

Supporting Information

Unified Synthesis and Biological Evaluation of Makaluvamine J and its Analogs

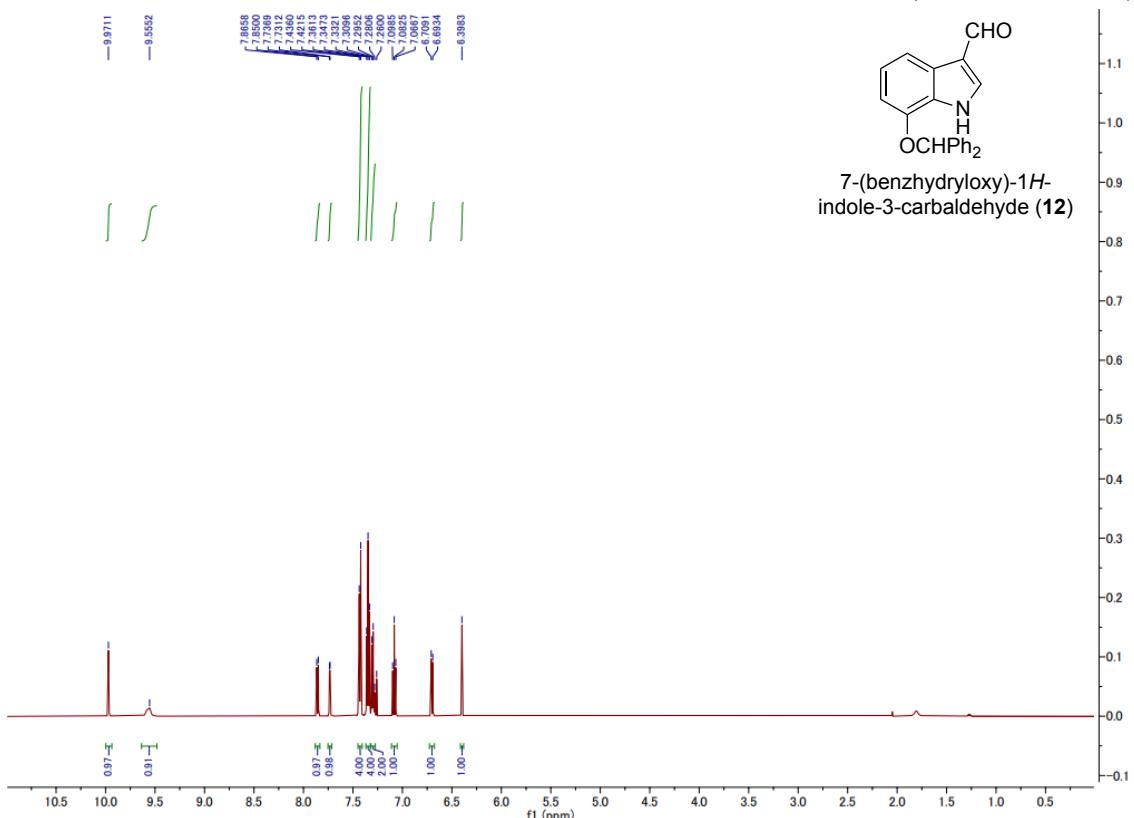
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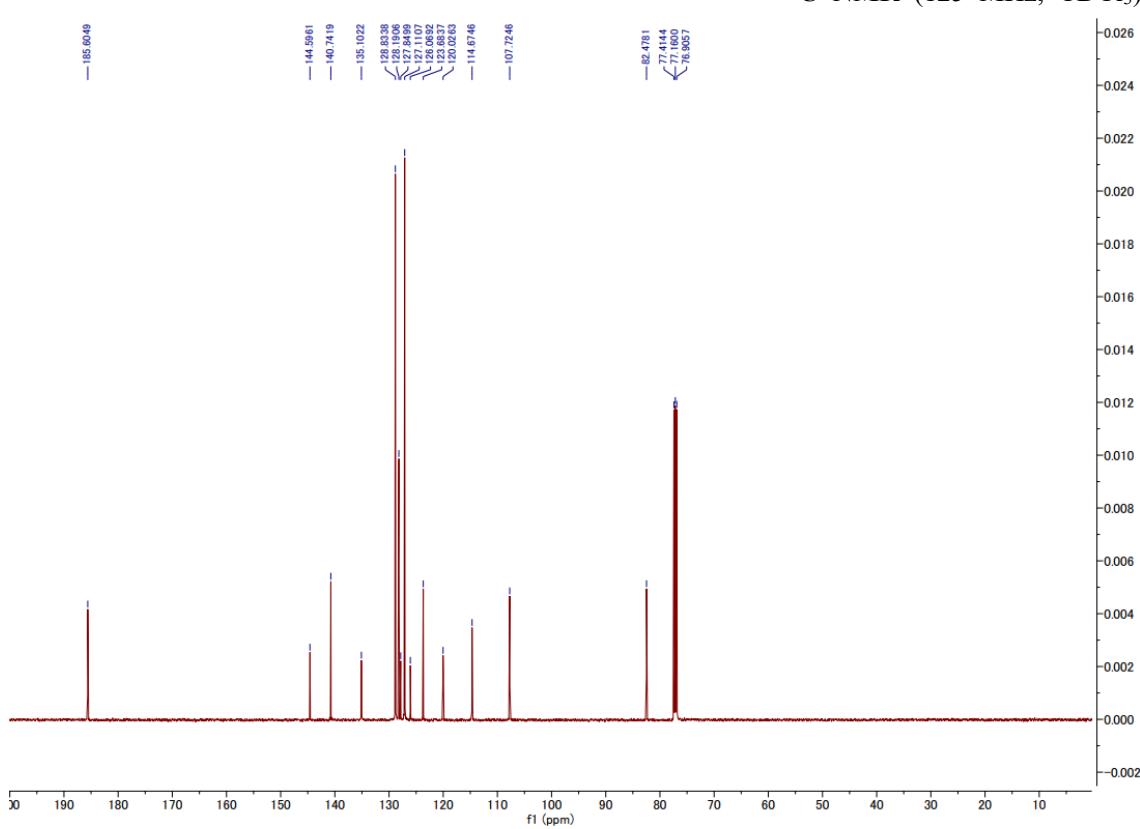
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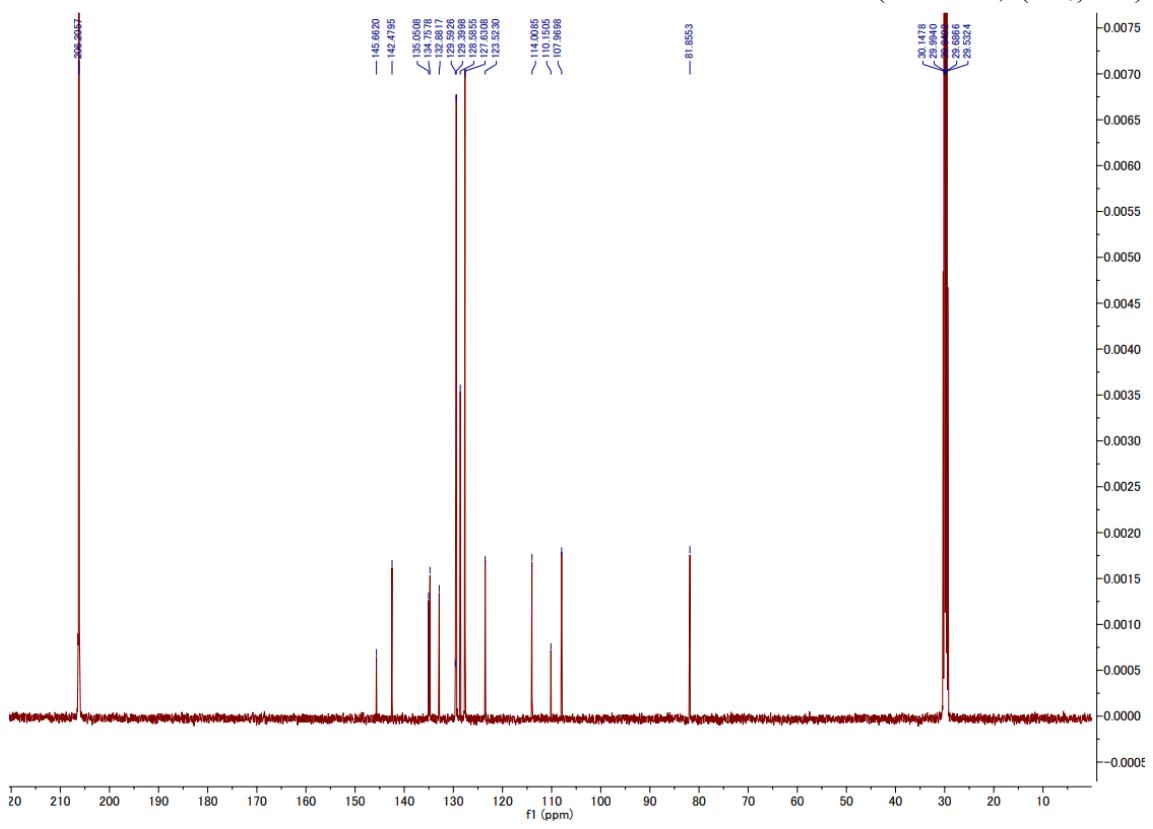
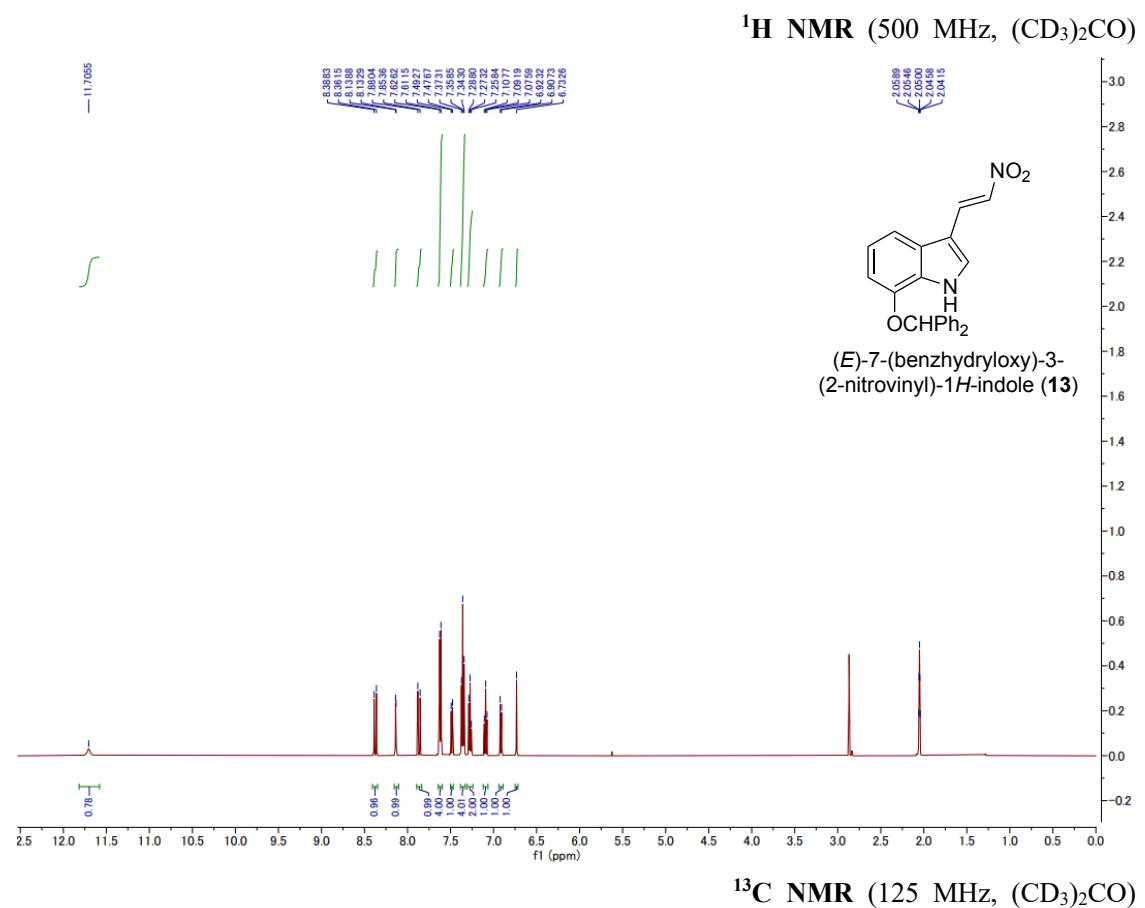
¹H- and ¹³C-NMR spectra for new compounds: pages S2-S11

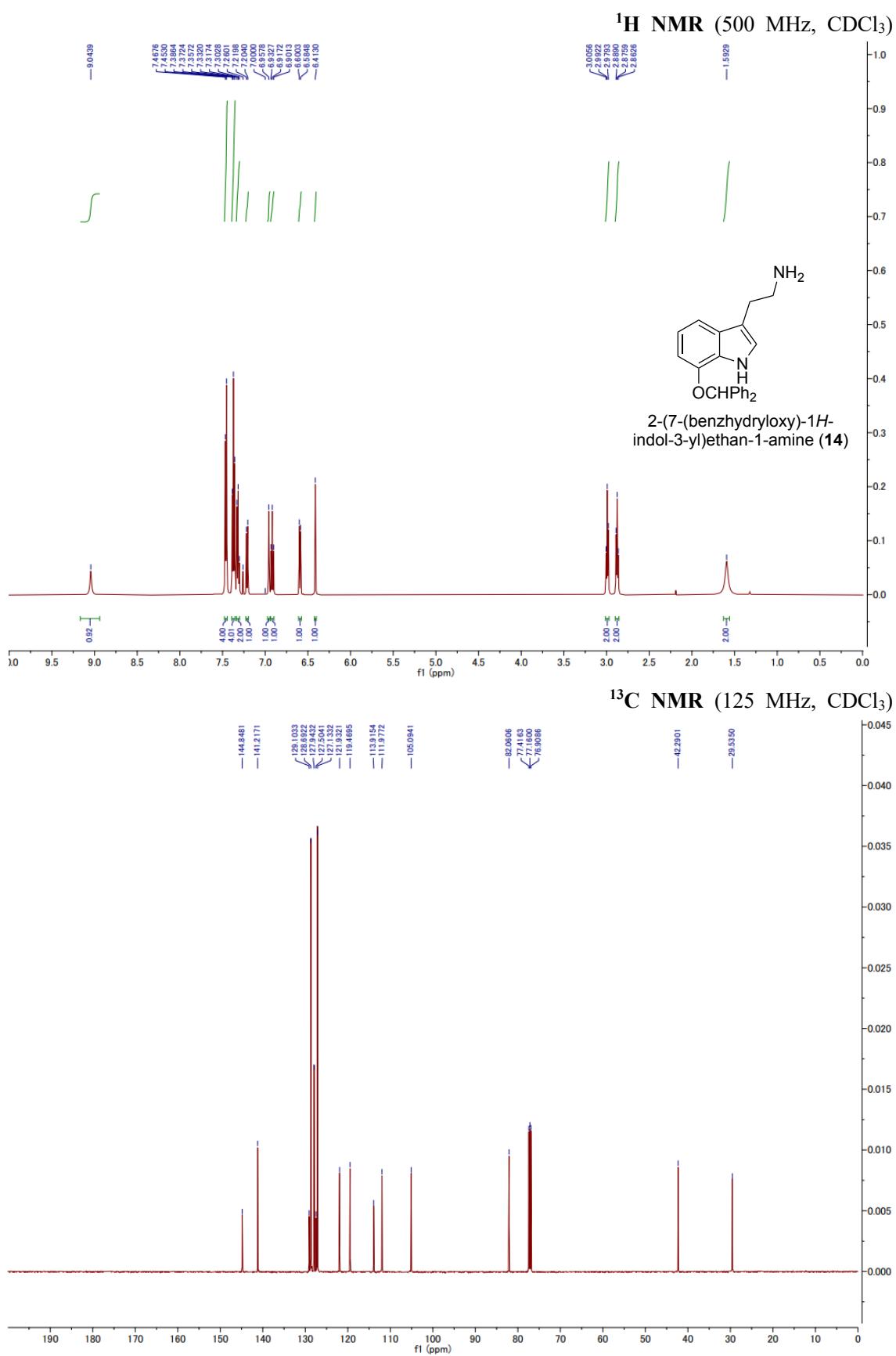
¹H NMR (500 MHz, CDCl₃)

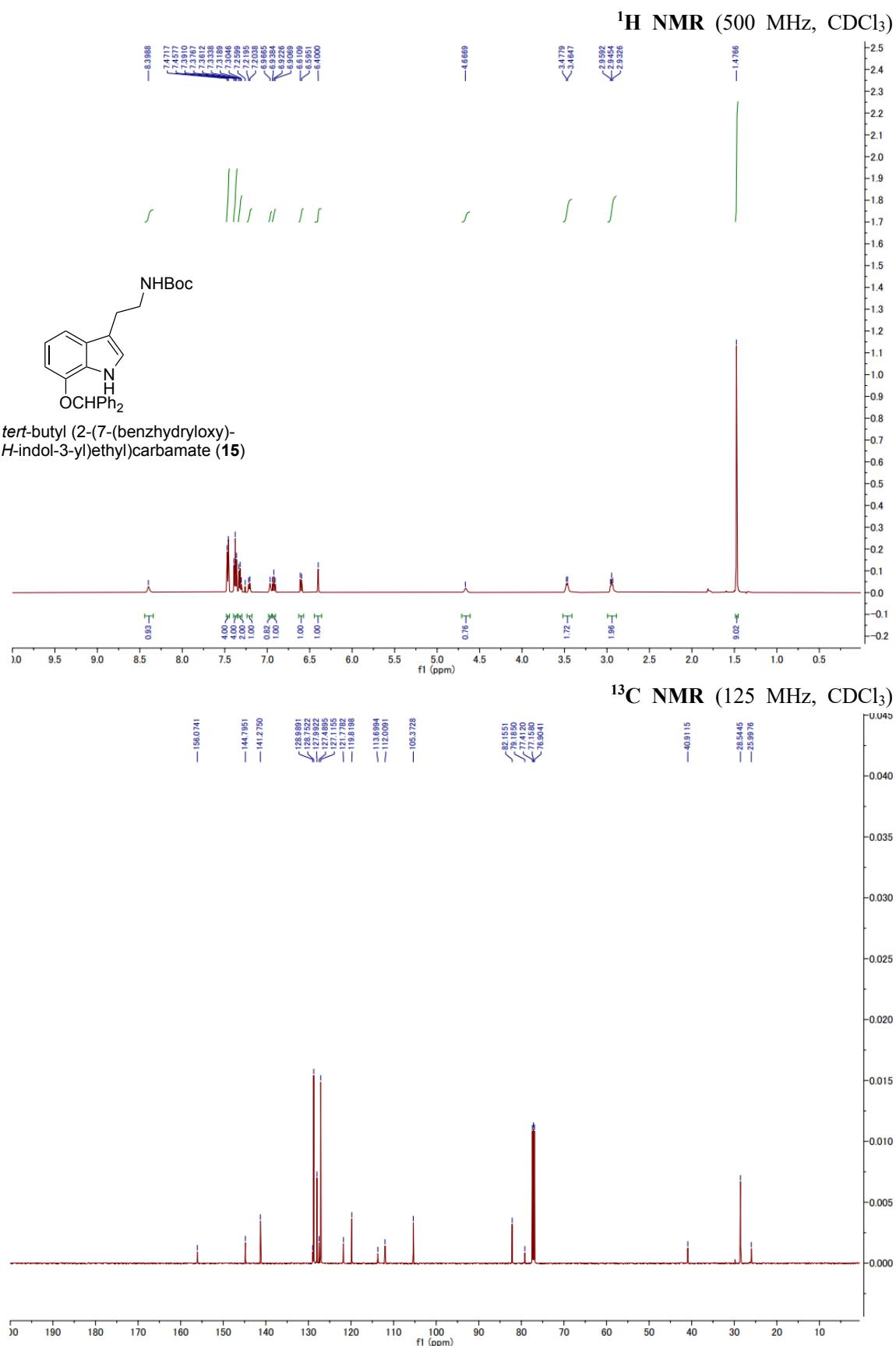


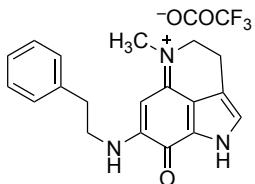
¹³C NMR (125 MHz, CDCl₃)



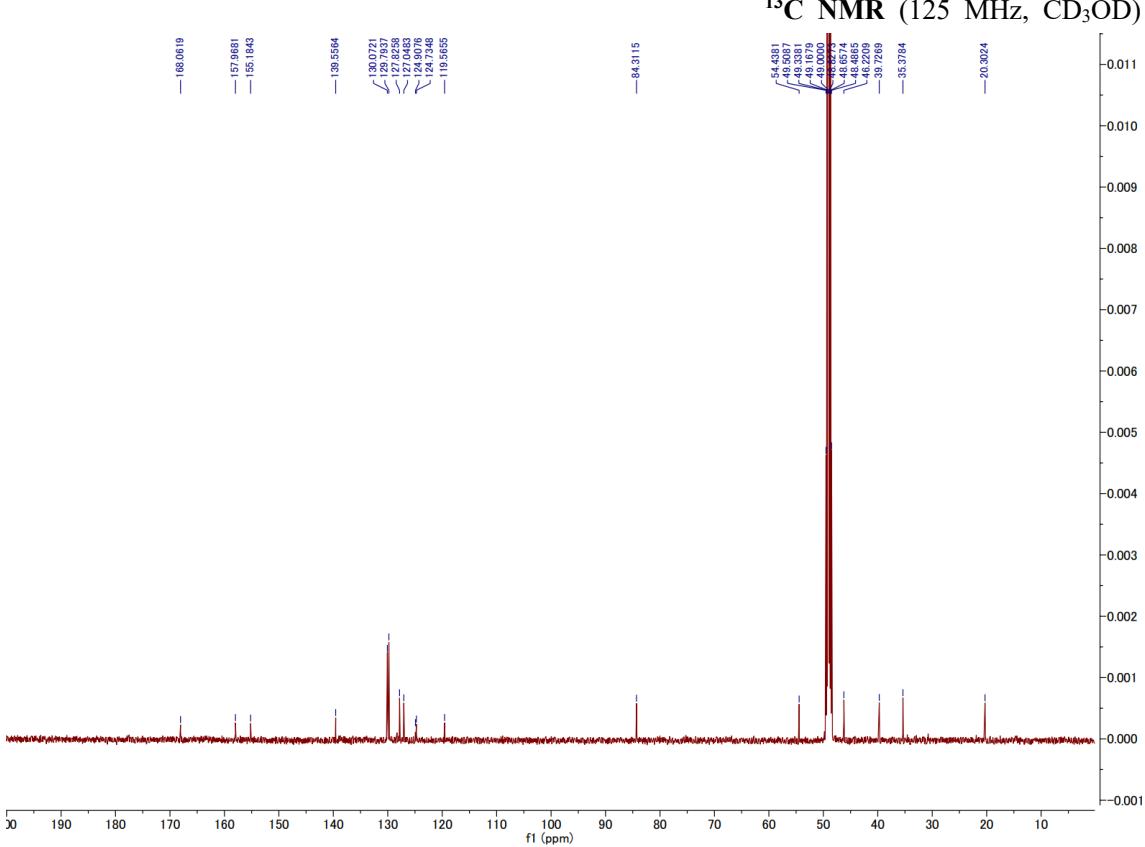
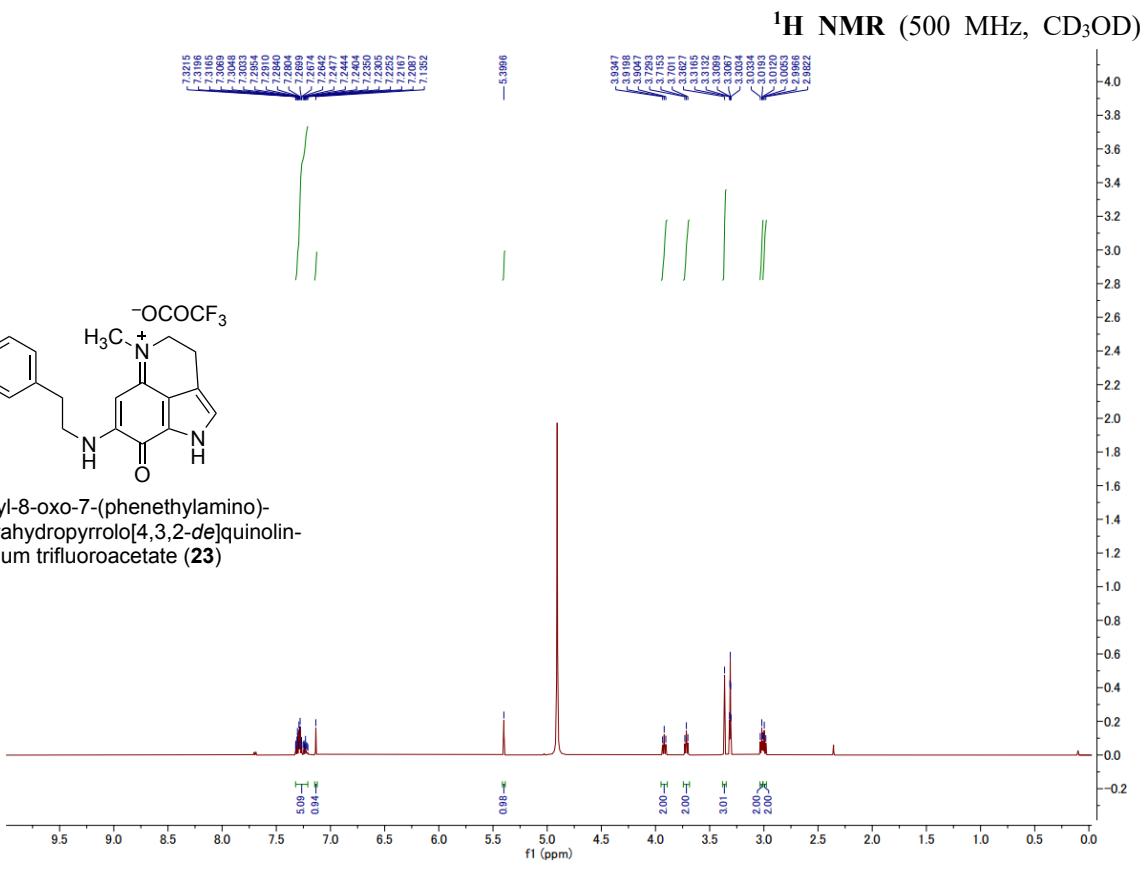


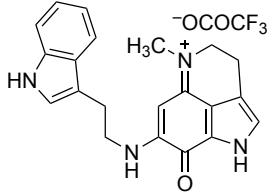




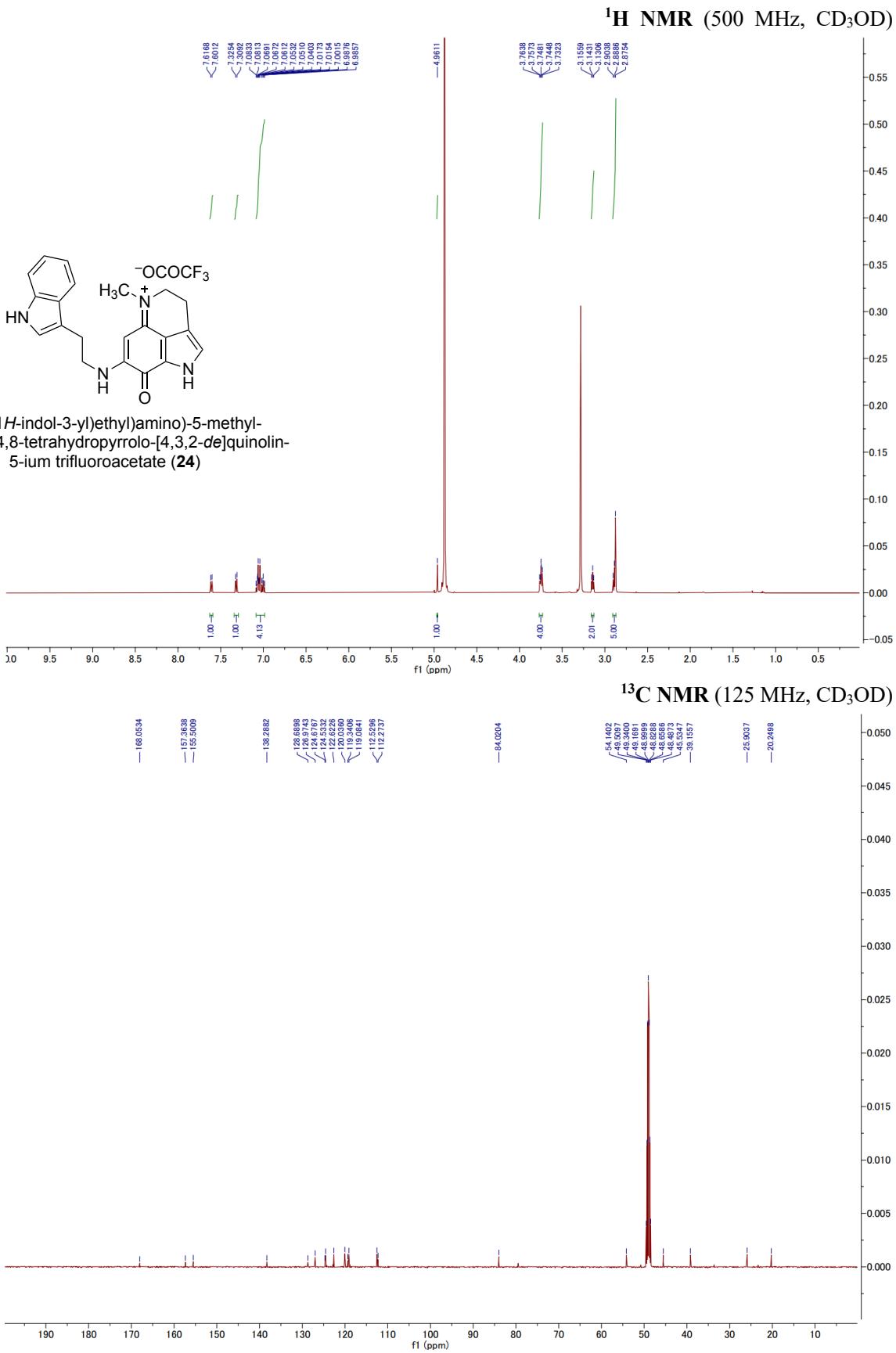


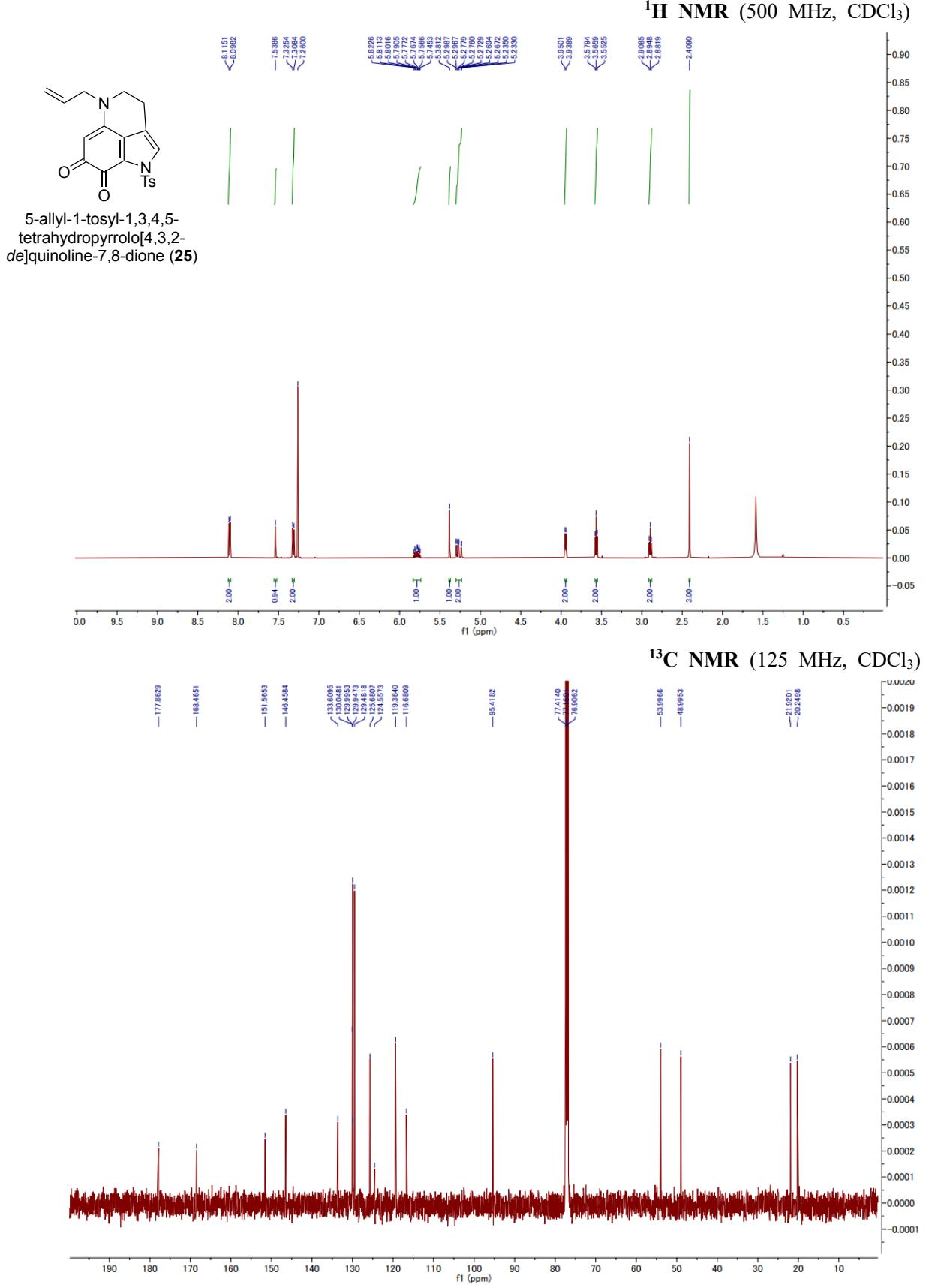
5-methyl-8-oxo-7-(phenethylamino)-1,3,4,8-tetrahydropyrrolo[4,3,2-de]quinolin-5-ium trifluoroacetate (**23**)

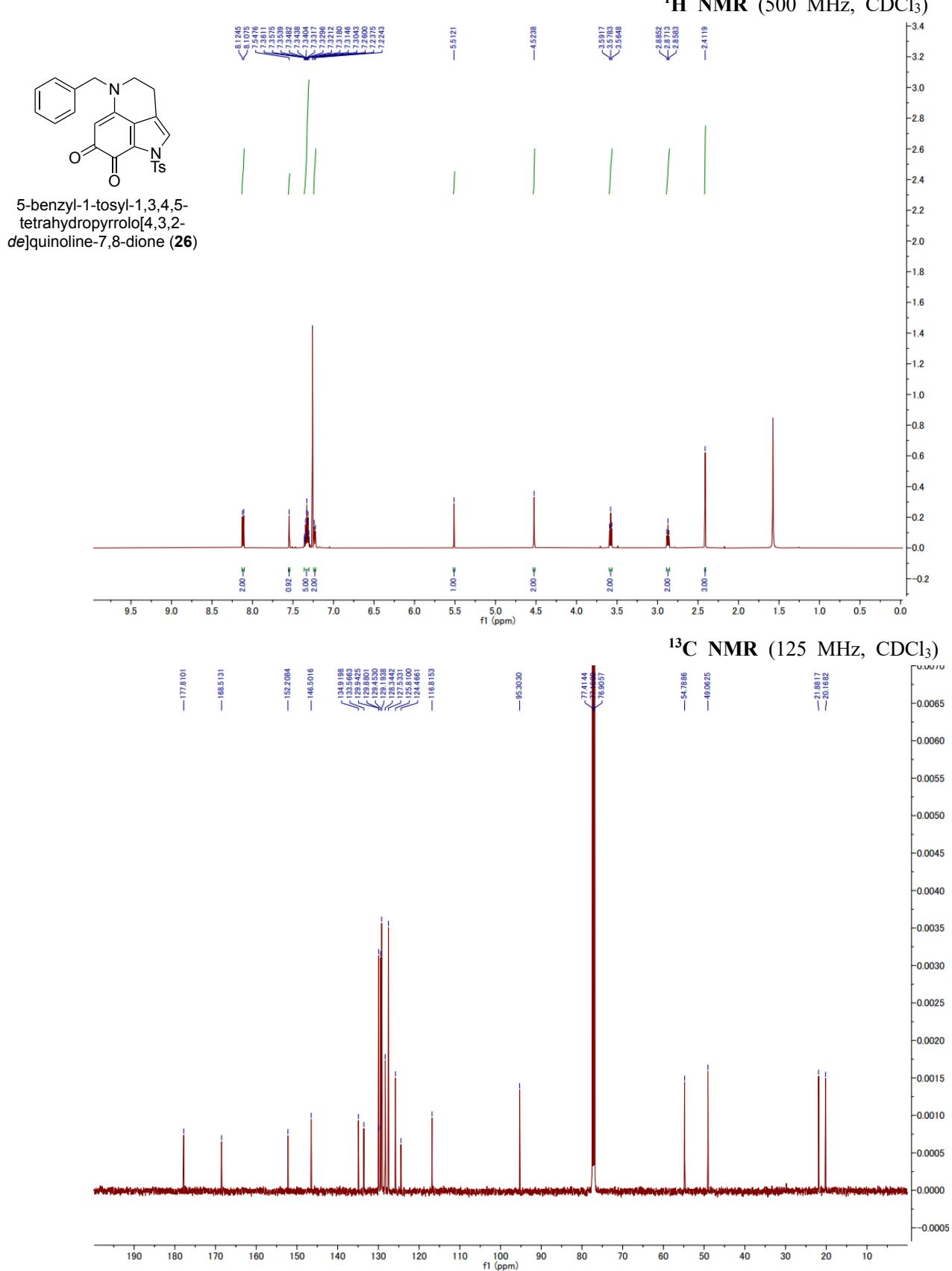


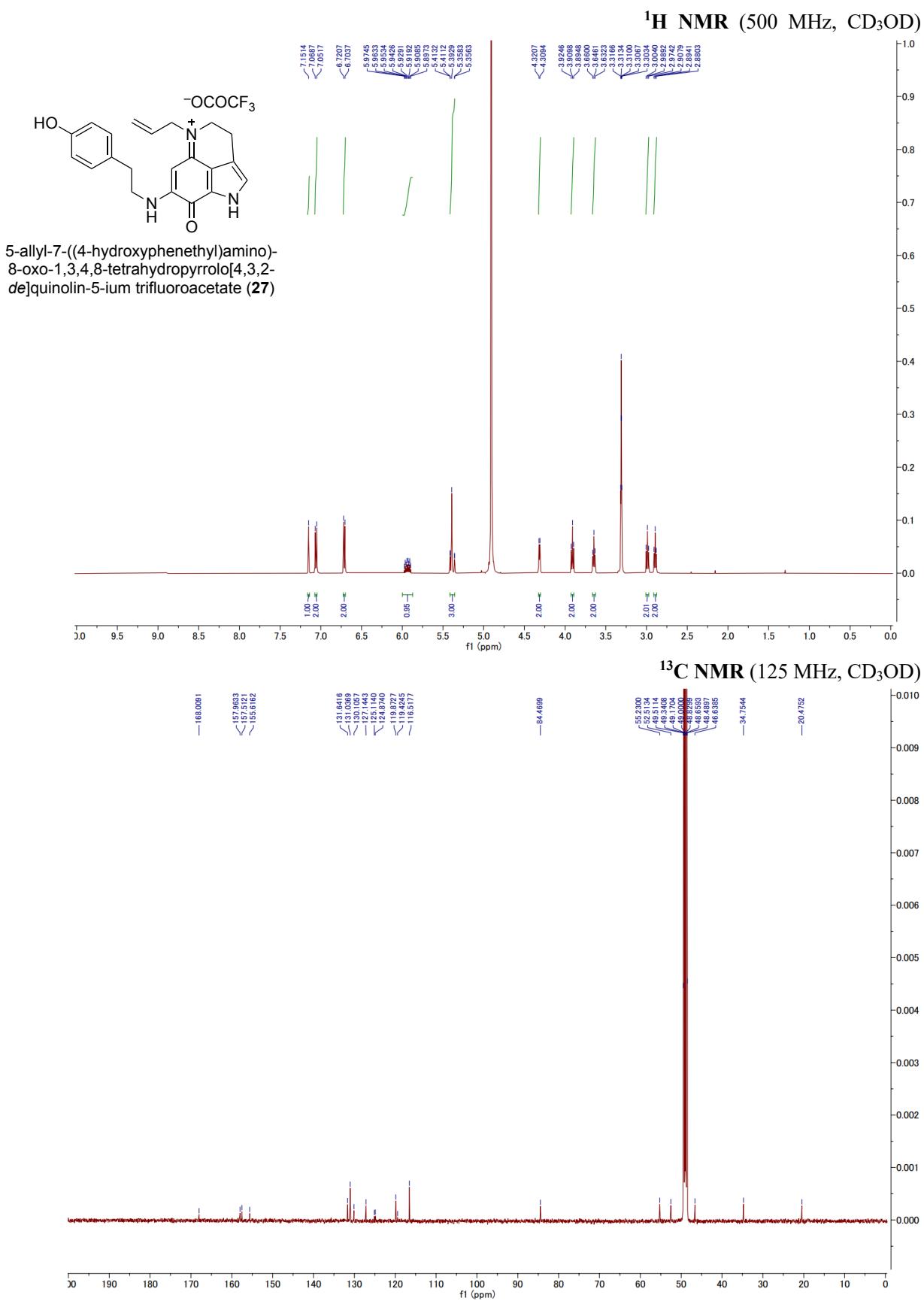


7-((2-(1*H*-indol-3-yl)ethyl)amino)-5-methyl-8-oxo-1,3,4,8-tetrahydropyrrolo-[4,3,2-*d*]quinolin-5-i^{um} trifluoroacetate (**24**)









¹H NMR (500 MHz, CD₃OD)

