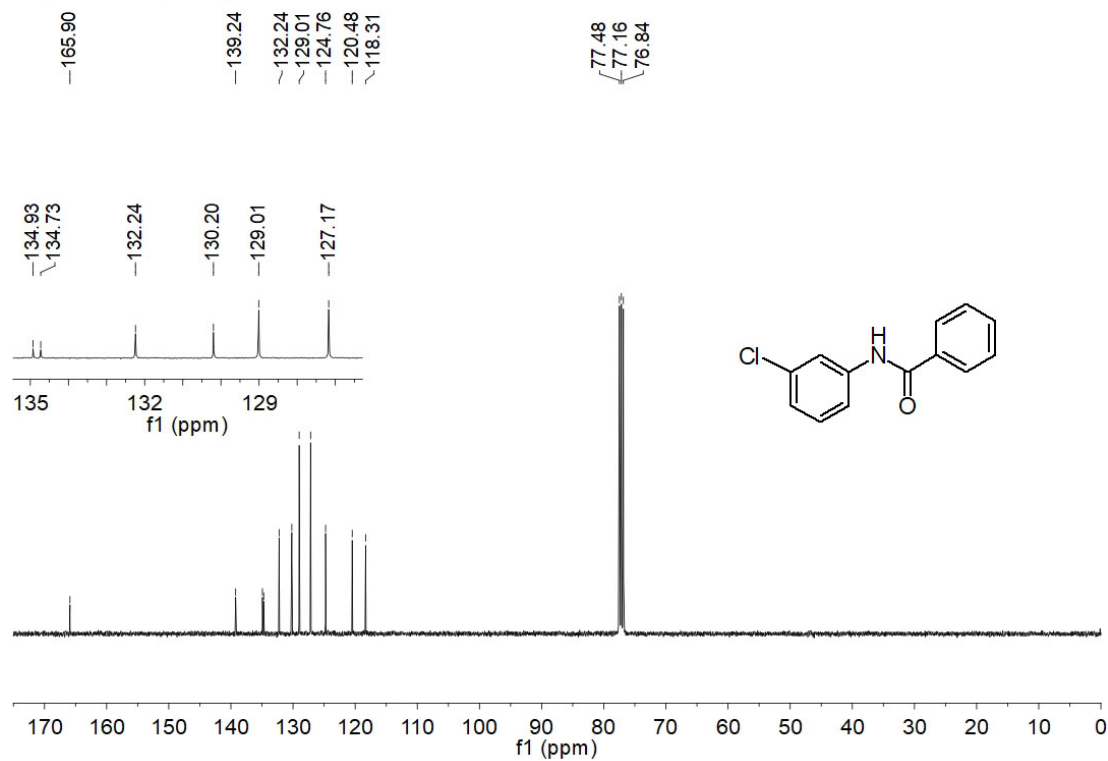
¹³C NMR of **3c**

yu20052926, CDCl₃, 400

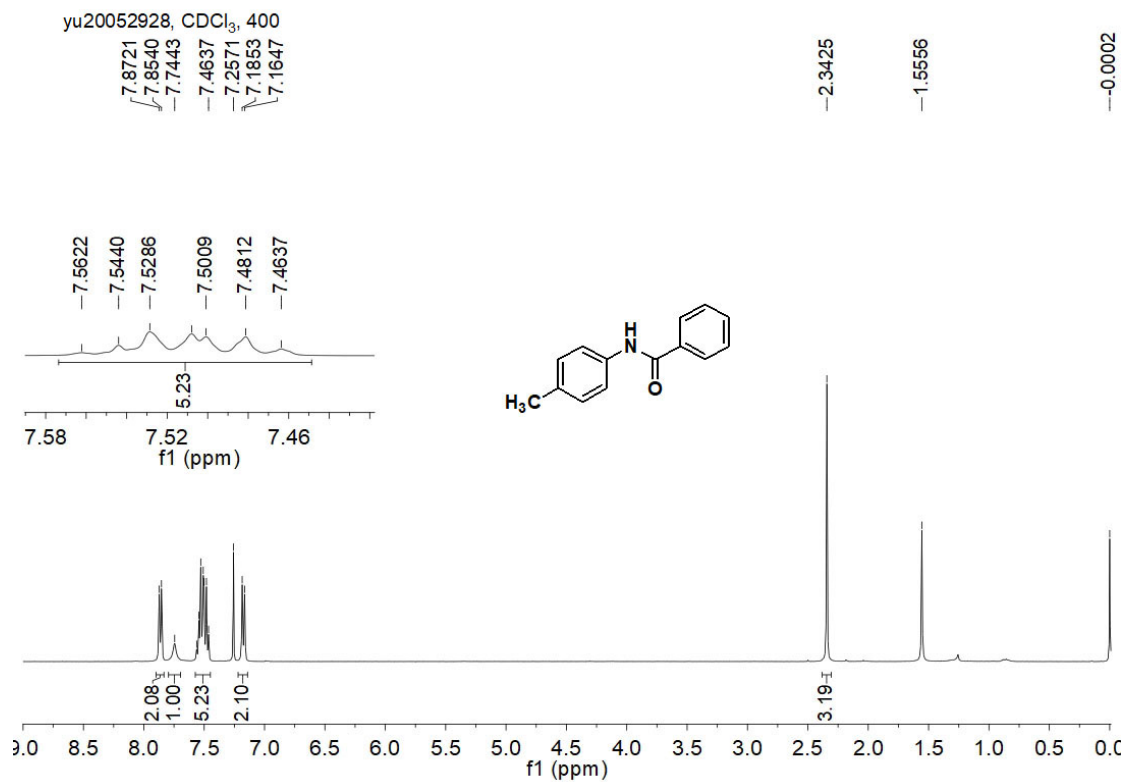


¹H NMR of **3d**

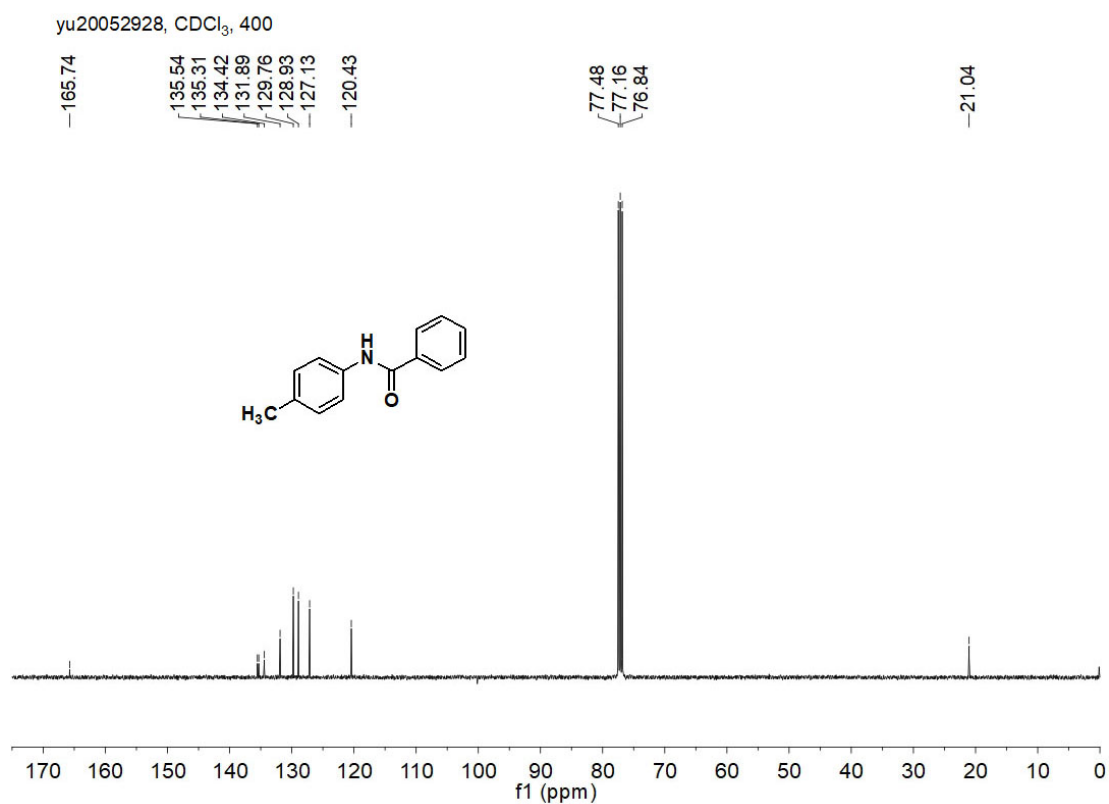
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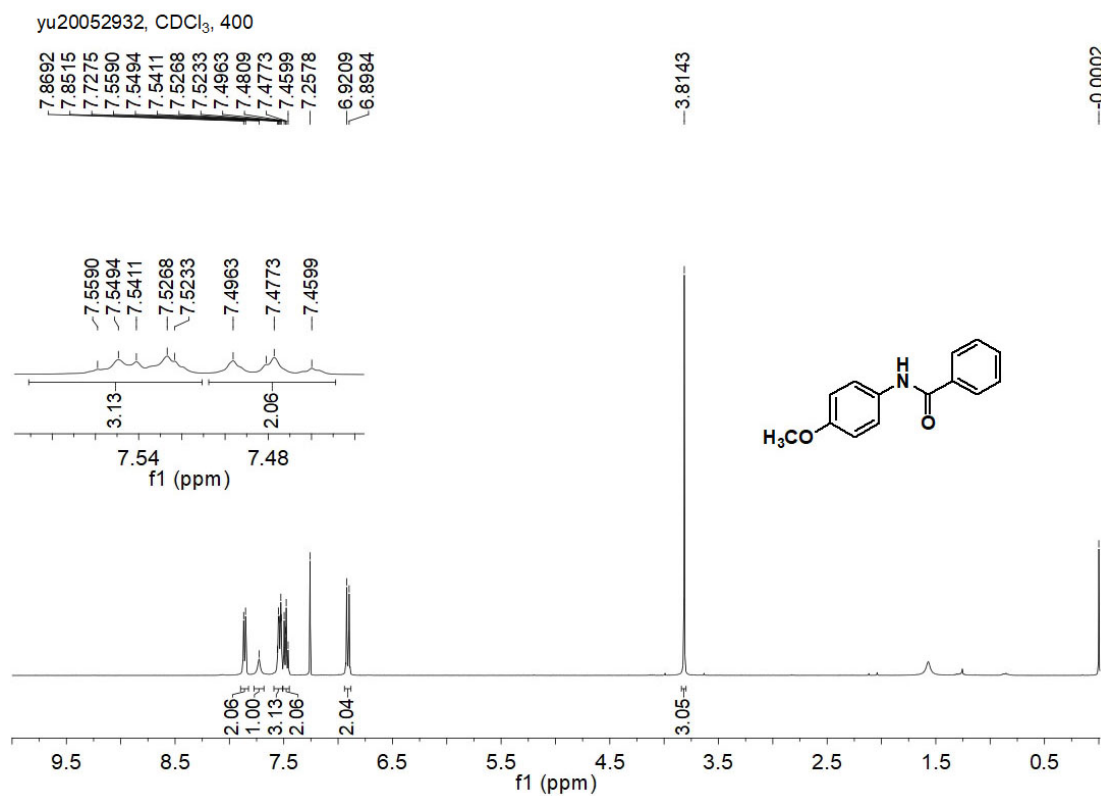
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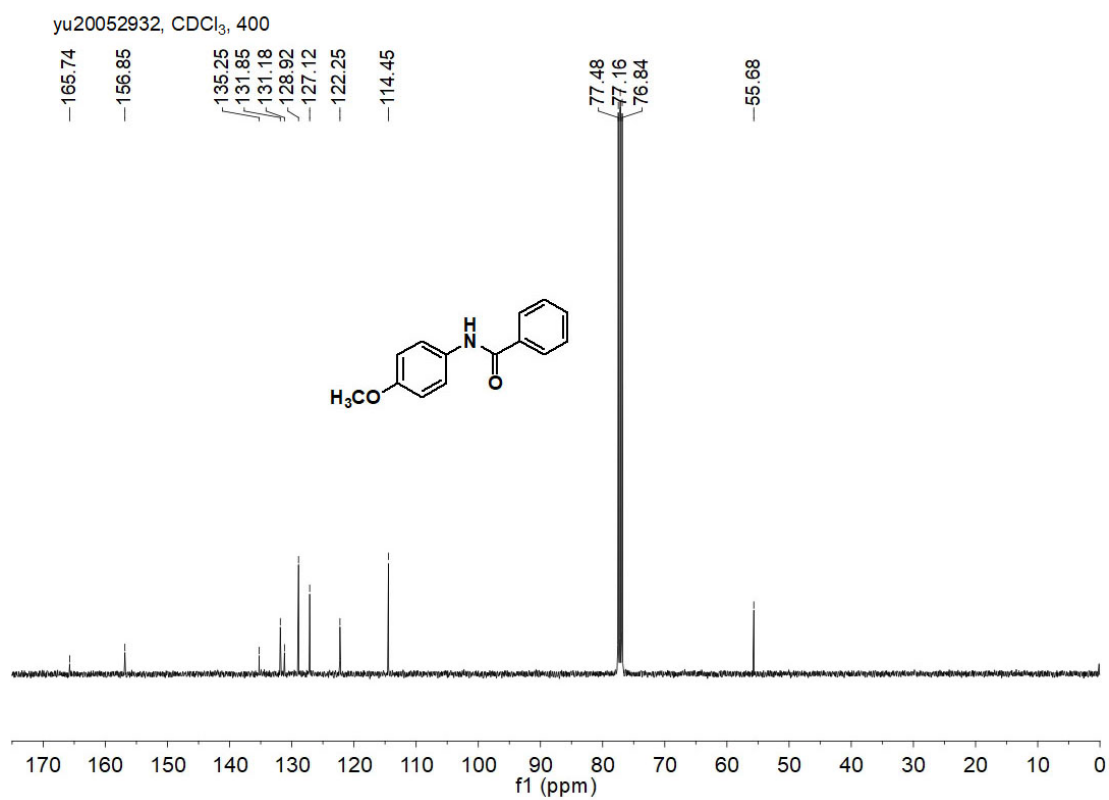
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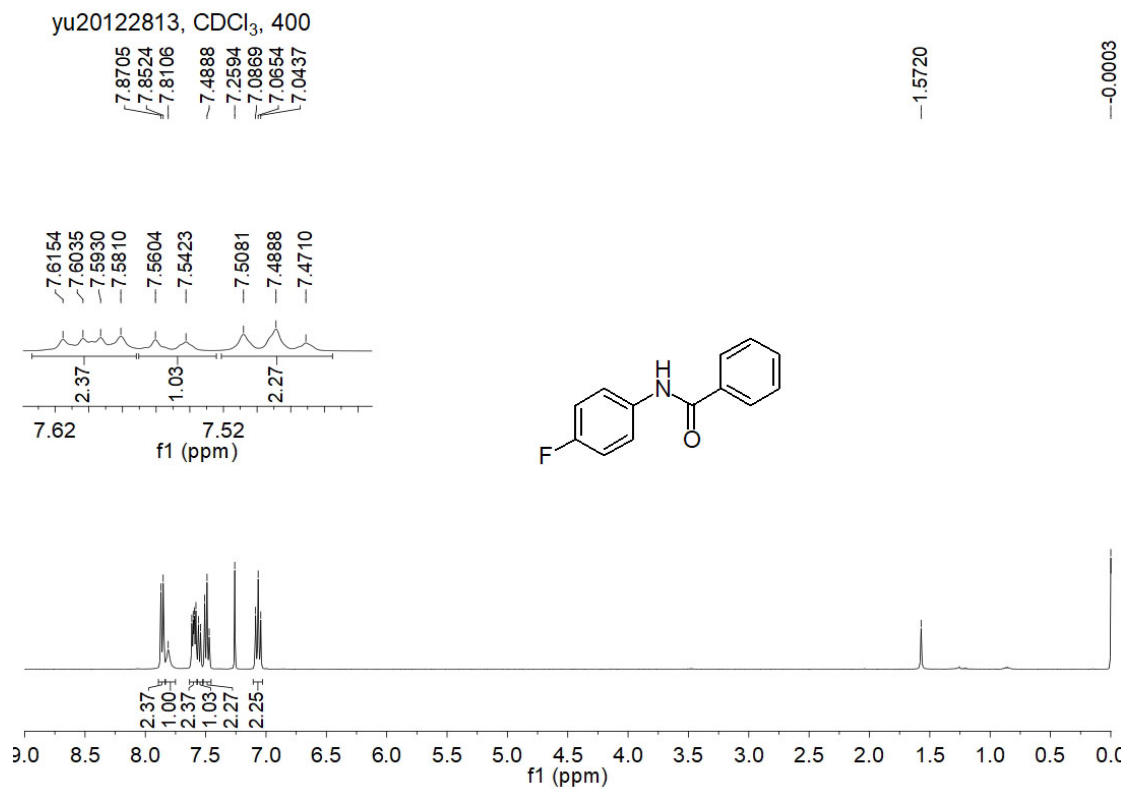
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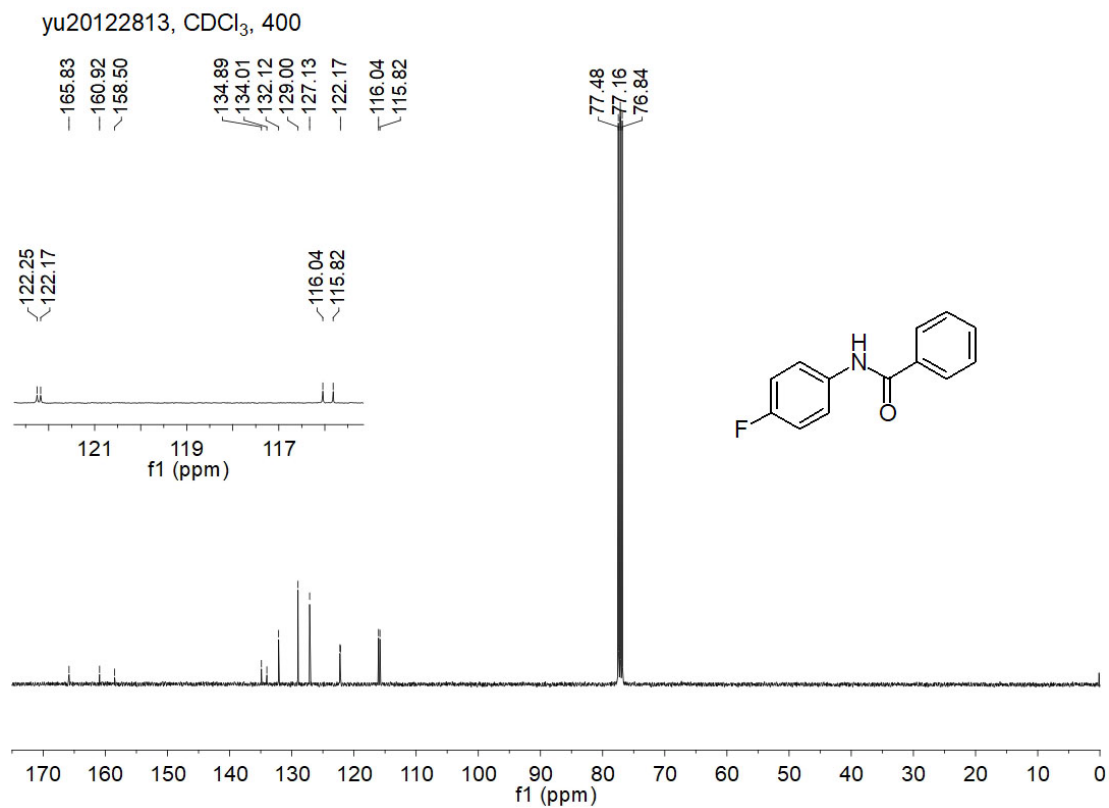
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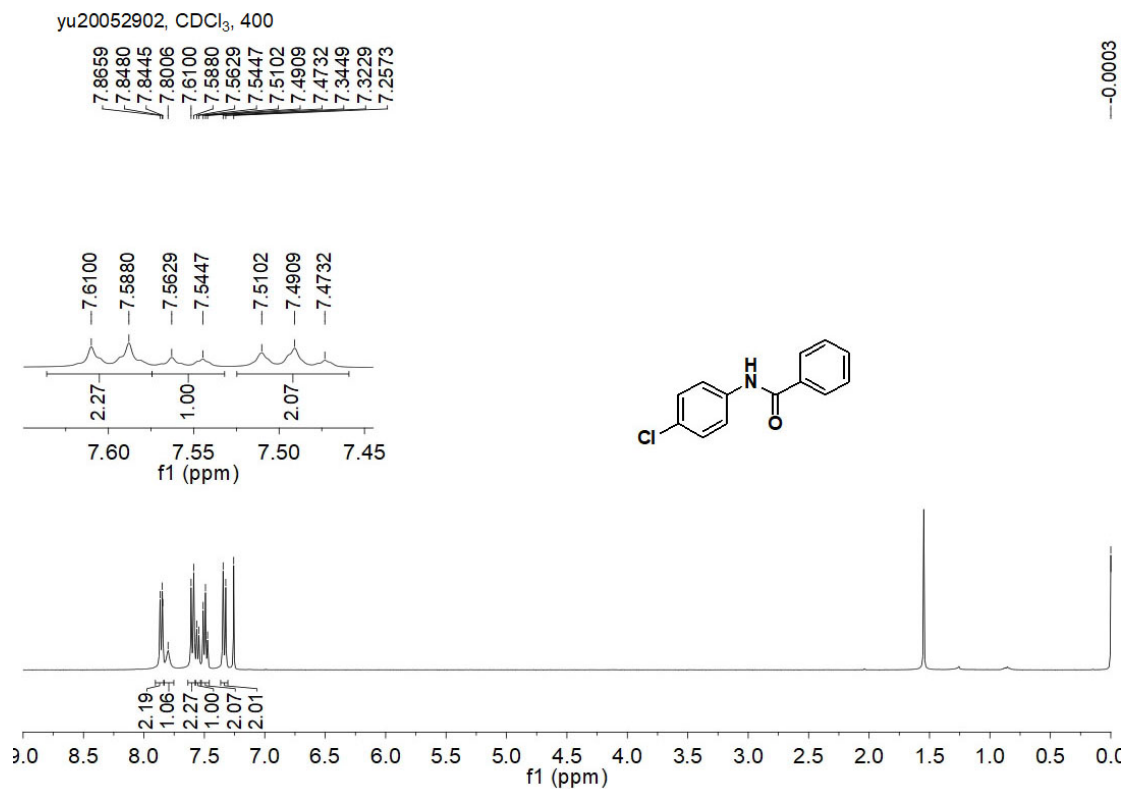
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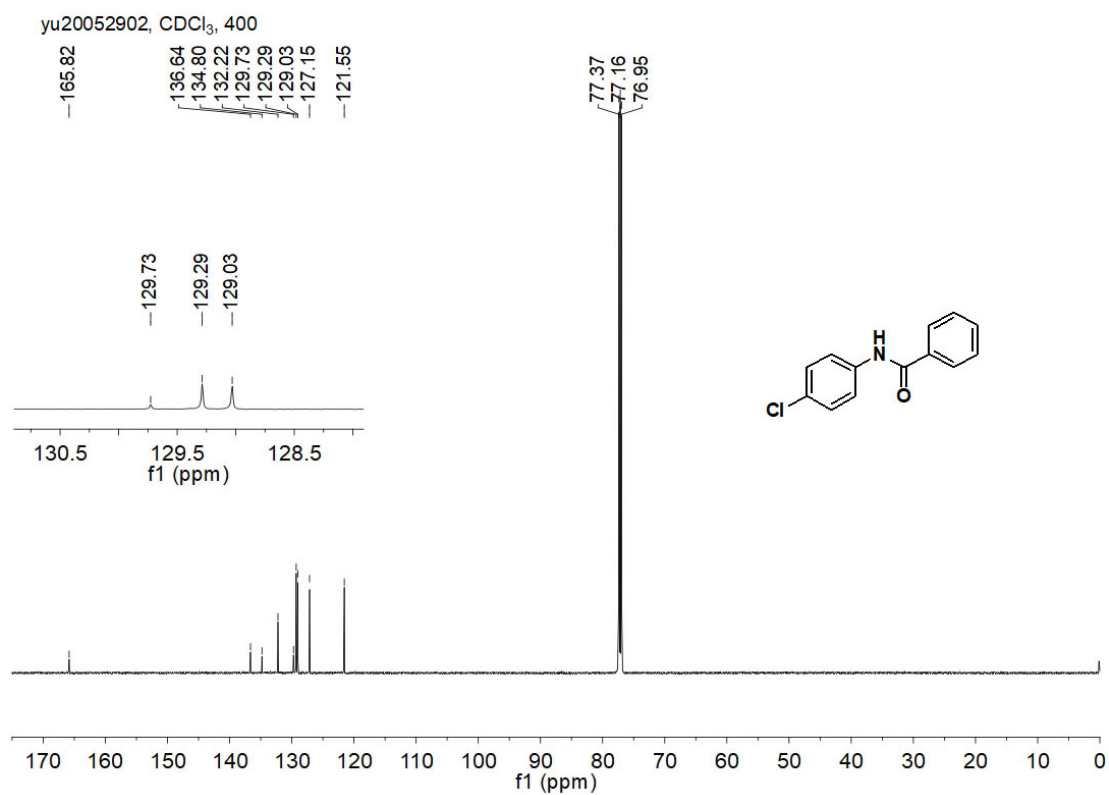
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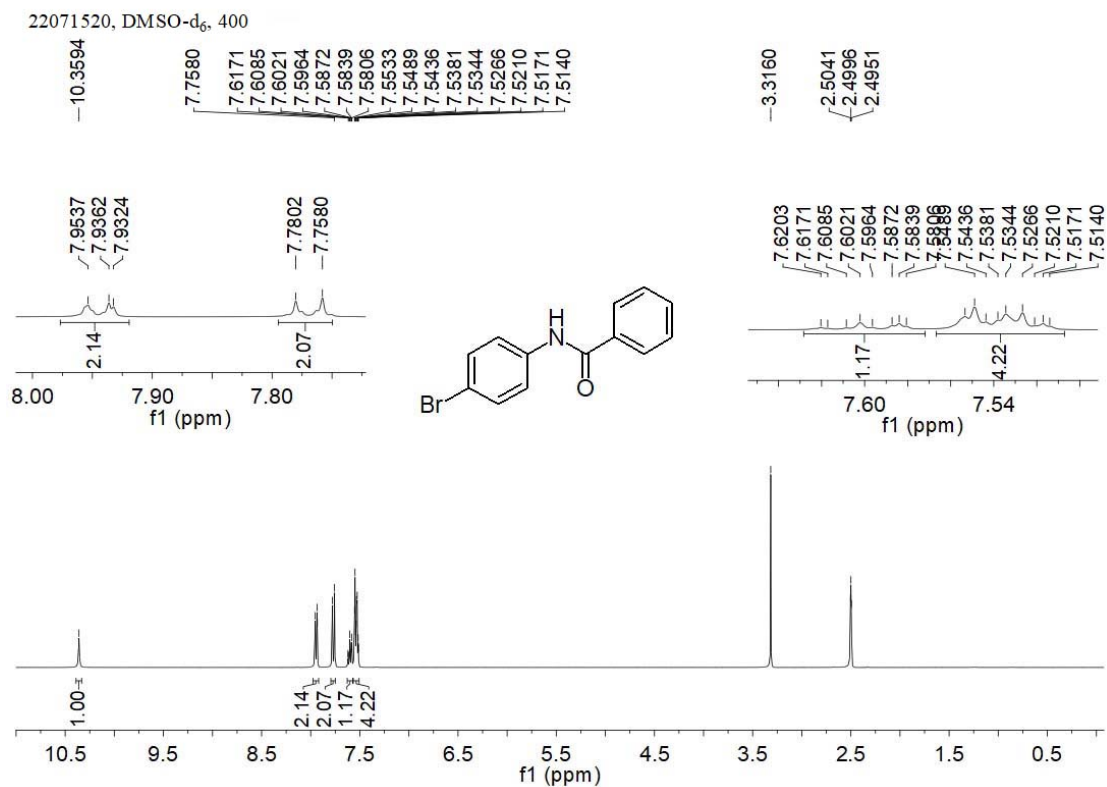
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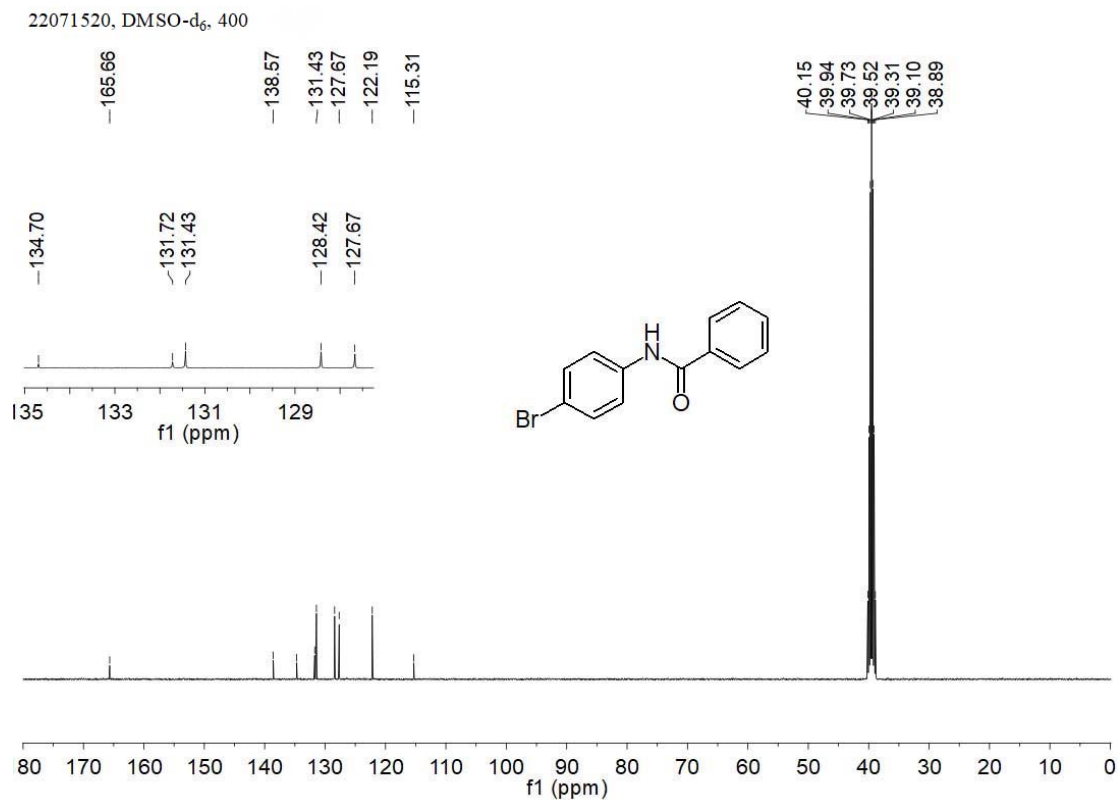
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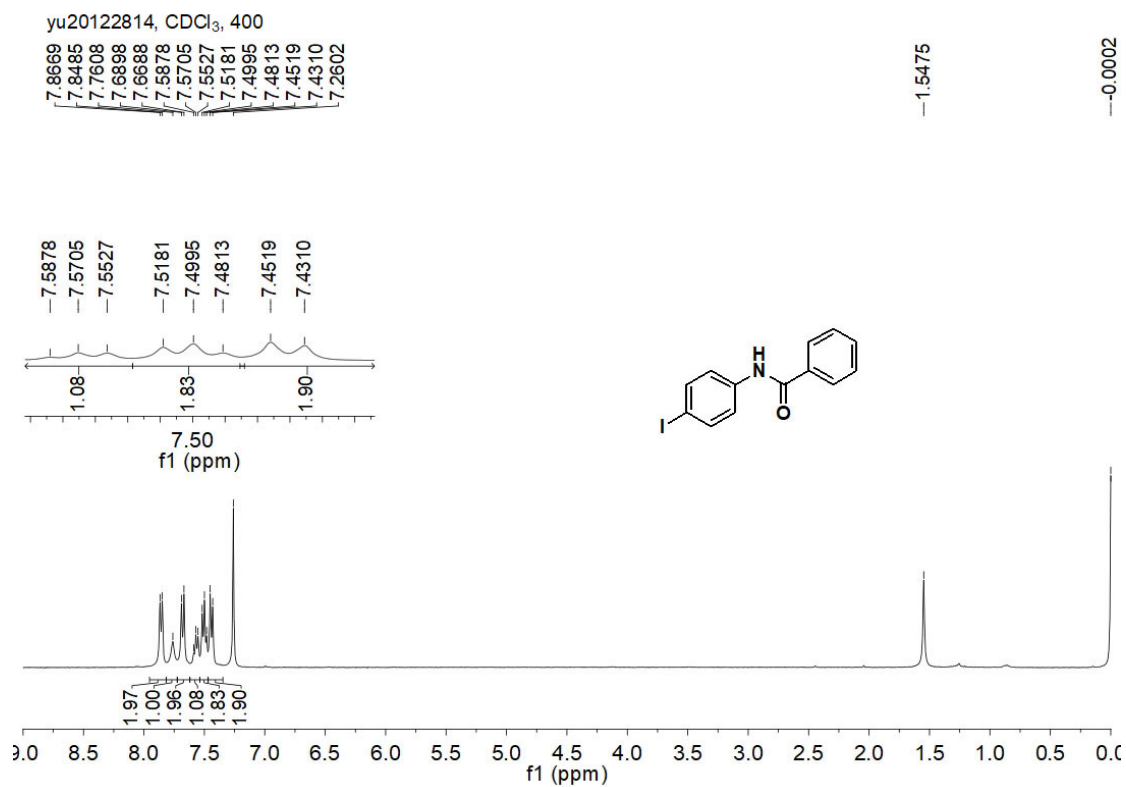
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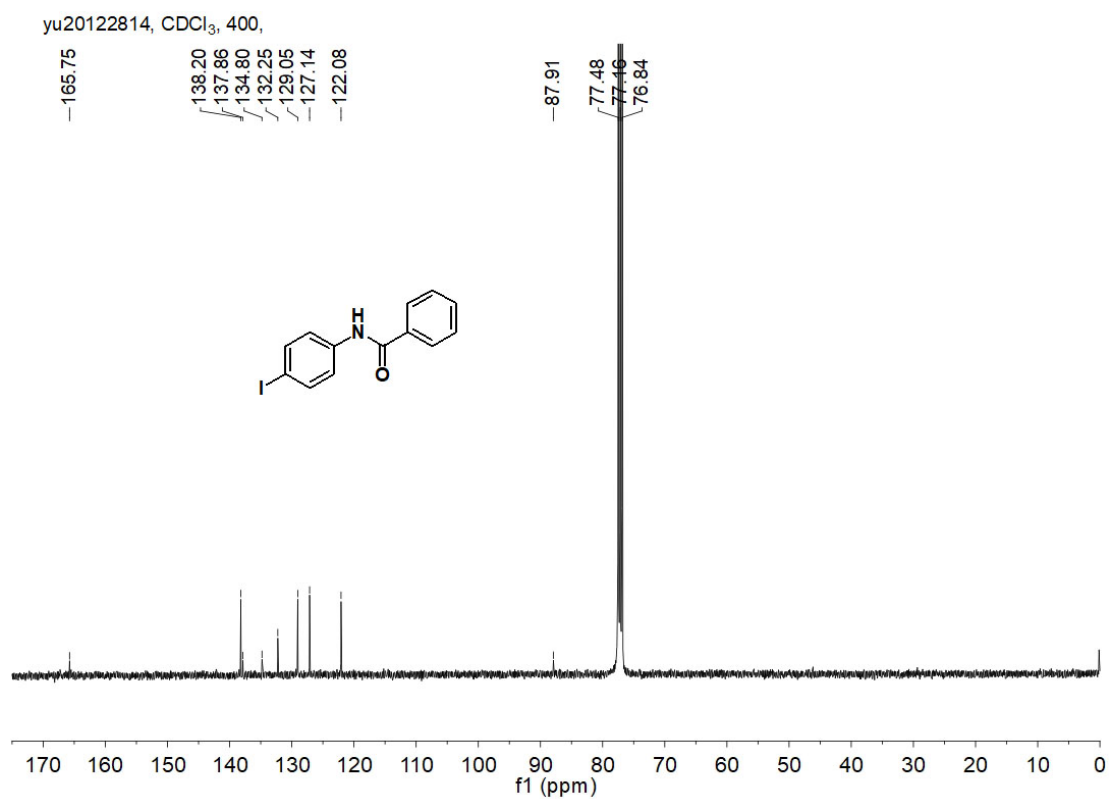
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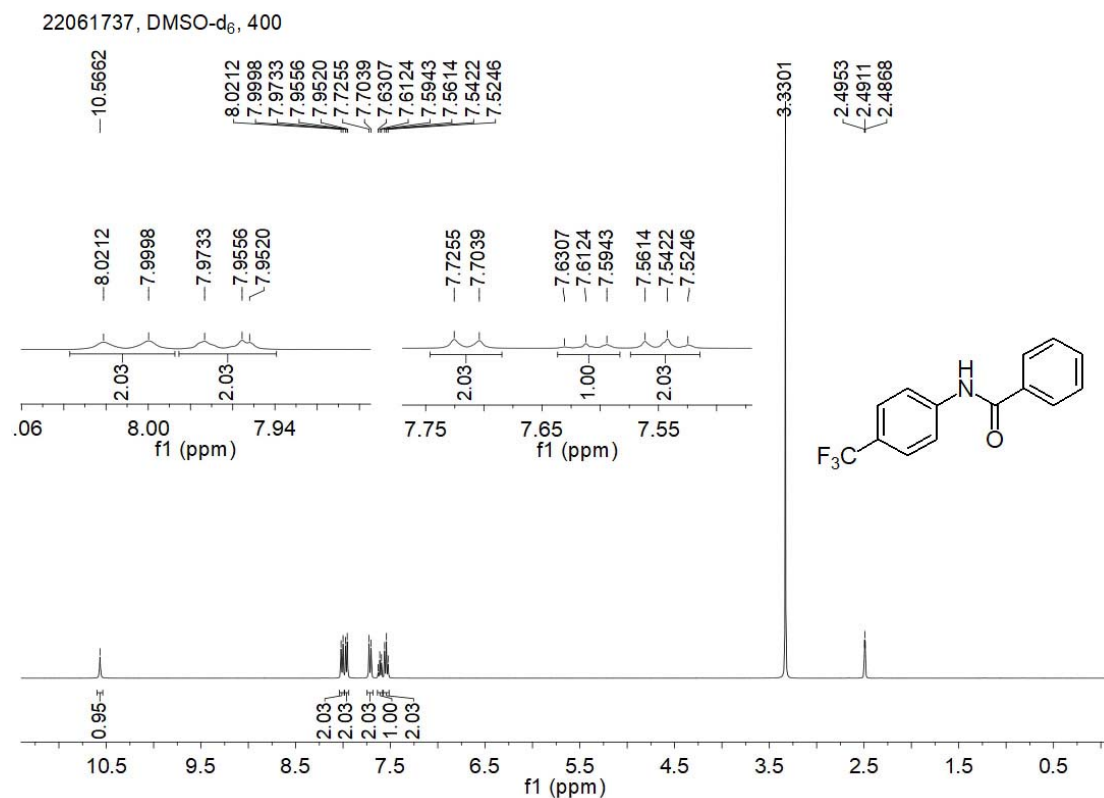
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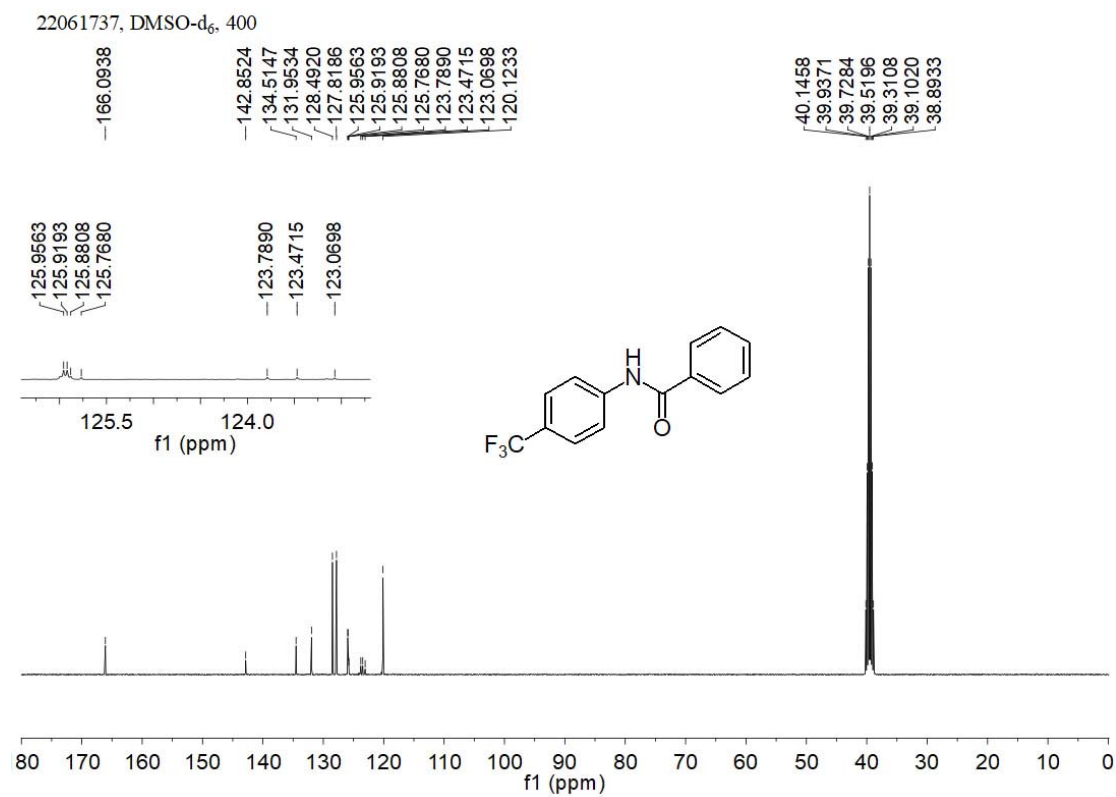
¹HNMR of **3j**



¹³CNMR of **3j**

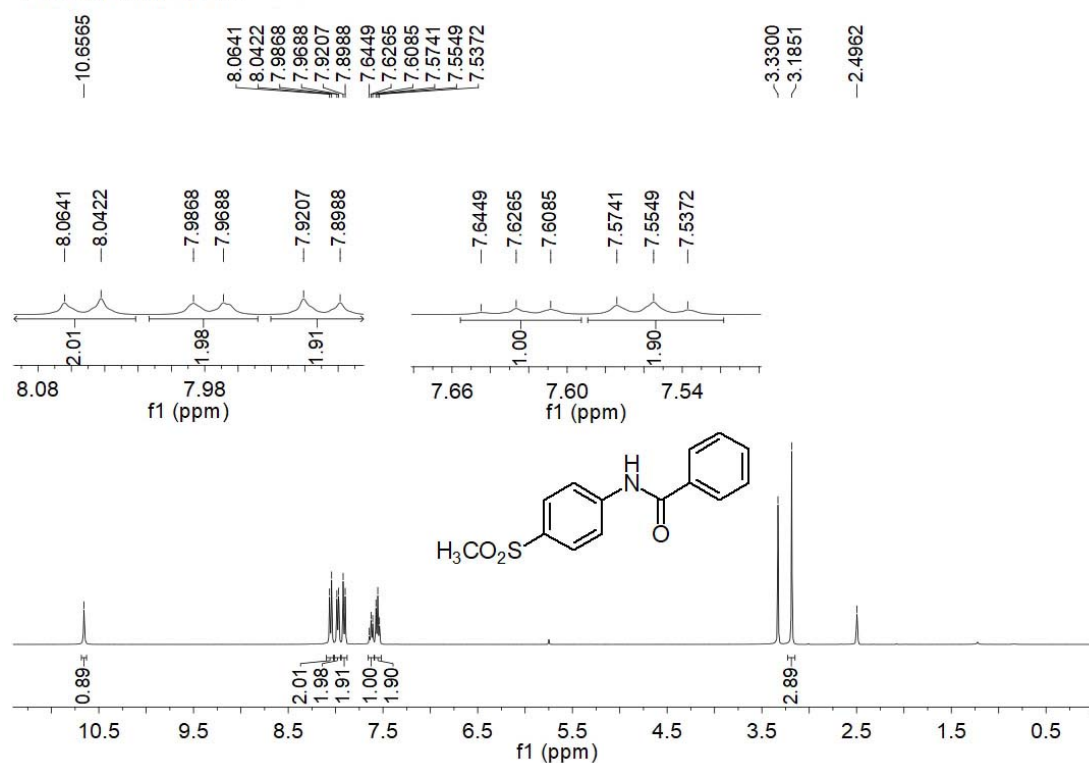


¹HNMR of **3k**



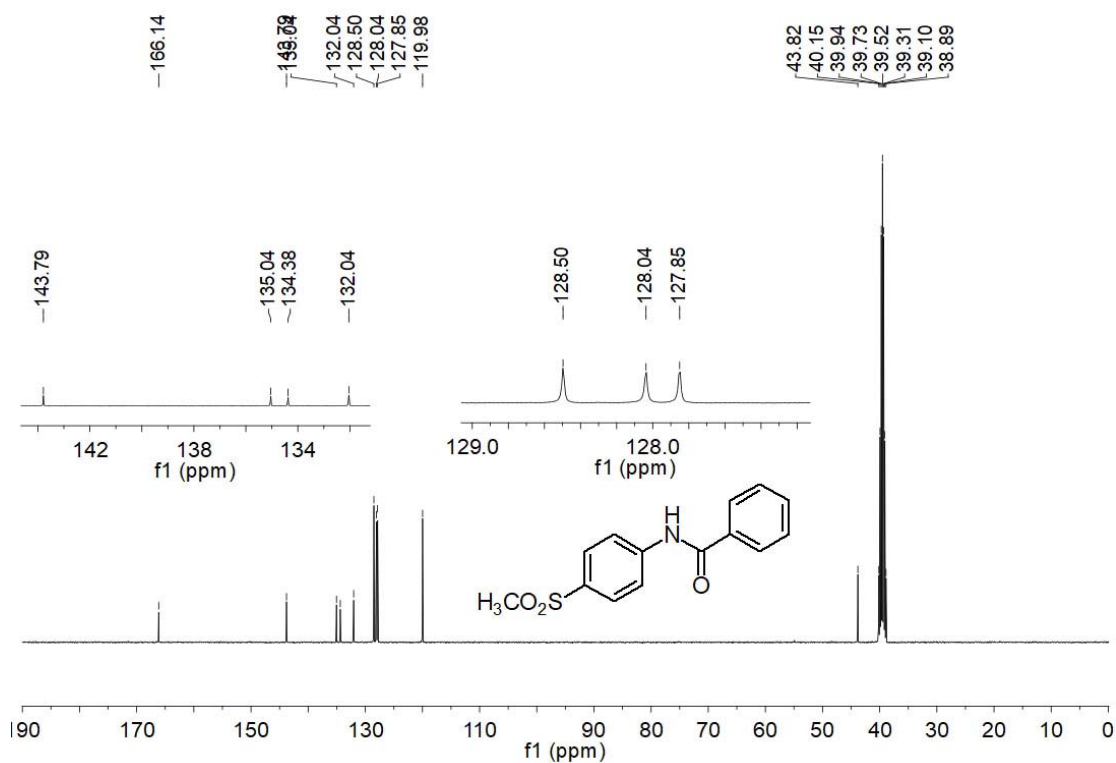
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22061734, DMSO-d₆, 400

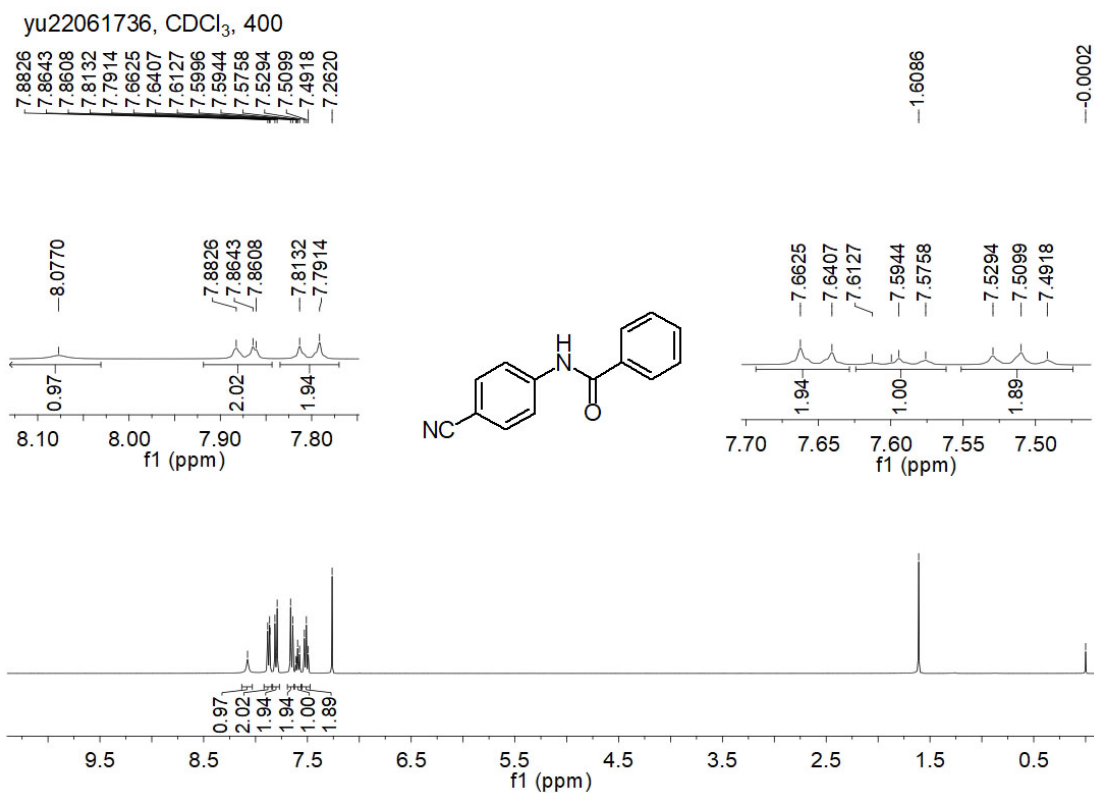


¹HNMR of **31**

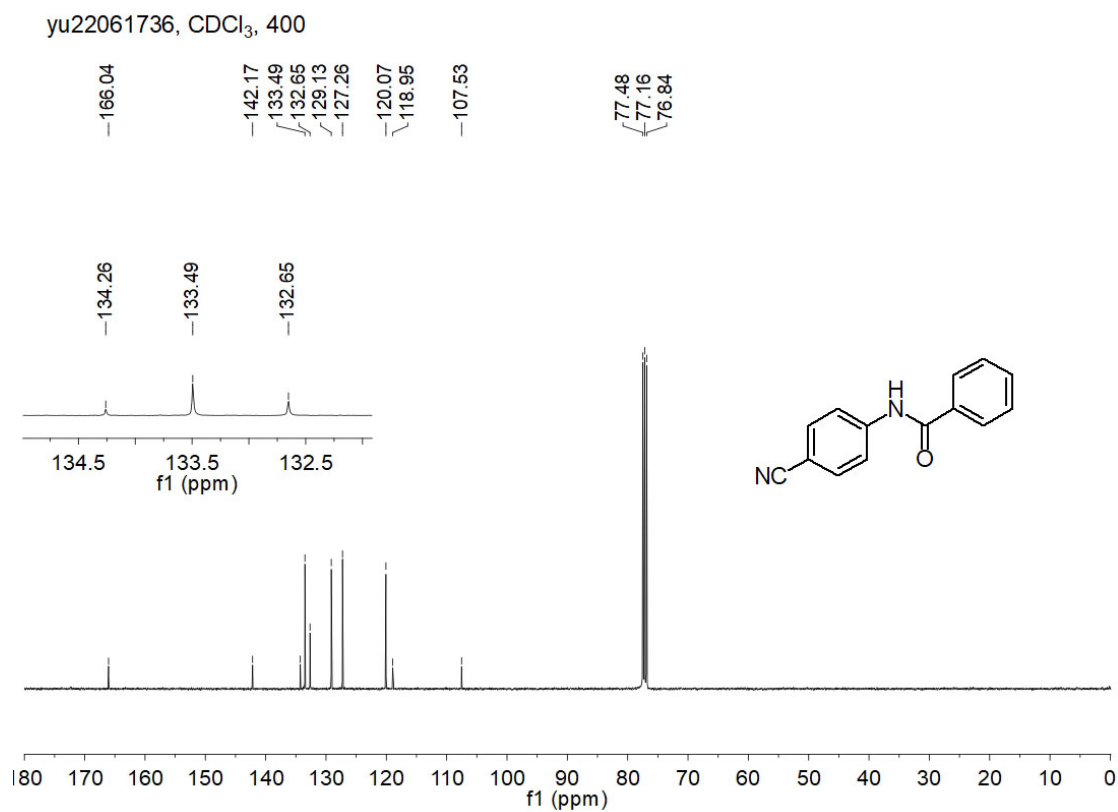
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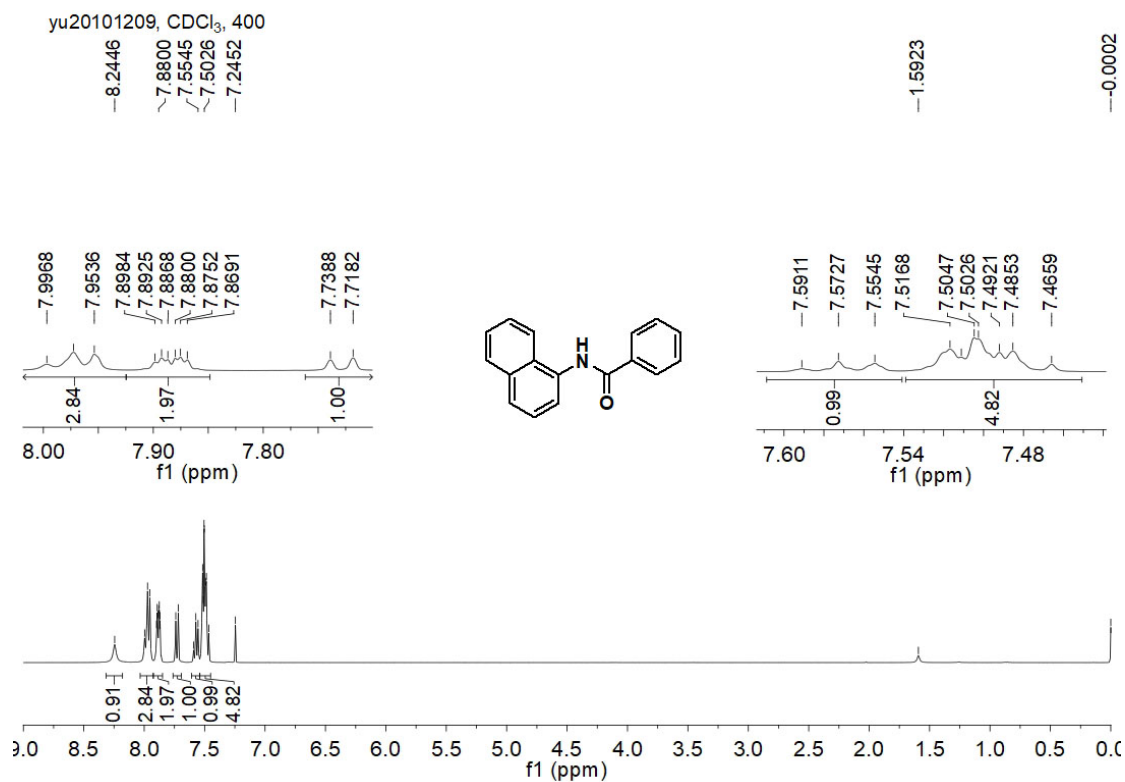
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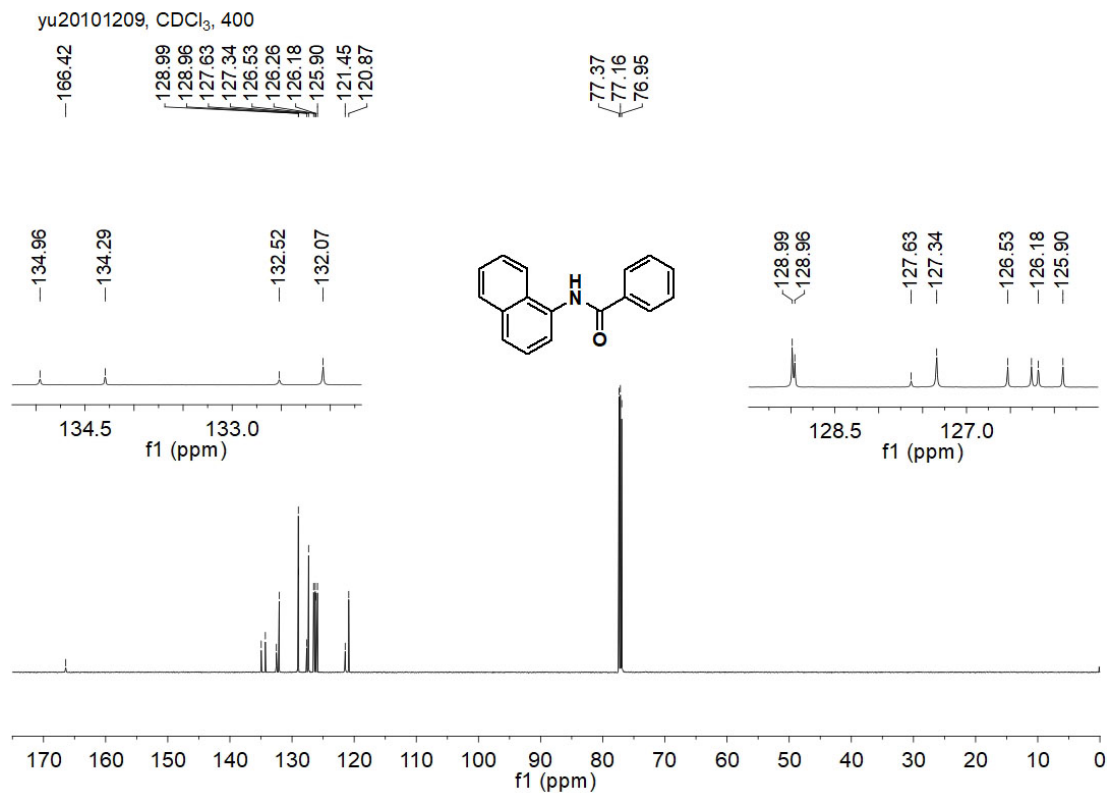
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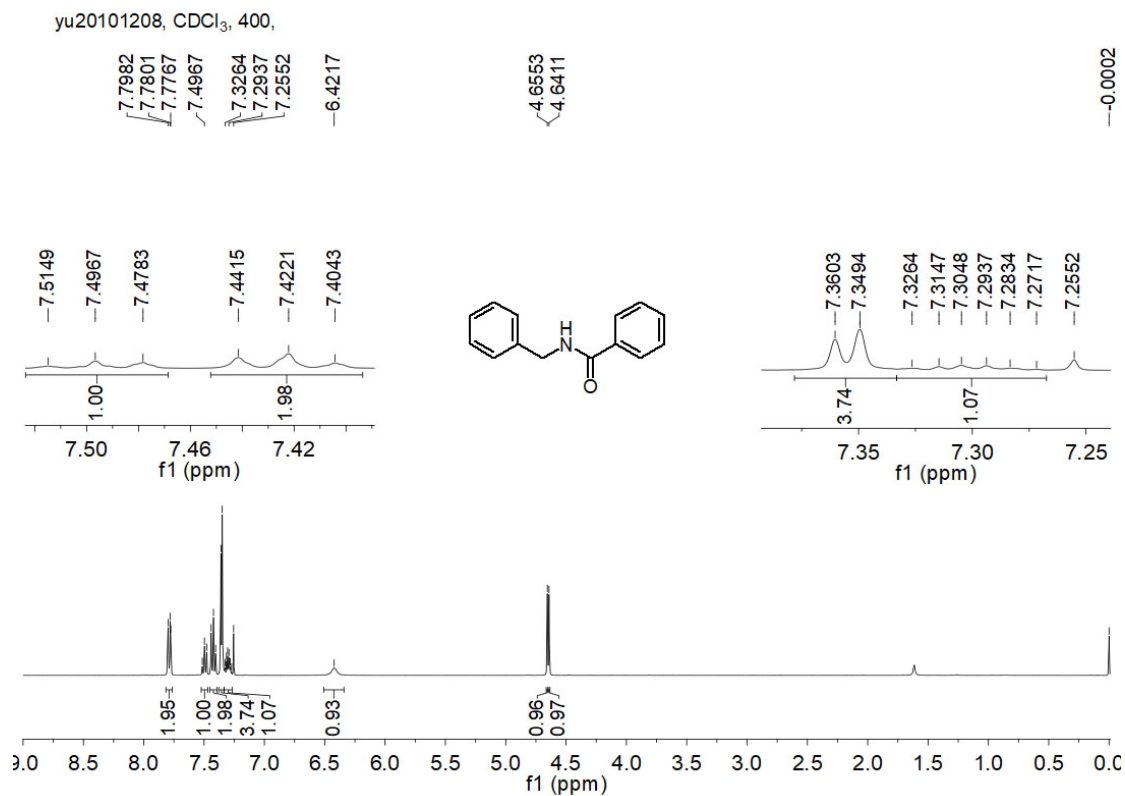
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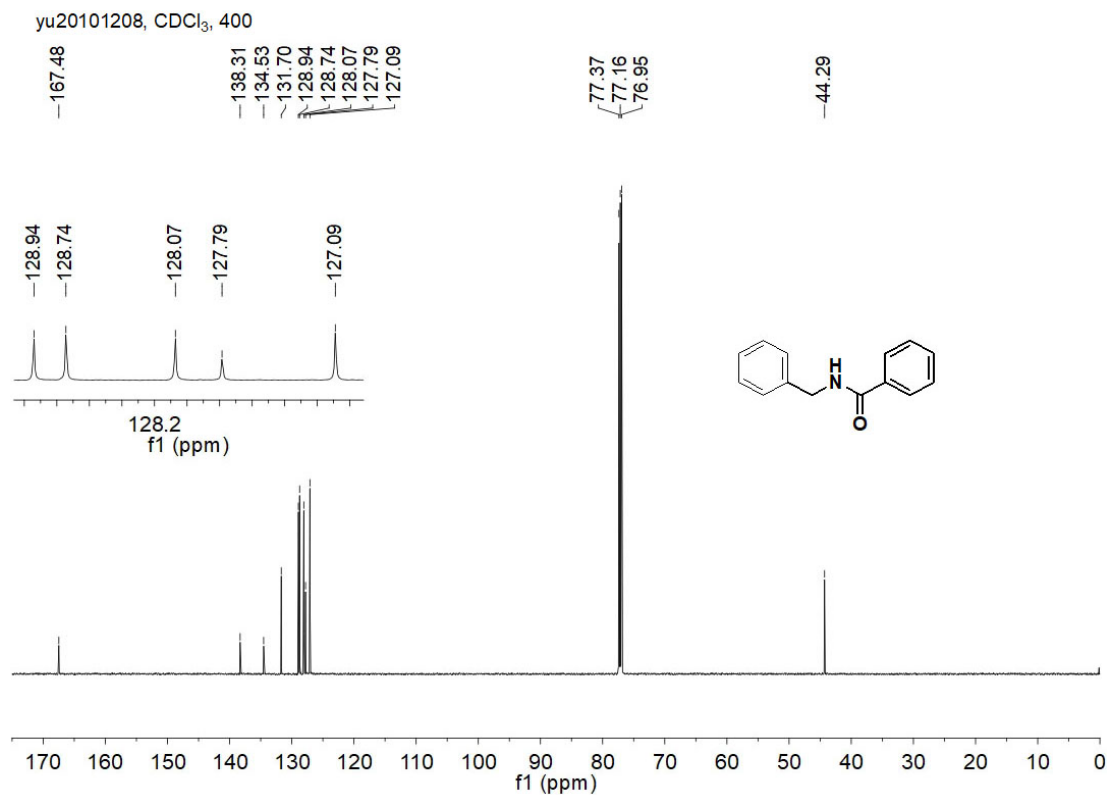
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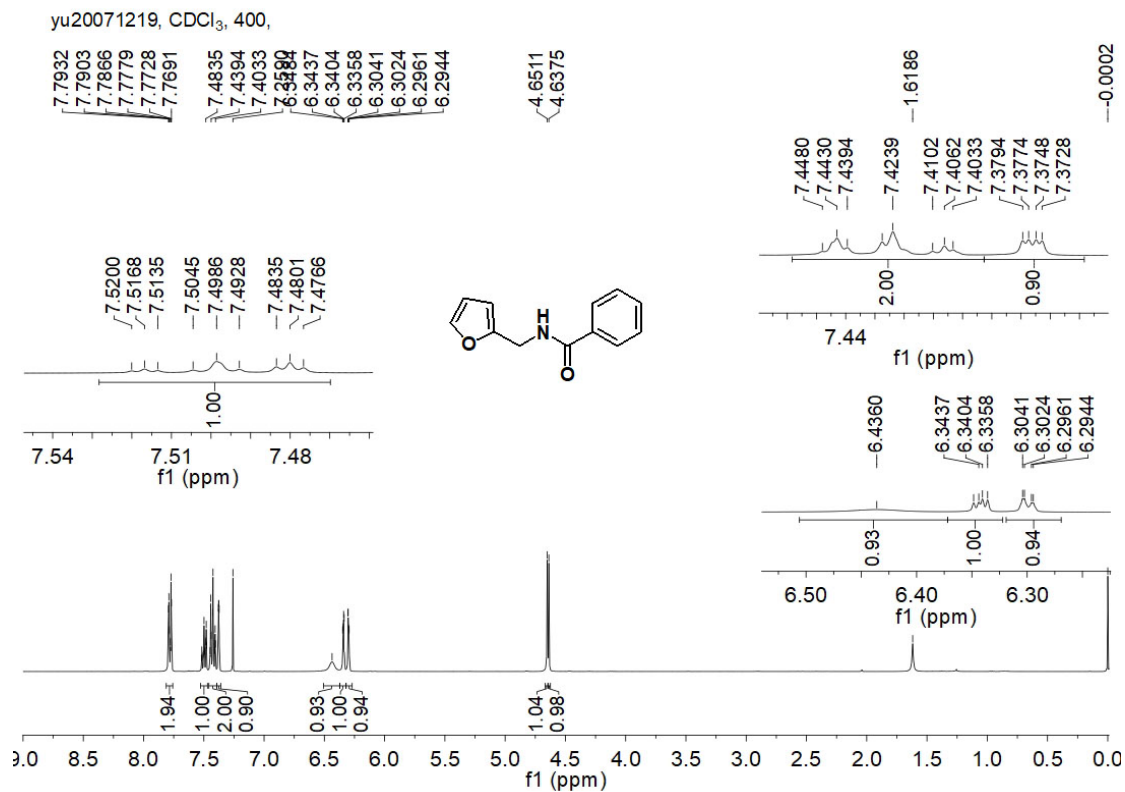
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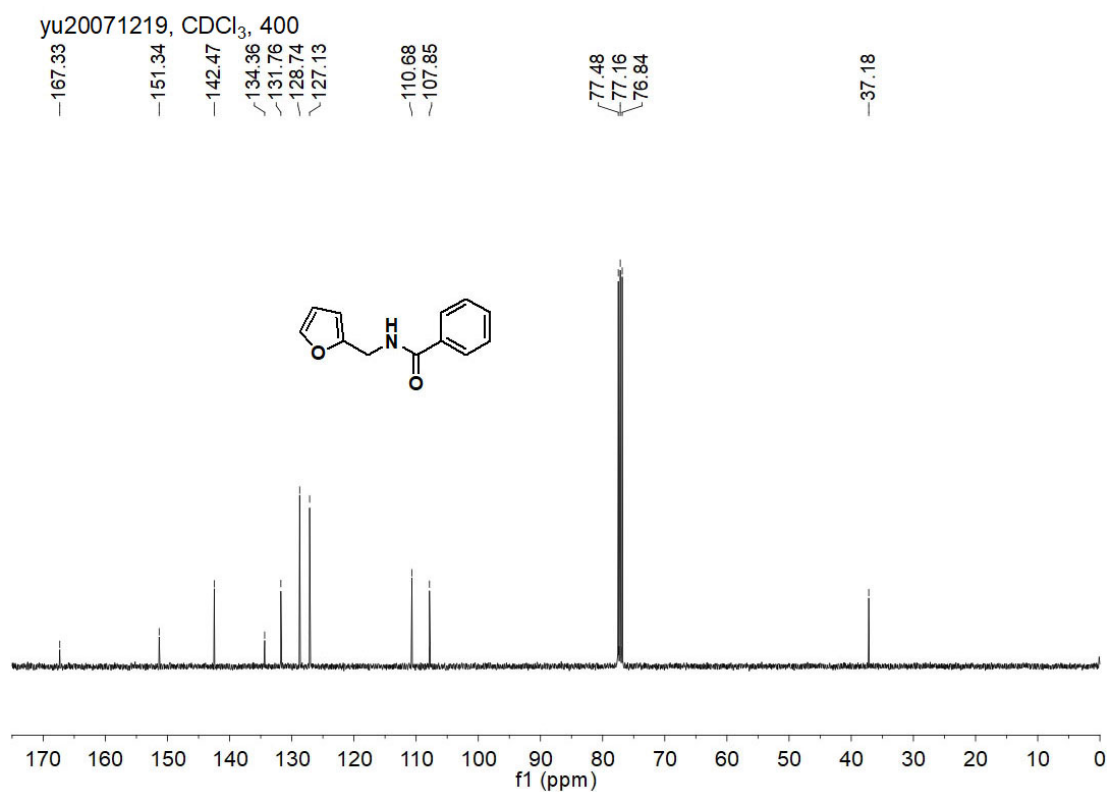
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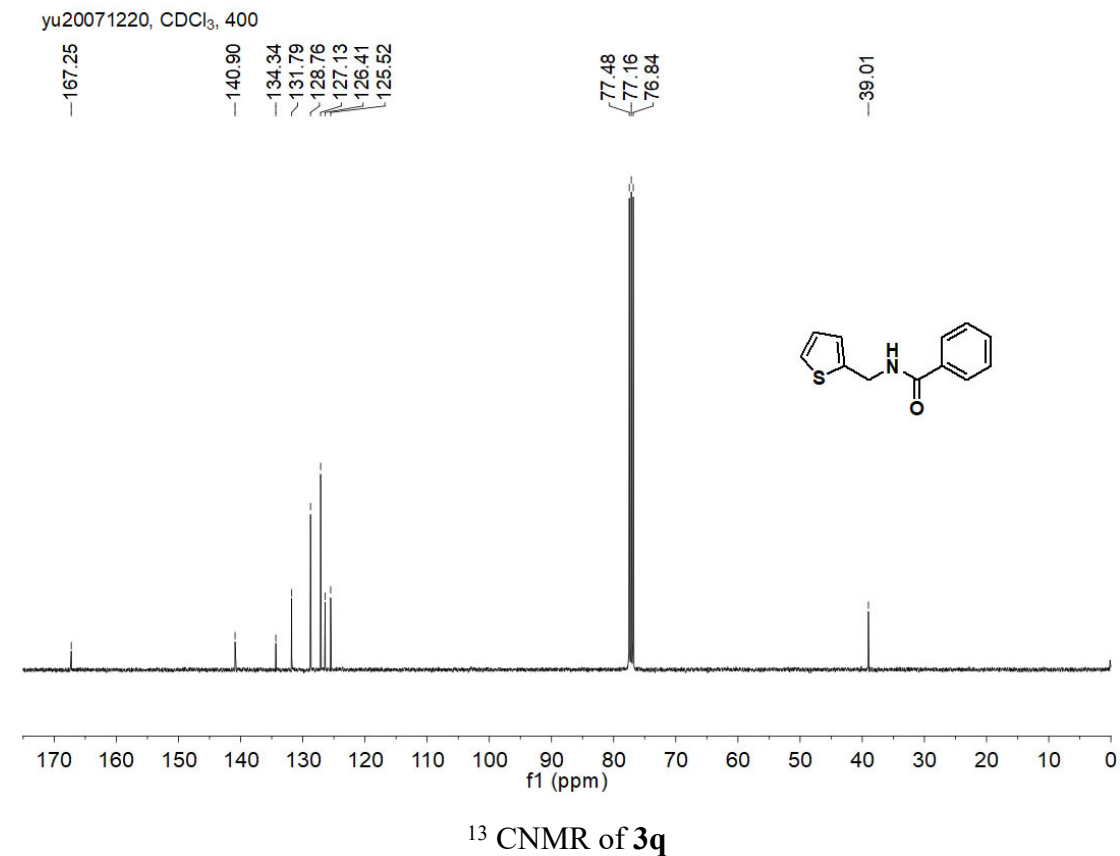
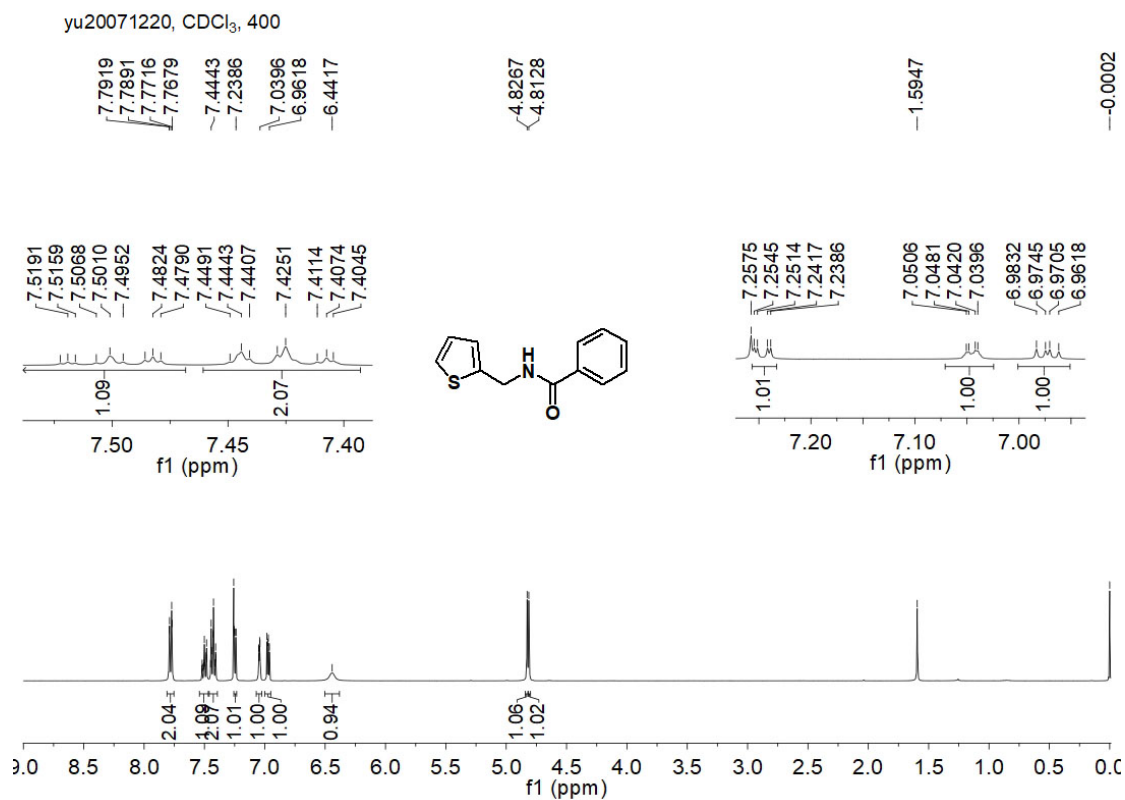
¹³CNMR of **3o**

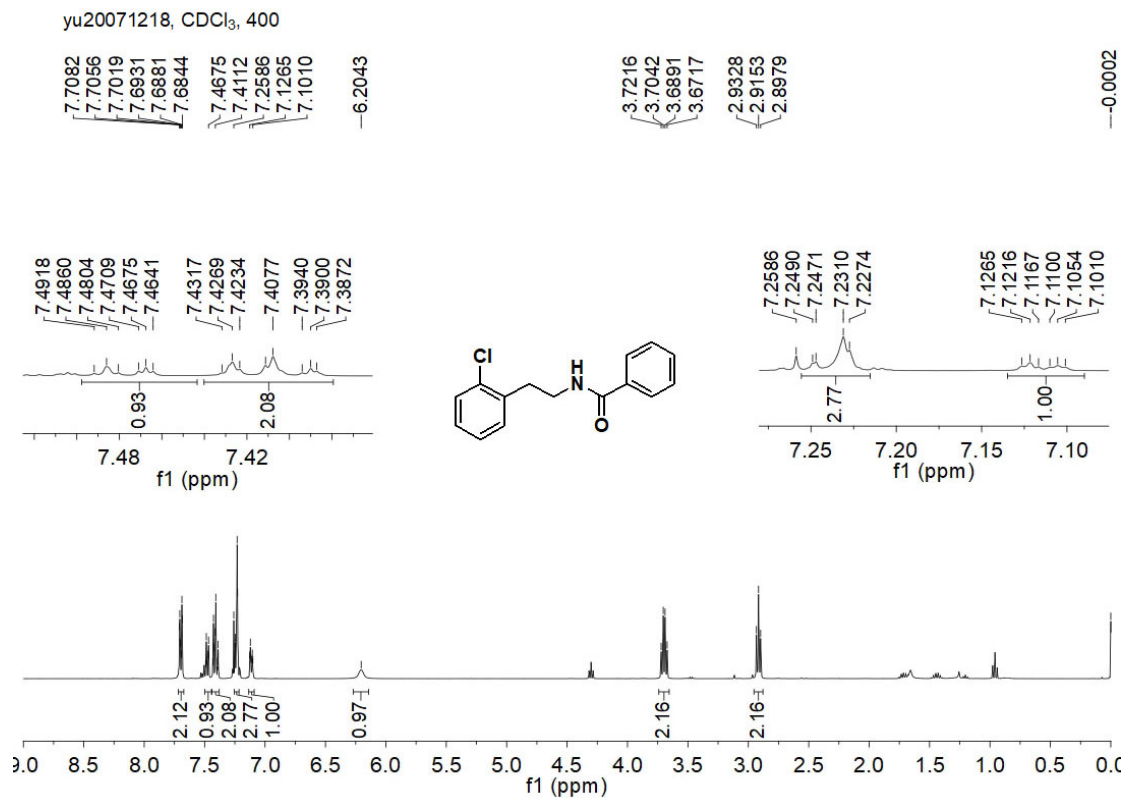


¹HNMR of **3p**

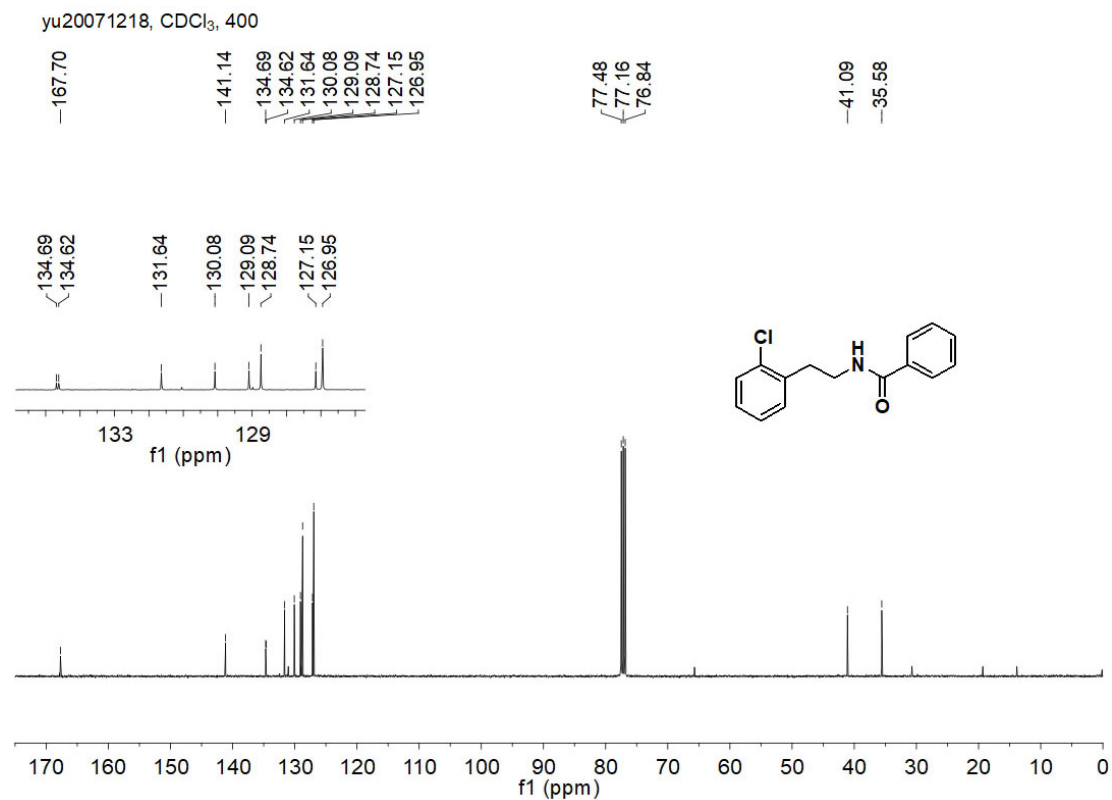


¹³CNMR of **3p**

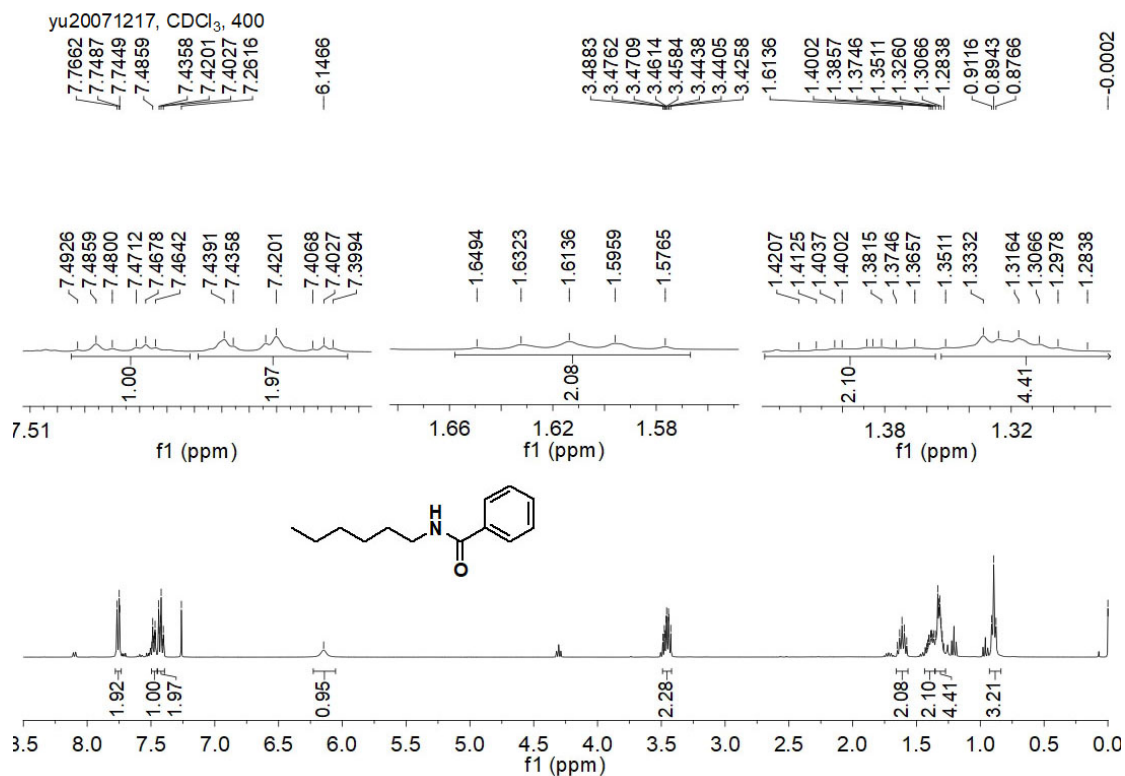




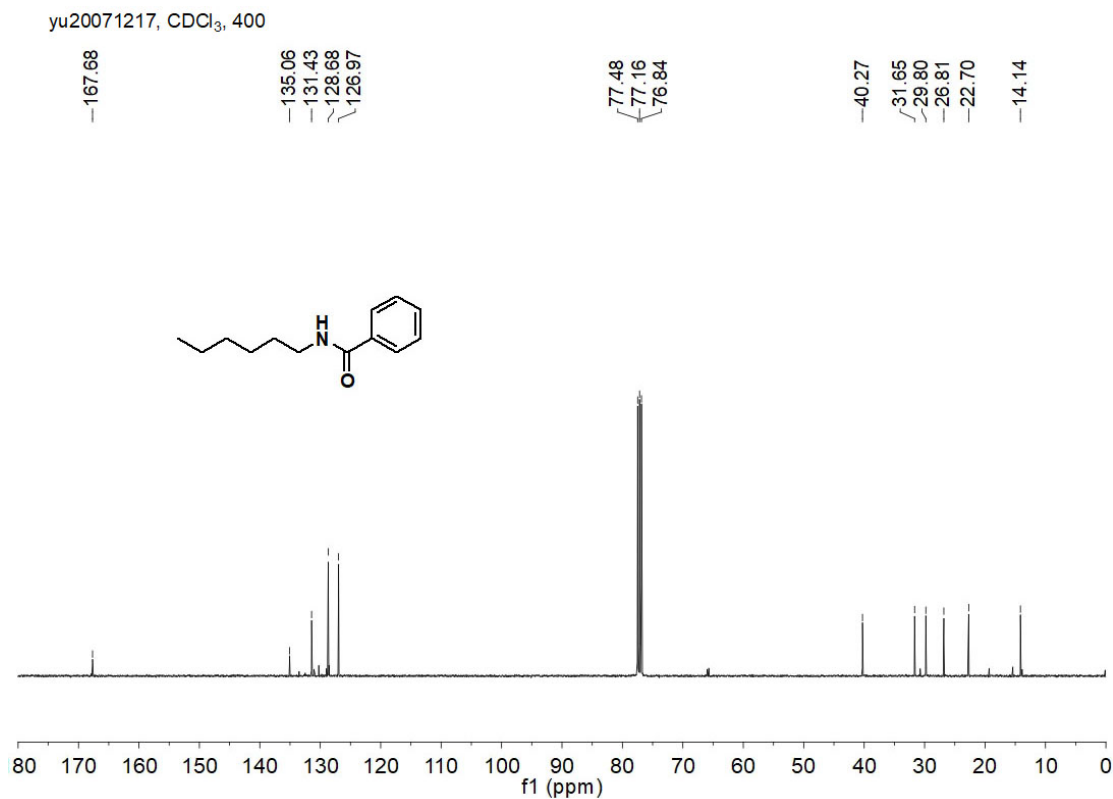
¹HNMR of **3r**



¹³CNMR of **3r**

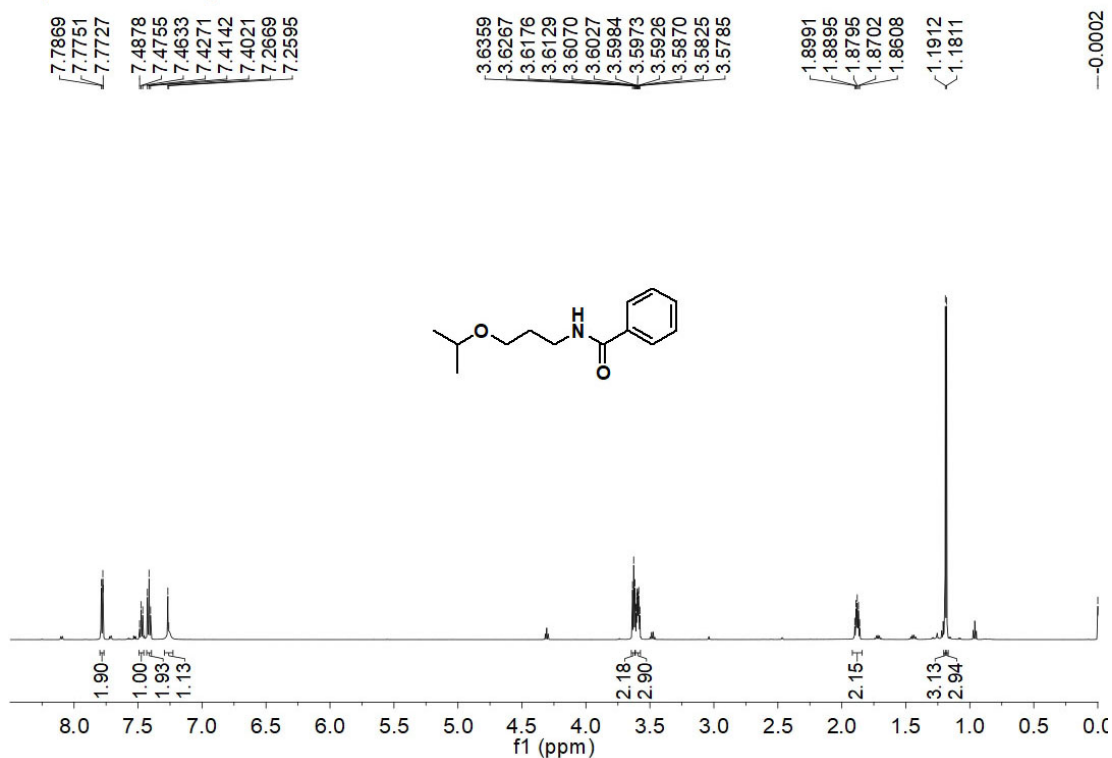


¹HNMR of **3s**

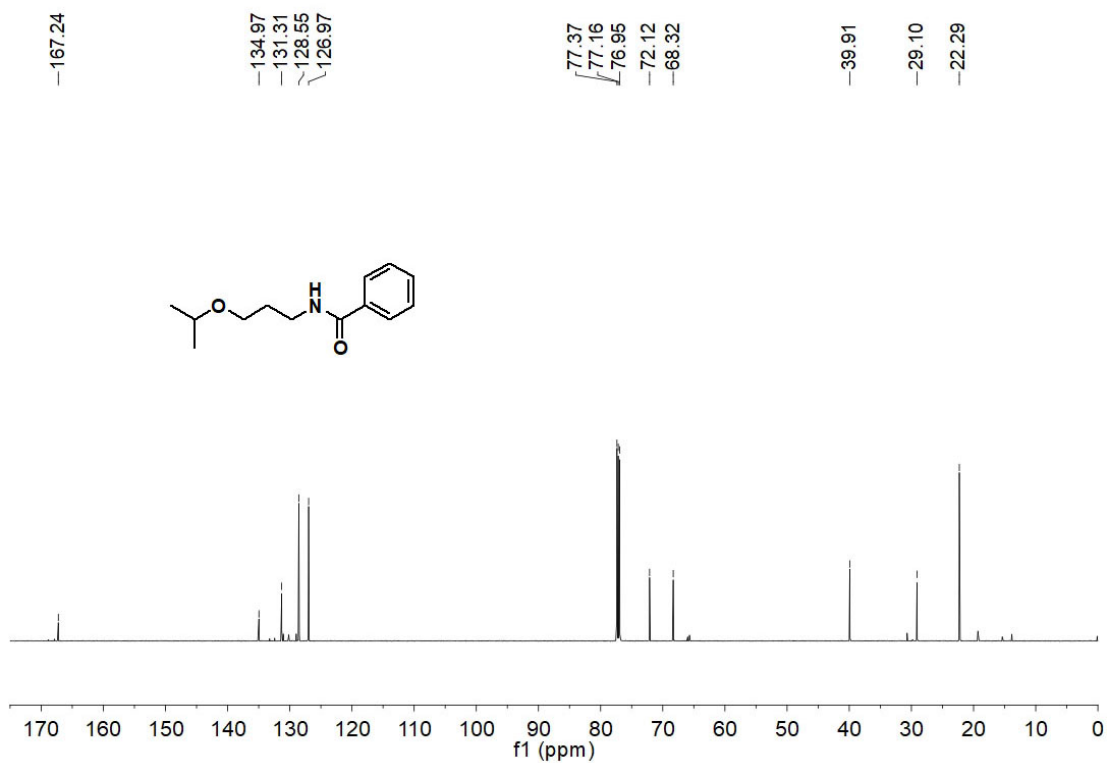


¹³CNMR of **3s**

yu20071216, CDCl₃, 400

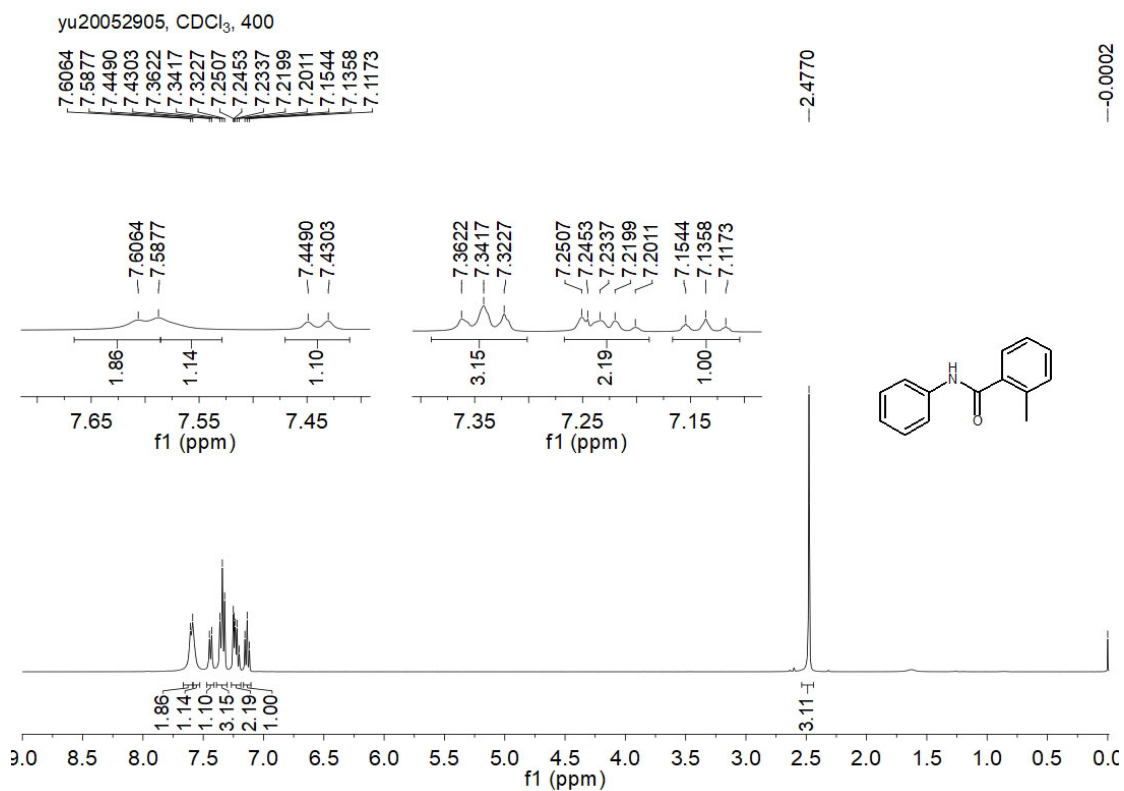


yu20071216, CDCl₃, 400

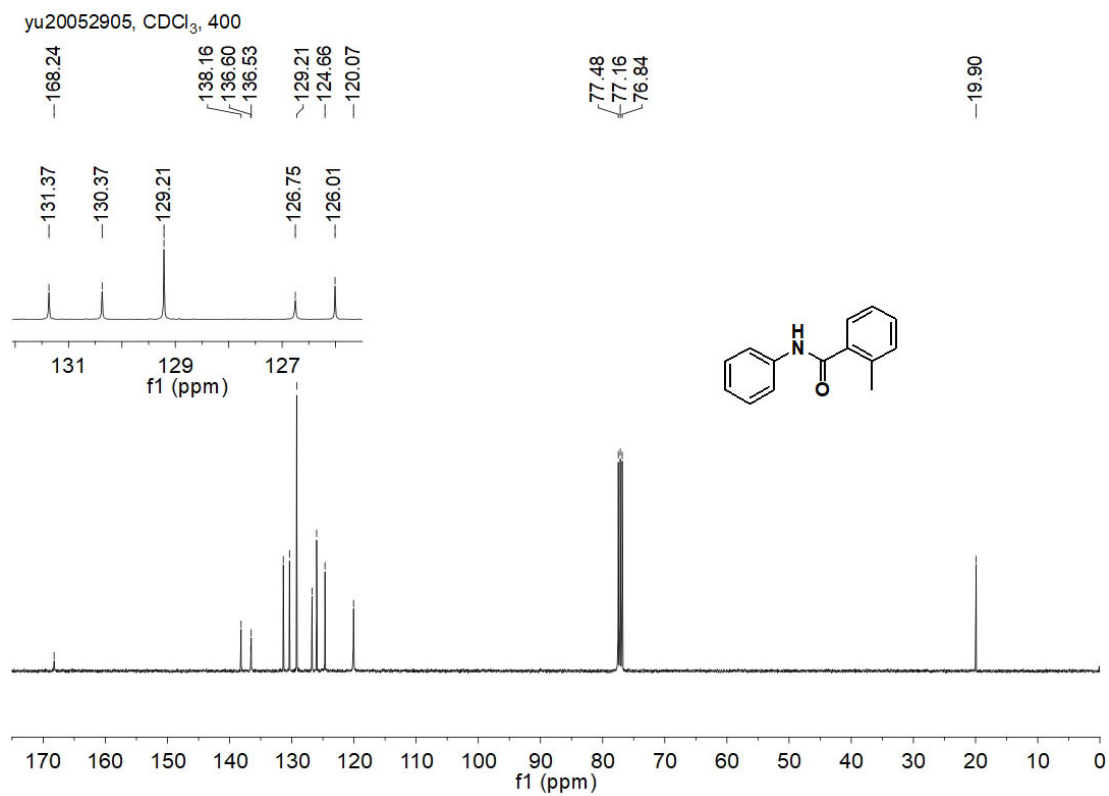


¹³C NMR of 3t

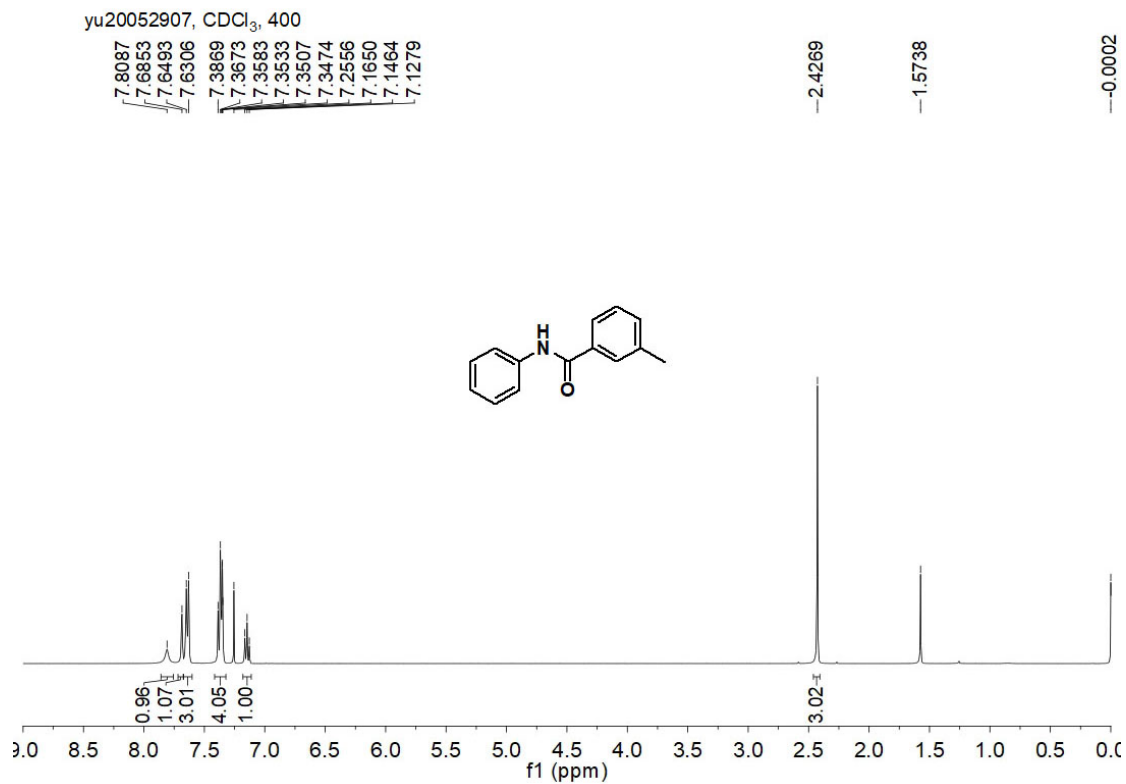
IV. Copies of NMR spectra for 4



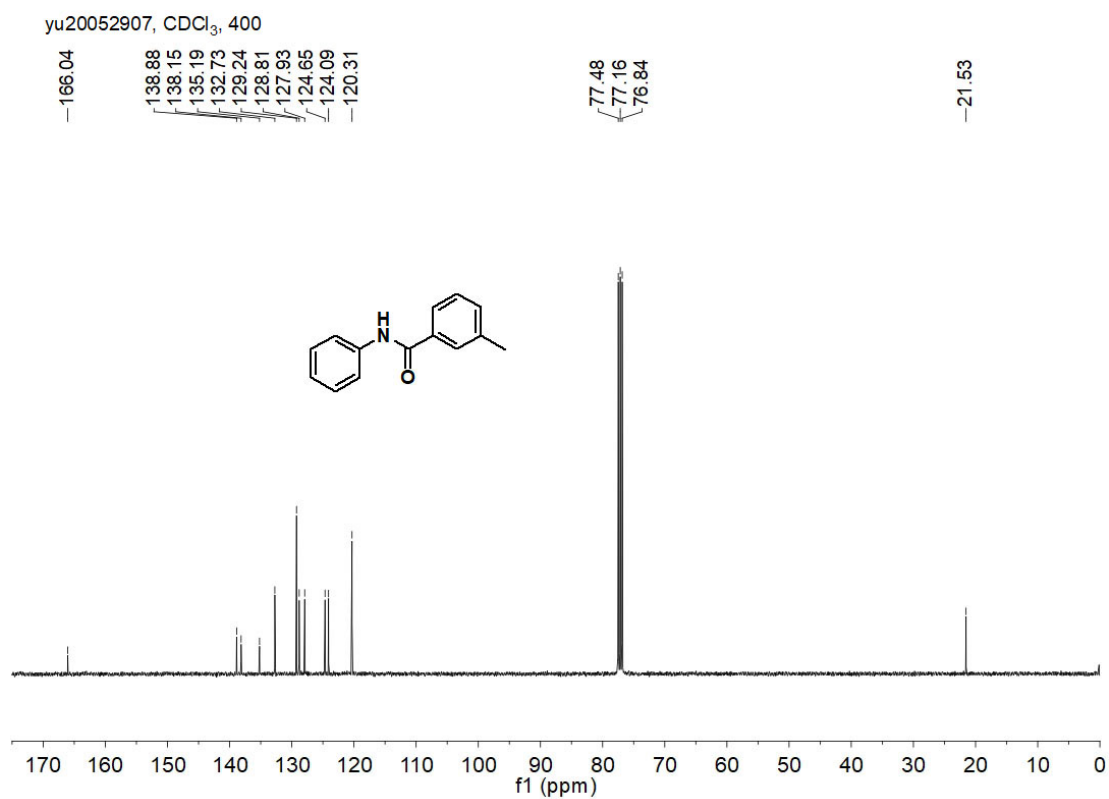
¹H NMR of 4a



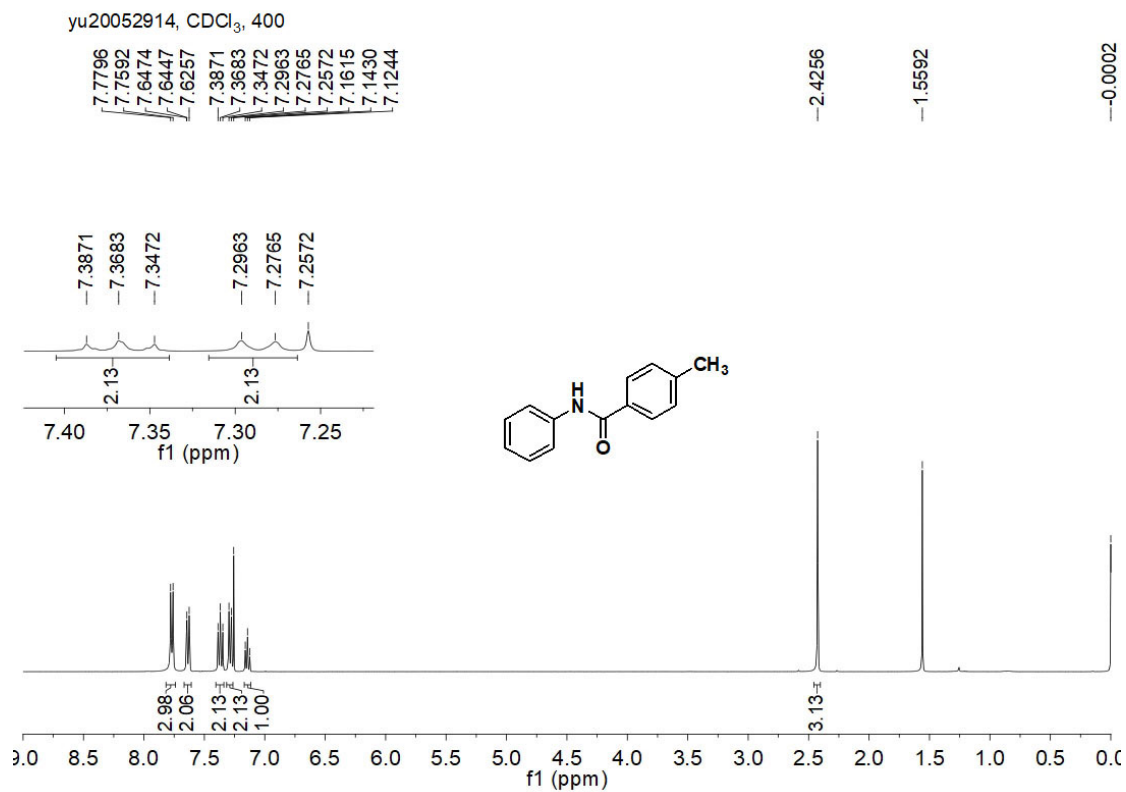
¹³C NMR of 4a



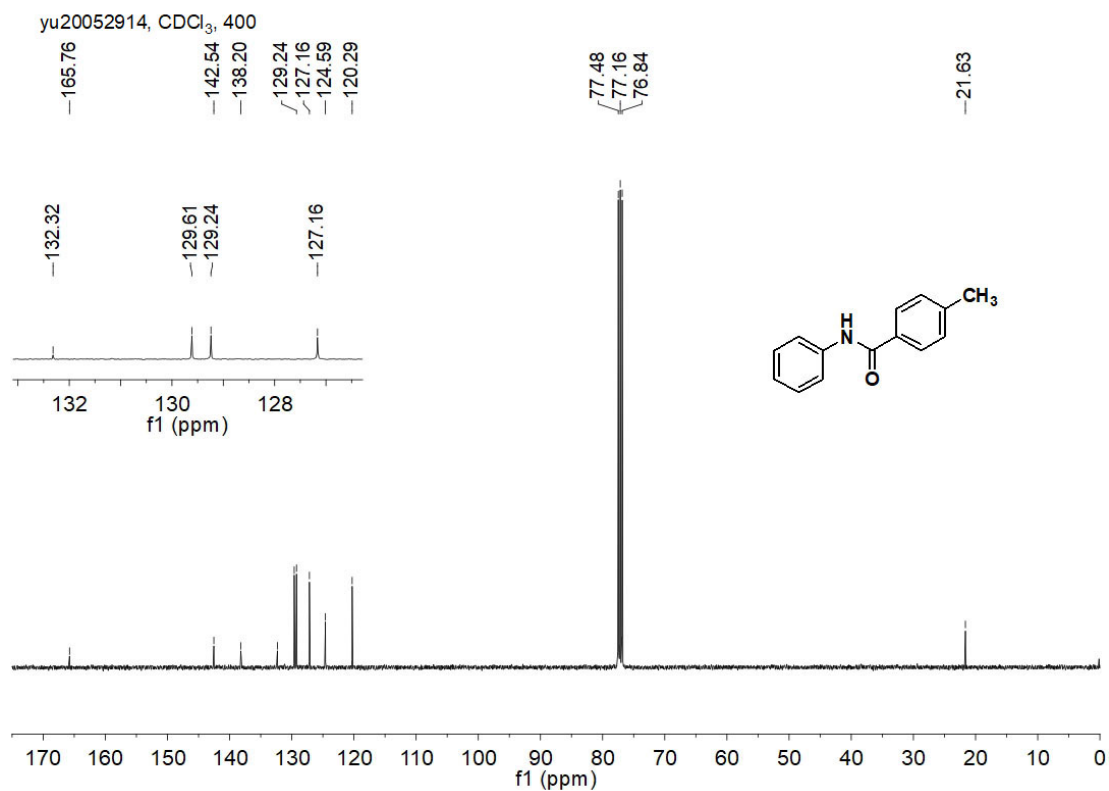
¹H NMR of **4b**



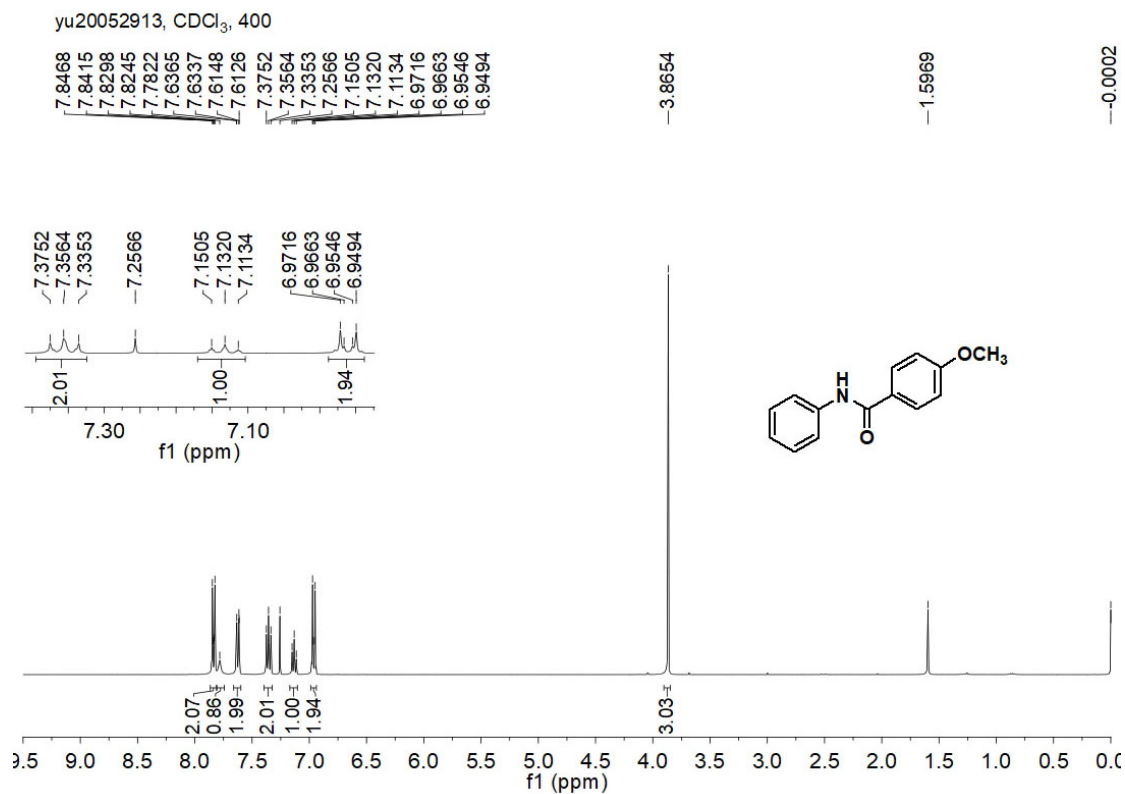
¹³C NMR of **4b**



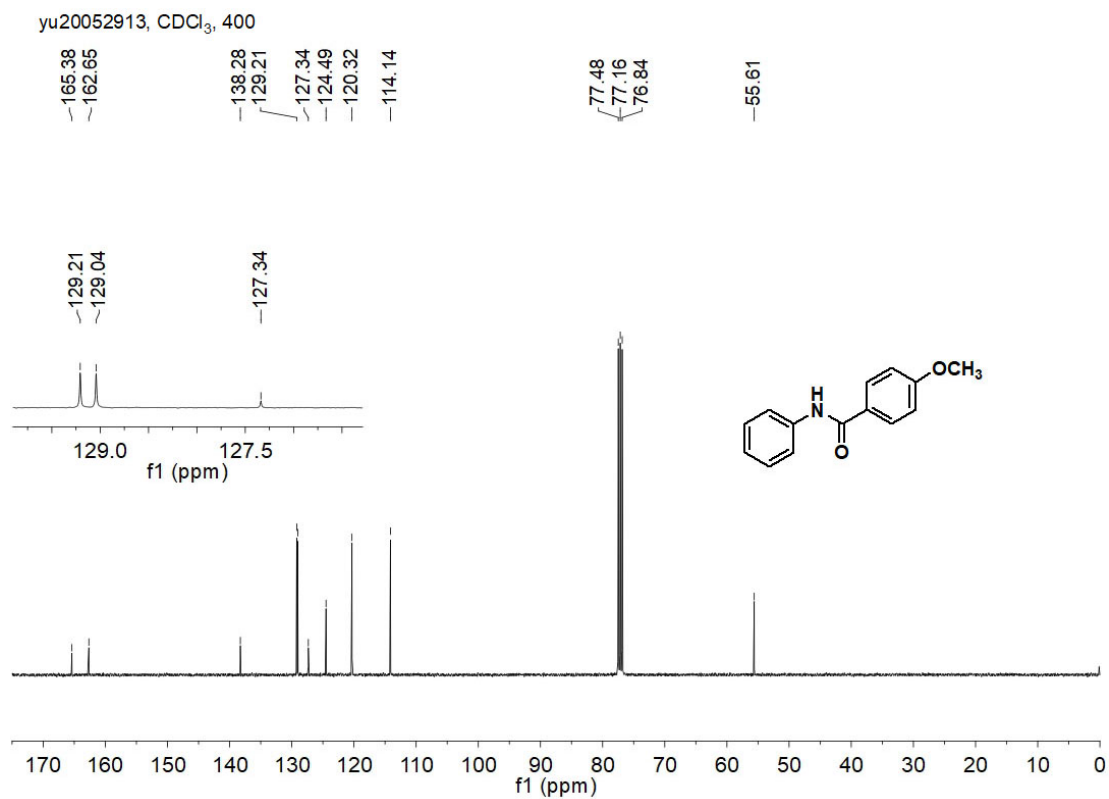
¹HNMR of **4c**



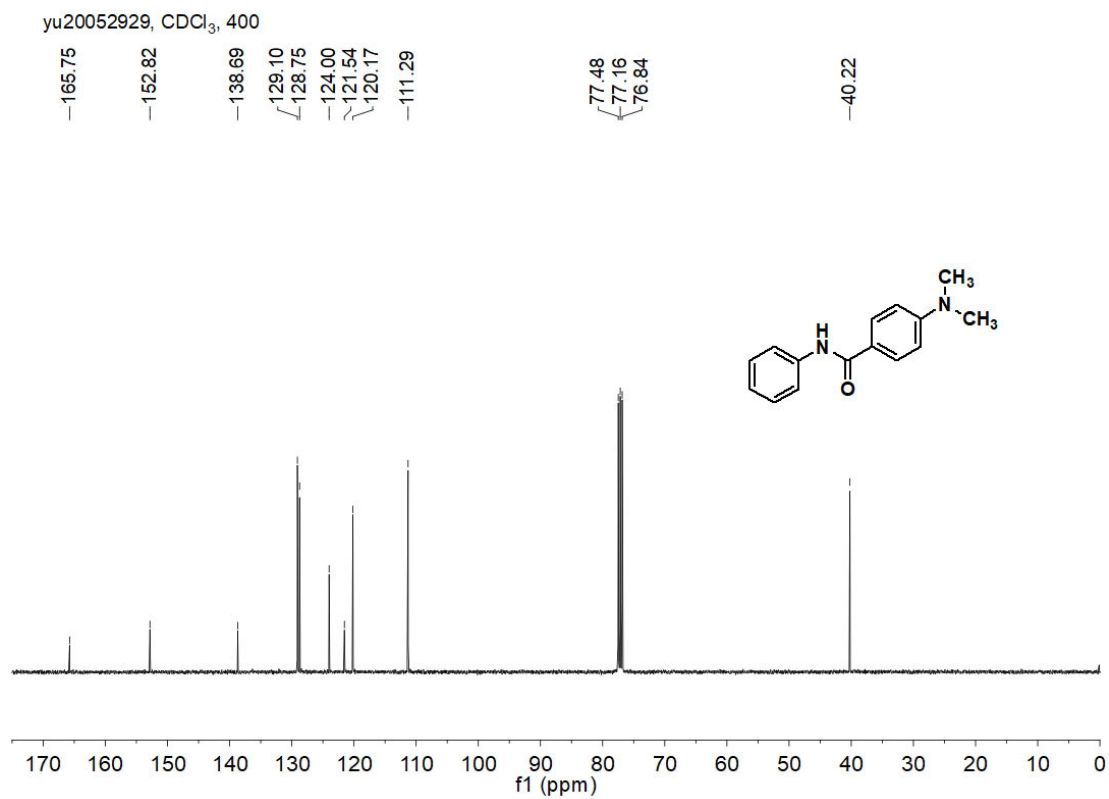
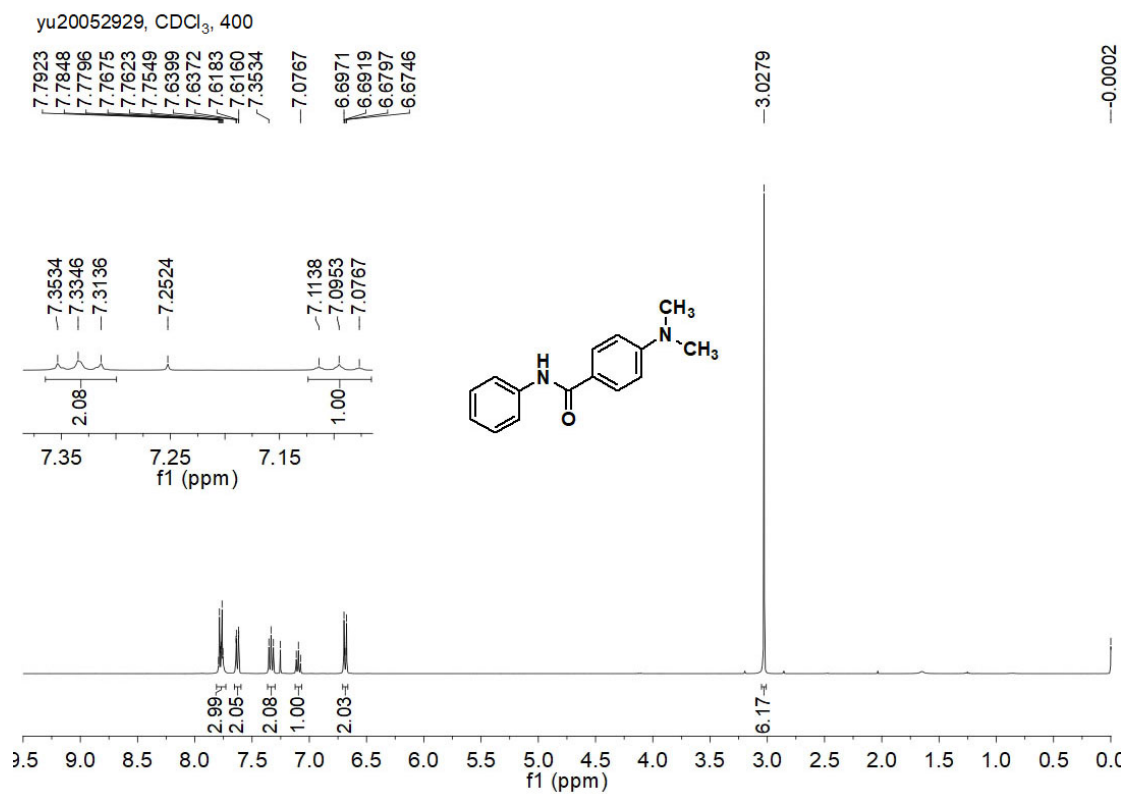
¹³CNMR of **4c**

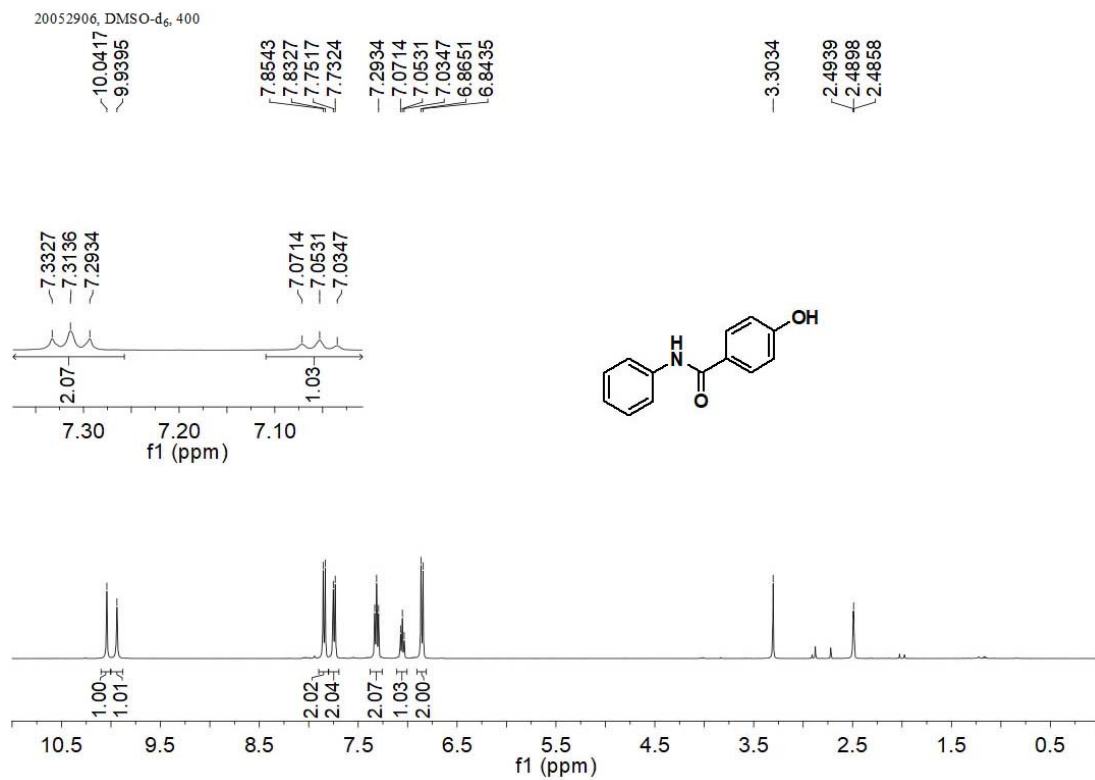


¹H NMR of **4d**

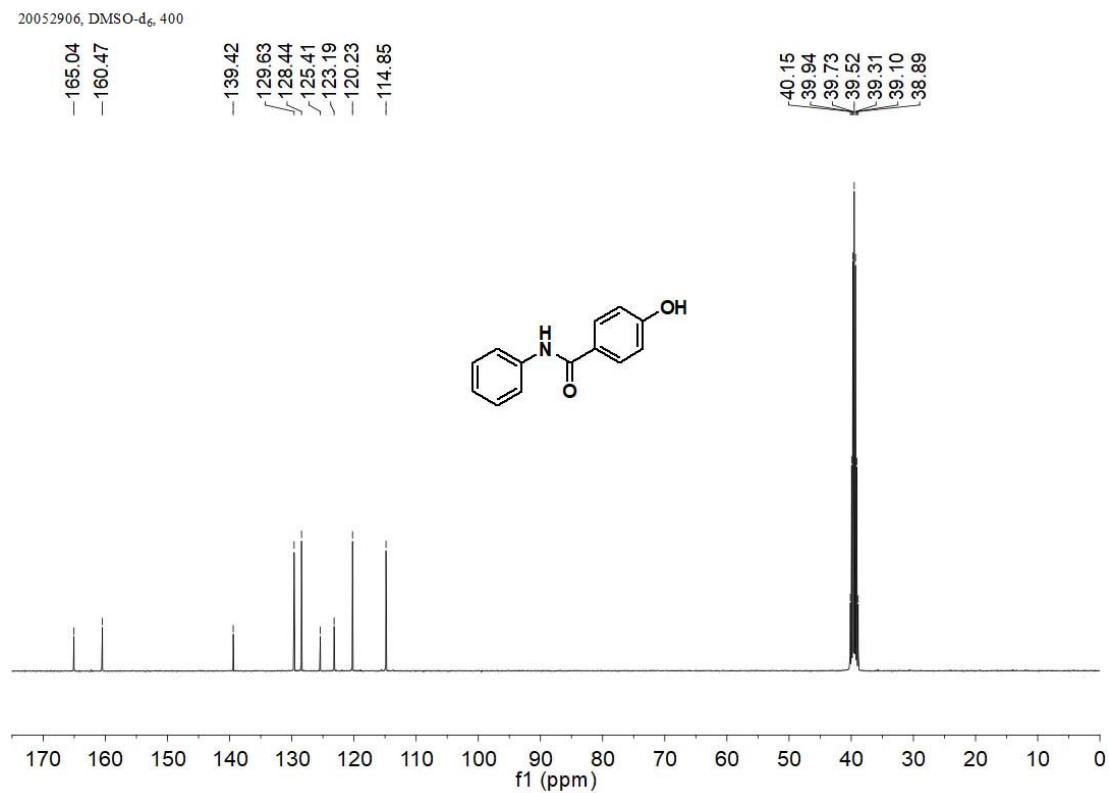


¹³C NMR of **4d**

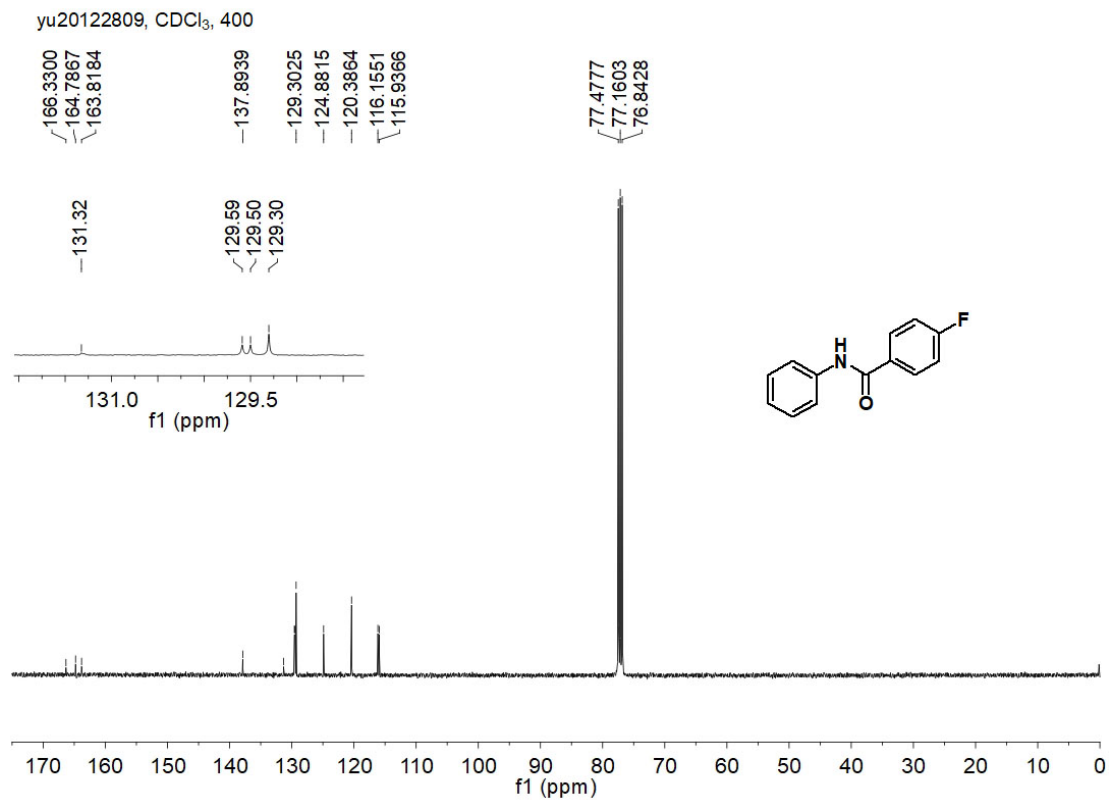
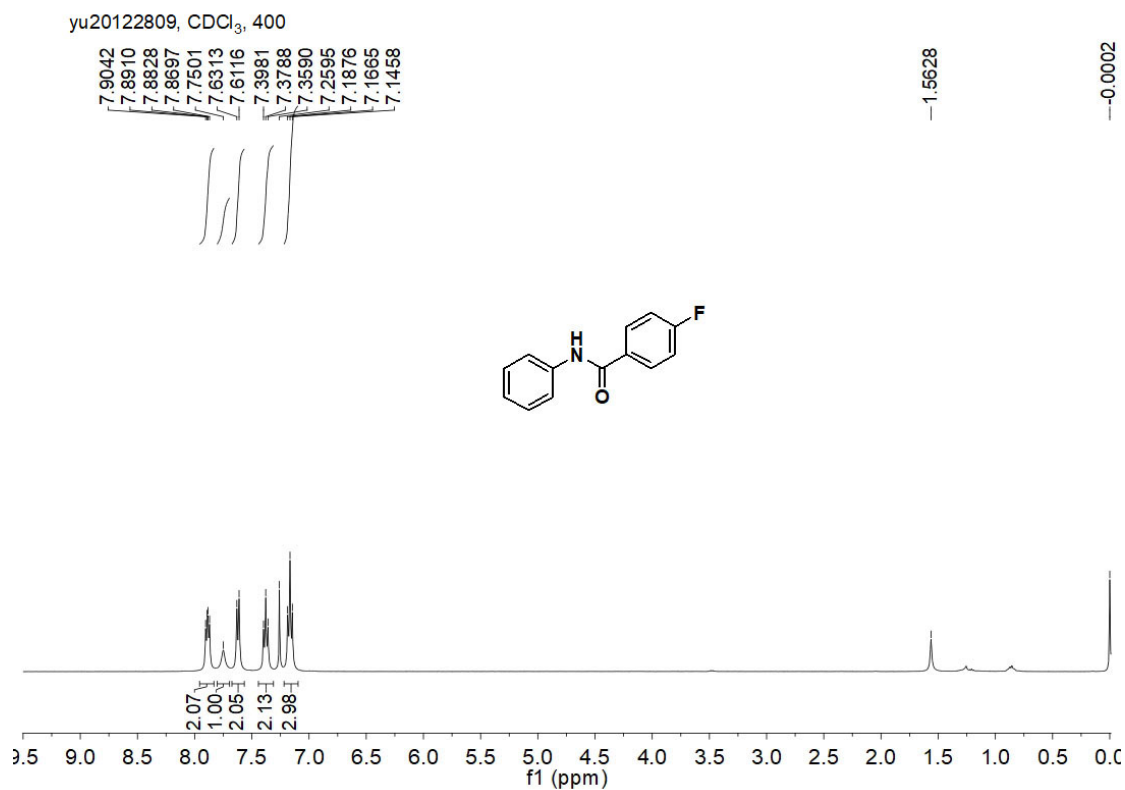


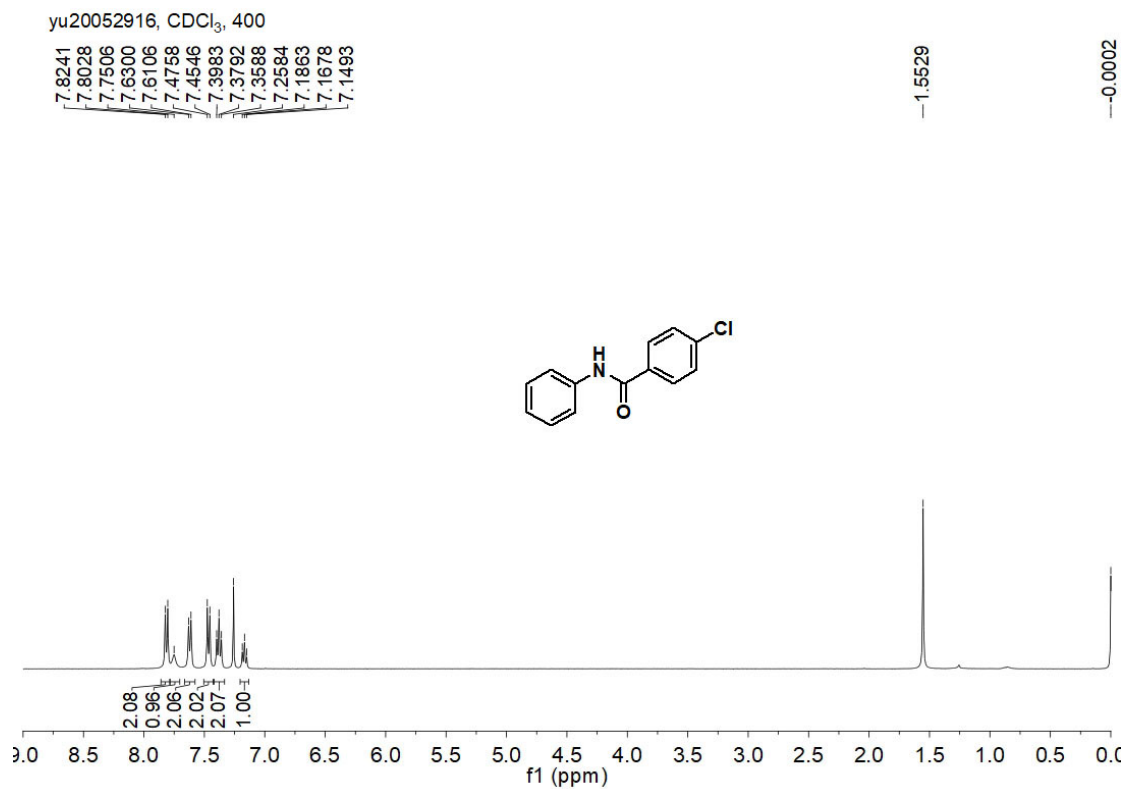


¹H NMR of **4f**

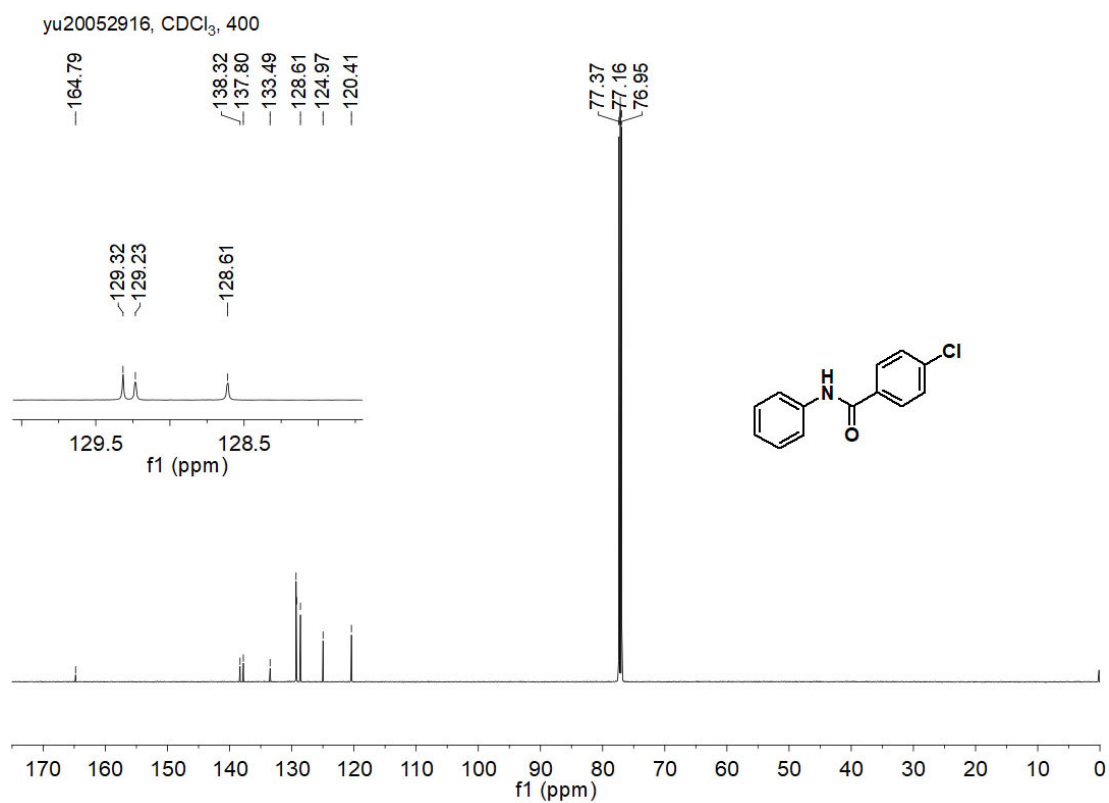


¹³C NMR of **4f**

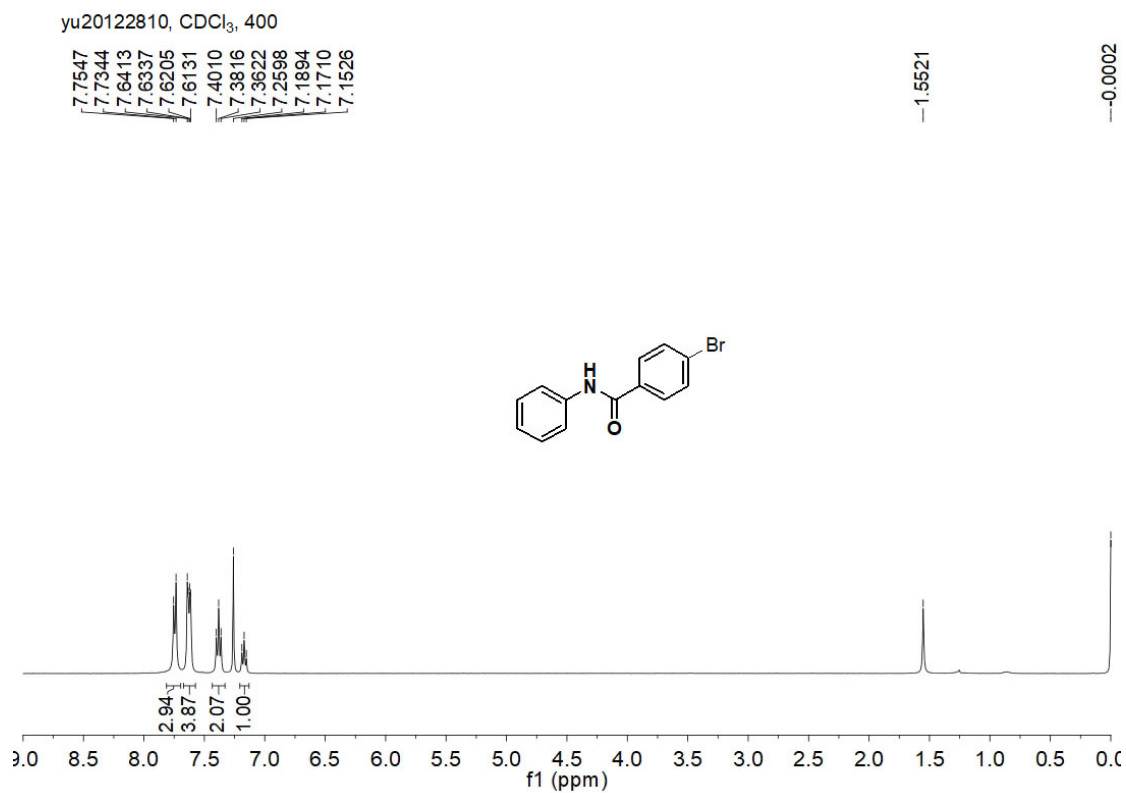




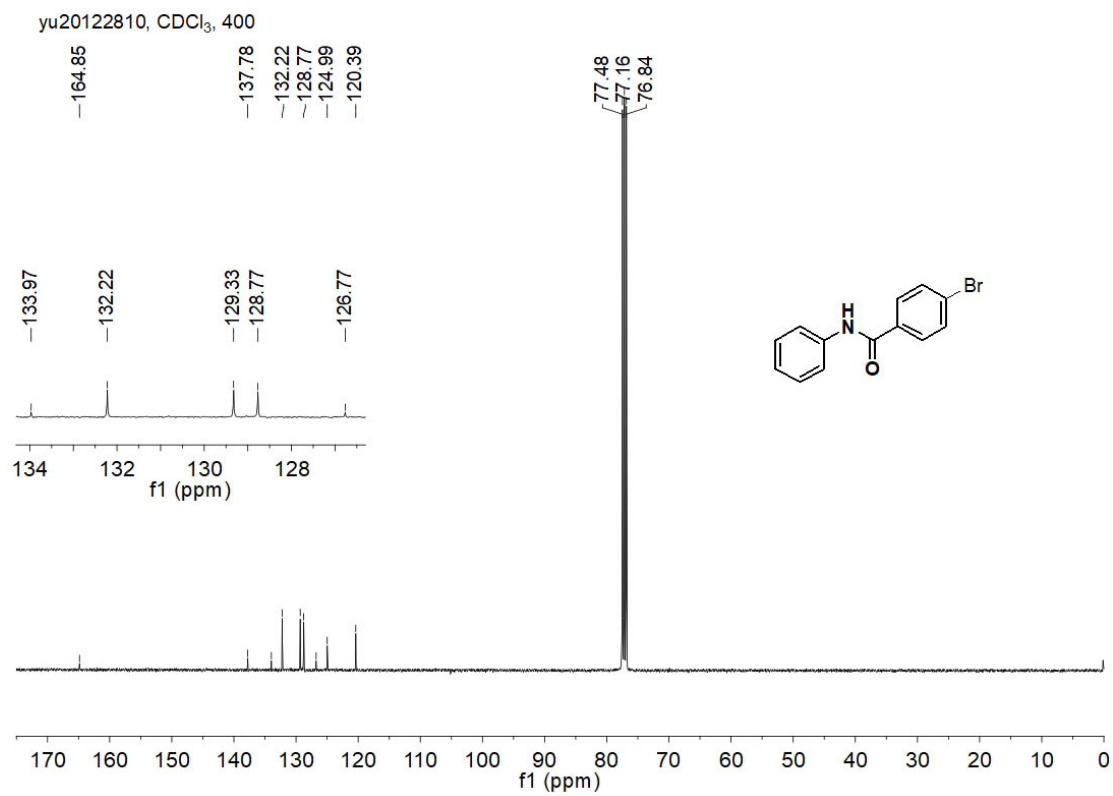
¹HNMR of **4h**



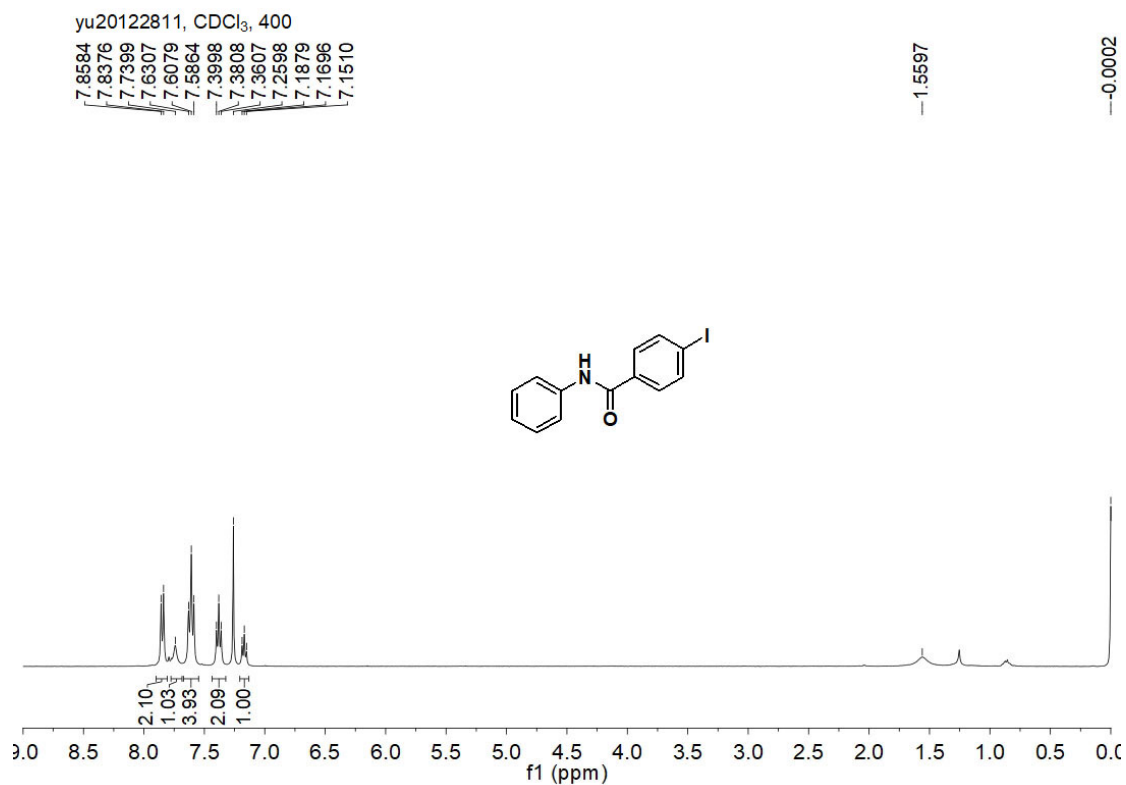
¹HNMR of **4h**



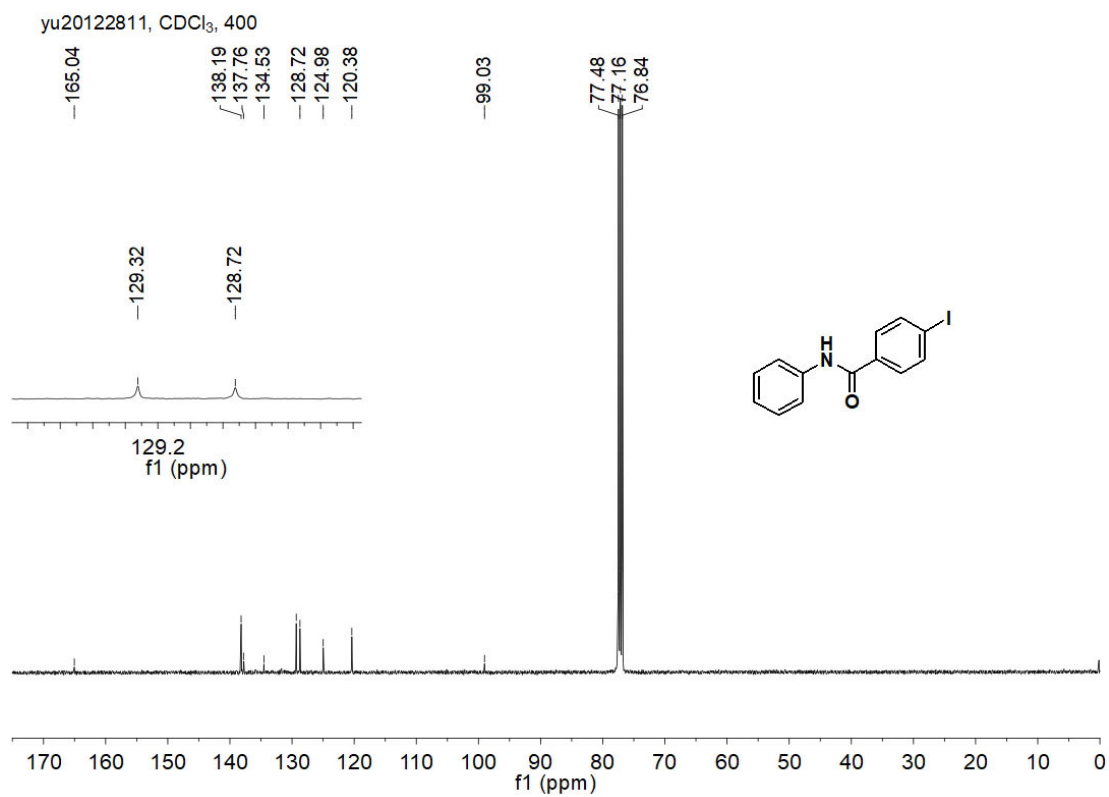
¹HNMR of **4i**



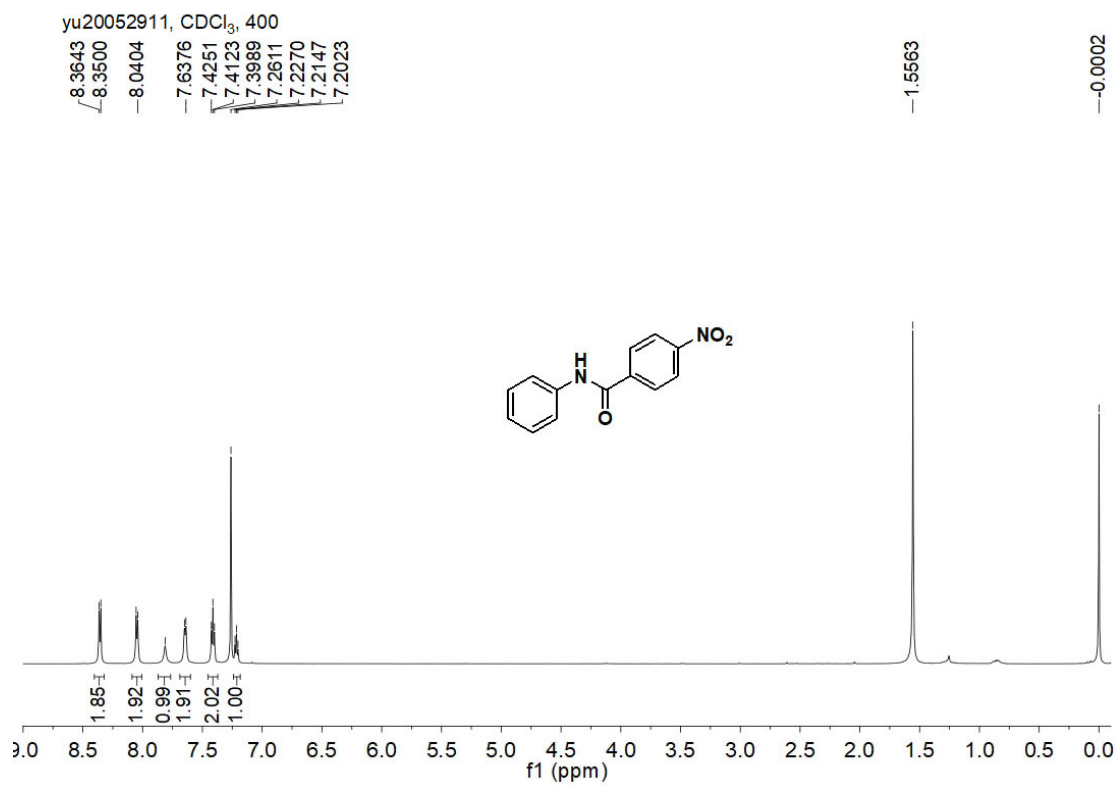
¹³CNMR of **4i**



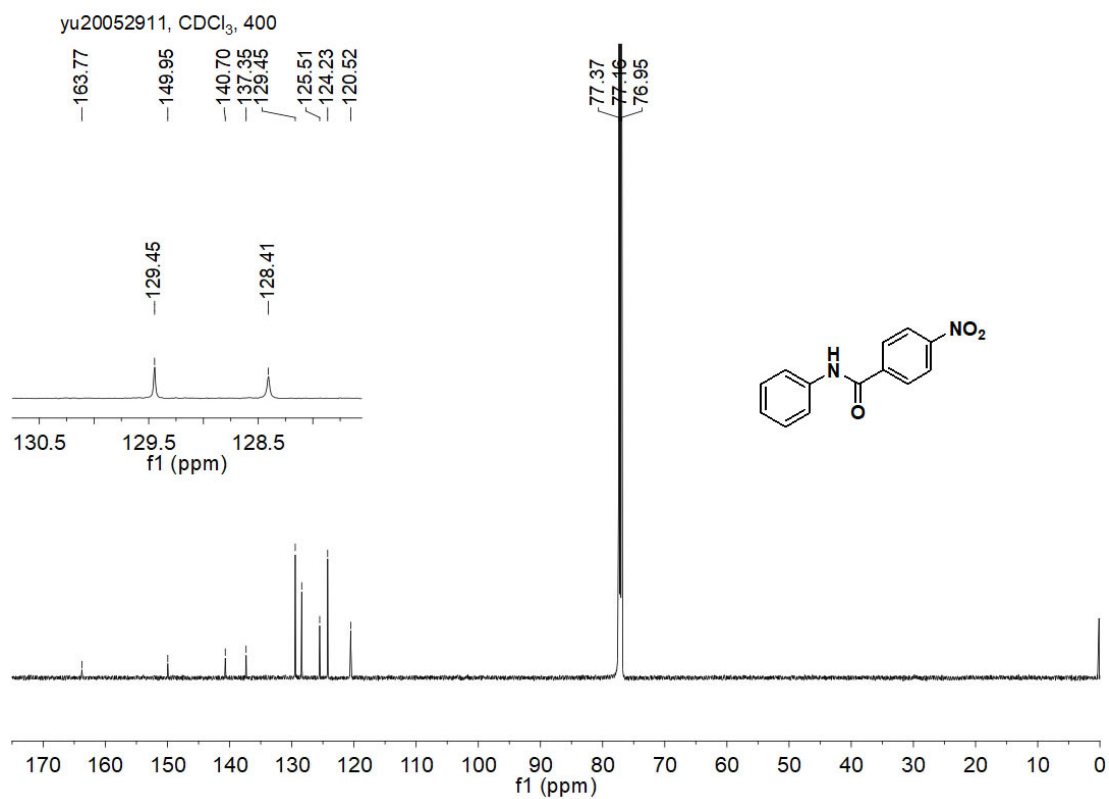
¹H NMR of **4j**



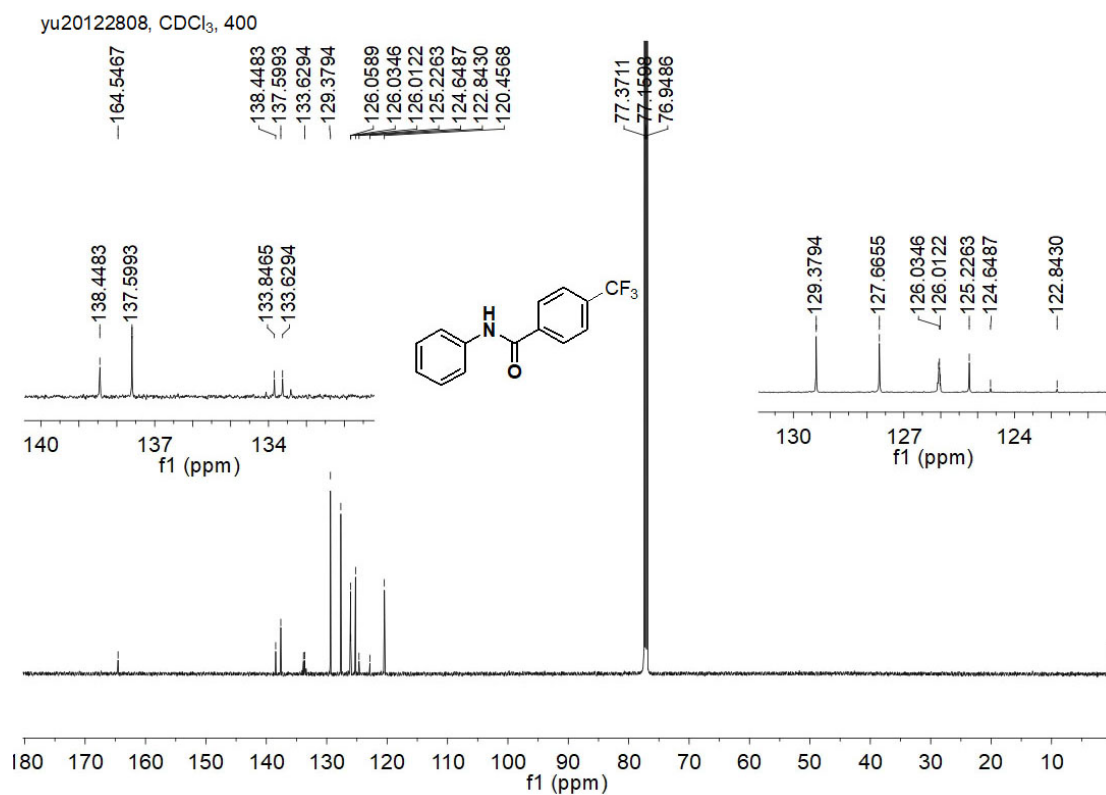
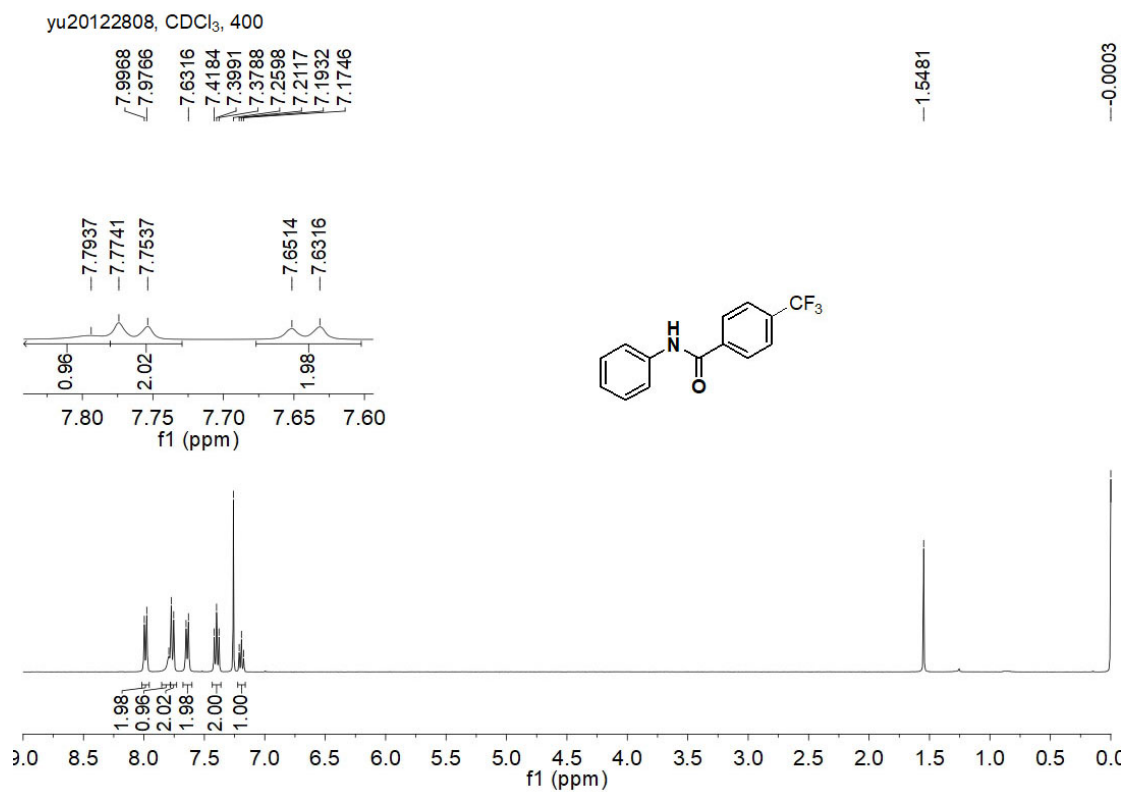
¹³C NMR of **4j**

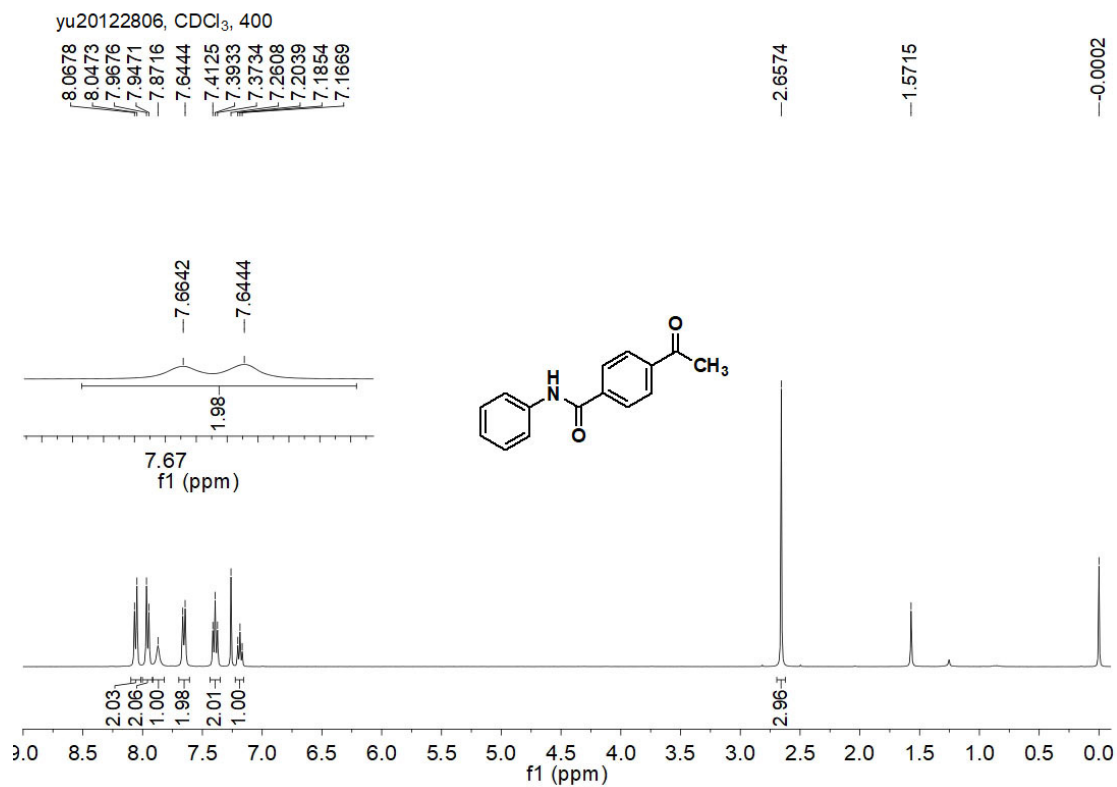


¹H NMR of **4k**

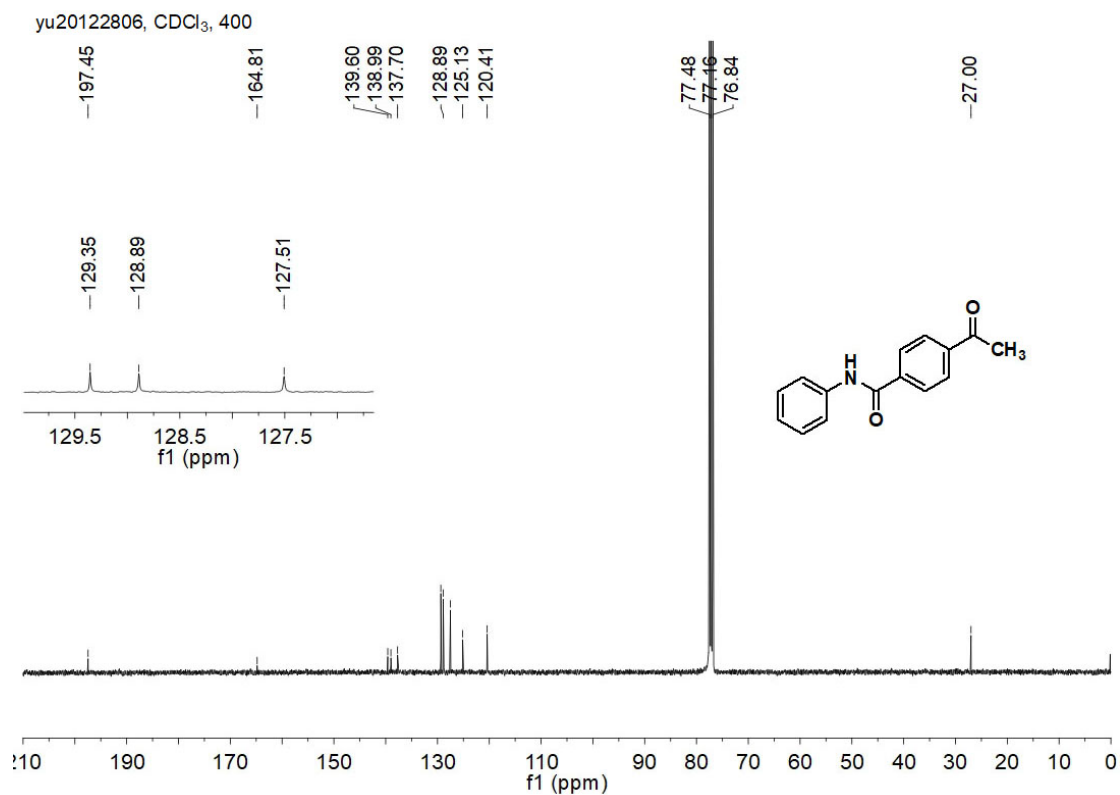


¹³C NMR of **4k**

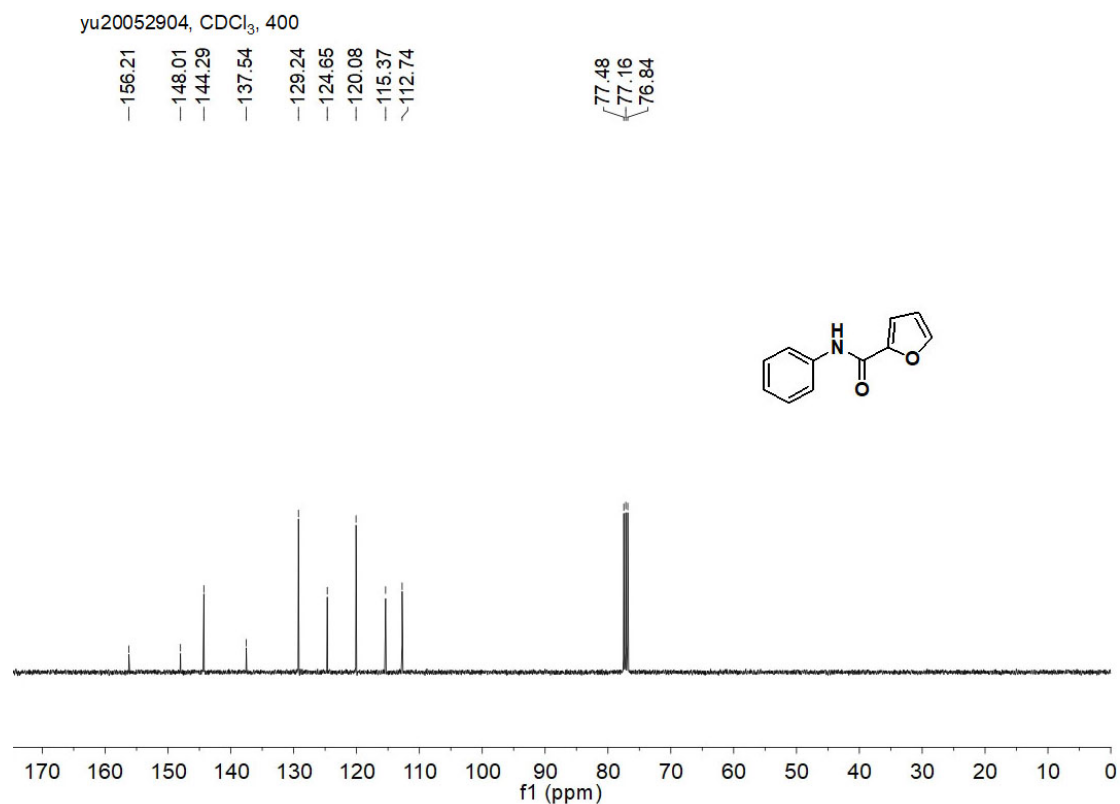
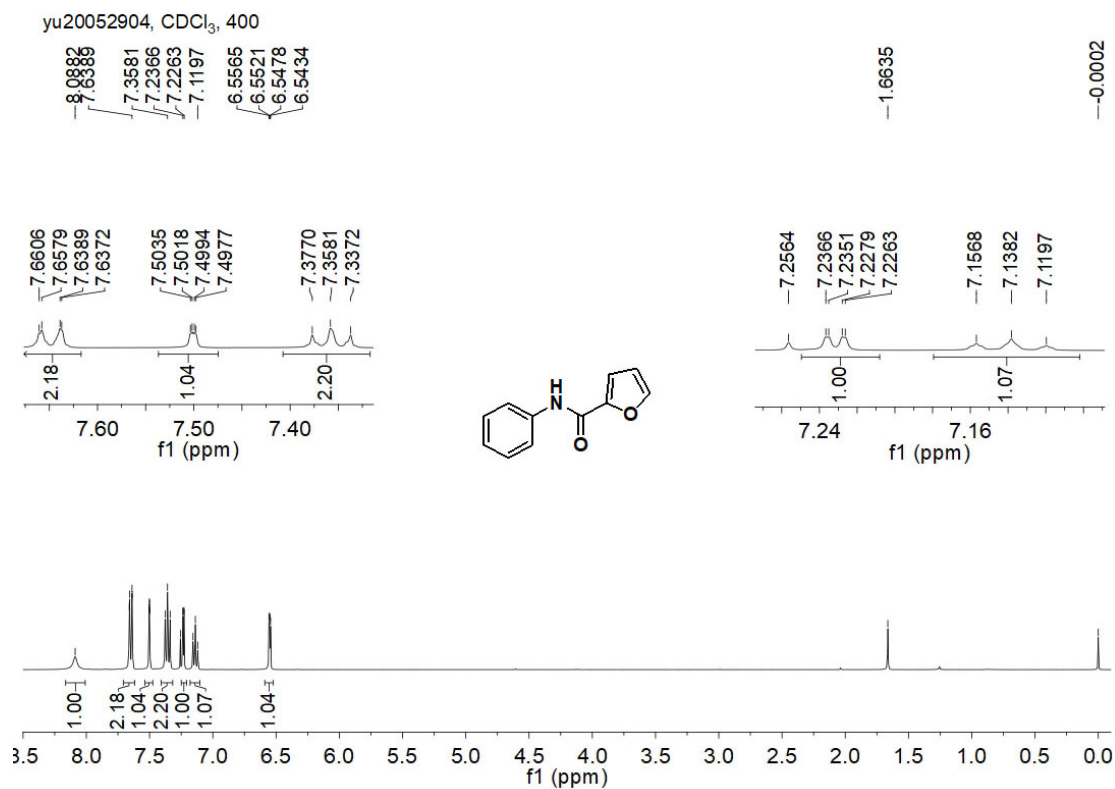




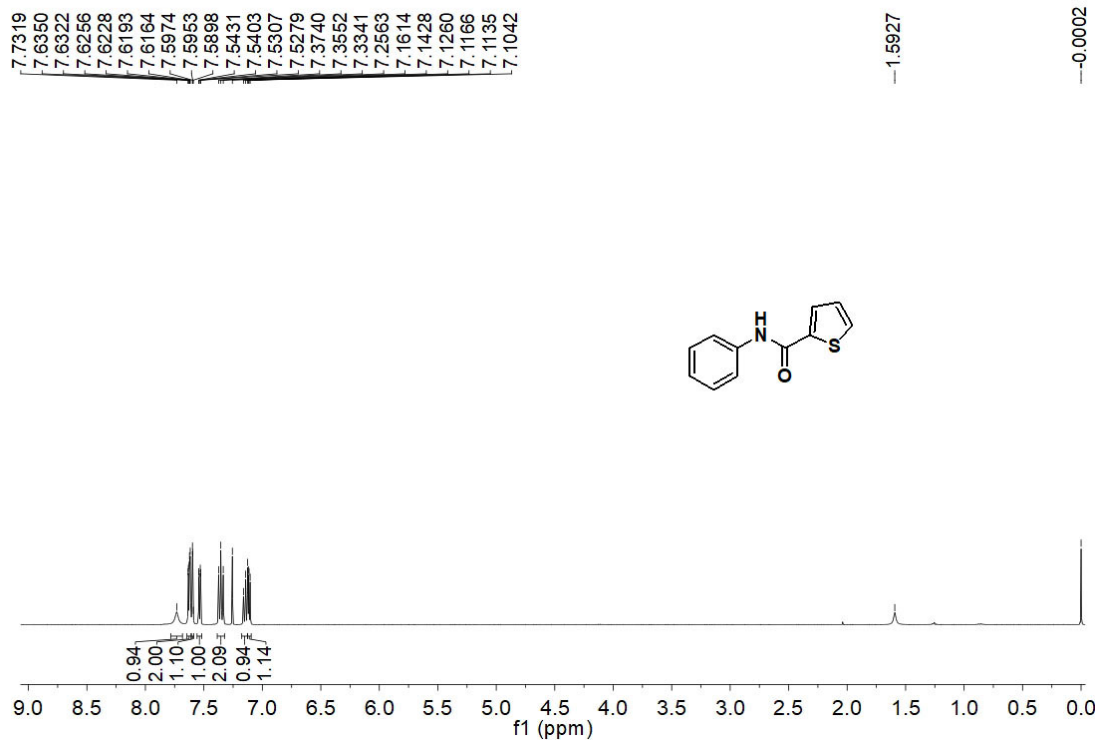
¹HNMR of **4m**



¹³CNMR of **4m**

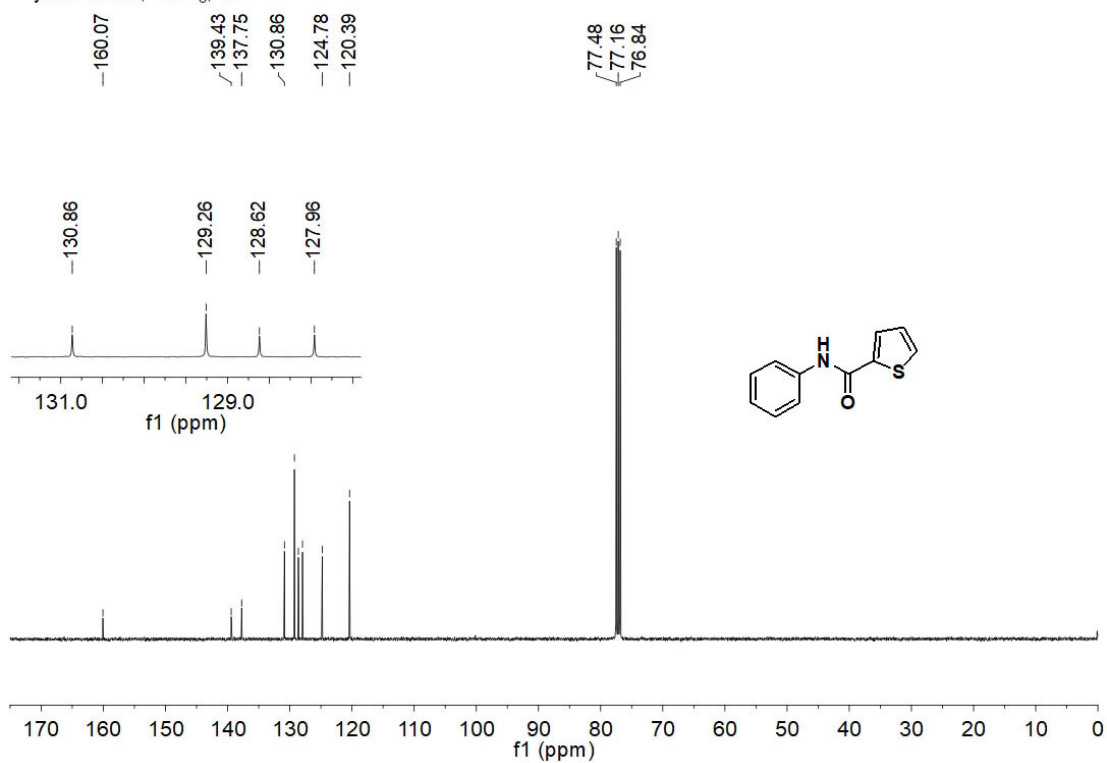


yu20052912, CDCl₃, 400

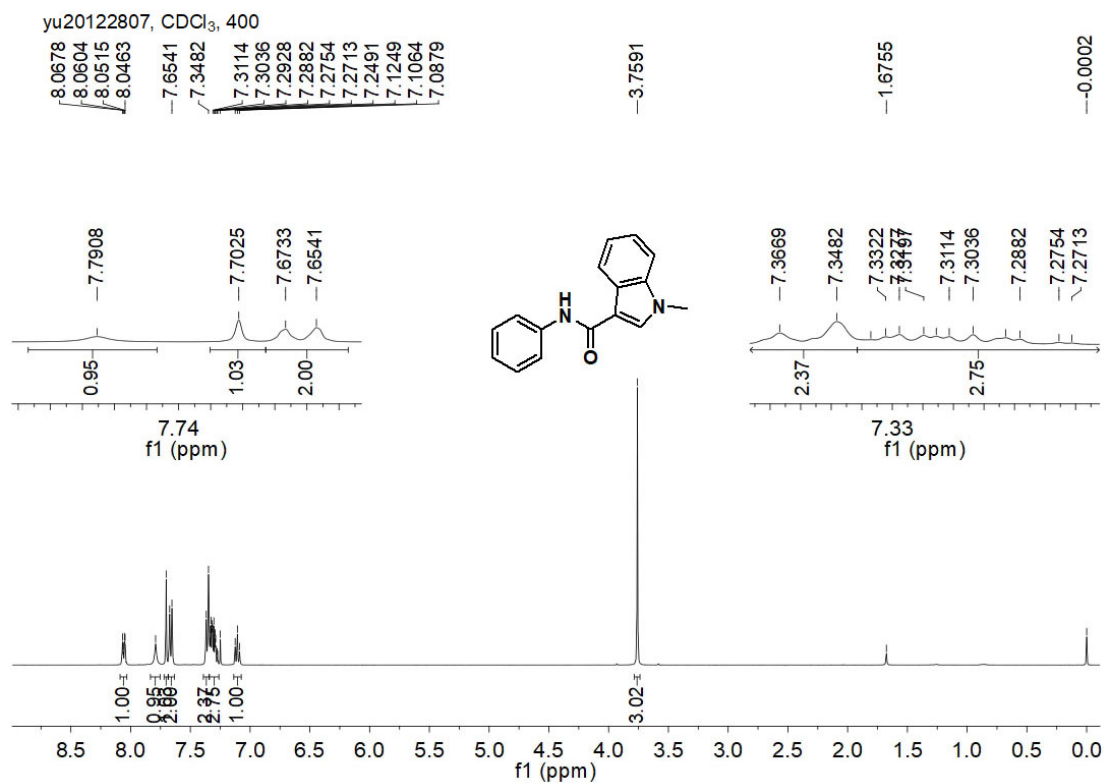


¹H NMR of **4o**

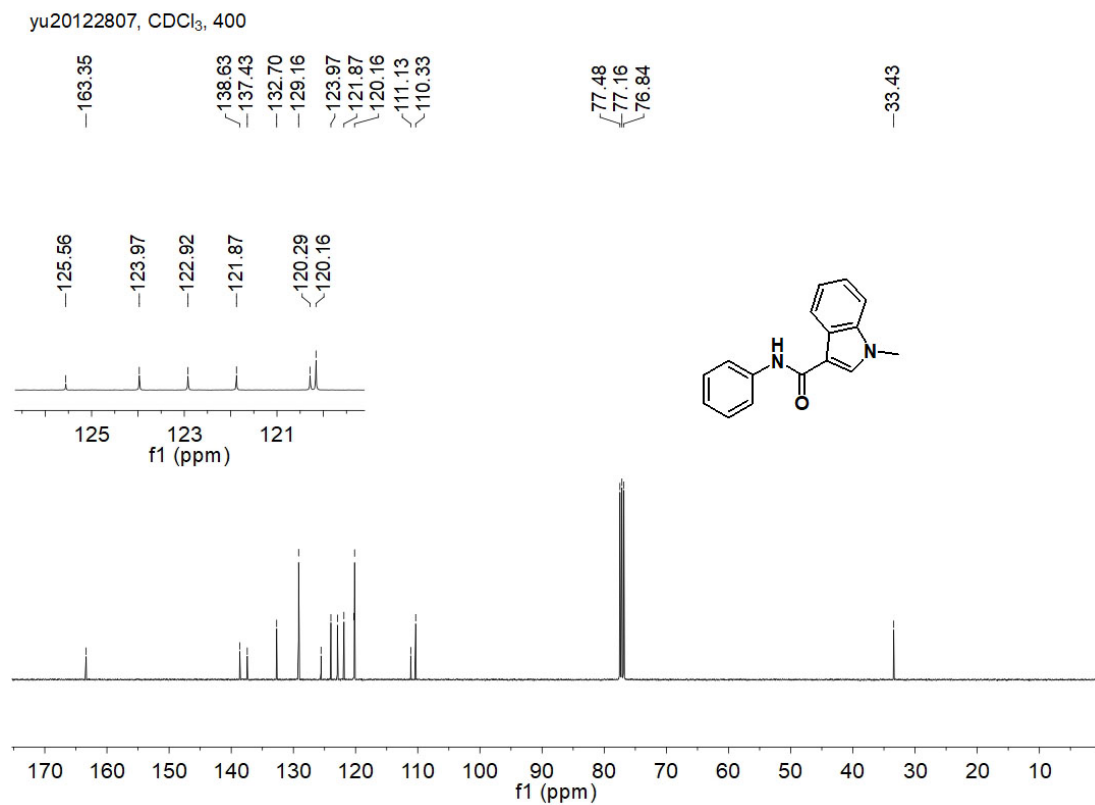
yu20052912, CDCl₃, 400



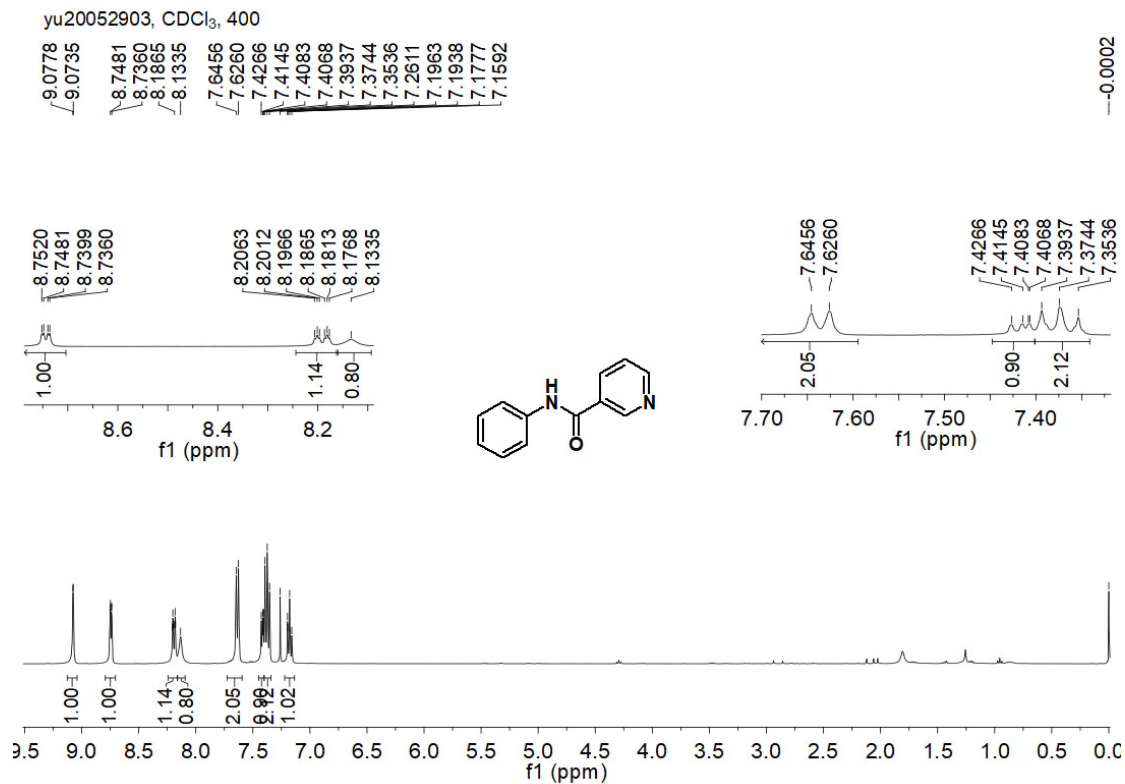
¹³C NMR of **4o**



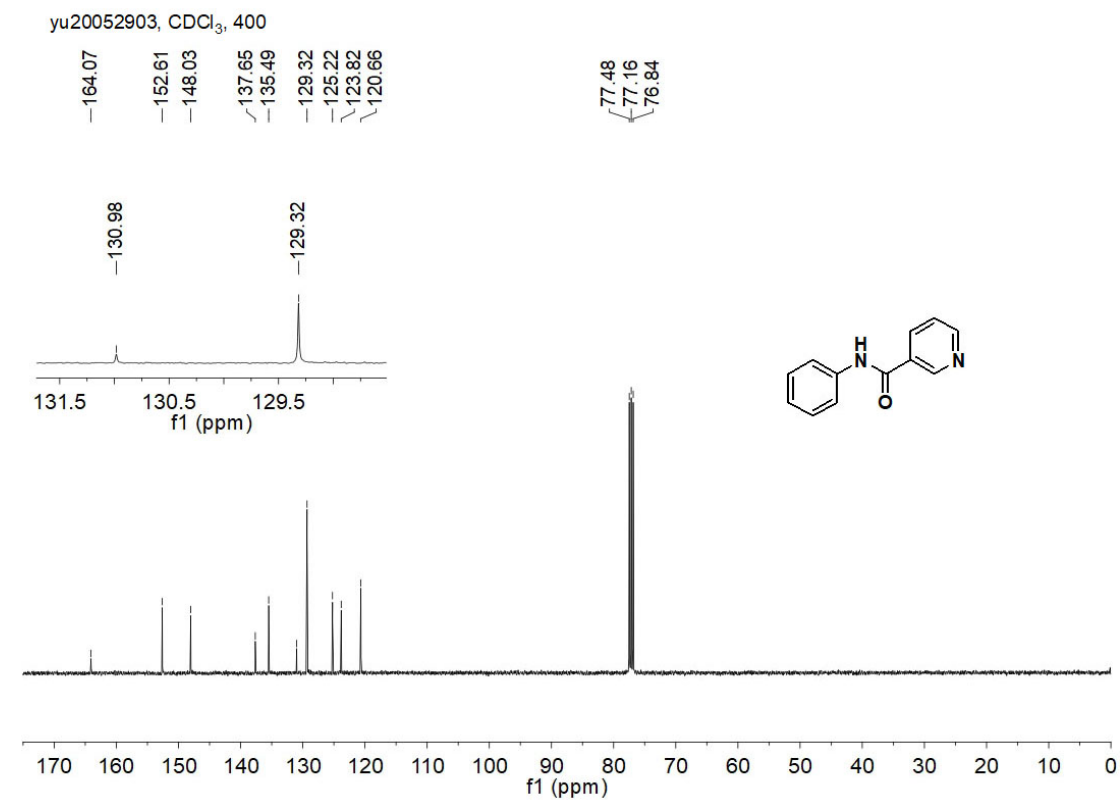
¹HNMR of **4p**



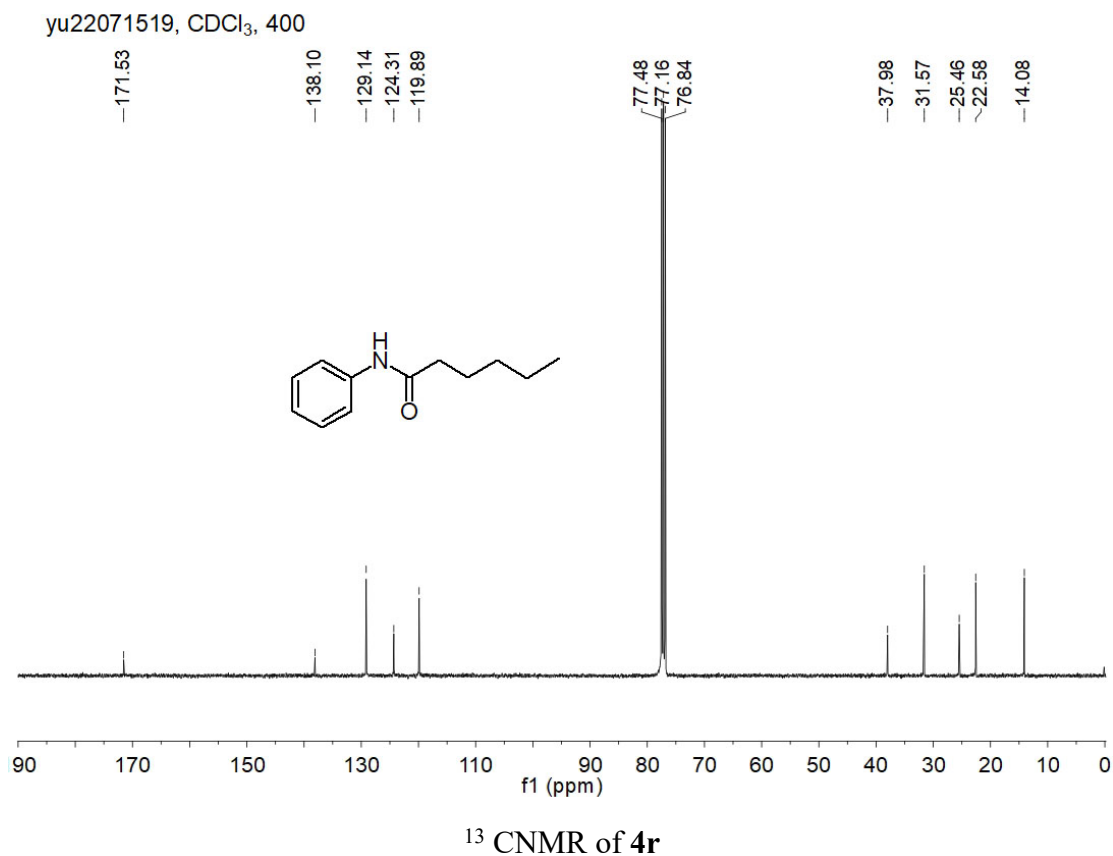
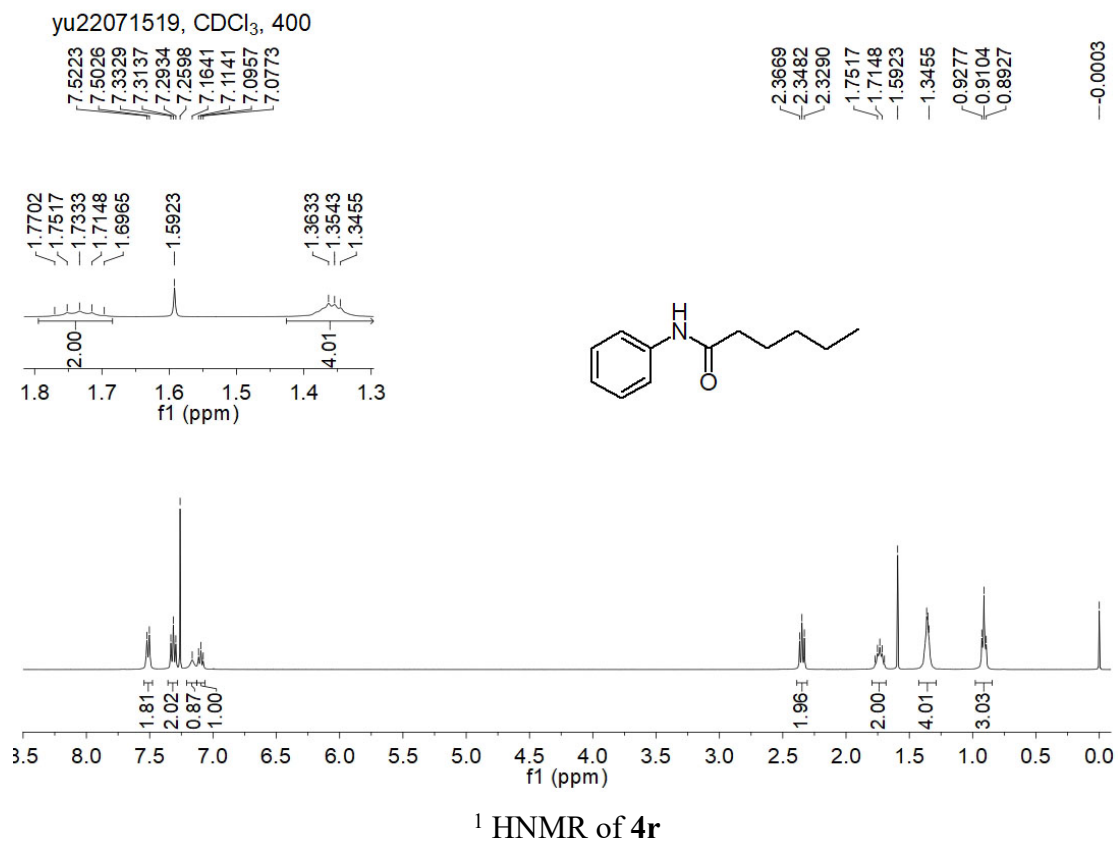
¹³CNMR of **4p**

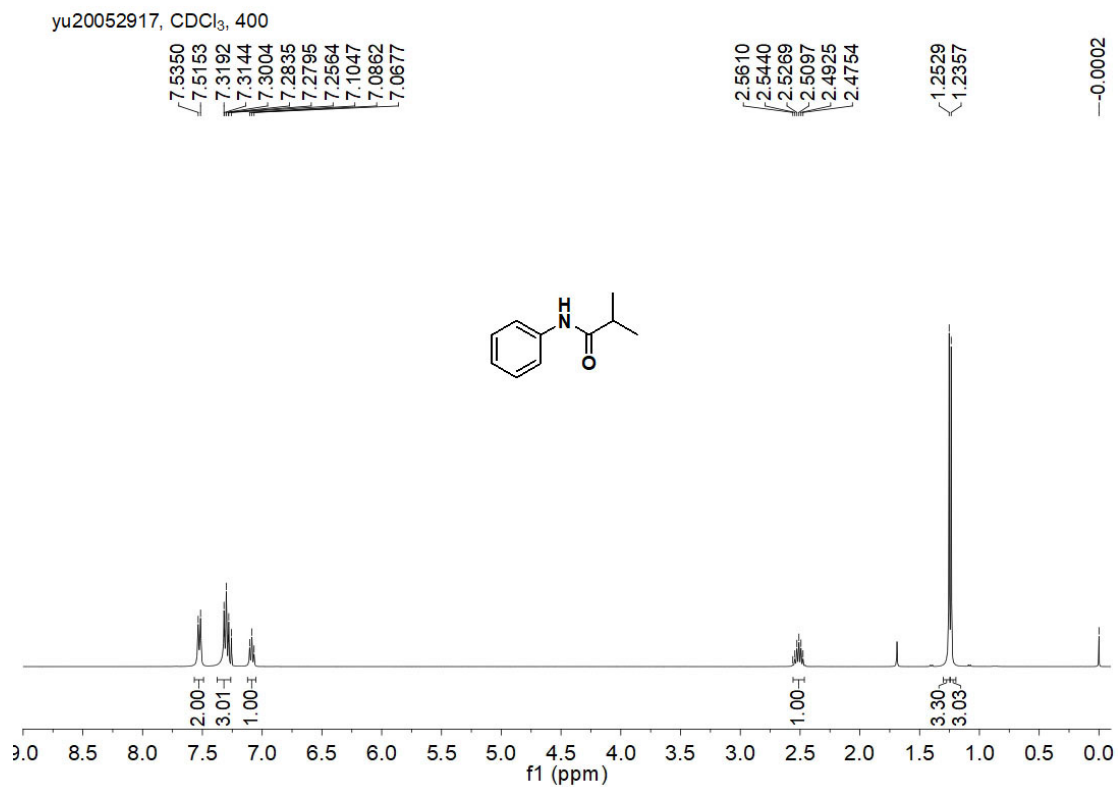


¹H NMR of **4q**

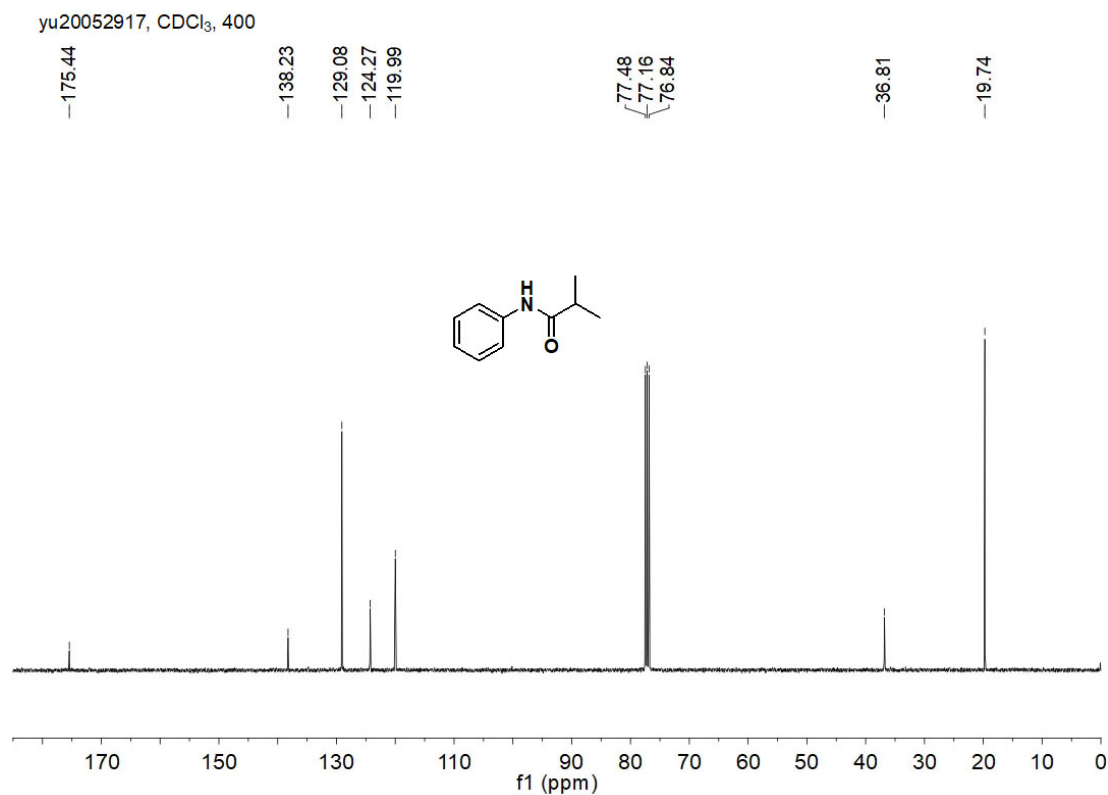


¹³C NMR of **4q**

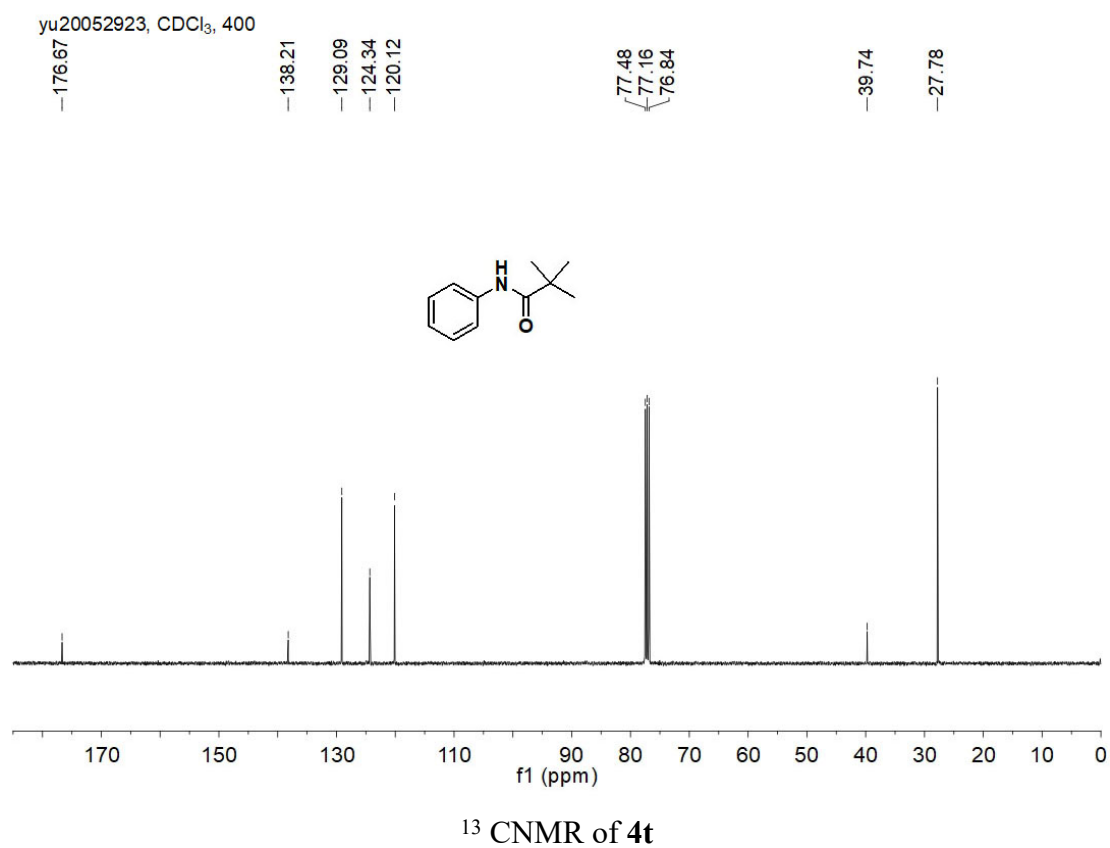
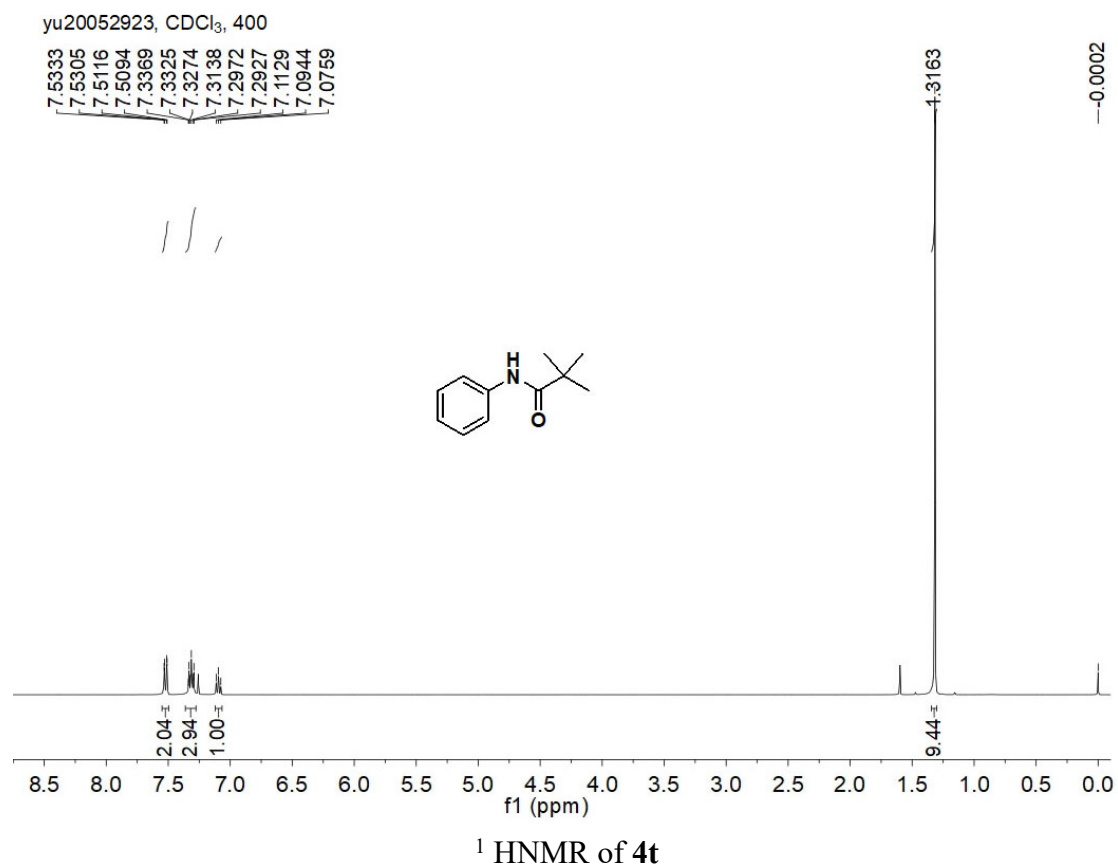




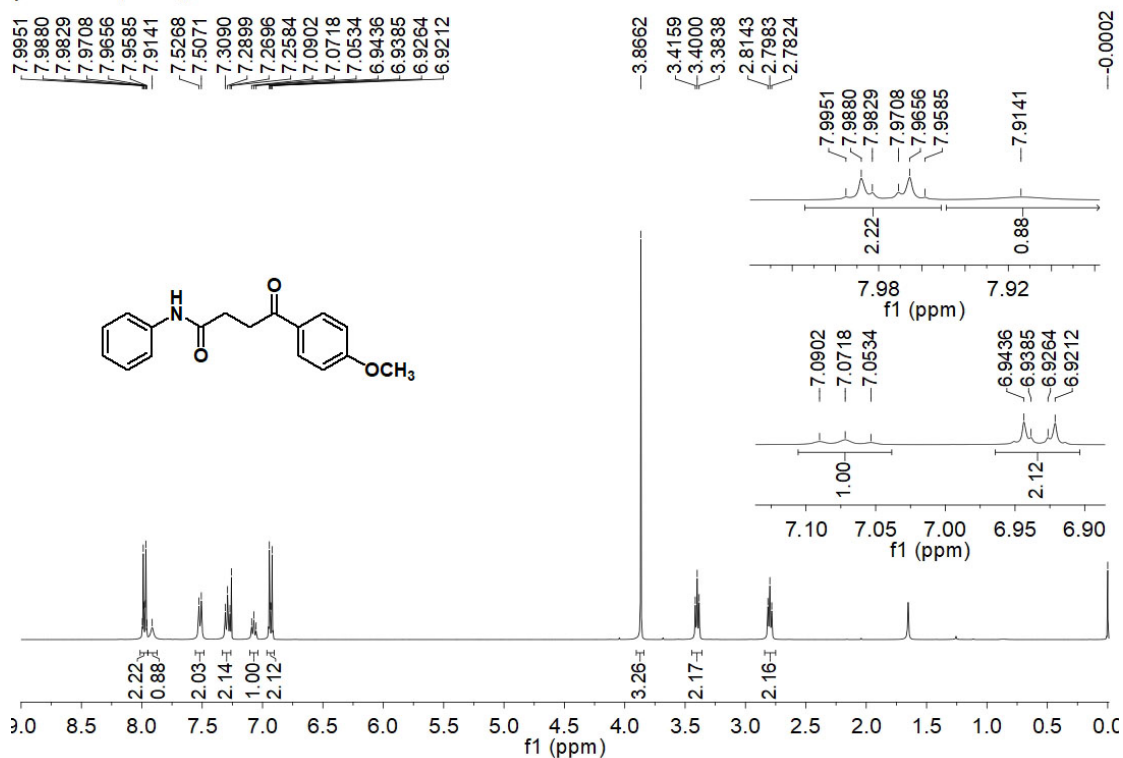
¹HNMR of **4s**



¹³CNMR of **4s**

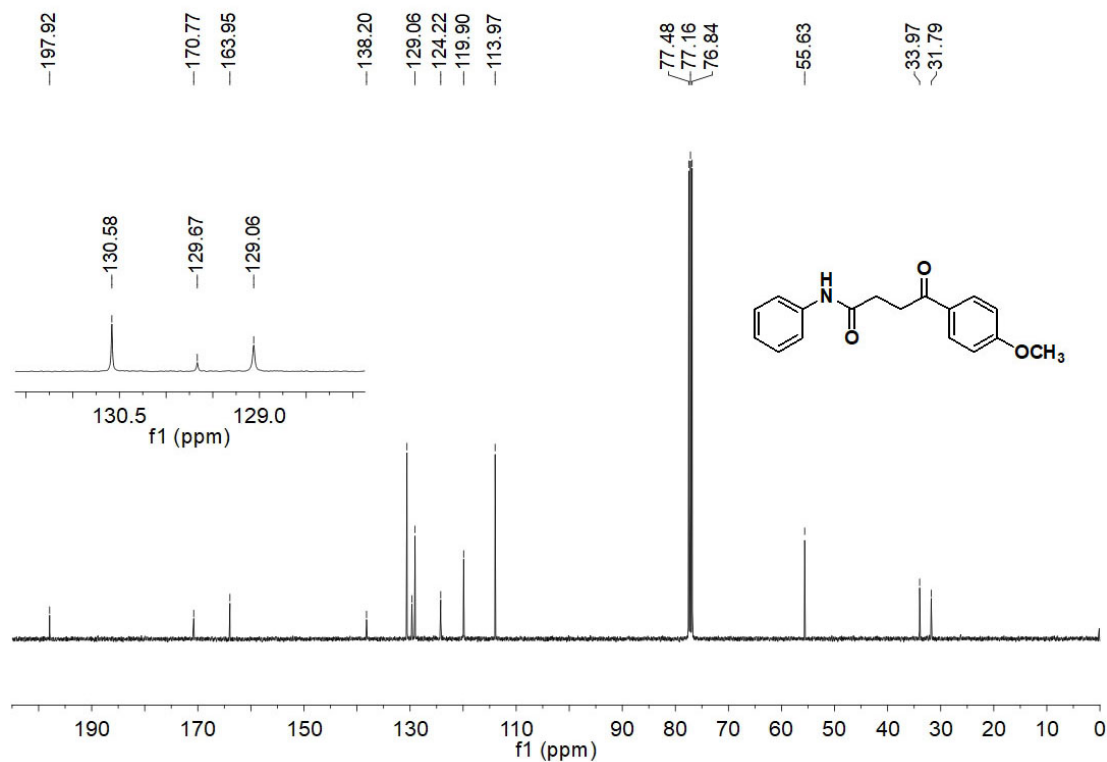


yu20052918, CDCl₃, 400

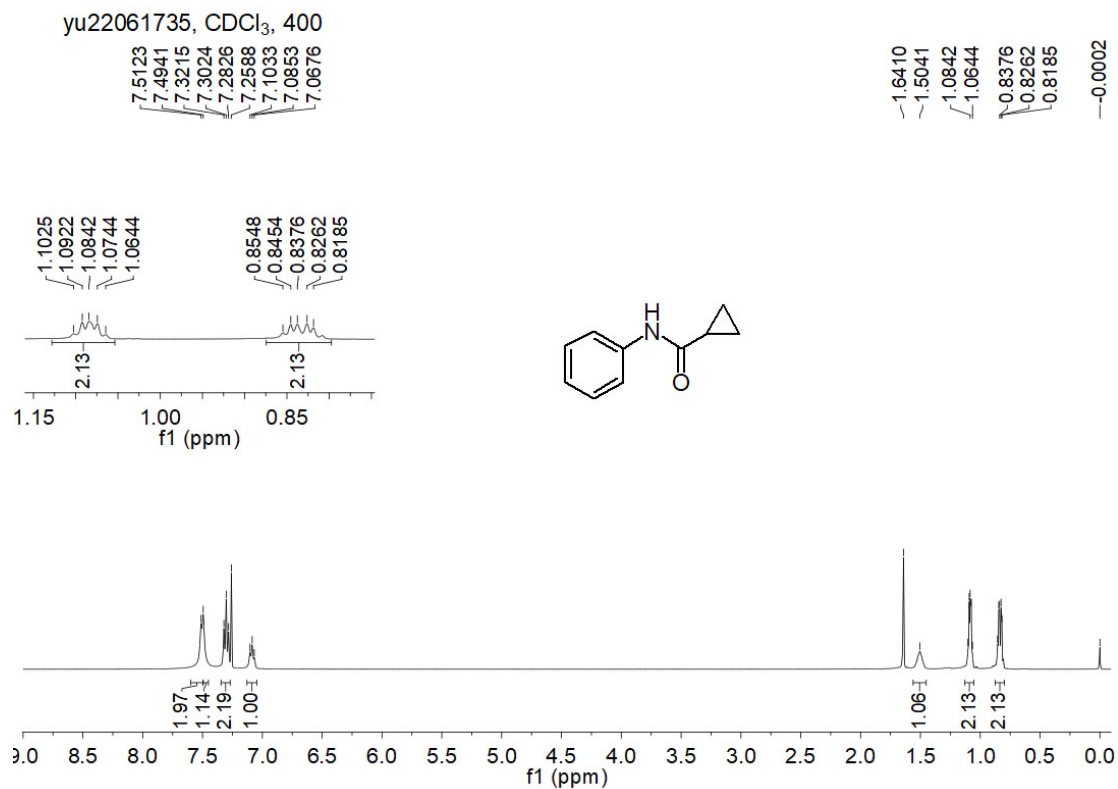


¹HNMR of **4u**

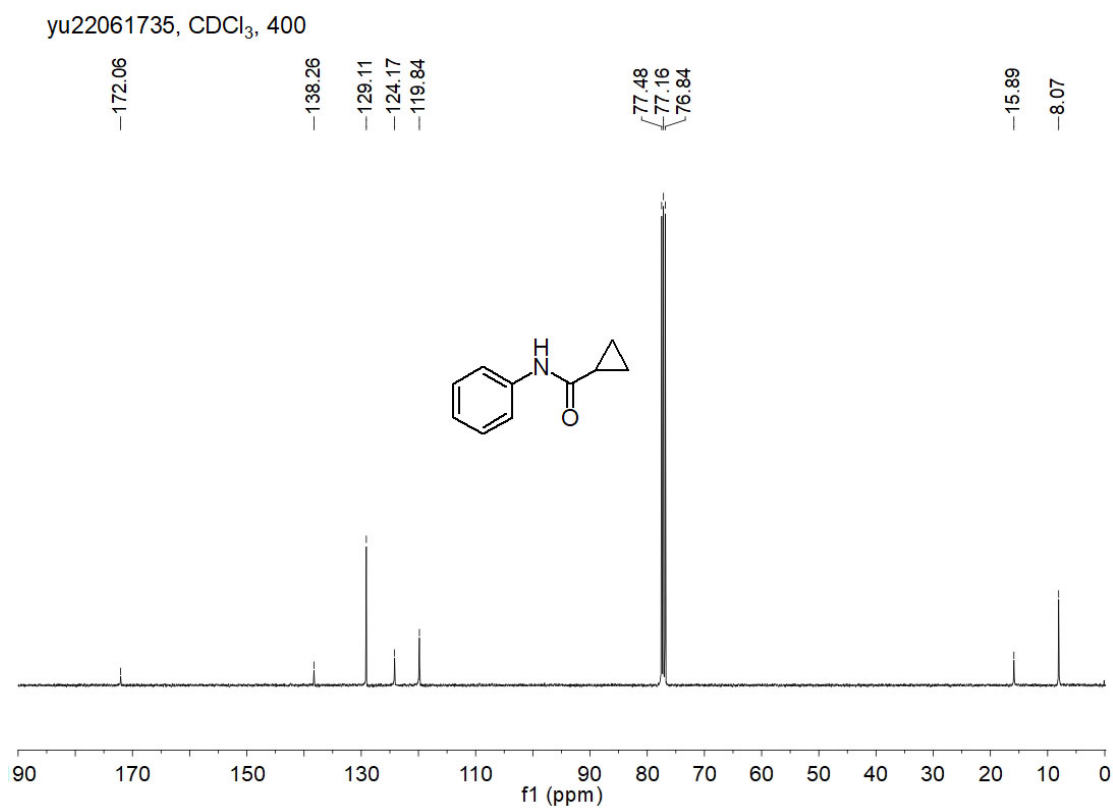
yu20052918, CDCl₃, 400



¹³CNMR of **4u**

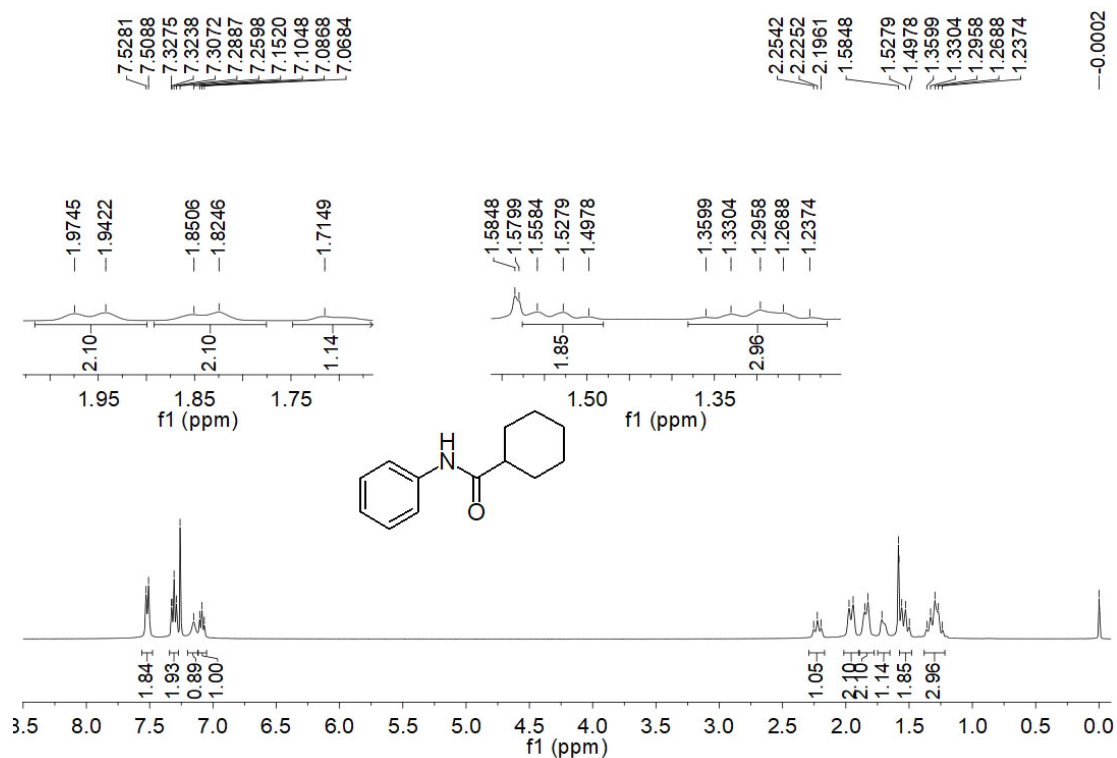


¹HNMR of **4v**

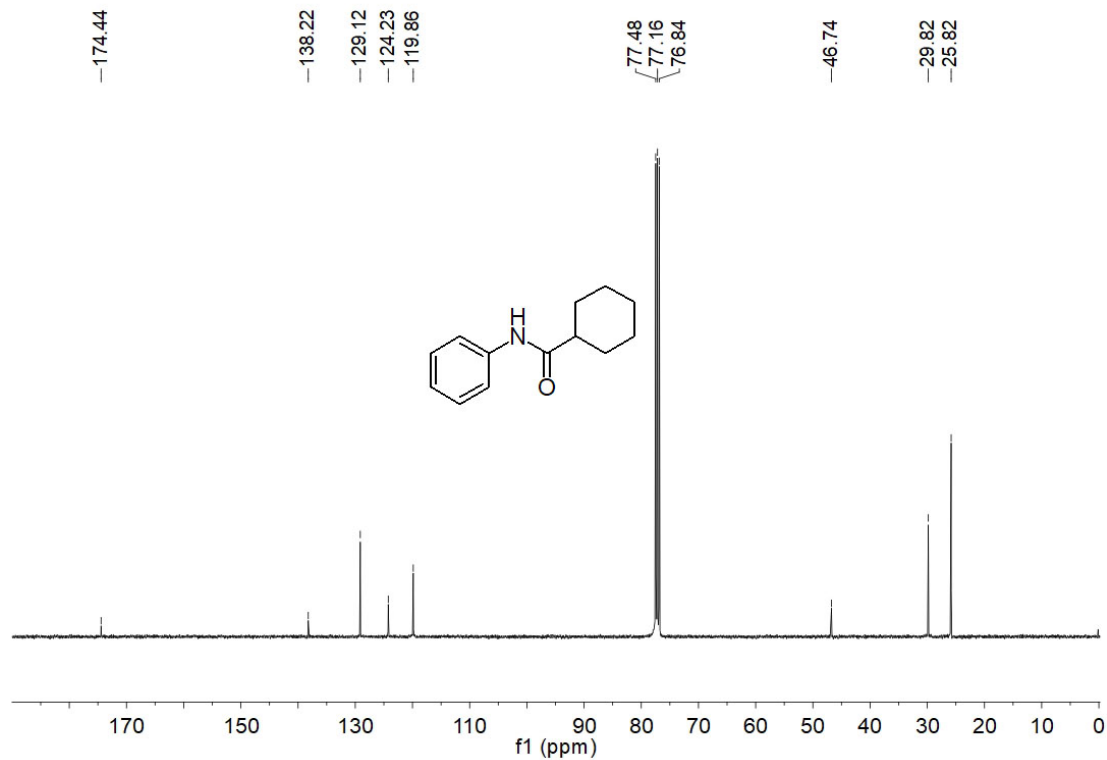


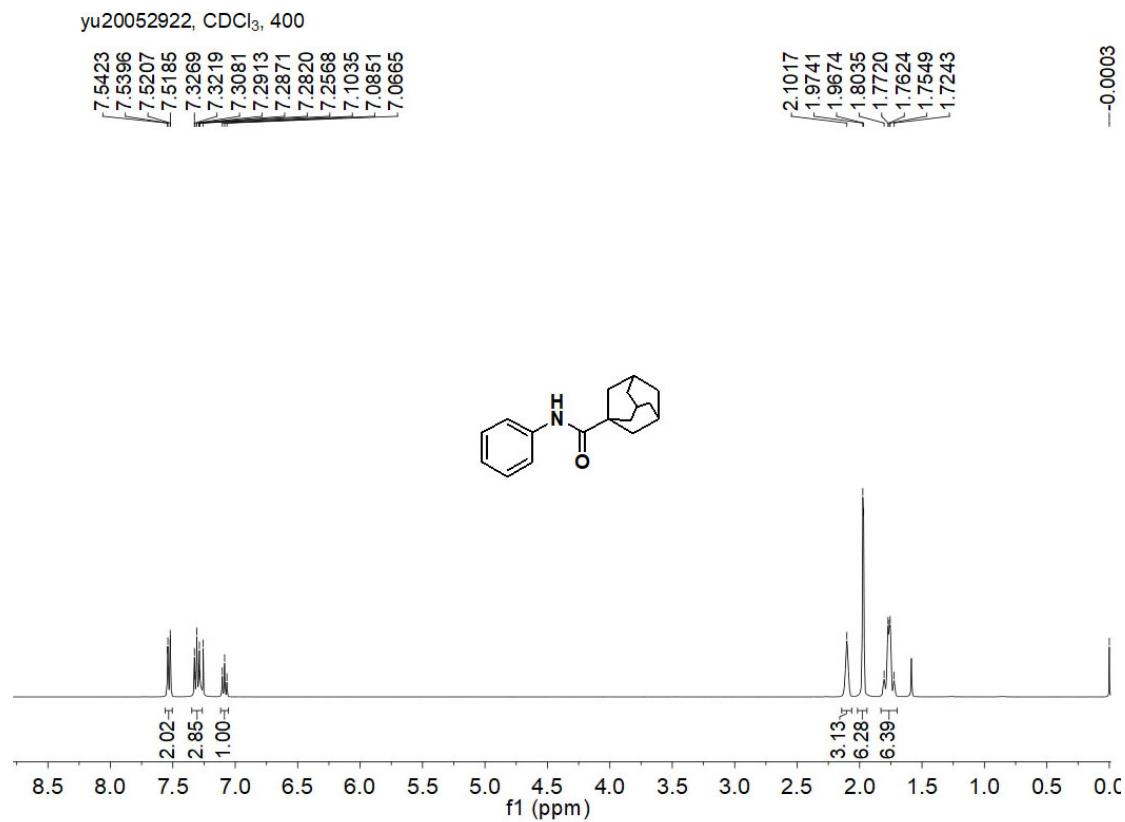
¹³CNMR of **4v**

yu22071518, CDCl₃, 400

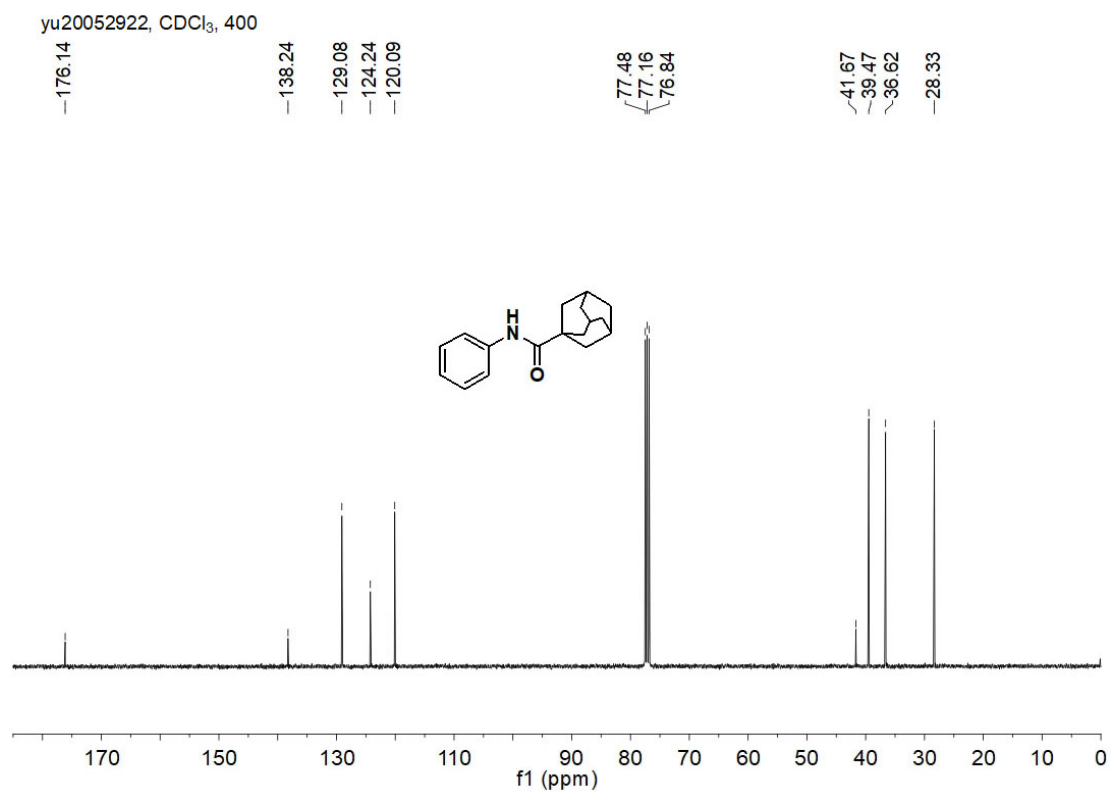


yu22071518, CDCl₃, 400

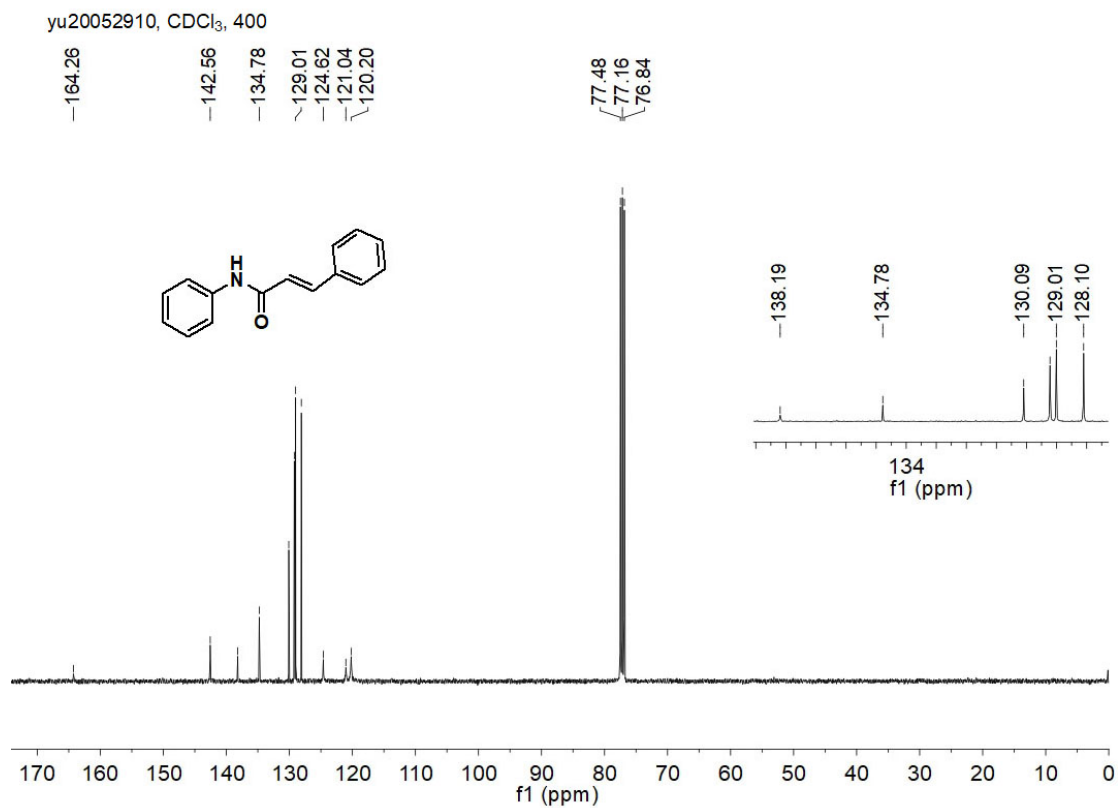
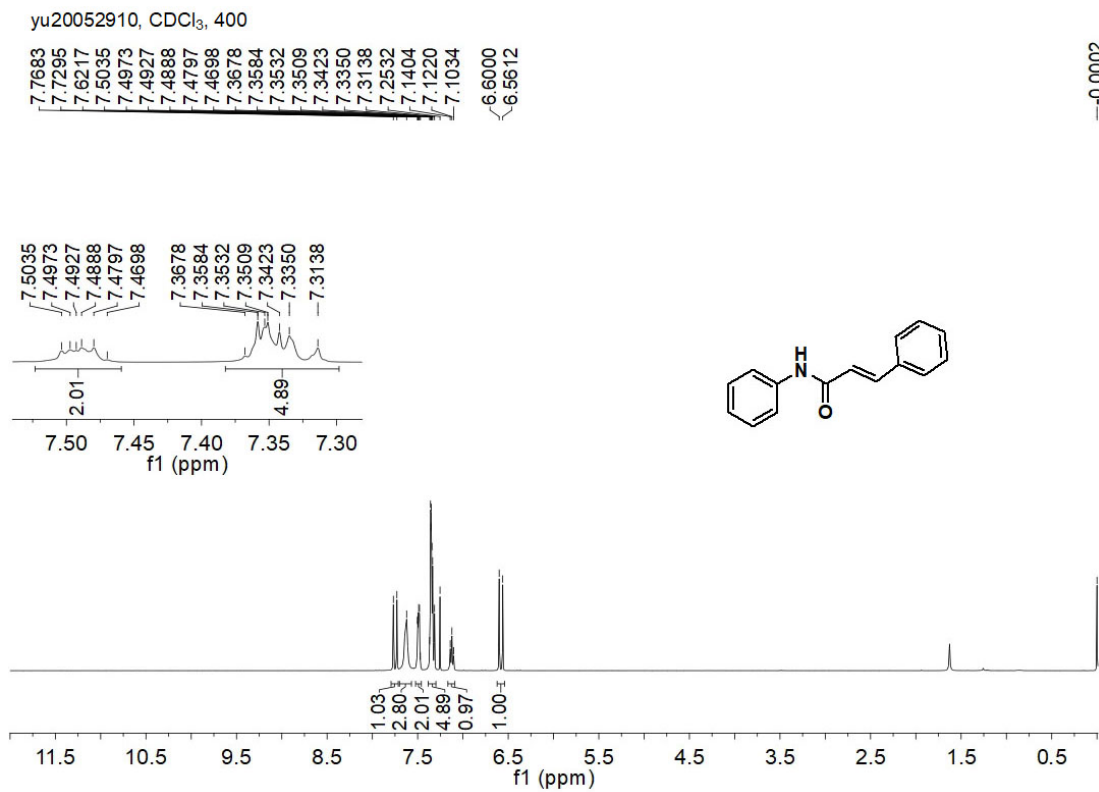




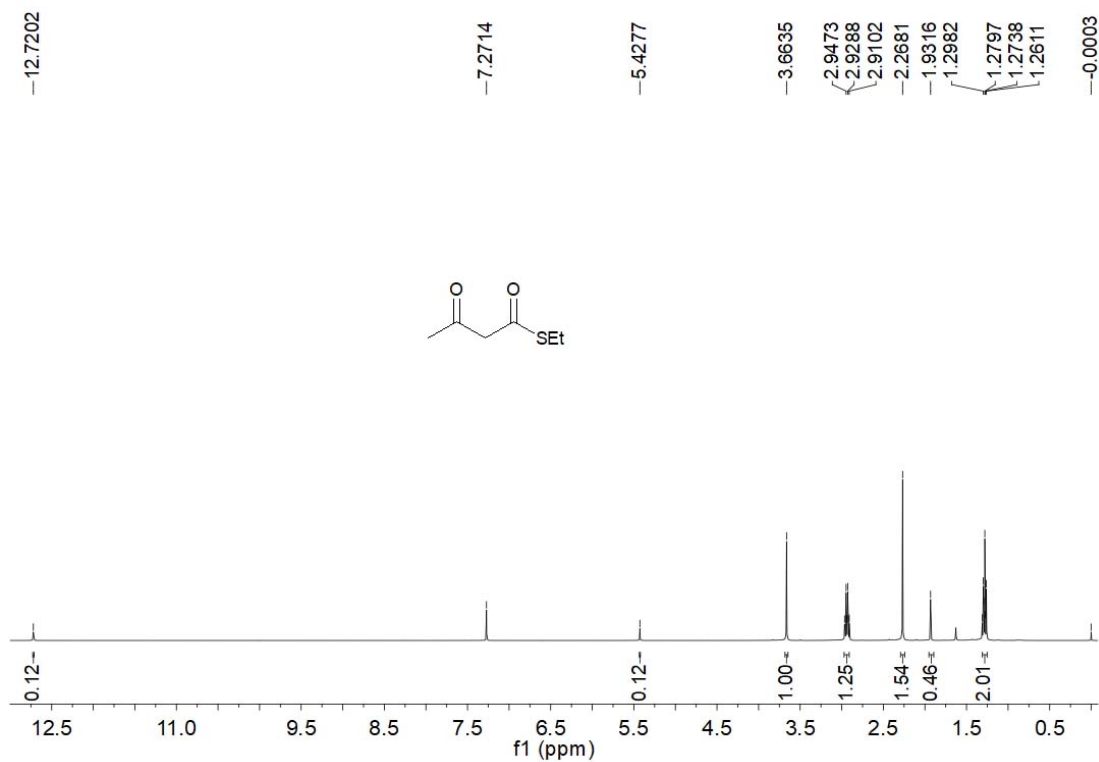
¹HNMR of **4x**



¹³CNMR of **4x**

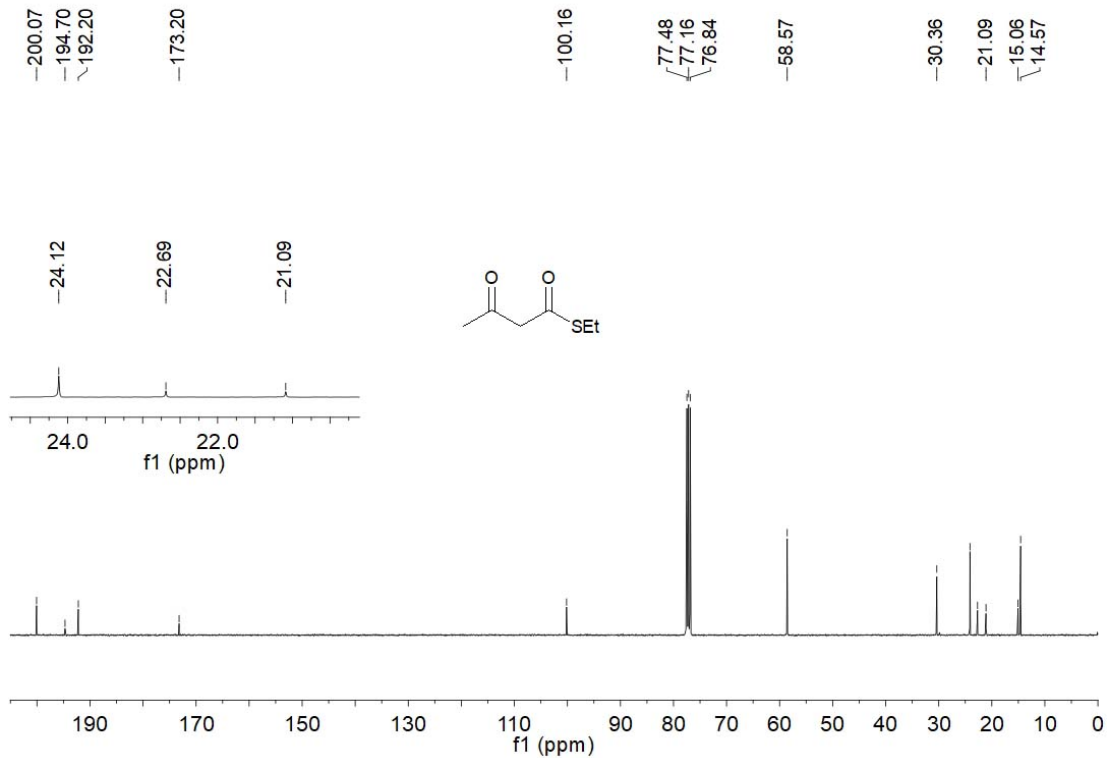


yu22112702, CDCl₃, 400



¹H NMR of A

yu22112702, CDCl₃, 400



¹³C NMR of A