

Supporting Information

Synthesis and pharmacological evaluation of thiophene bioisosteres of potent GluN2B ligands with a benzo[7]annulene-scaffold

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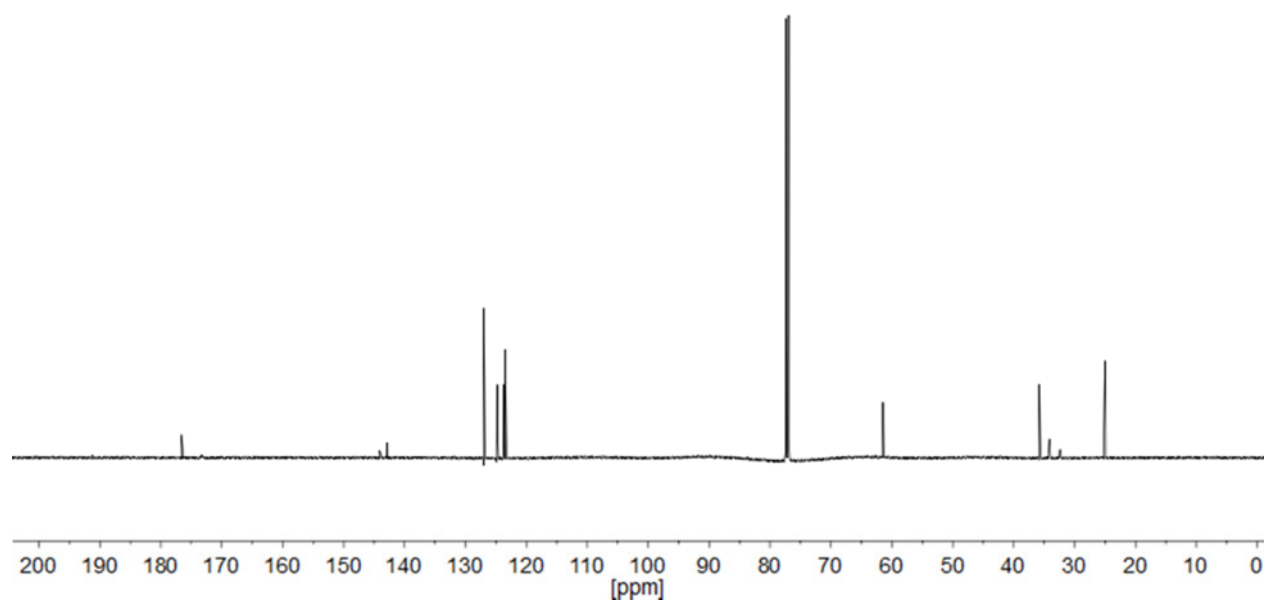
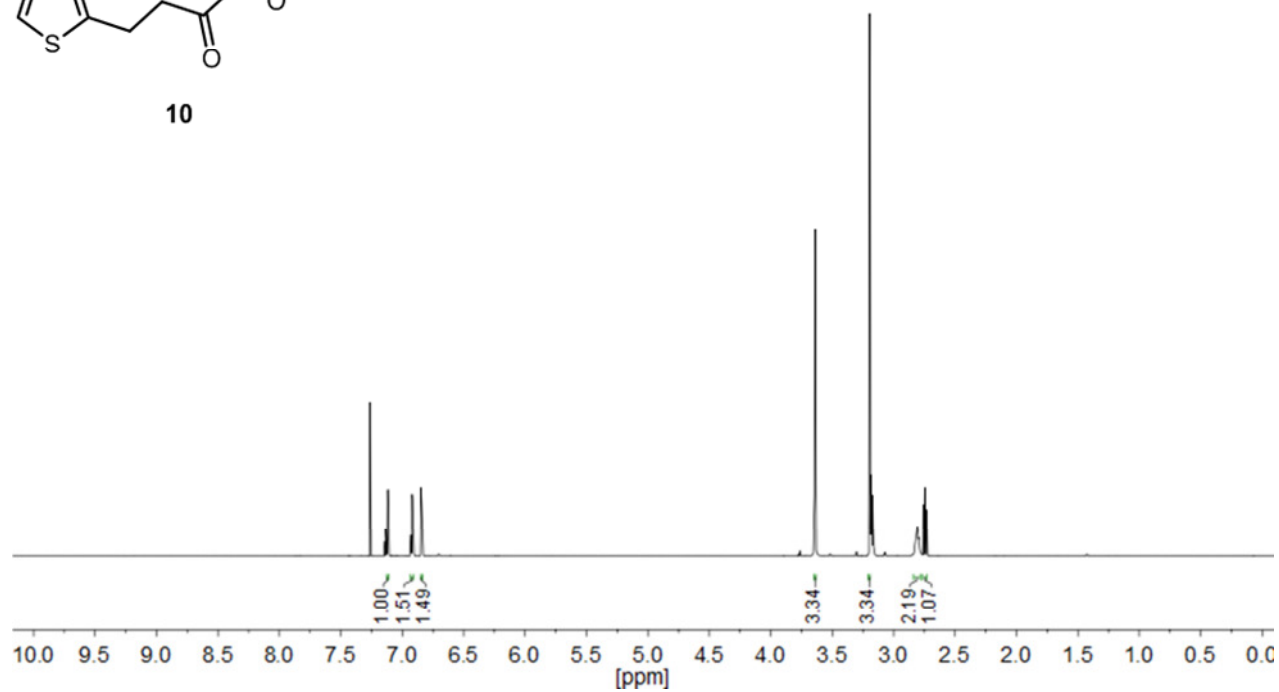
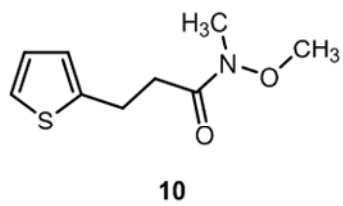
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Corrensstraße 48, D-48149 Münster, Germany.

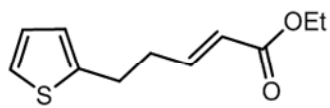
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Universität Münster, Germany.

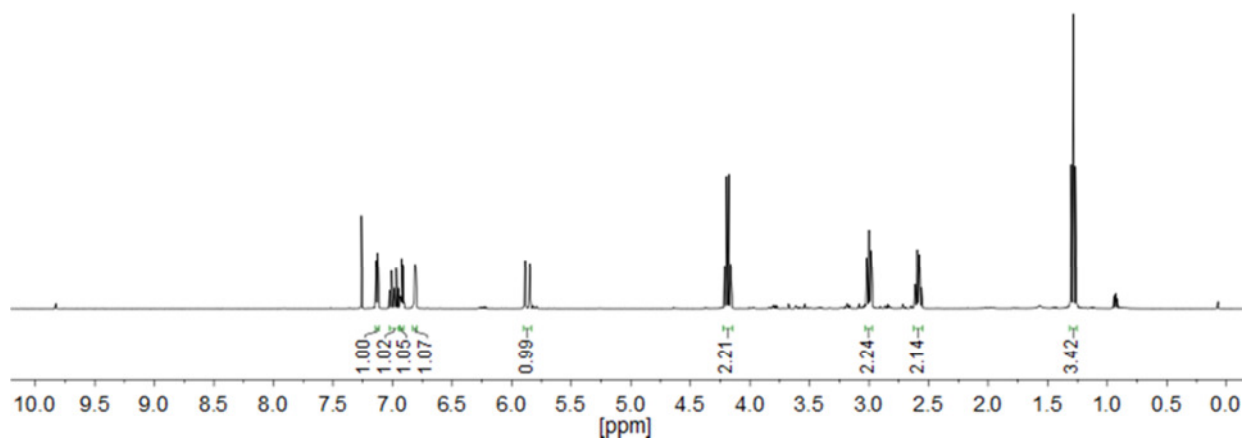
Contents

NMR spectra of prepared compounds

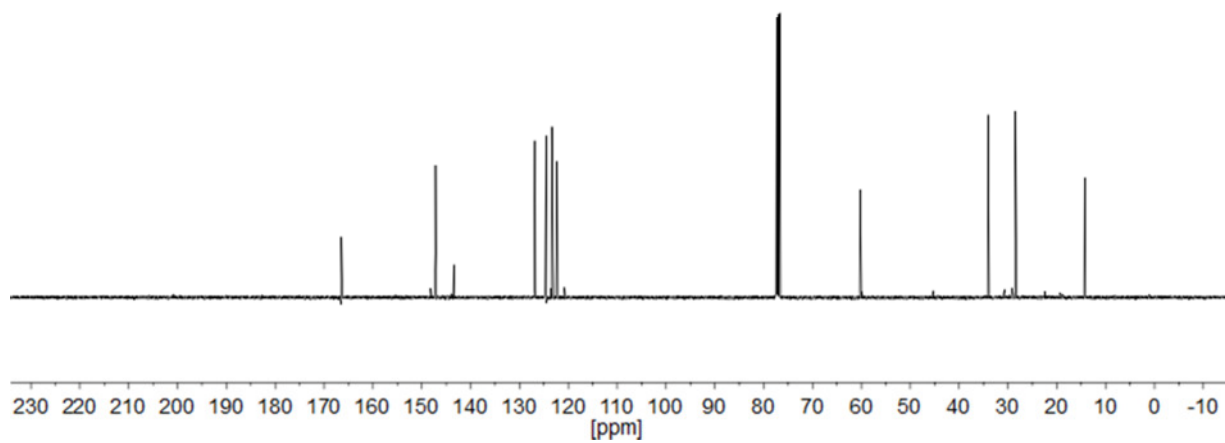




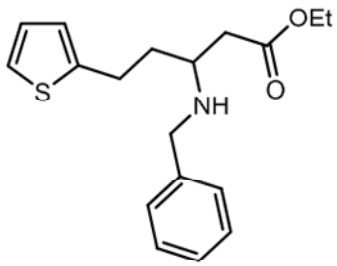
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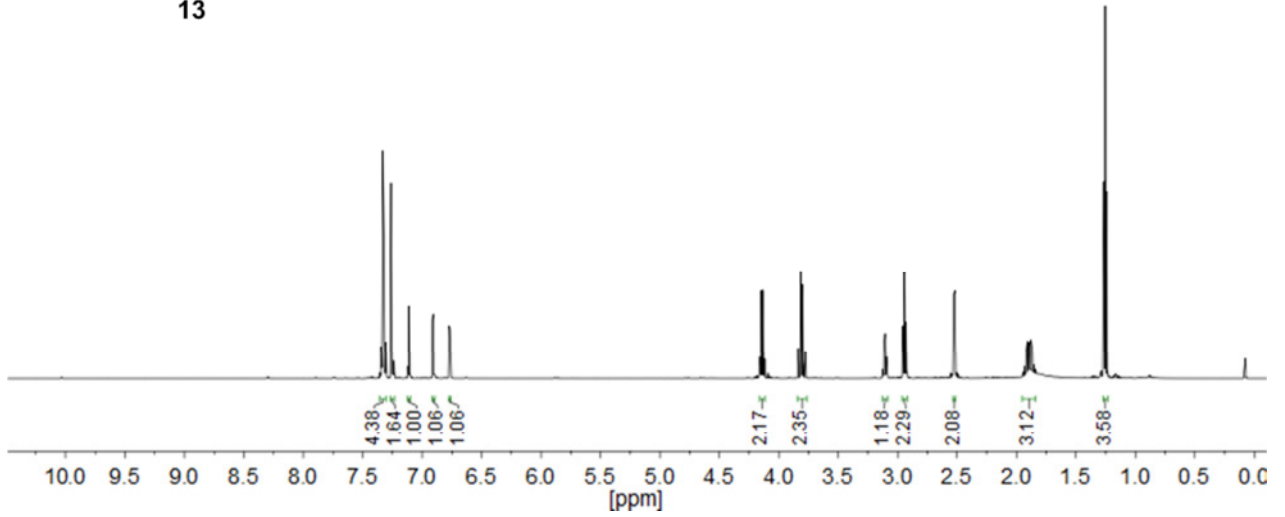
¹H NMR-Spectrum (CDCl₃) of **12**.



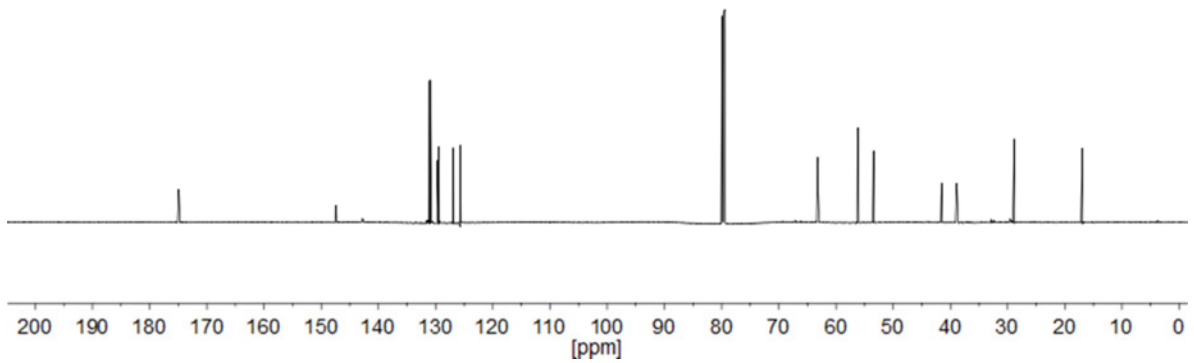
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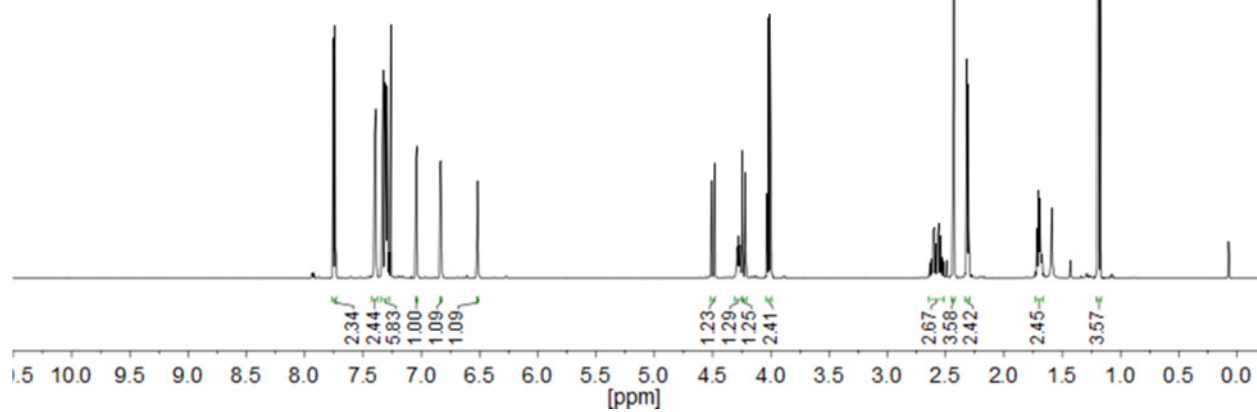
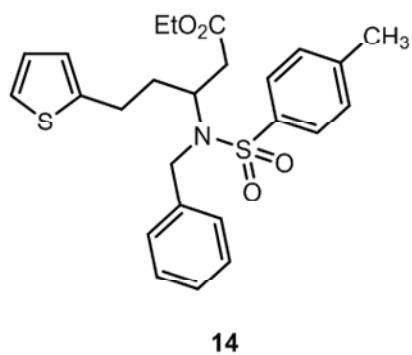
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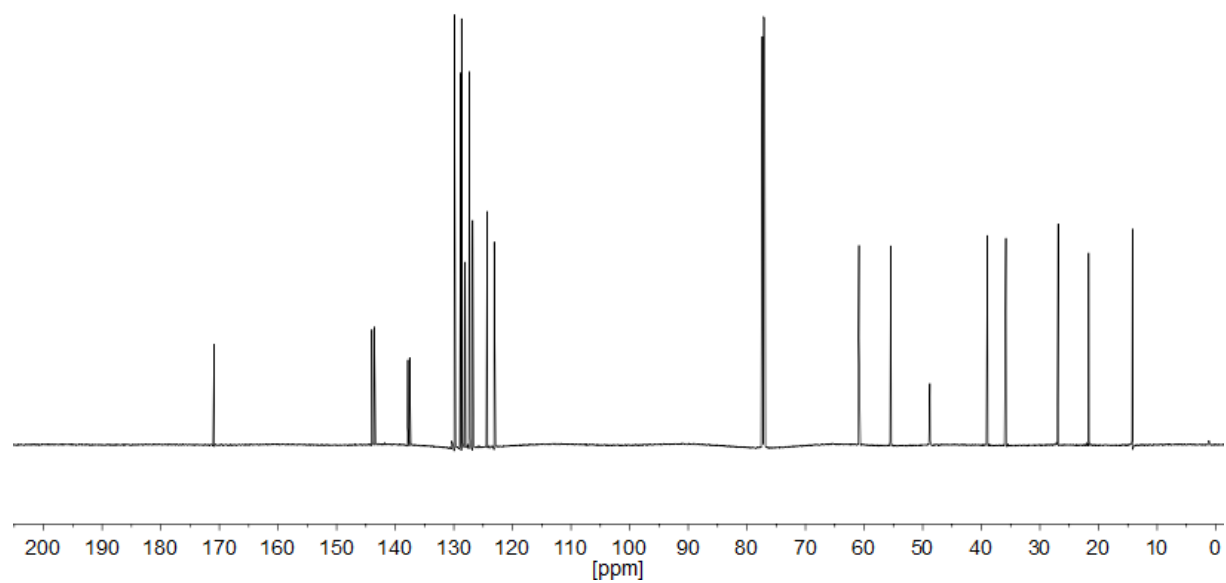
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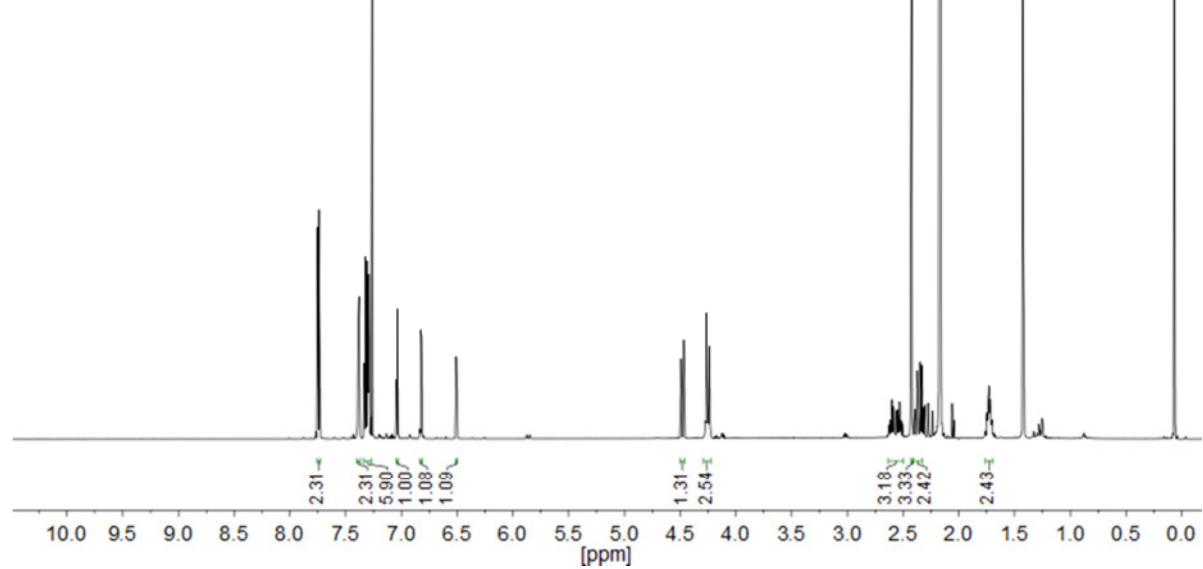
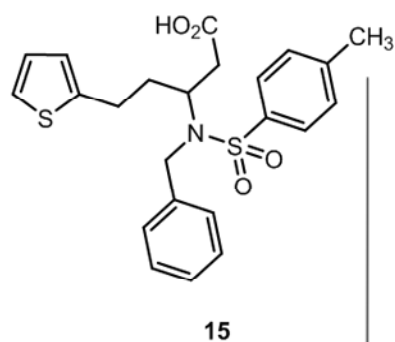
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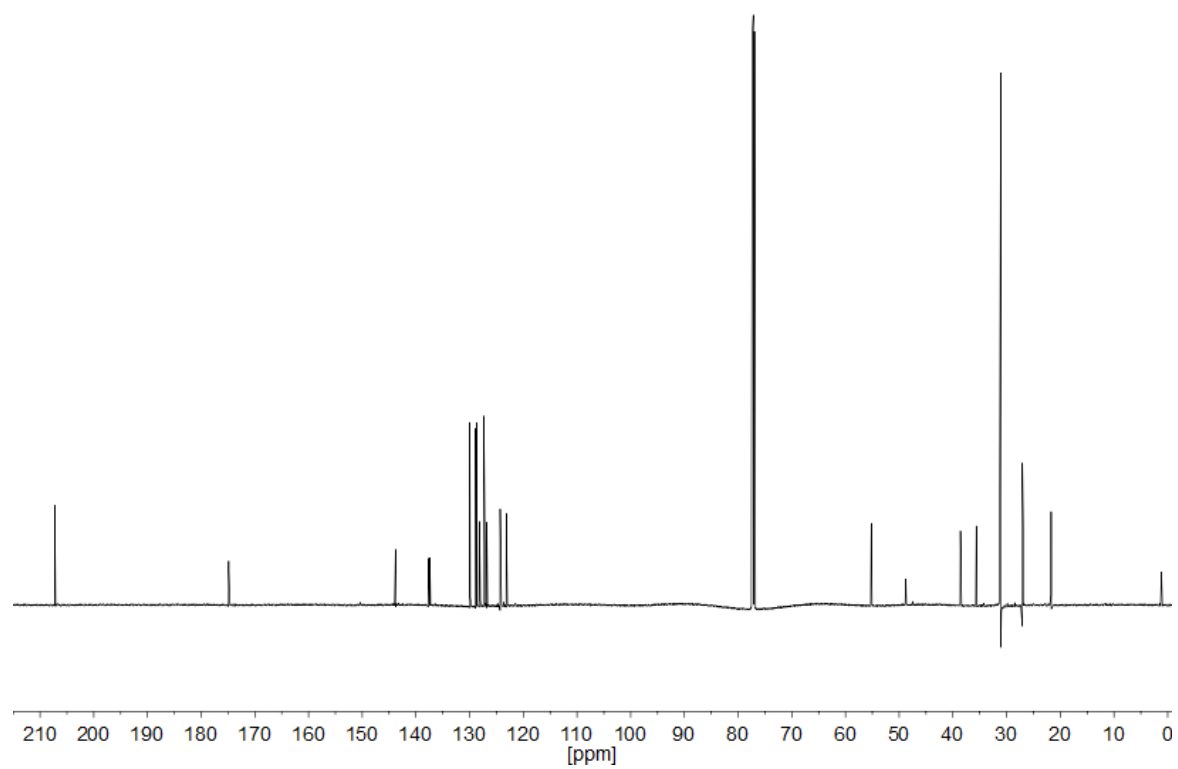
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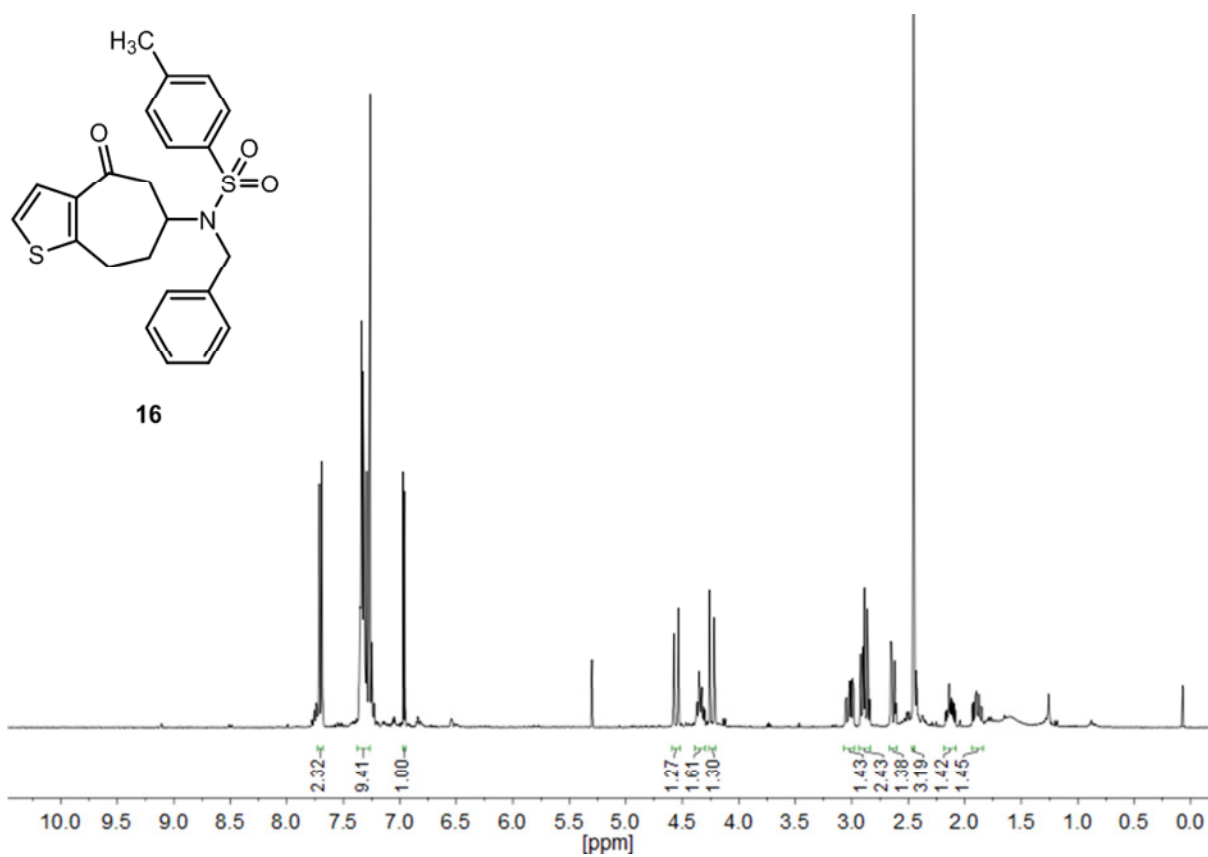
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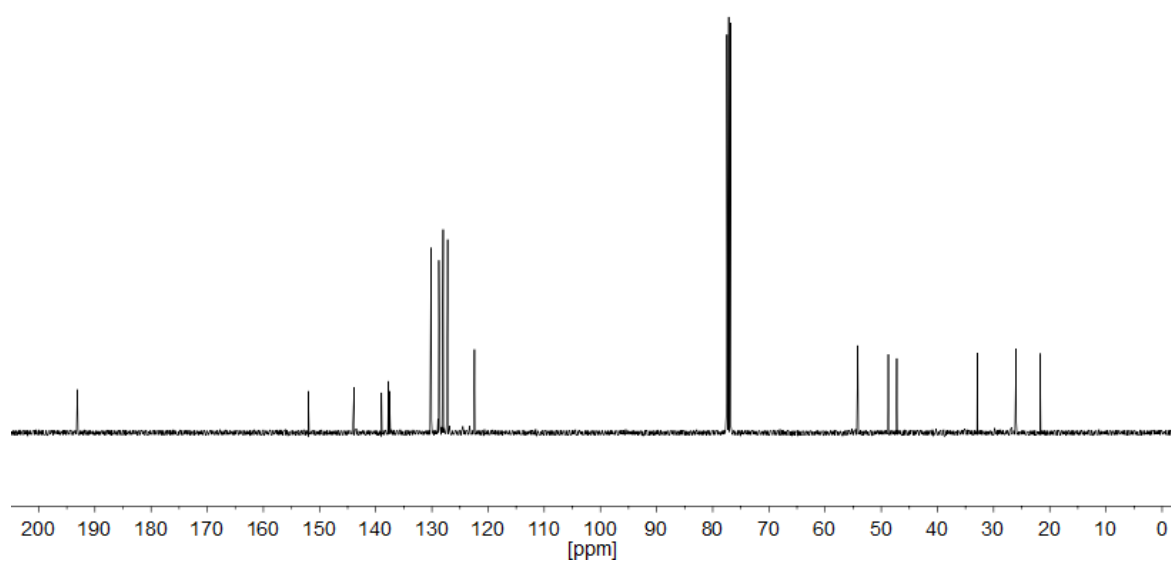
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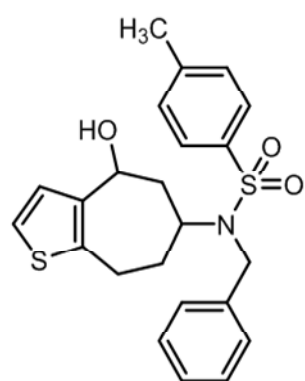
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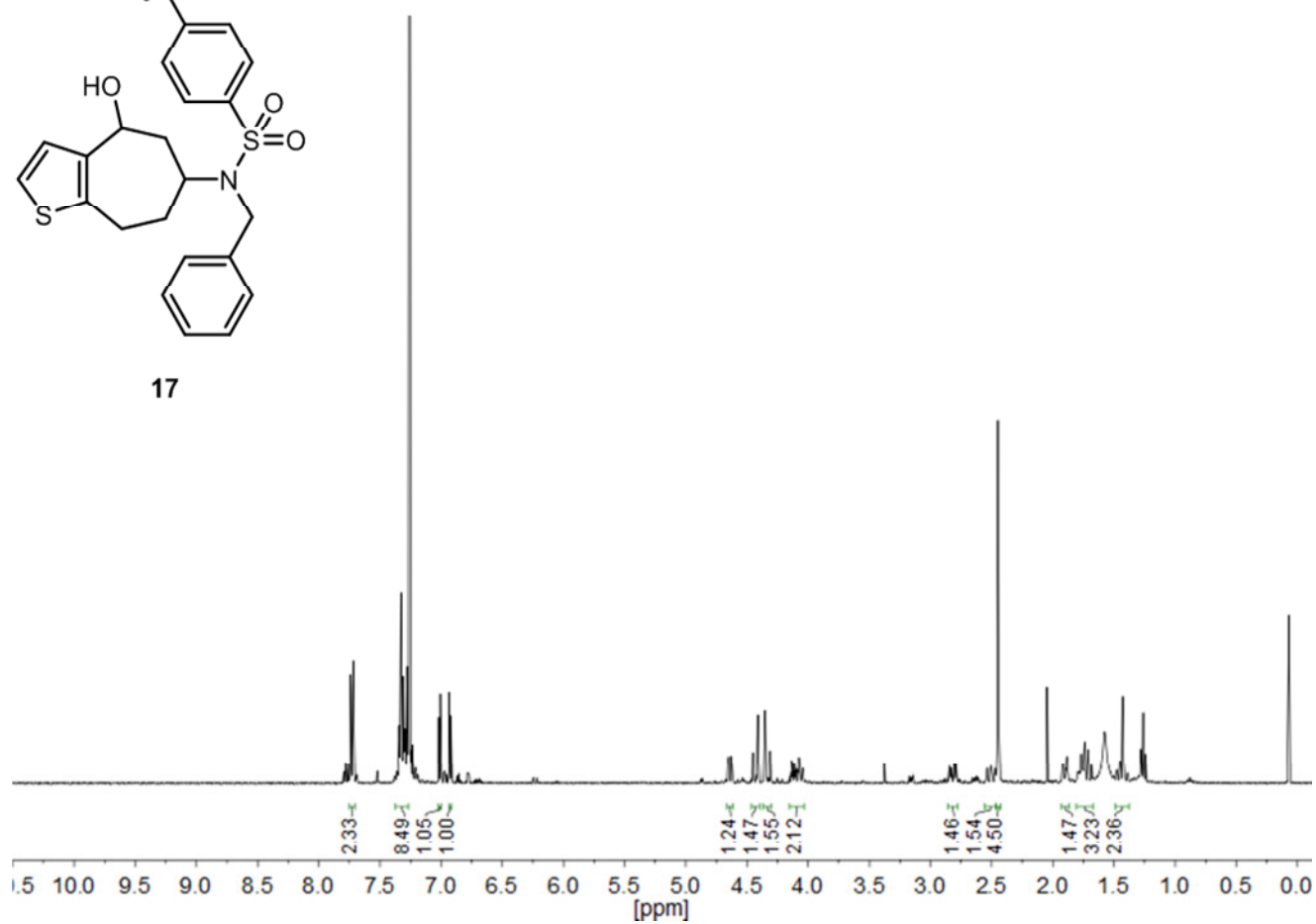
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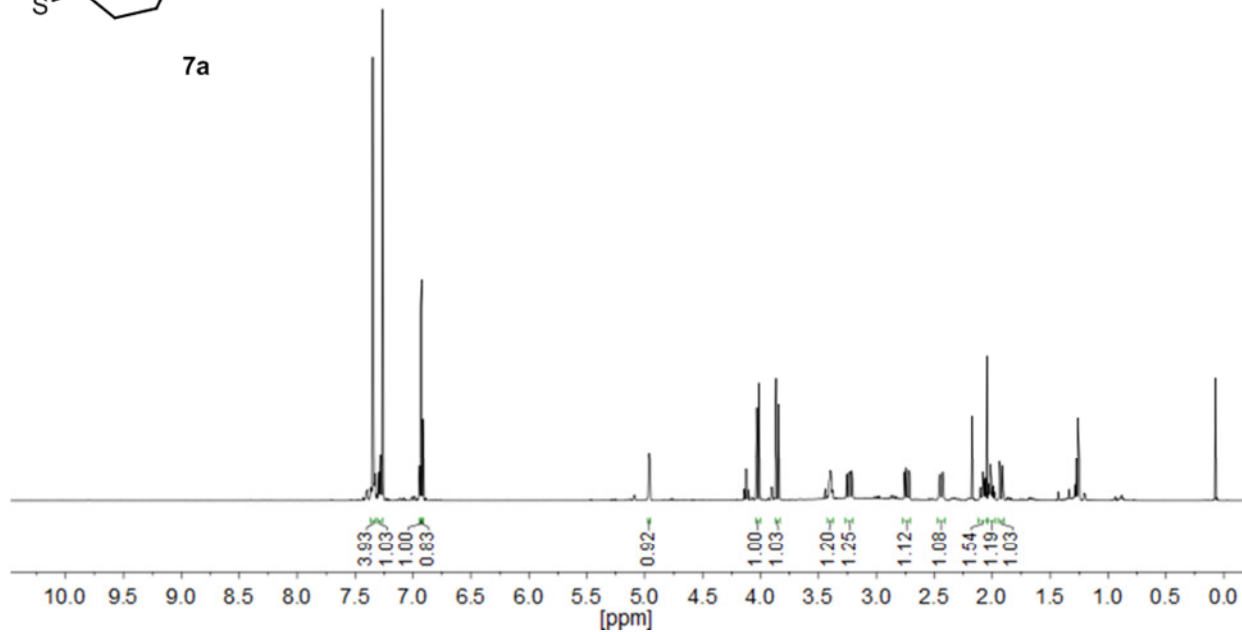
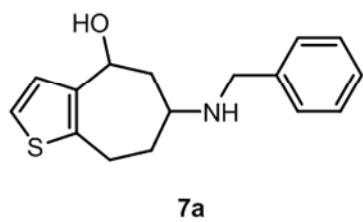
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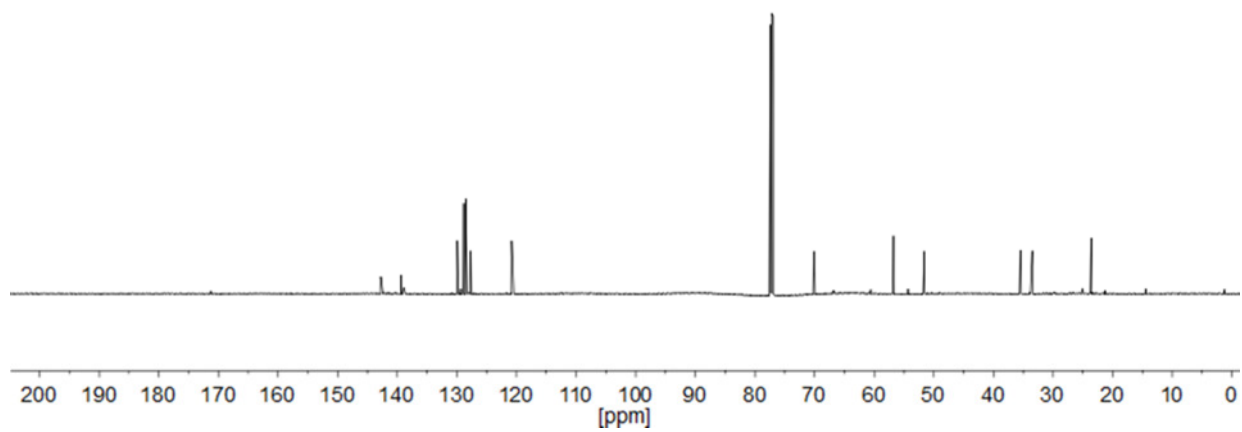
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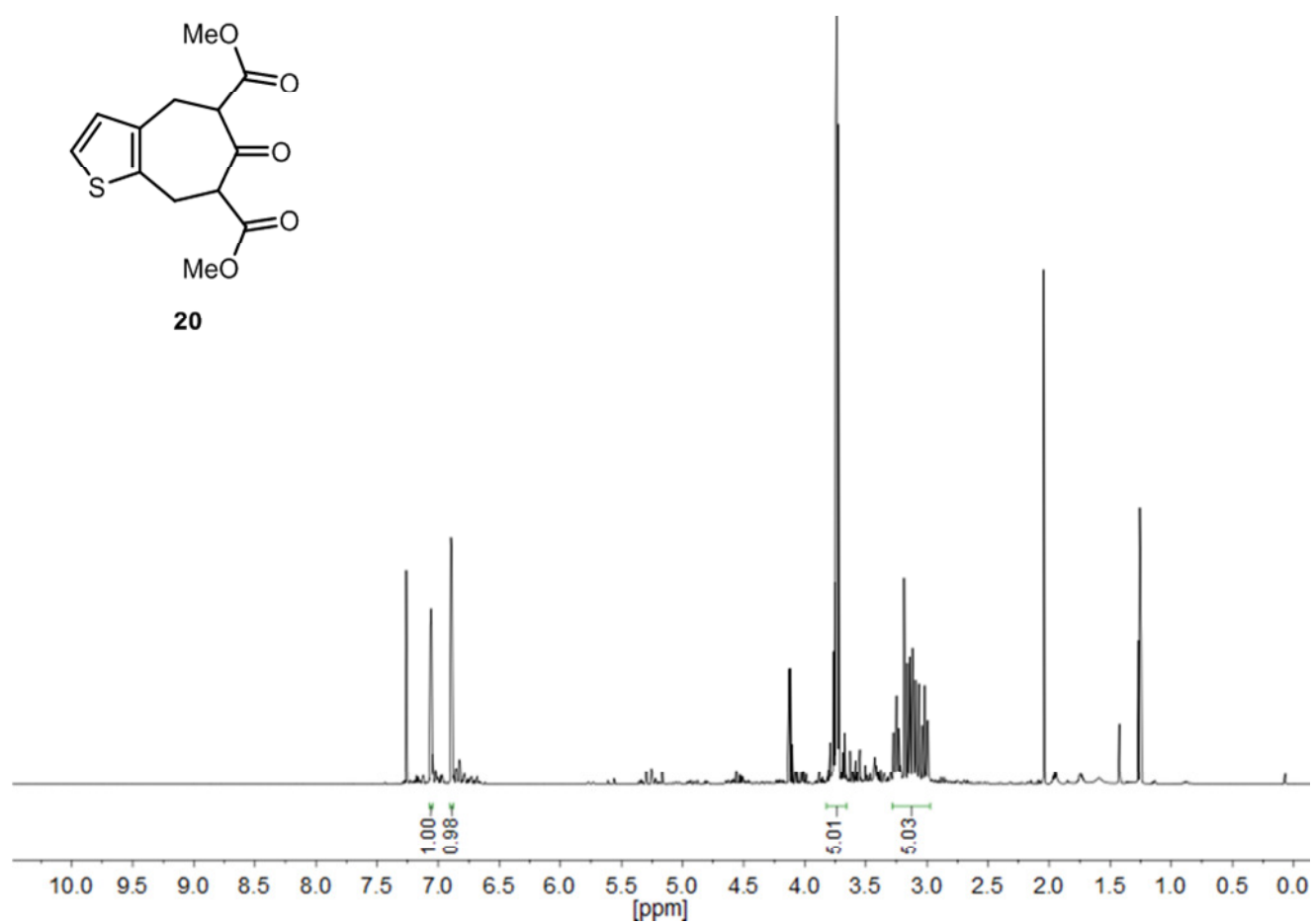
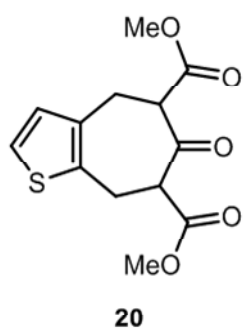
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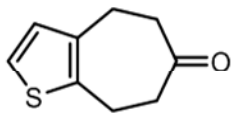
^1H NMR-Spectrum (CDCl_3) of **7a**.



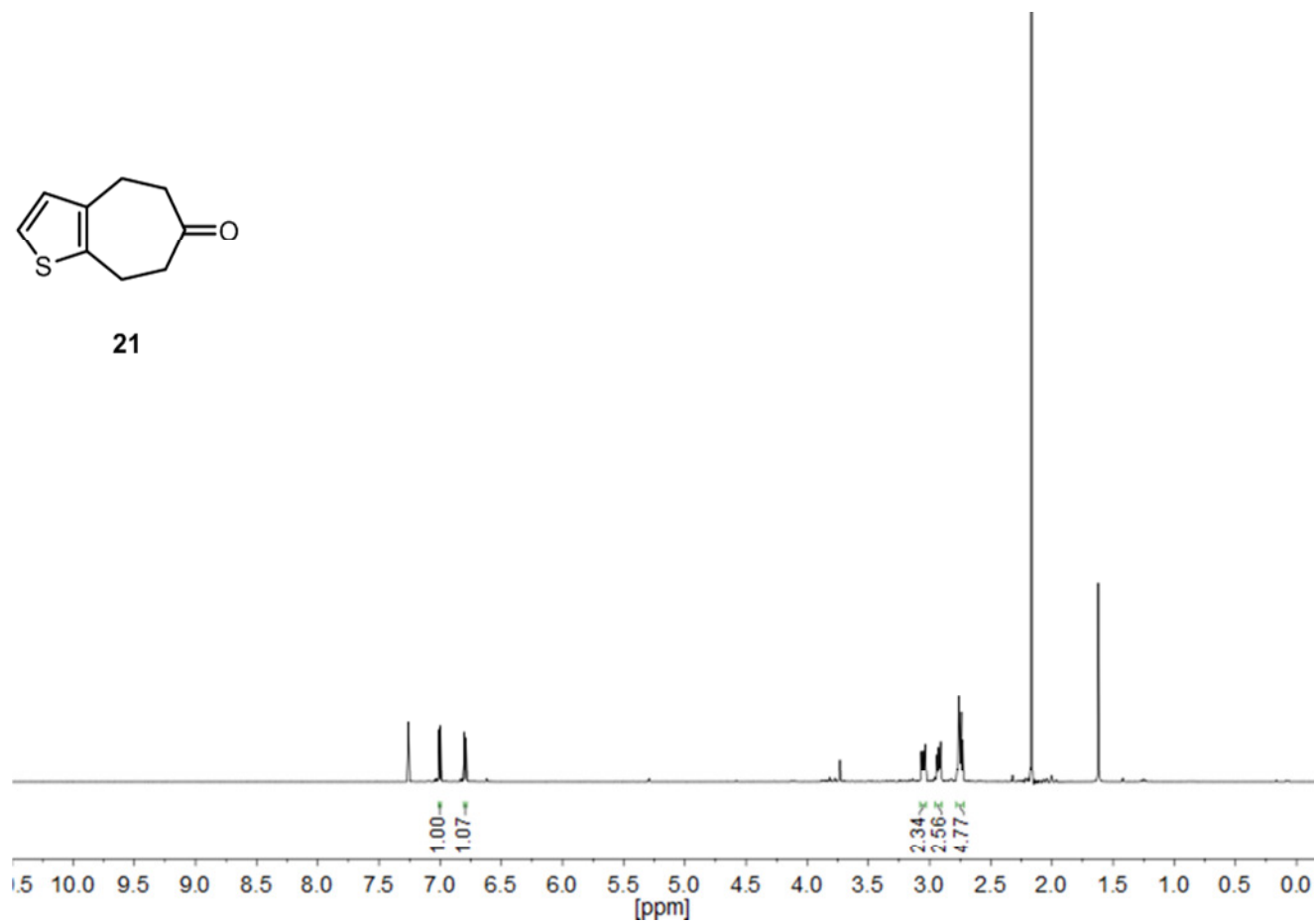
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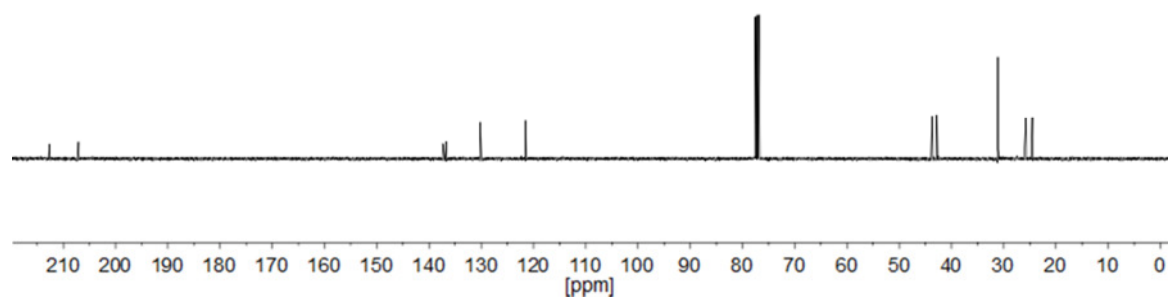
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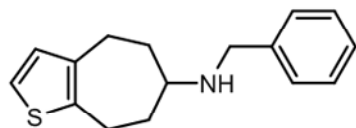
21



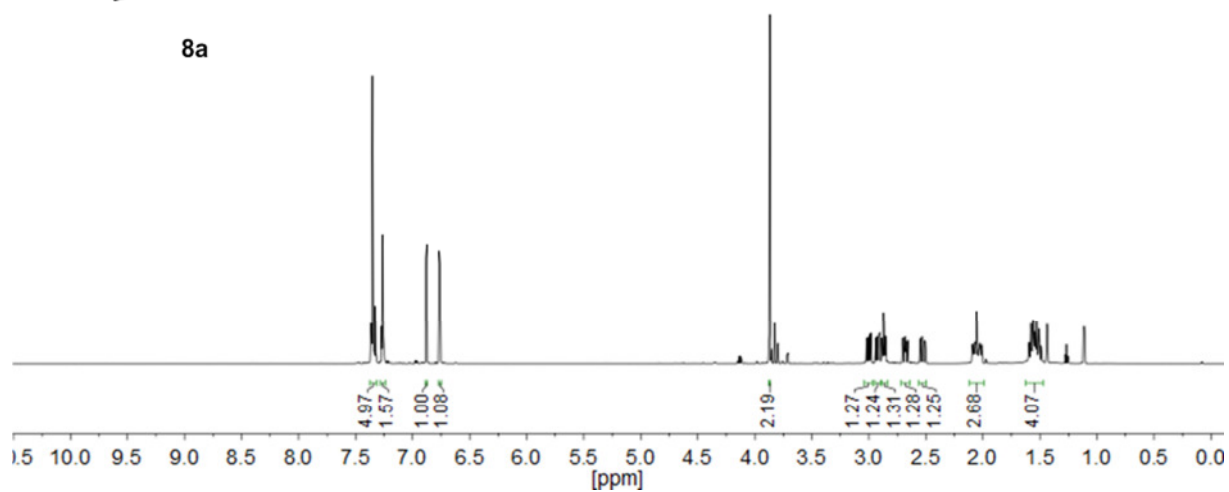
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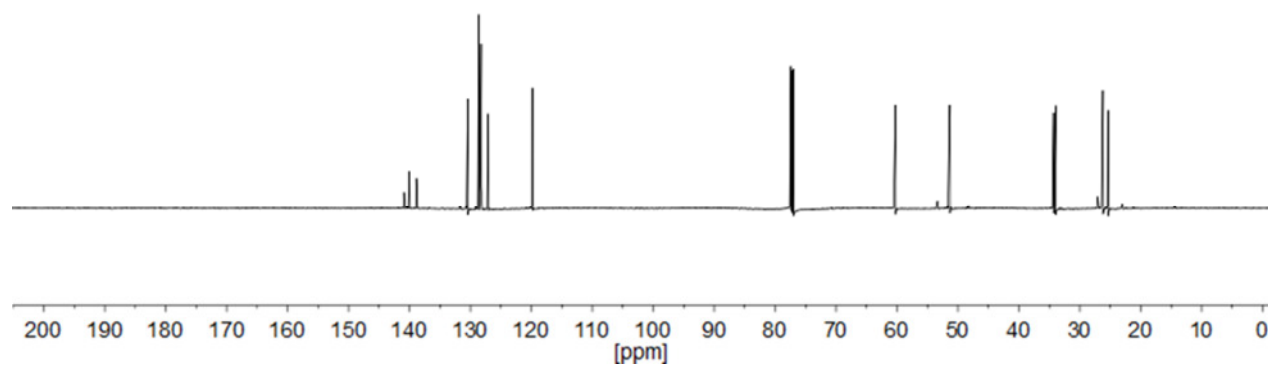
^{13}C NMR-Spectrum (CDCl_3) of **21**.



8a



^1H NMR-Spectrum (CDCl_3) of **8a**.



^{13}C NMR-Spectrum (CDCl_3) of **8a**.