

Supporting Information

Solvent-controlled regioselective C(5)–H/N(1)–H bond alkylations of indolines and C(6)–H bond alkylations of 1,2,3,4-tetrahydroquinolines with *para*-quinone methides

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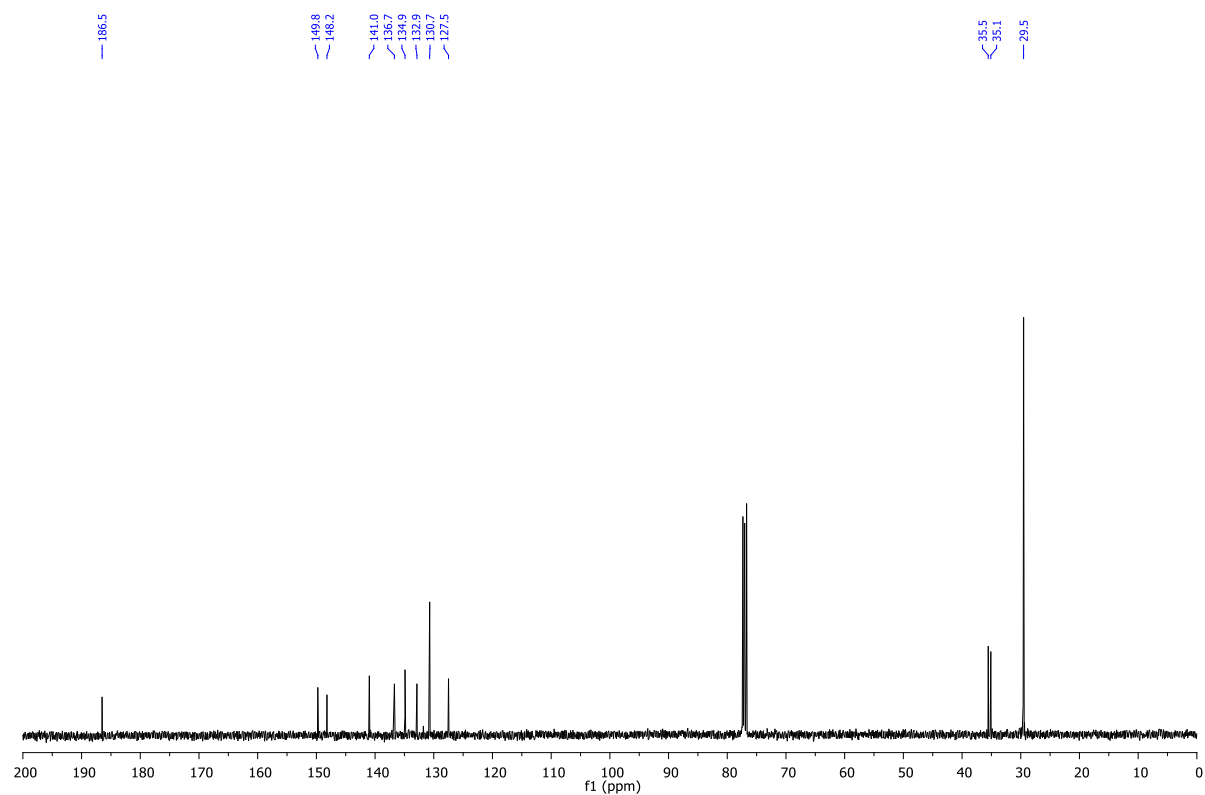
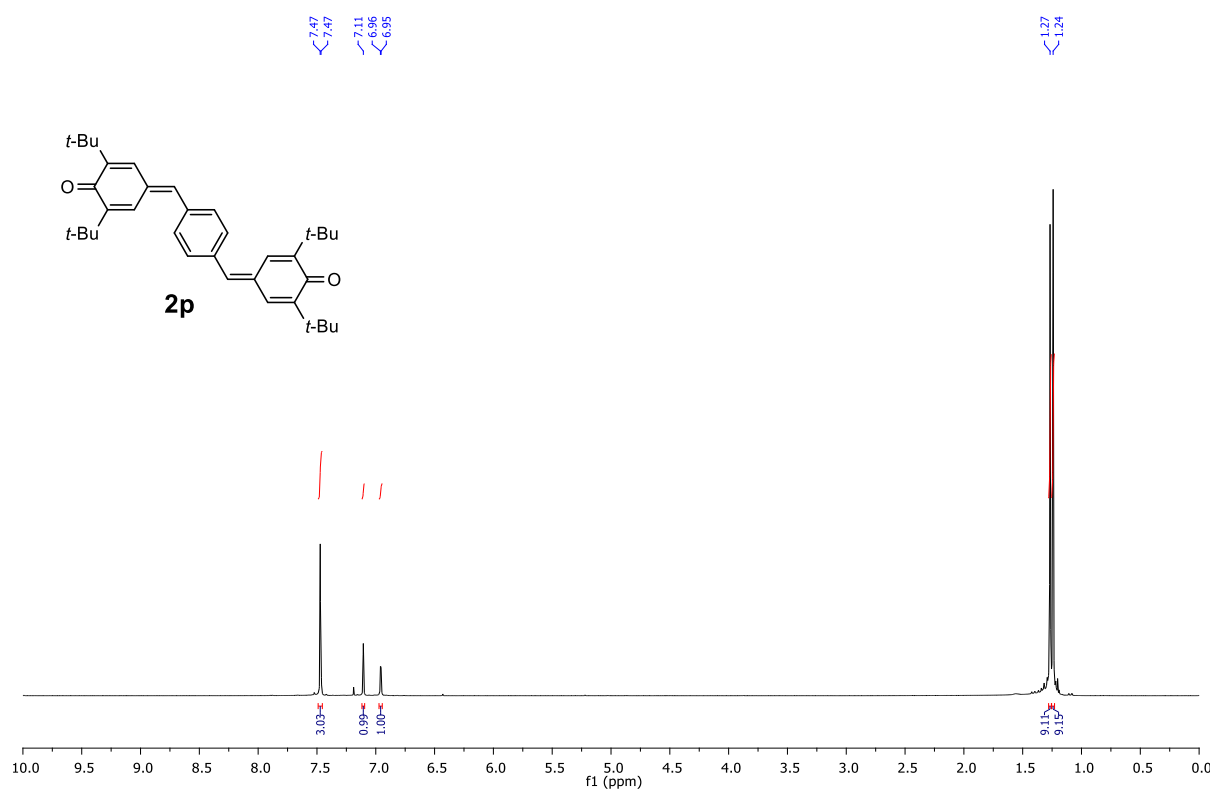
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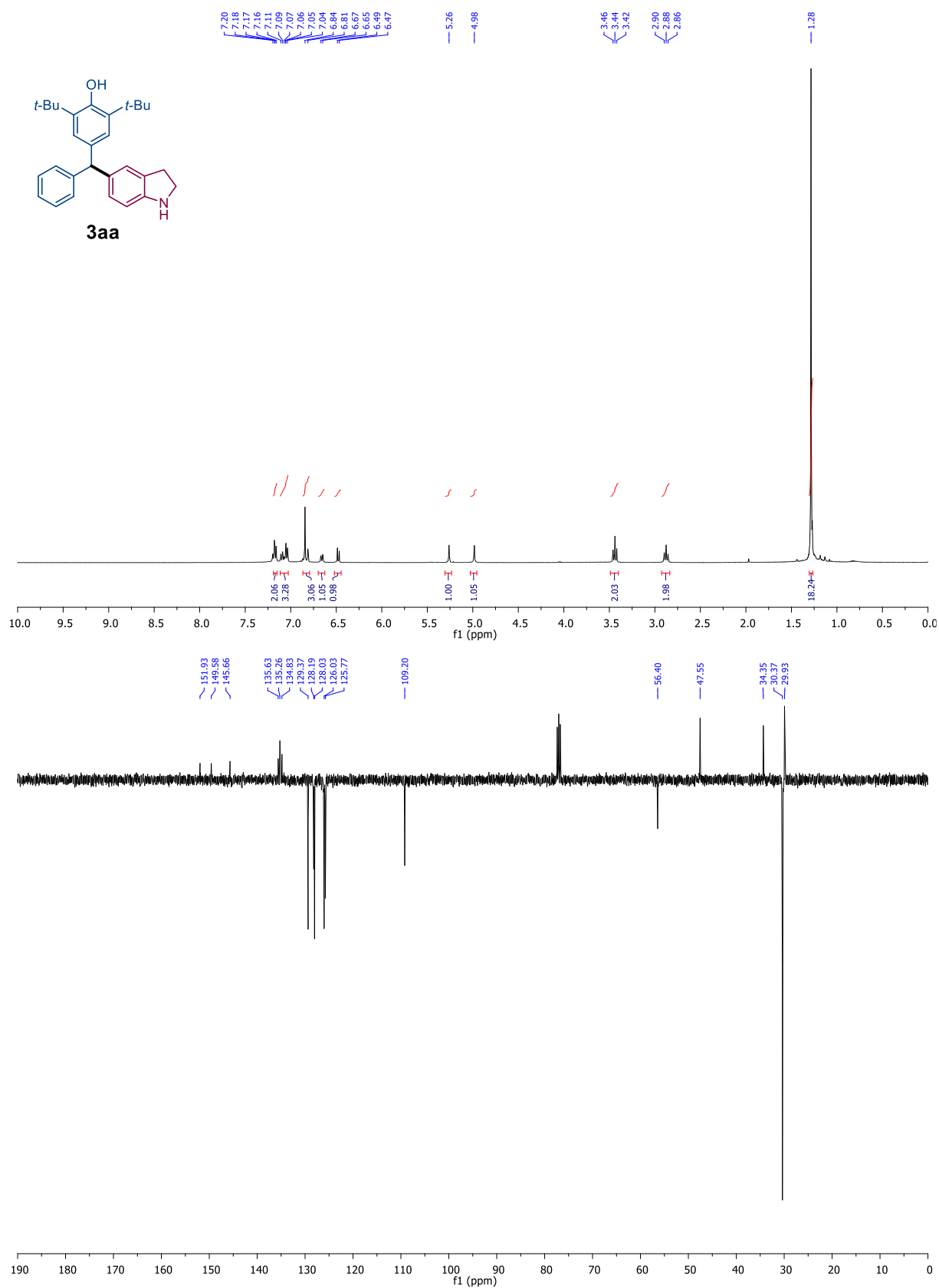
Email: nsarac@atauni.edu.tr

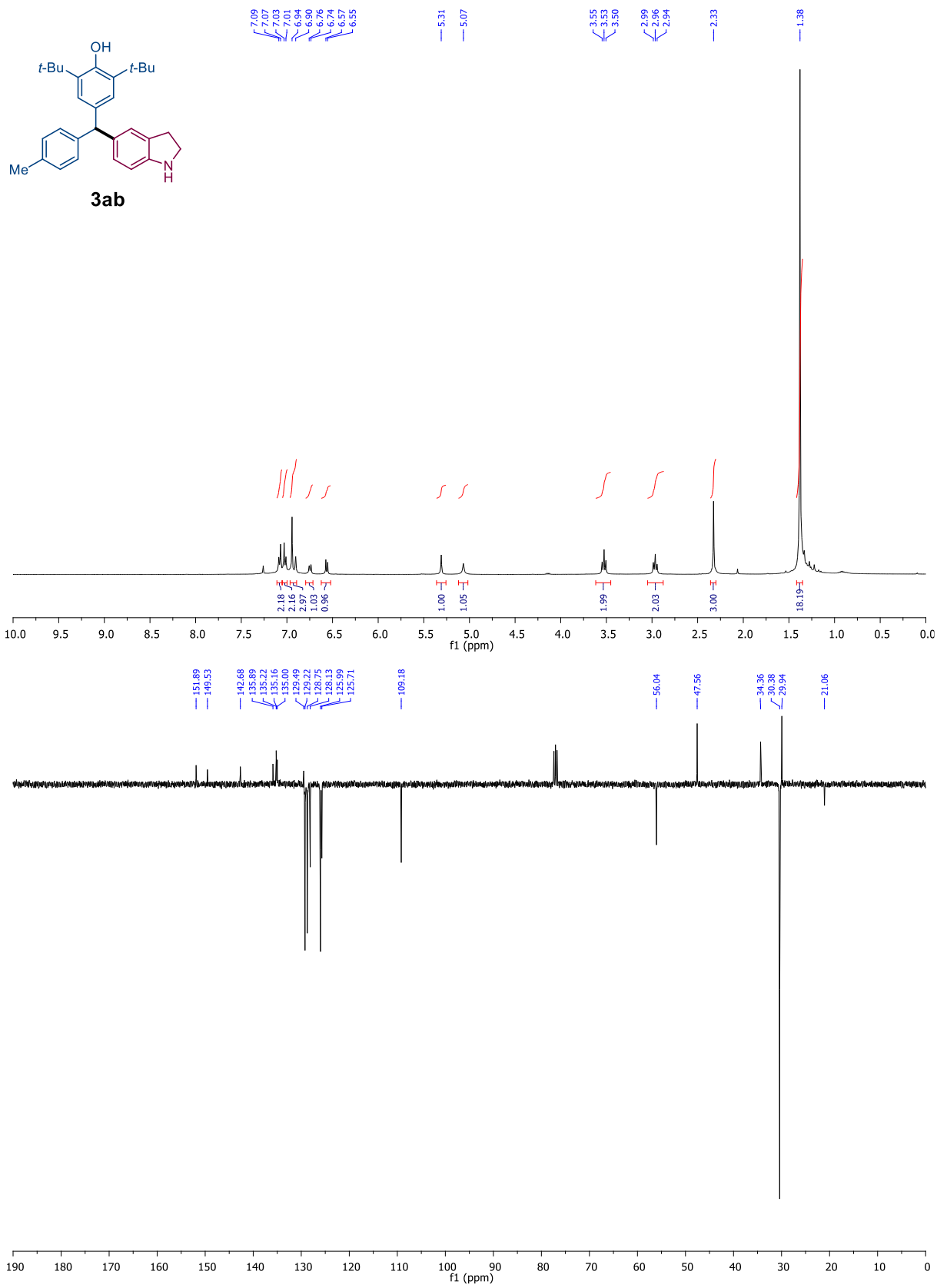
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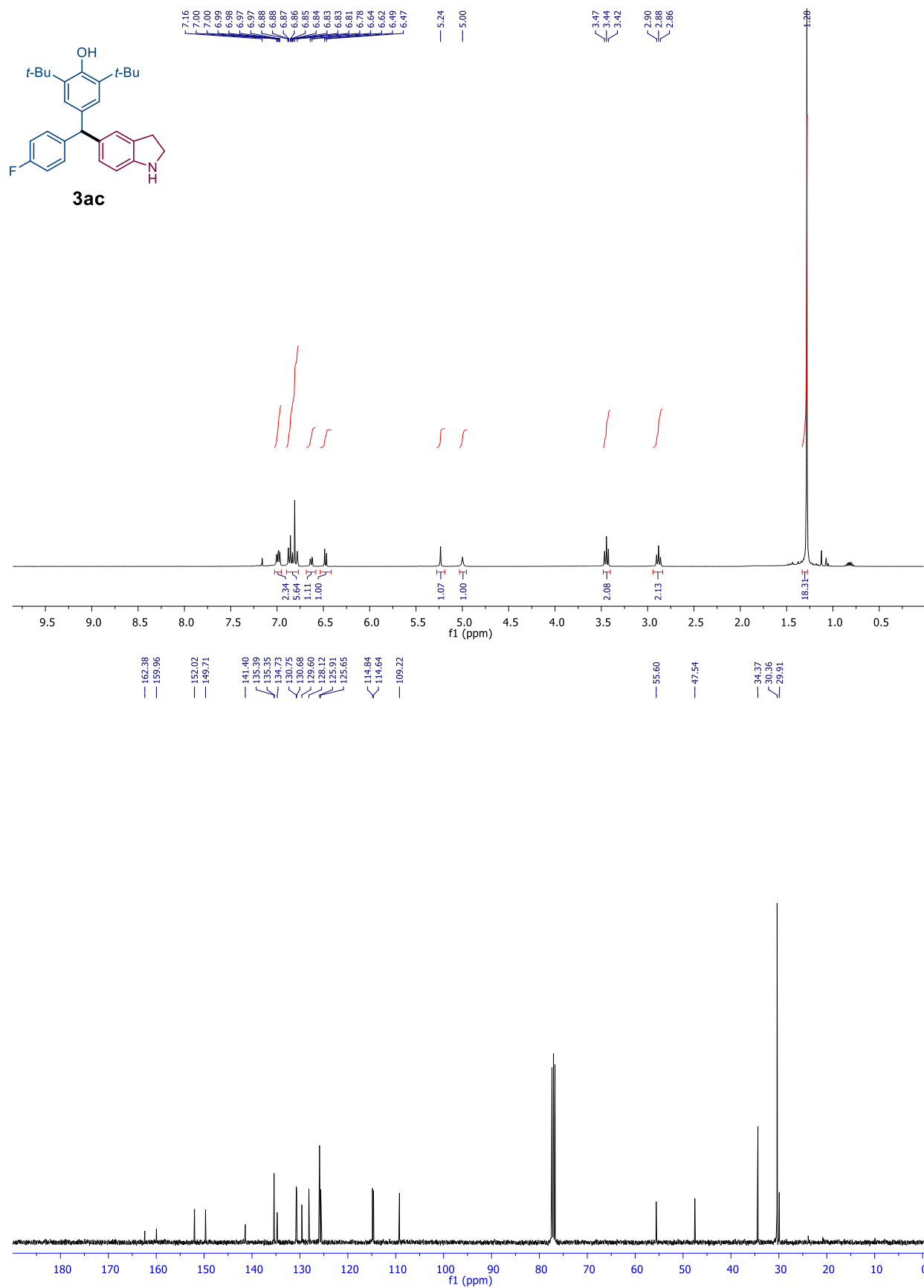
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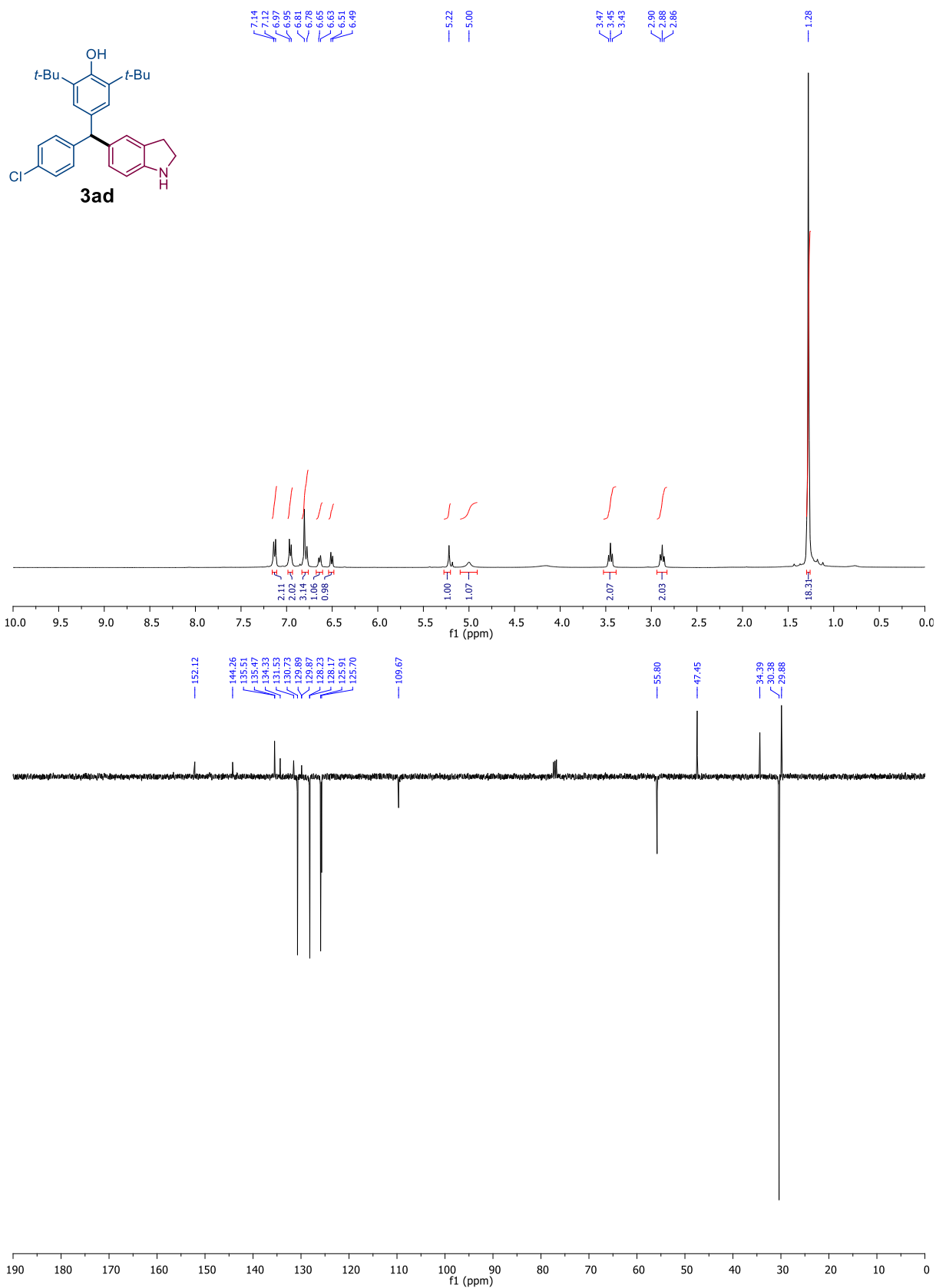
1. ^1H NMR (400 MHz) and ^{13}C NMR (100 MHz) spectra of compounds

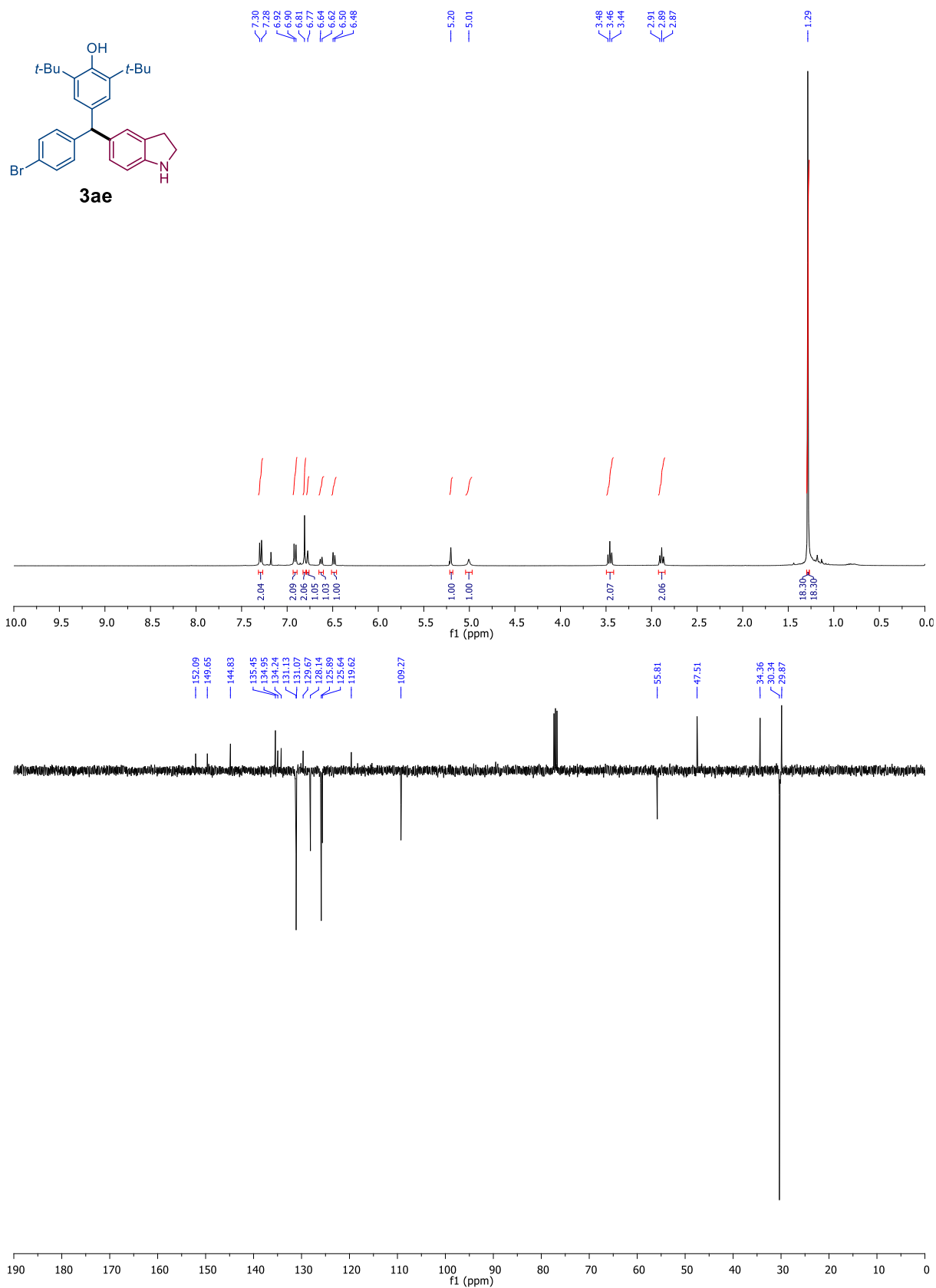


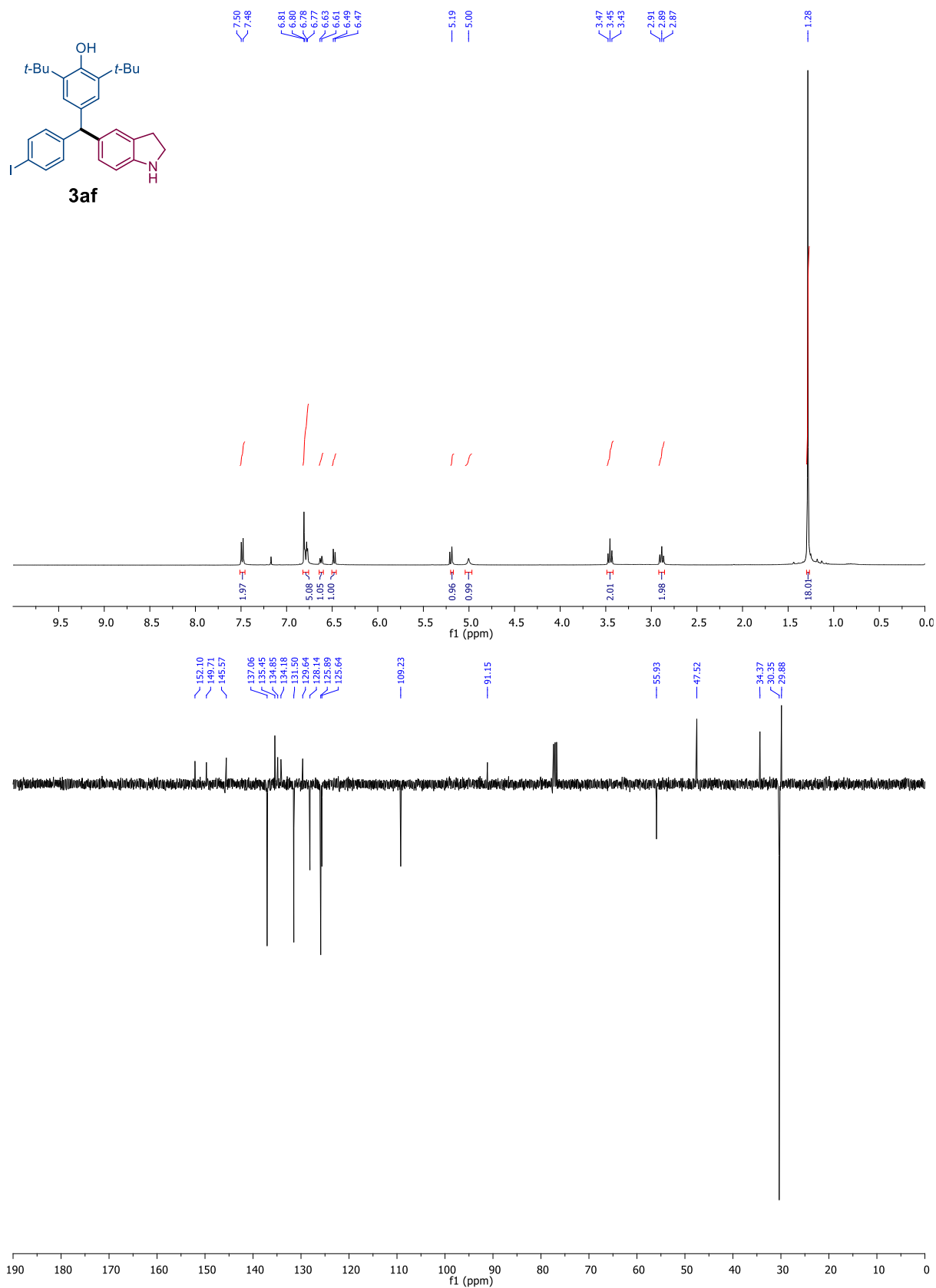


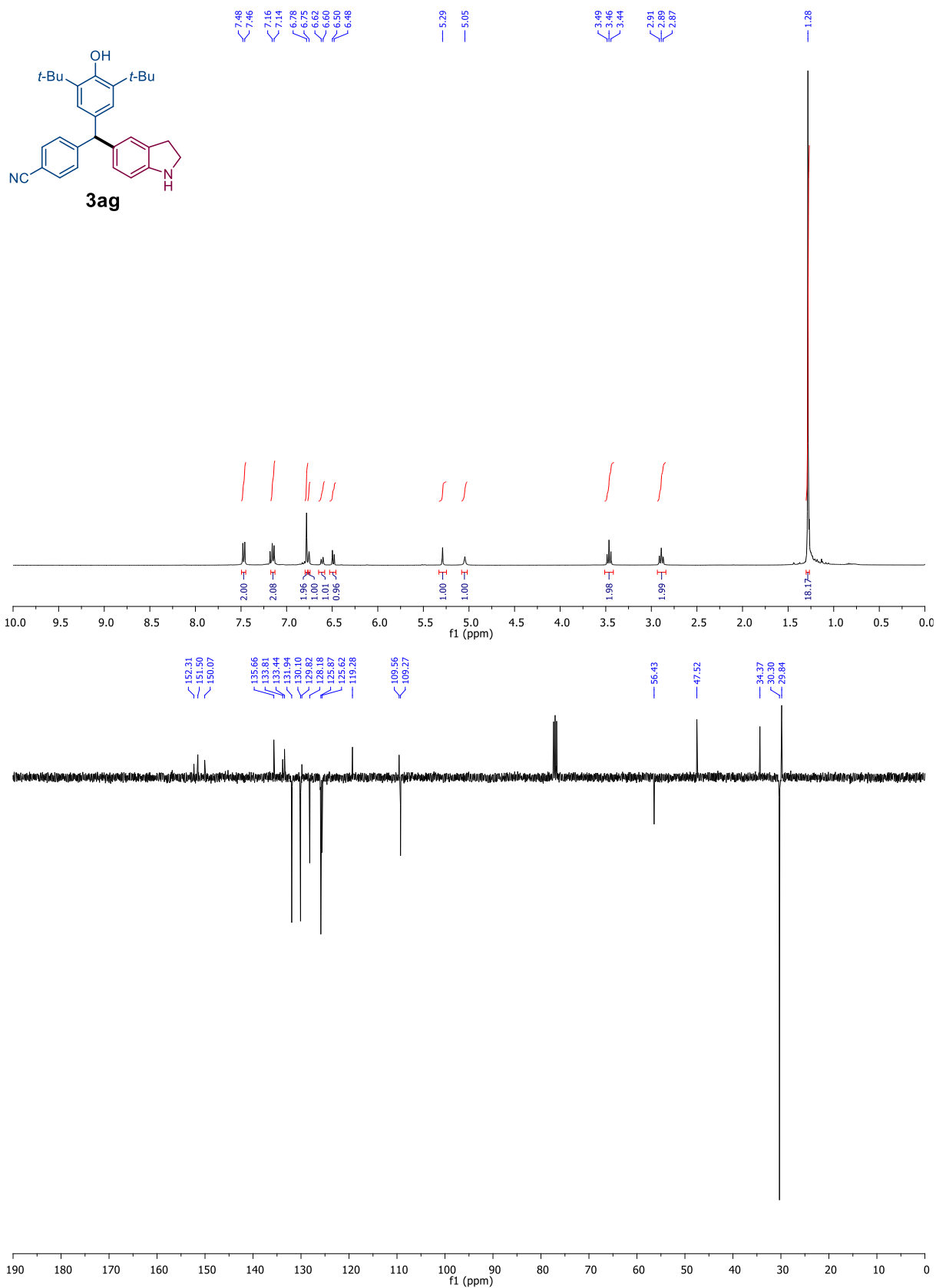


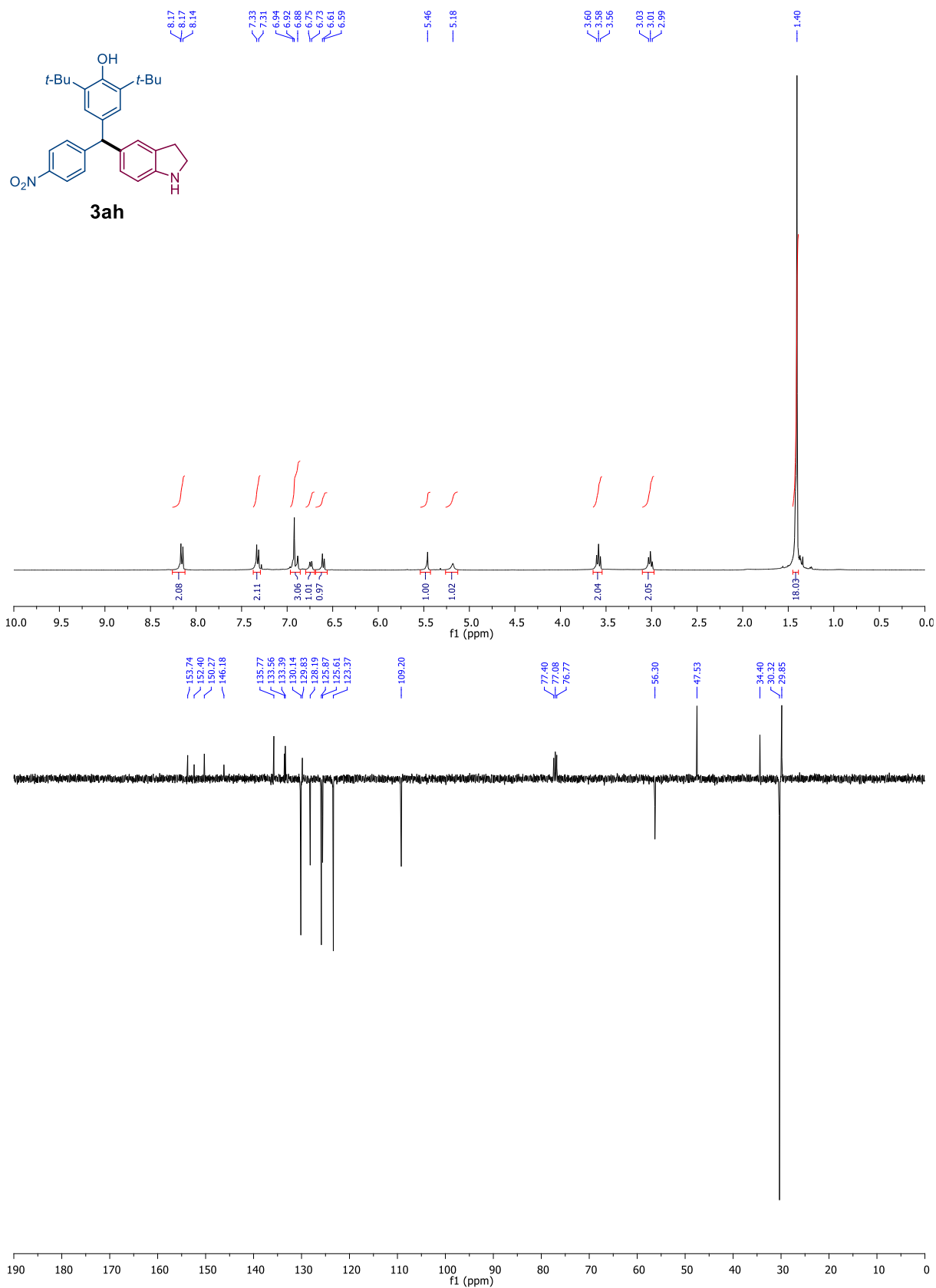


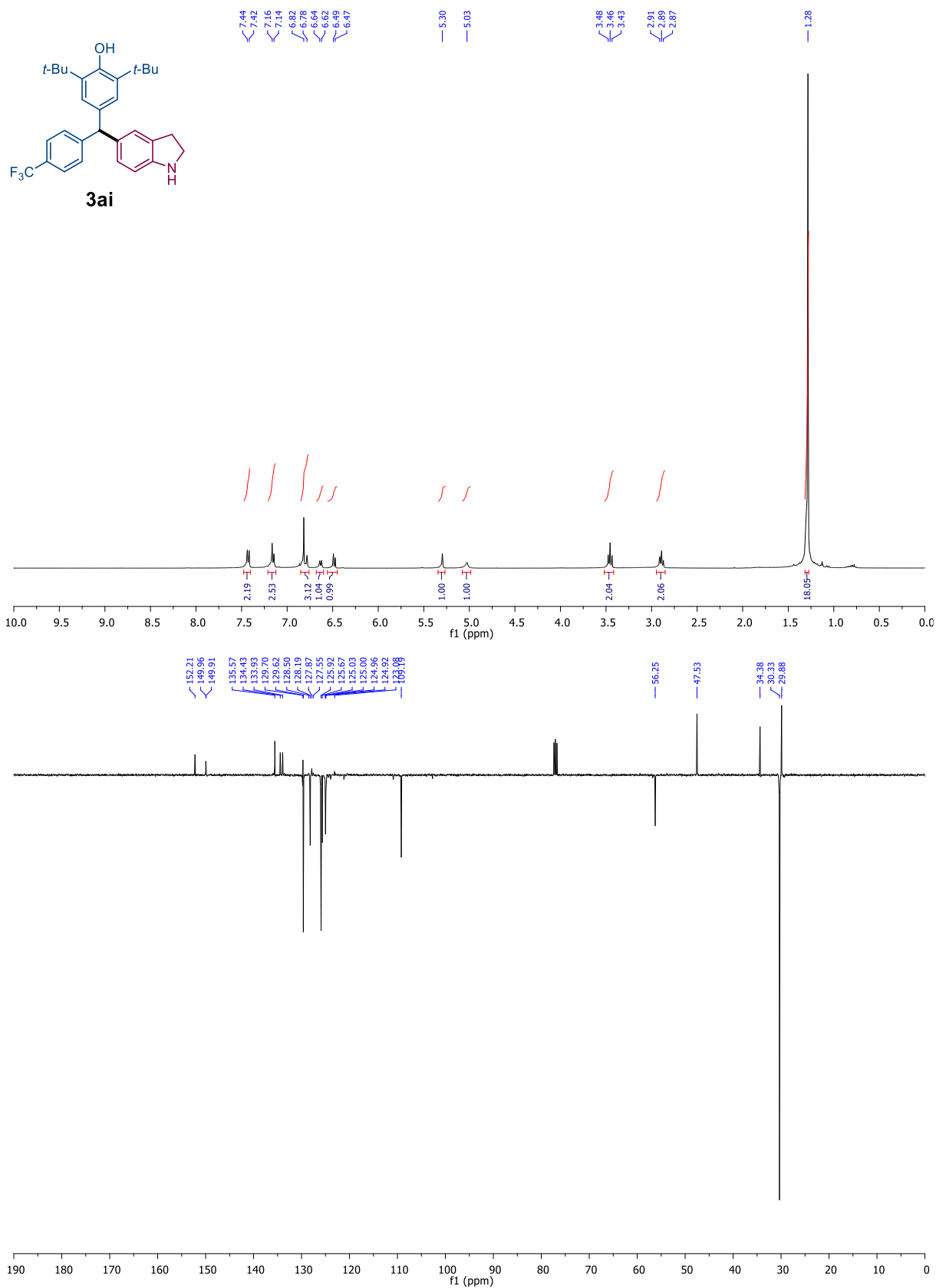


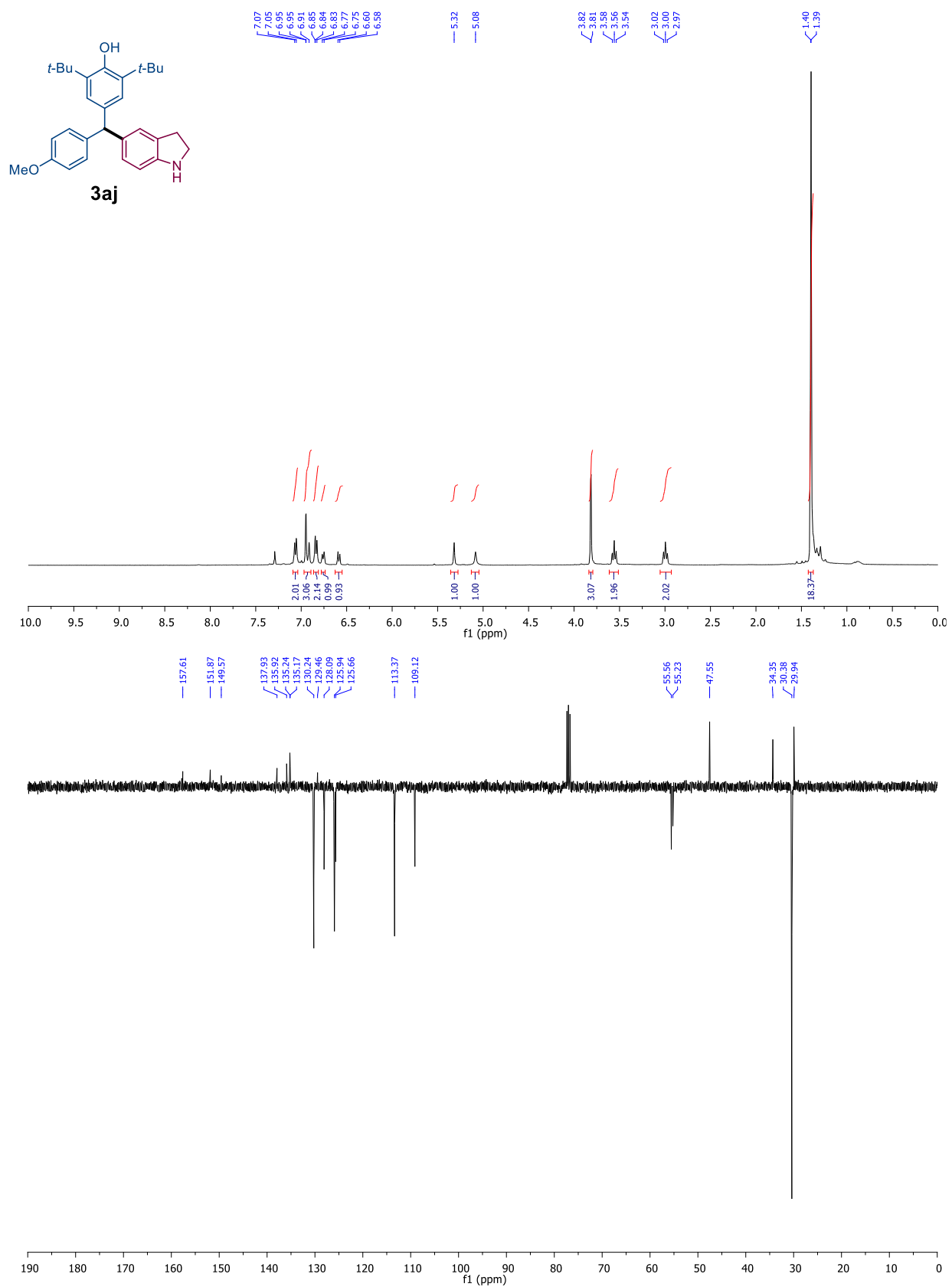


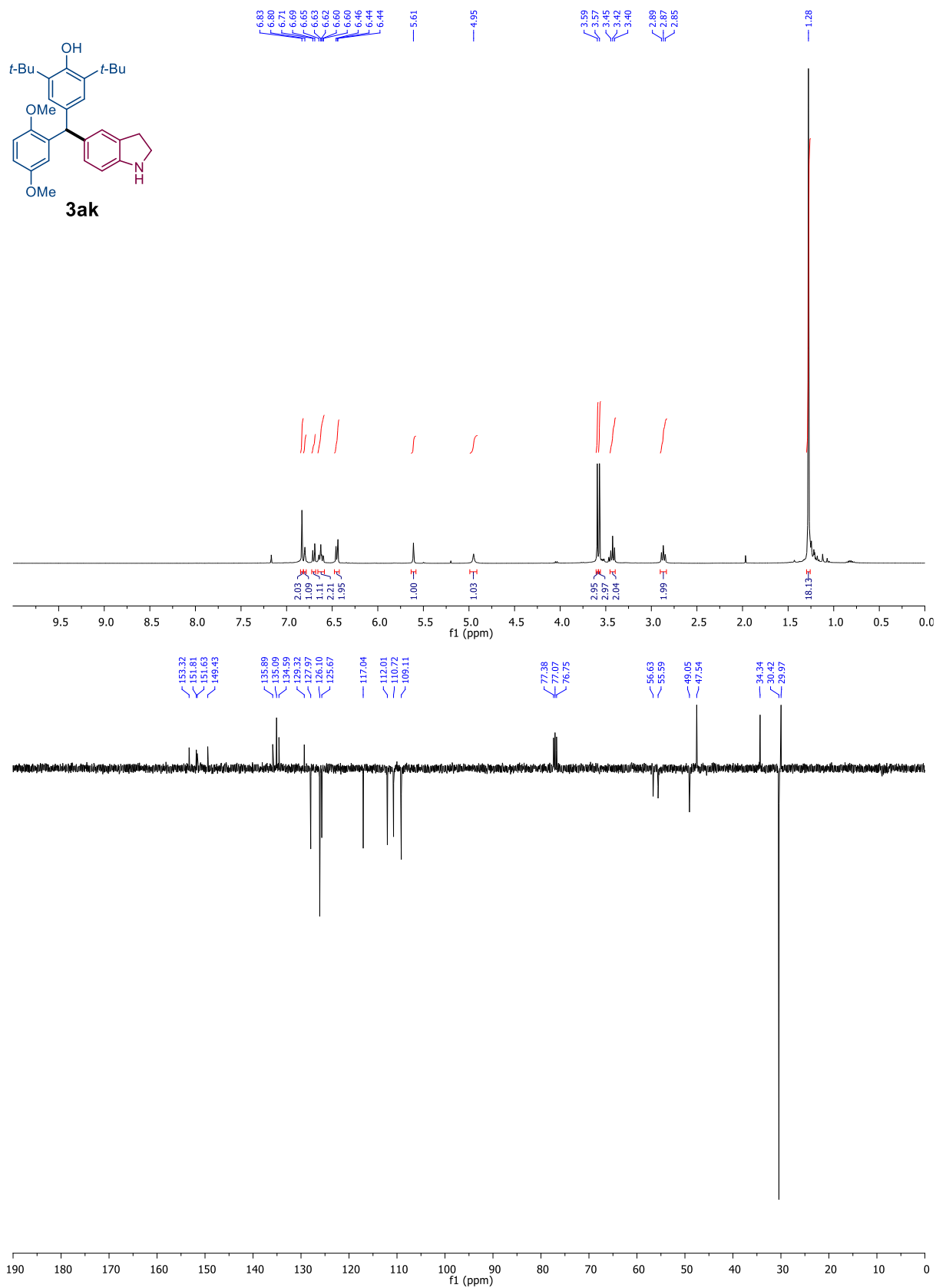


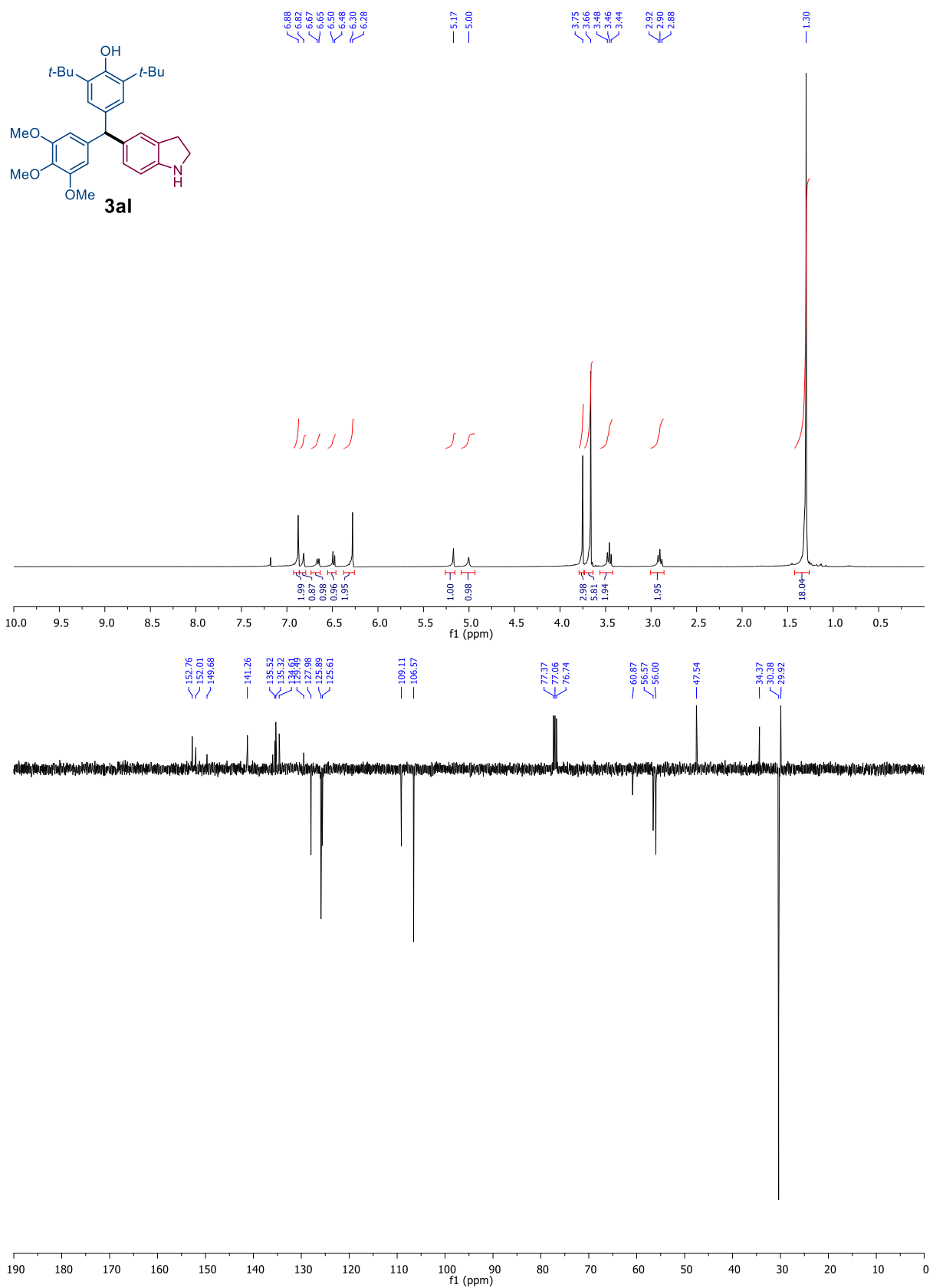


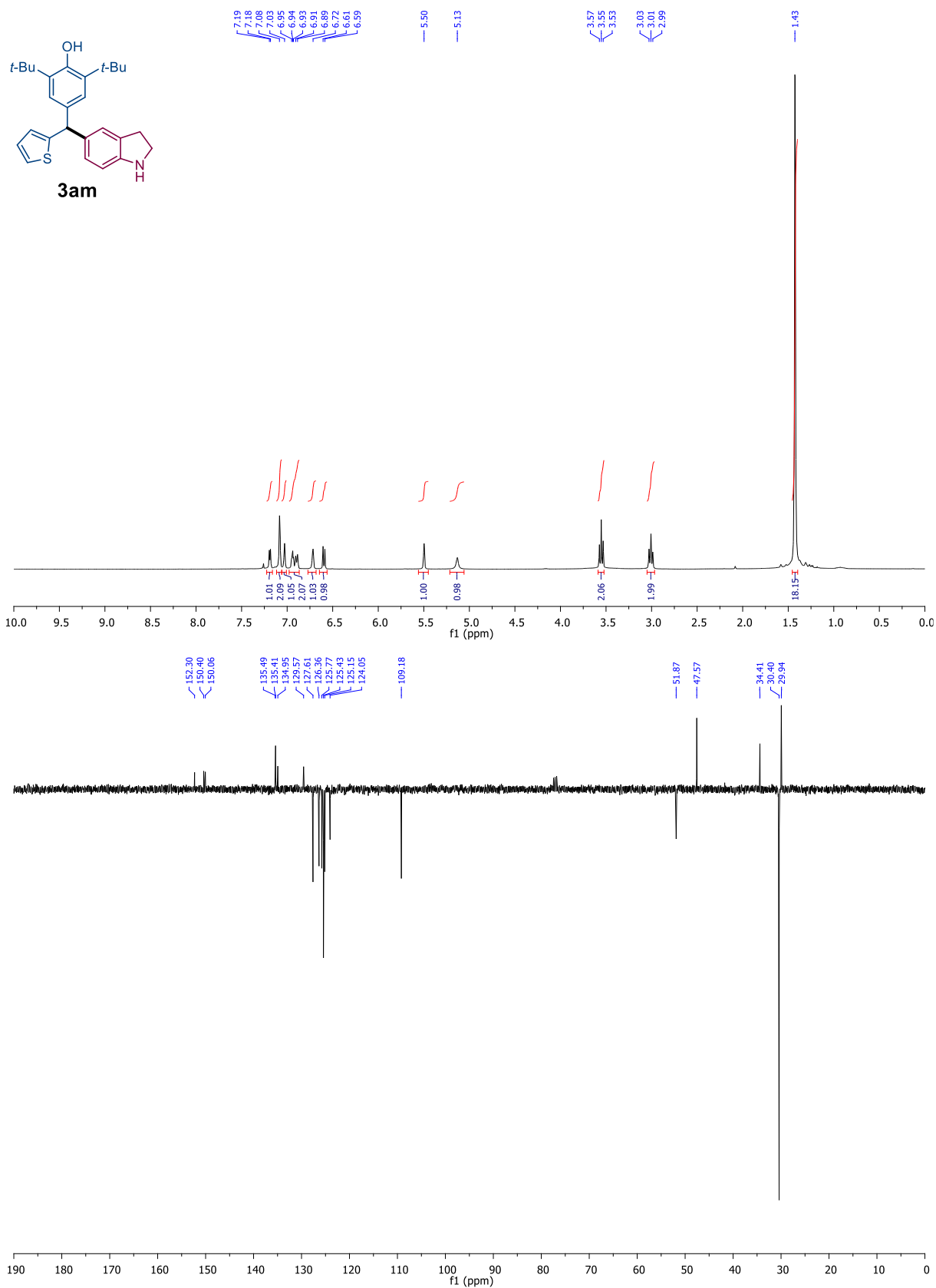




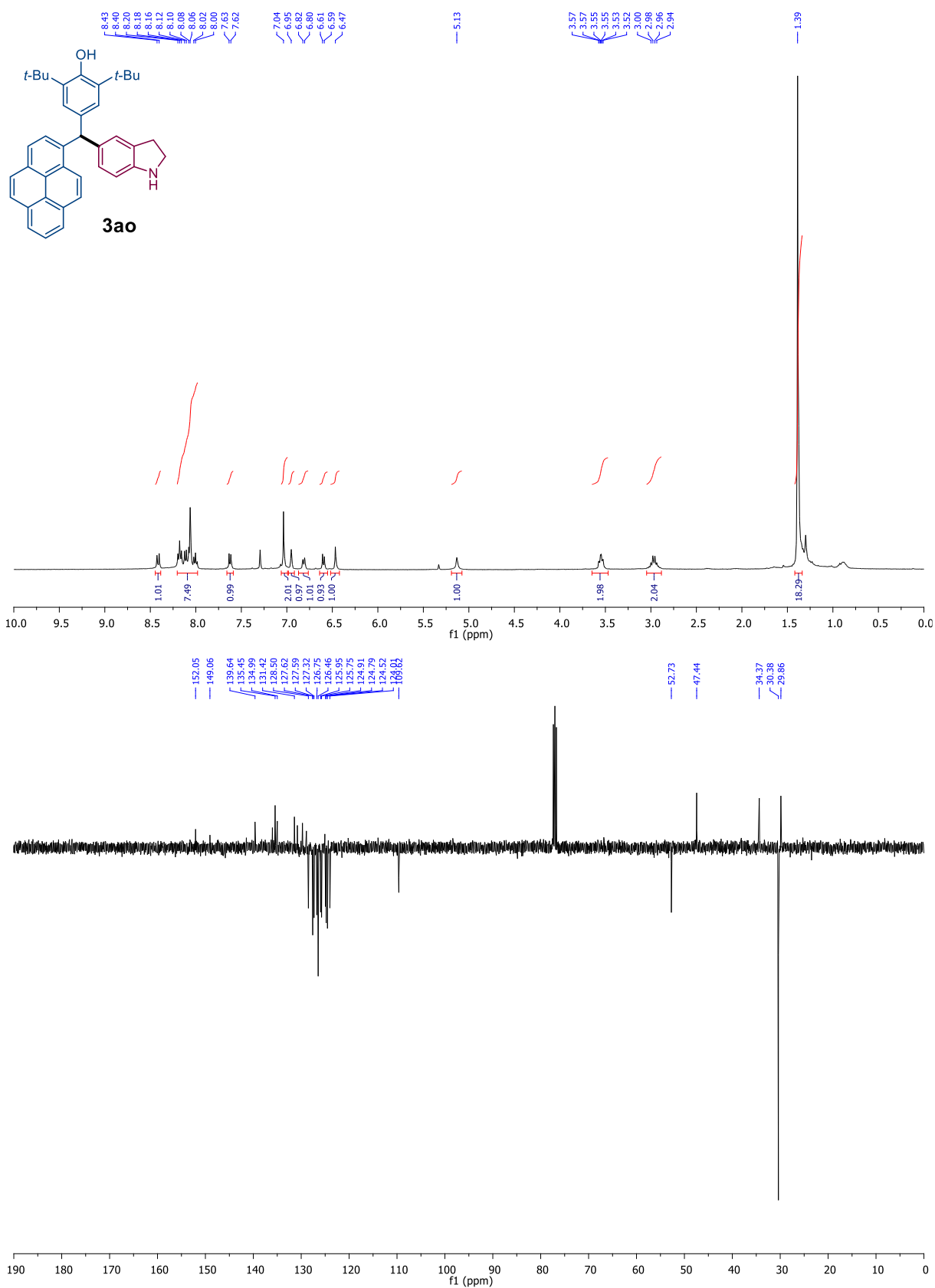


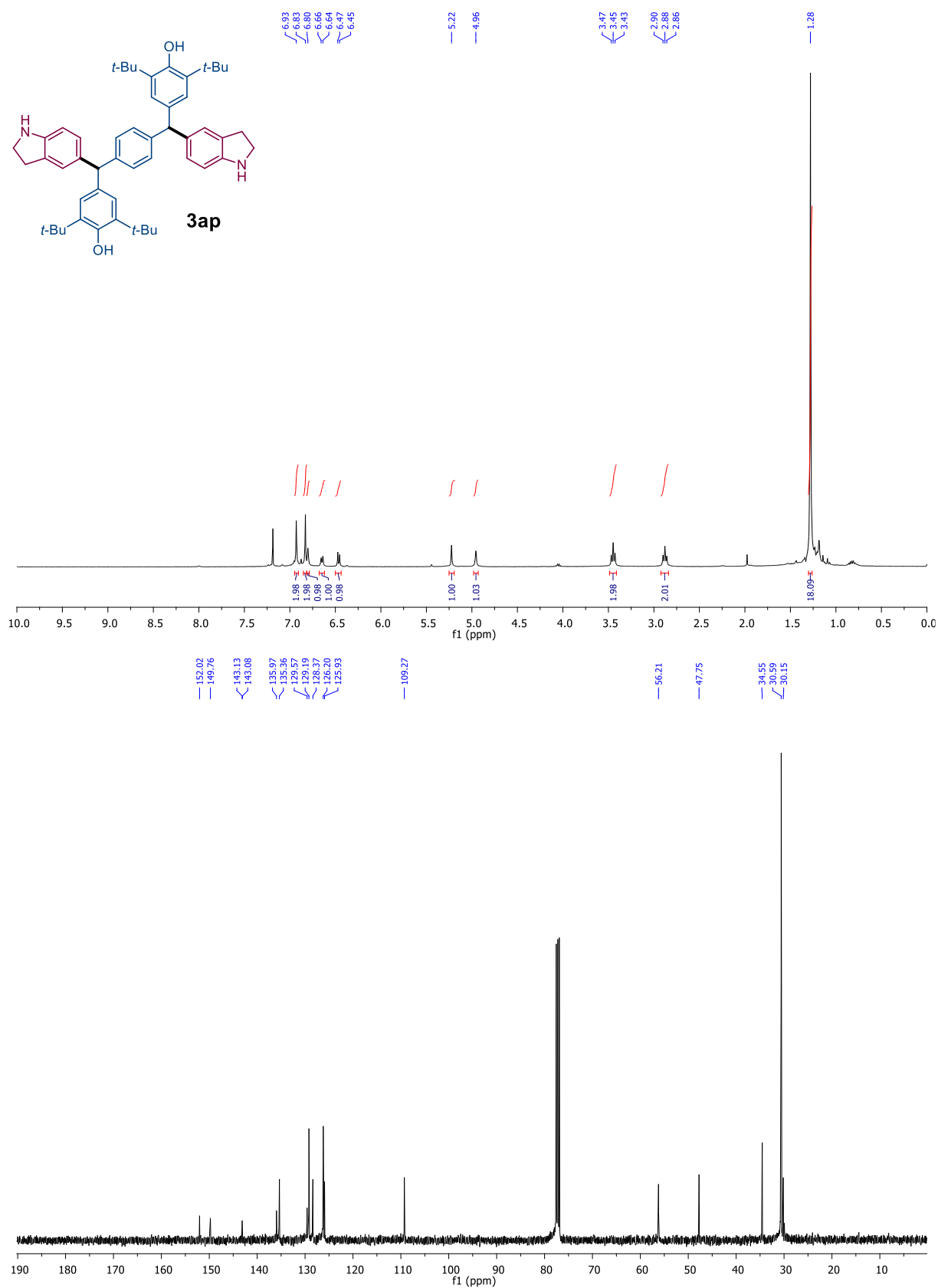


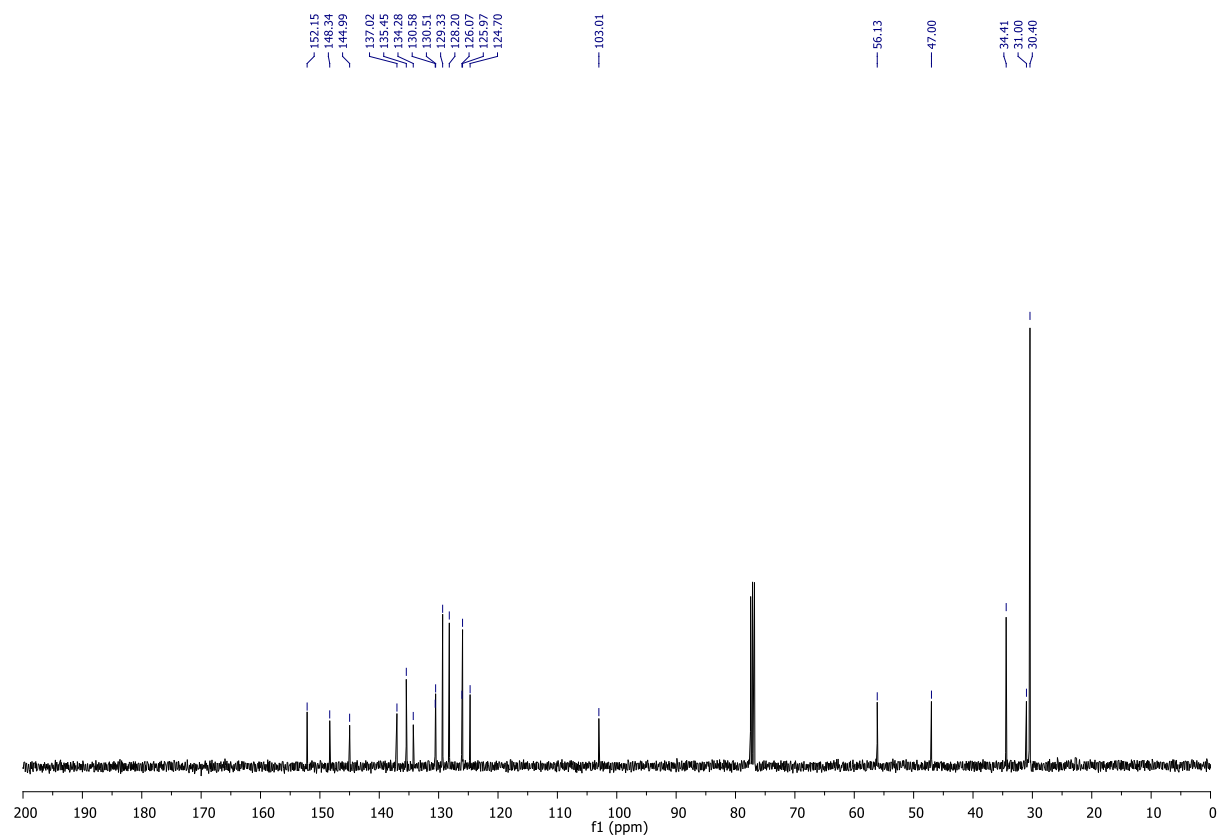
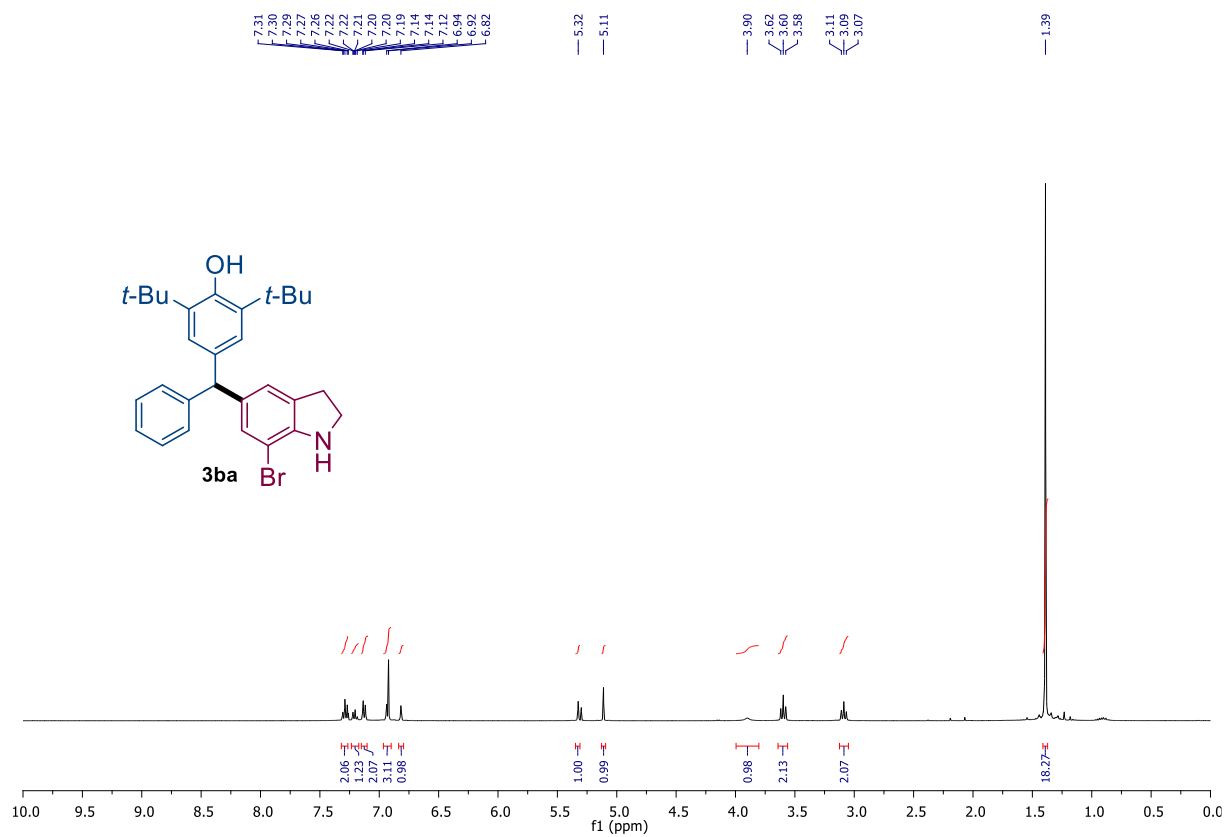


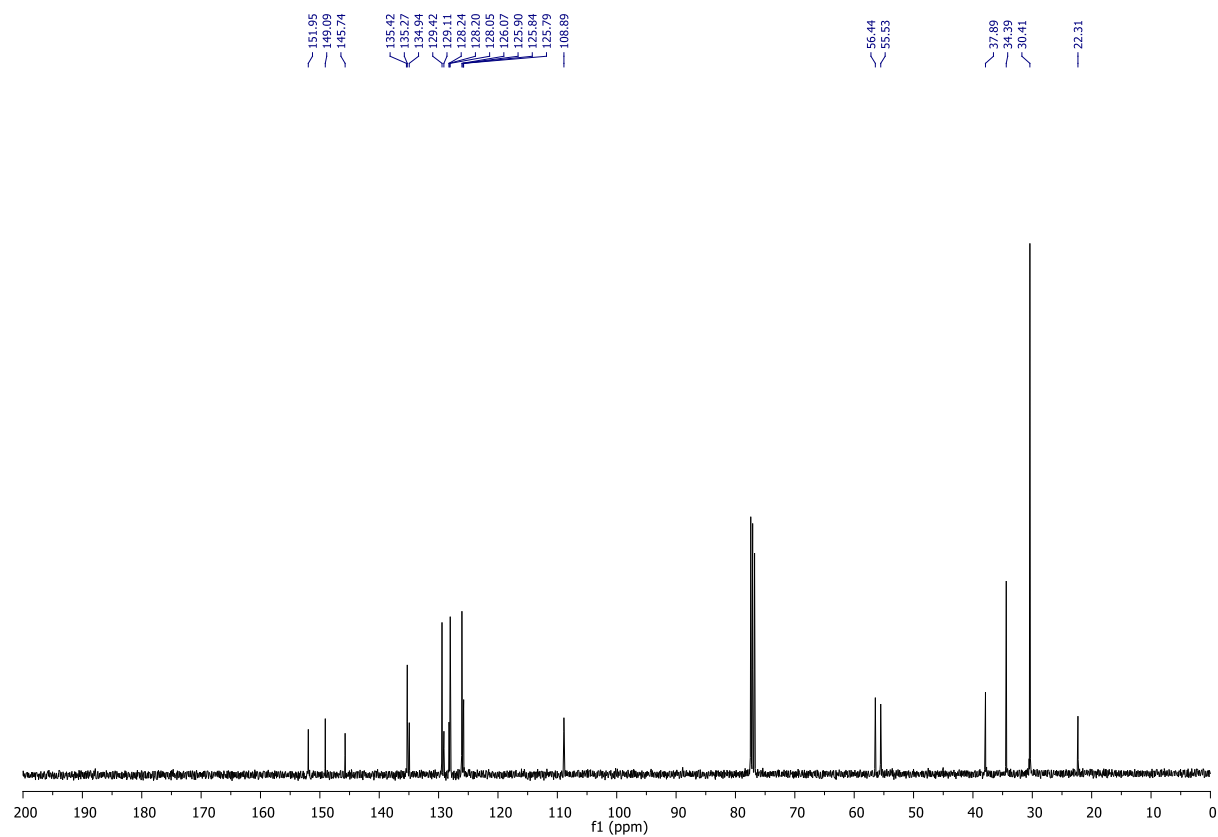
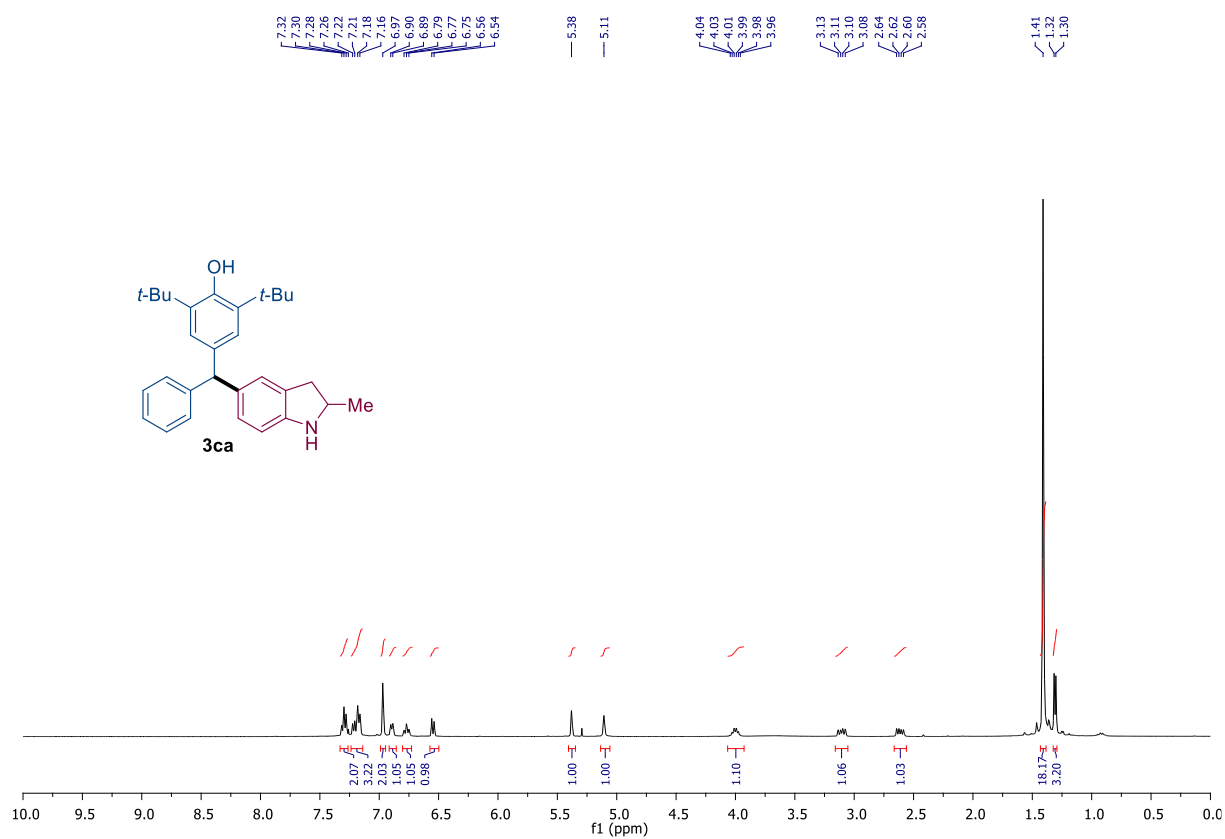


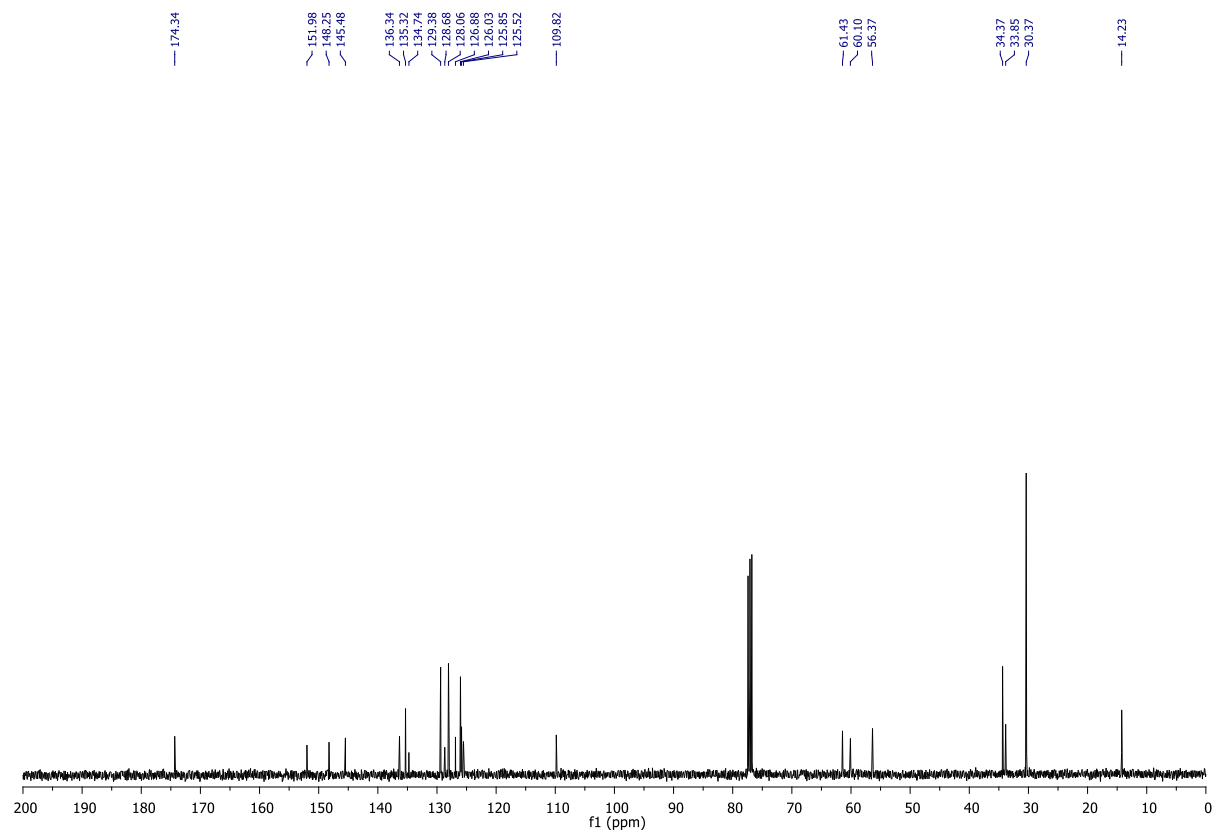
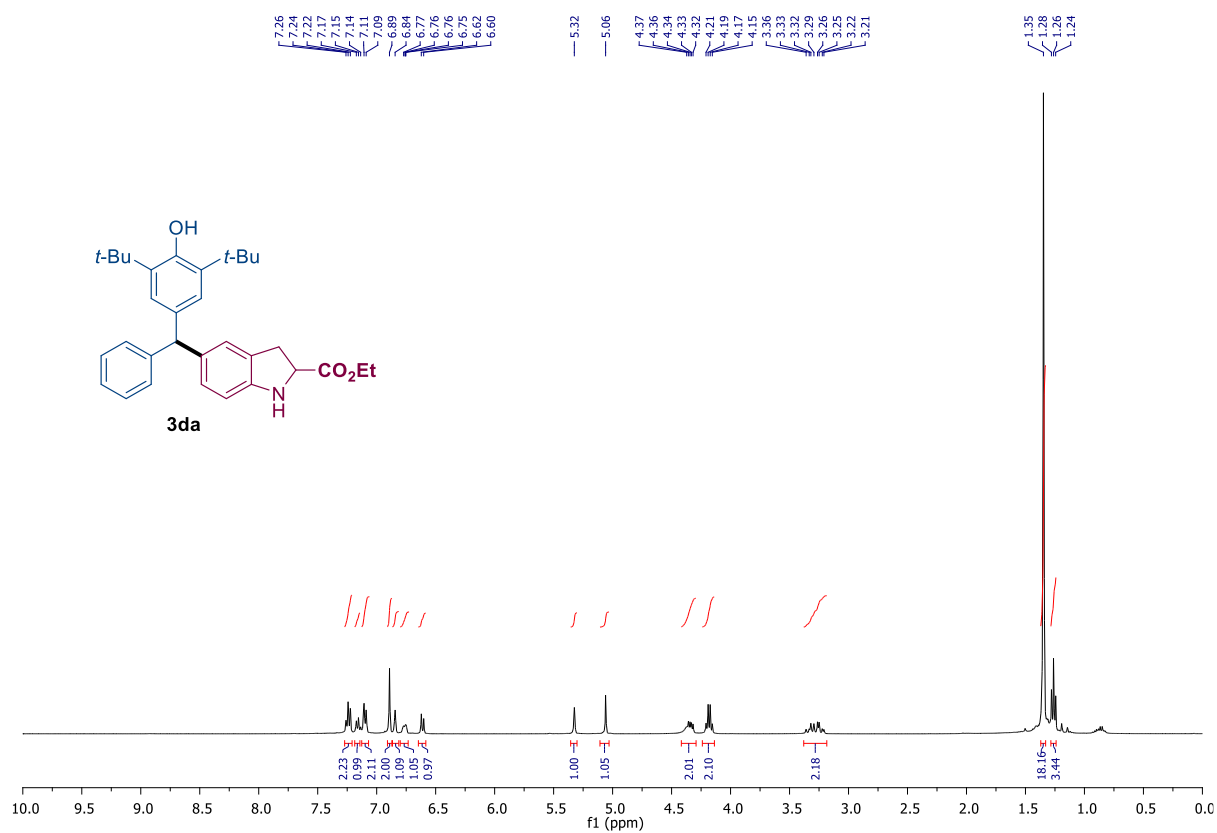


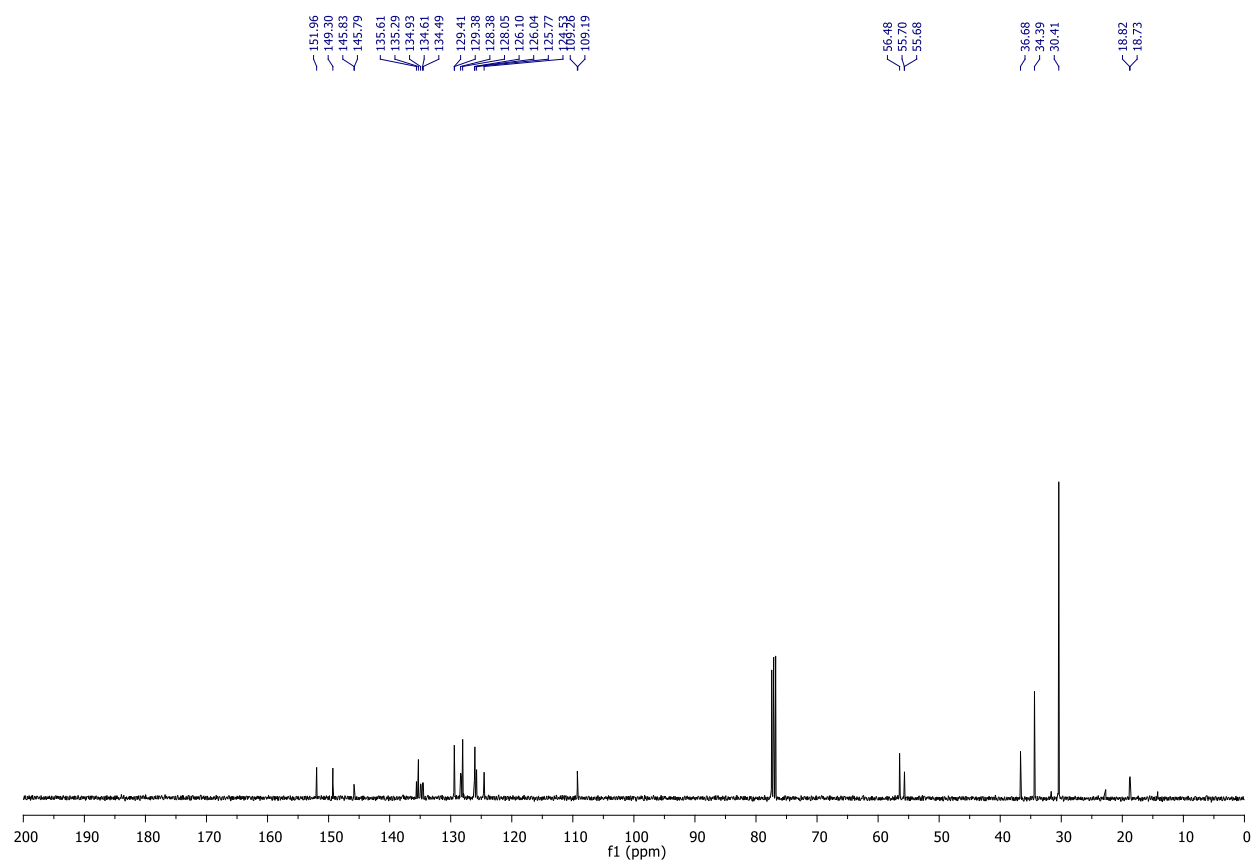
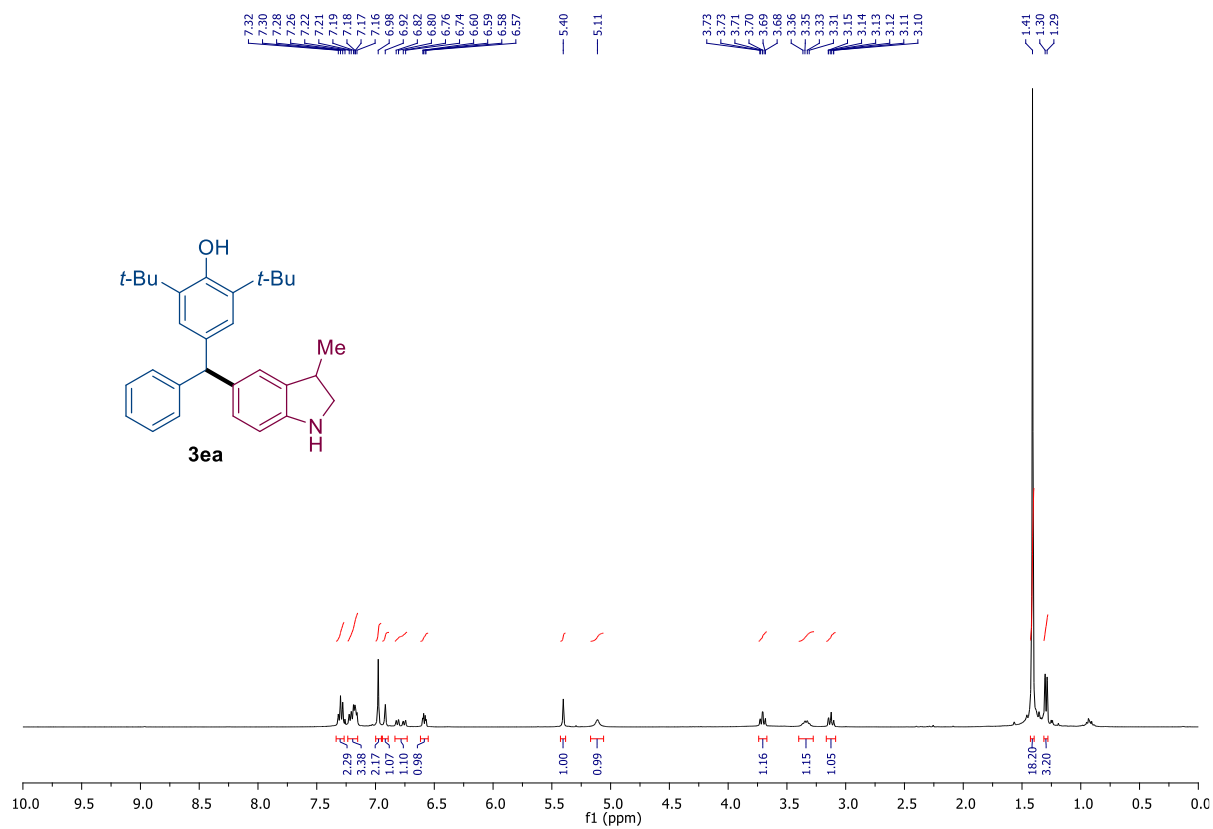


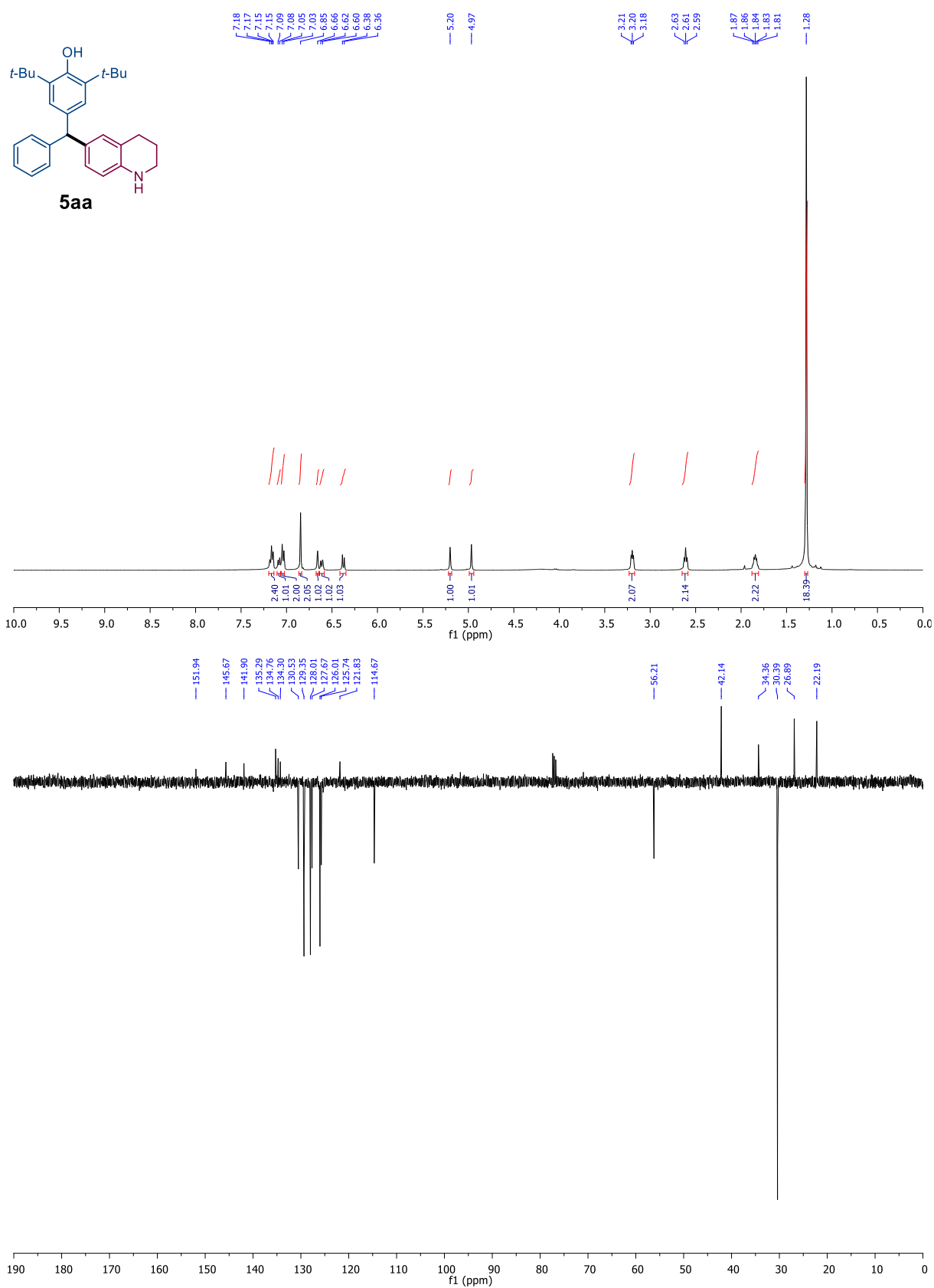


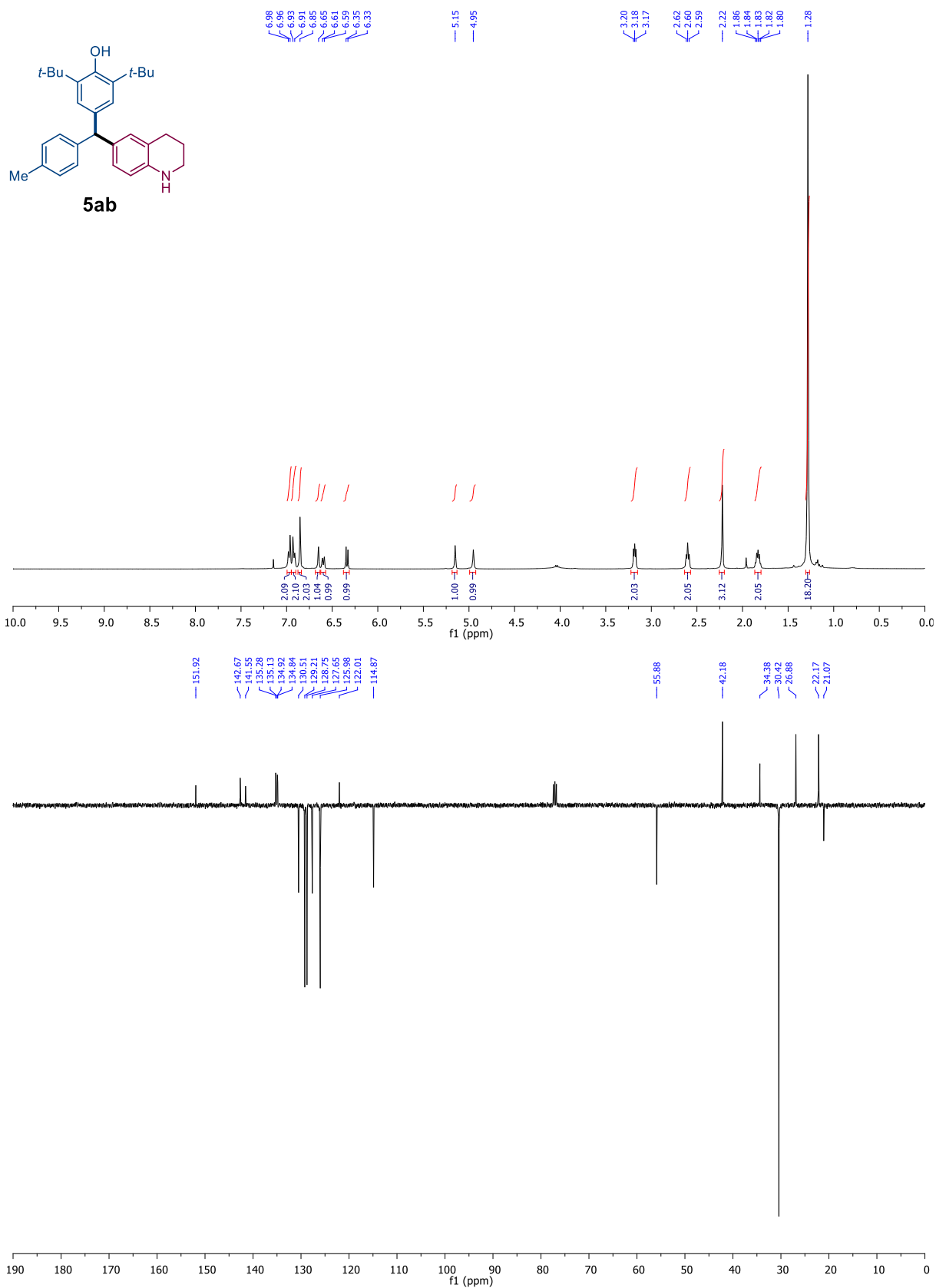


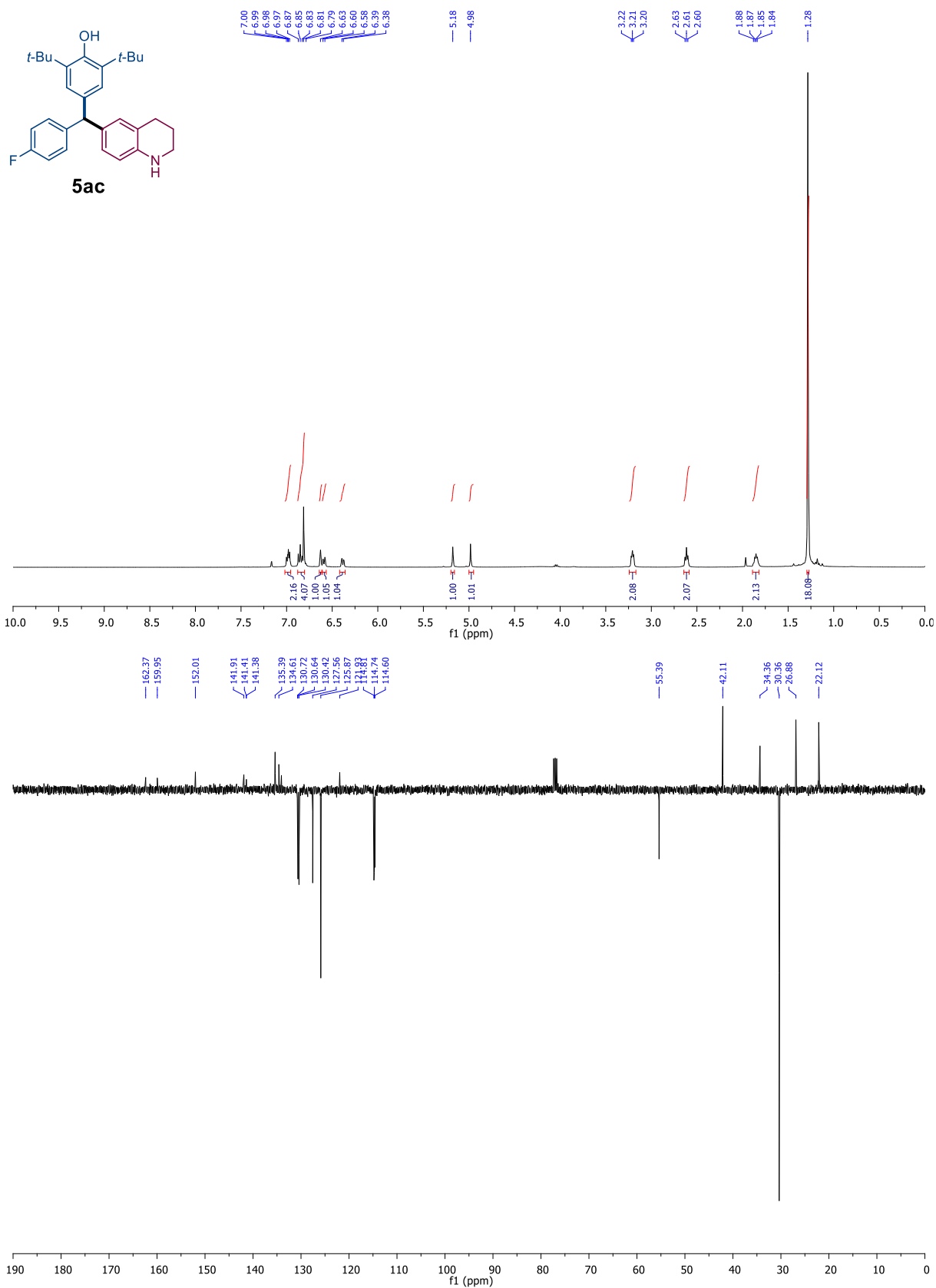


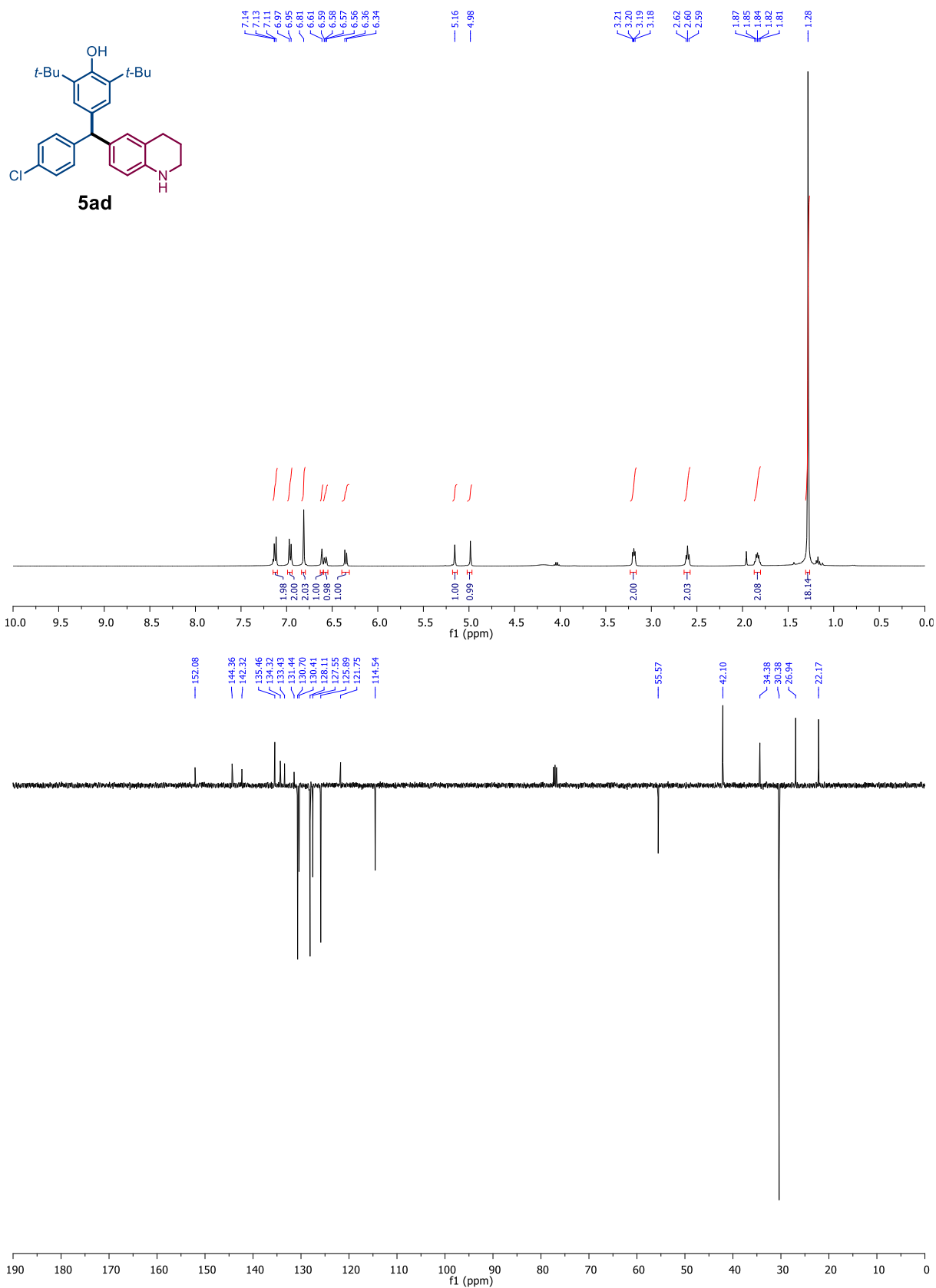


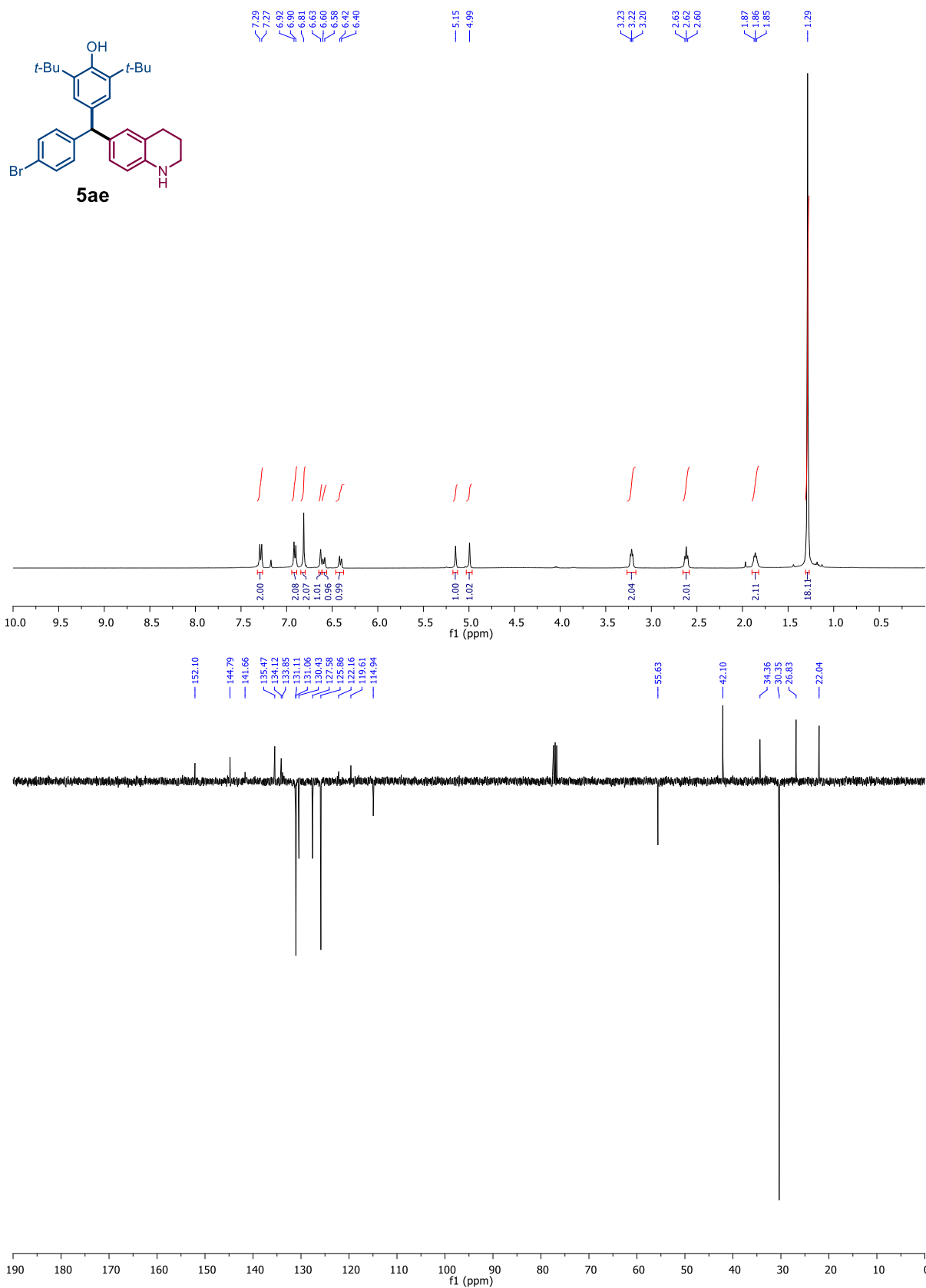


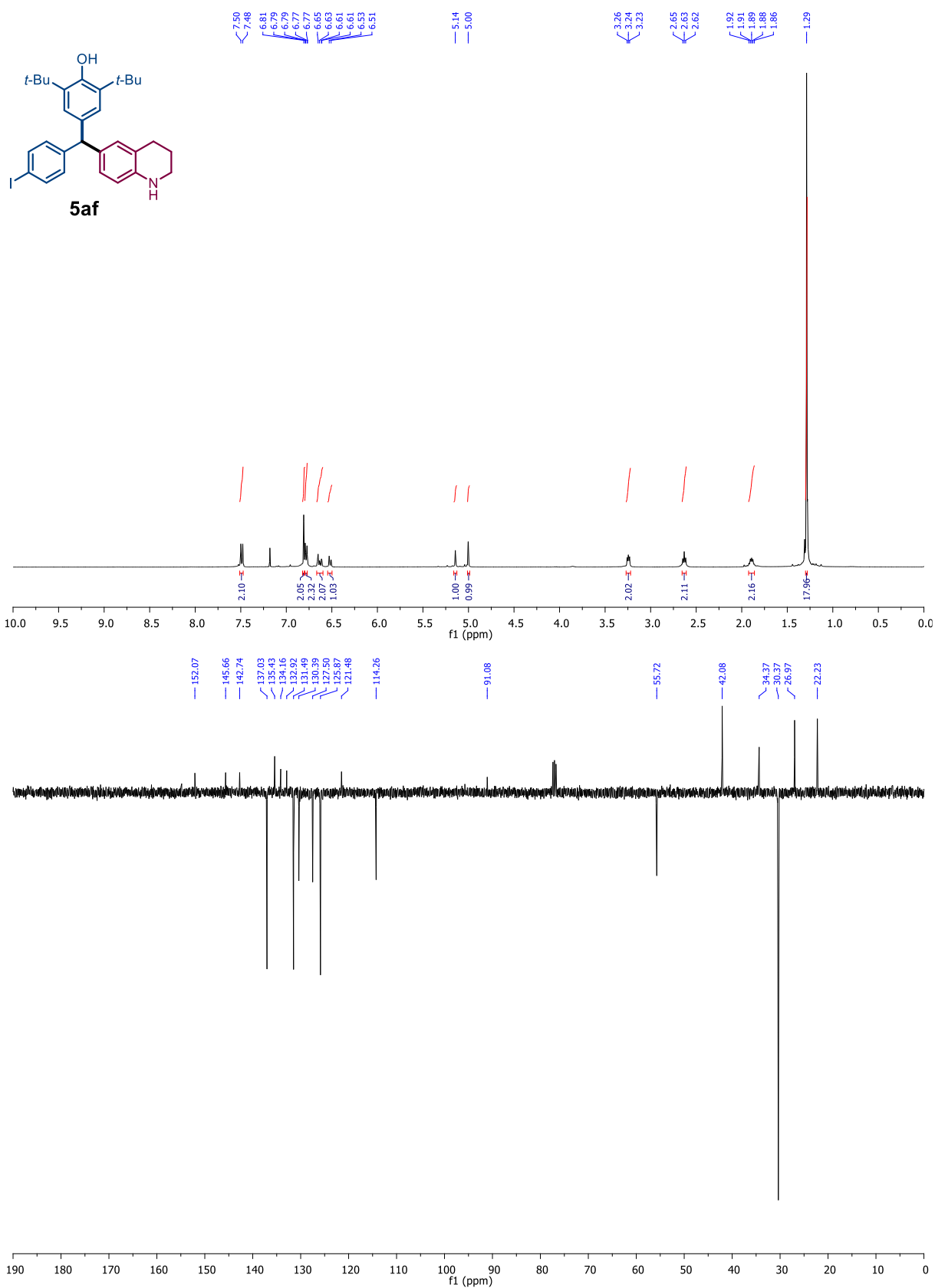


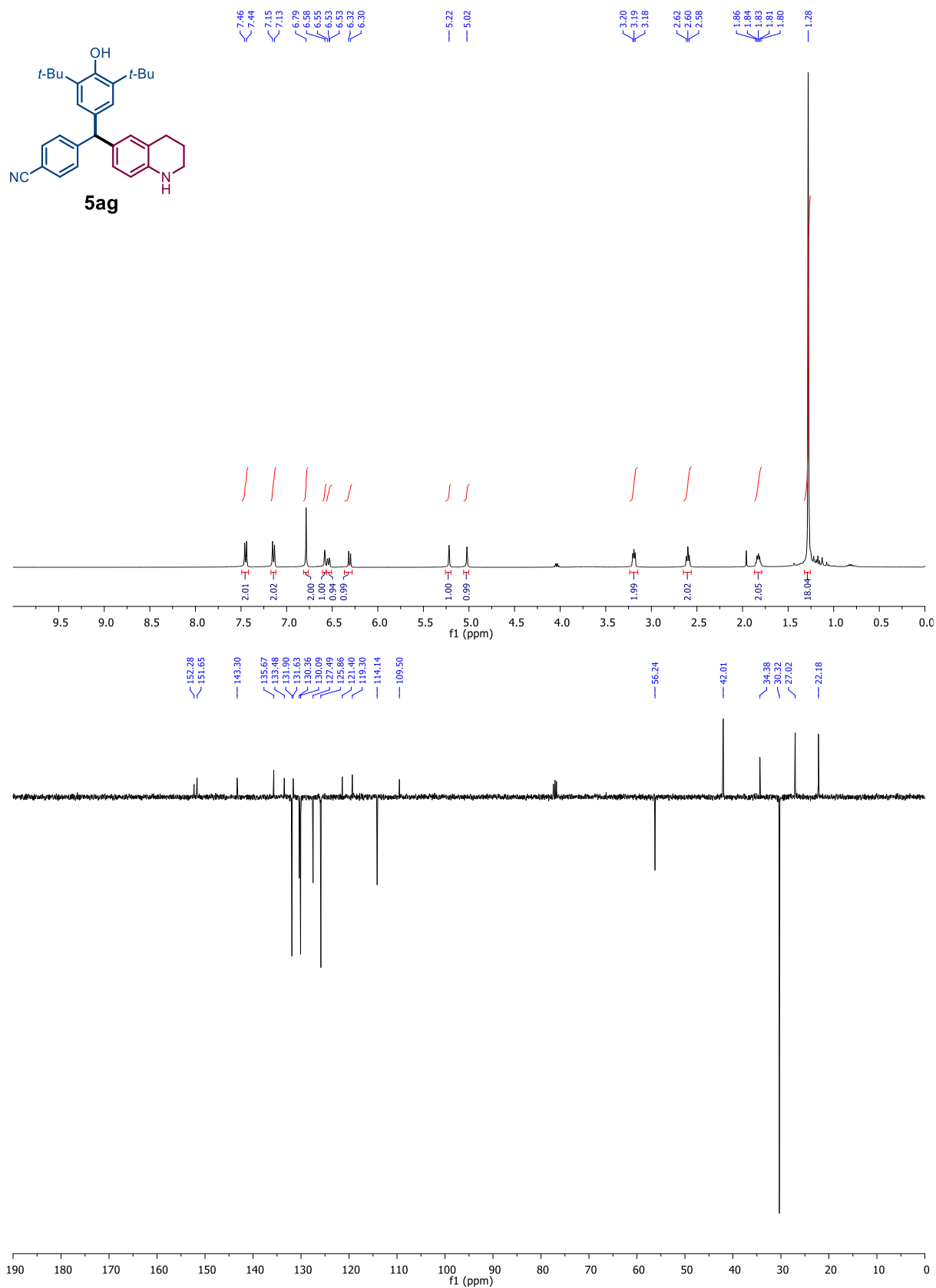


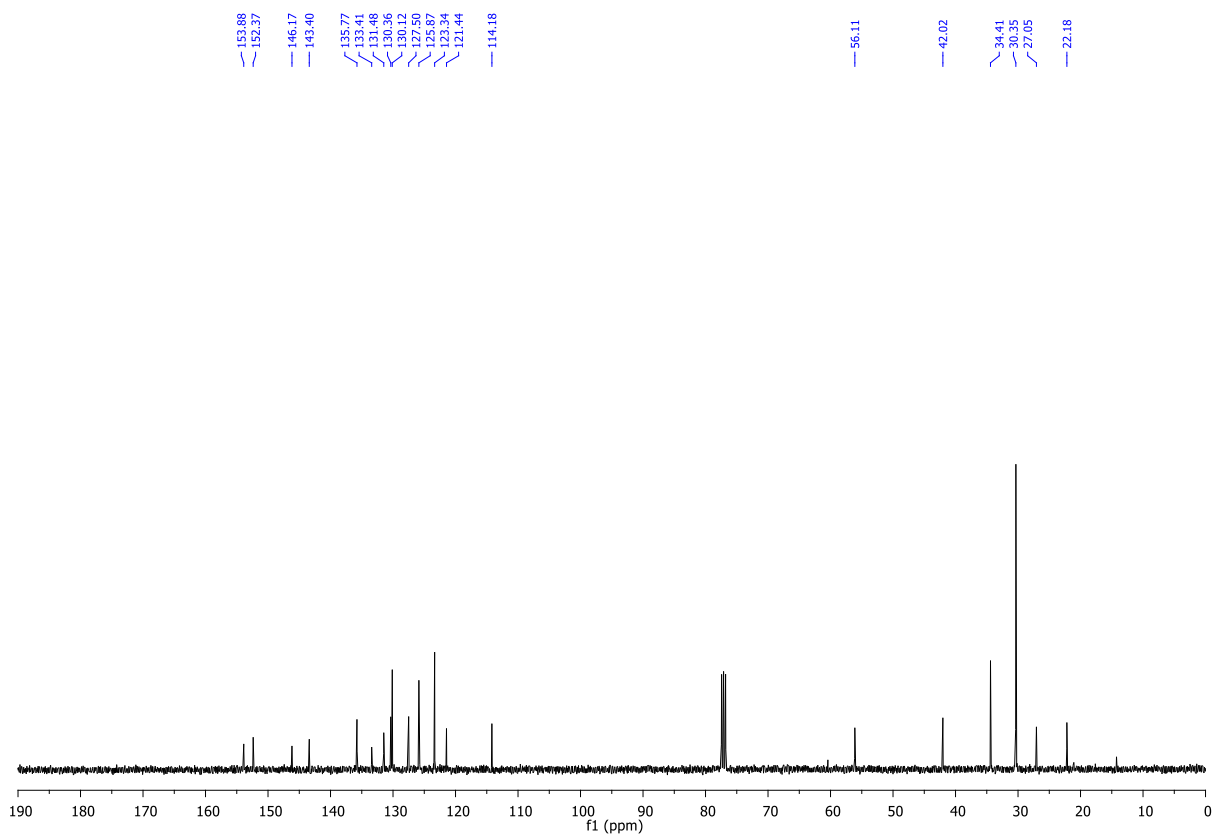
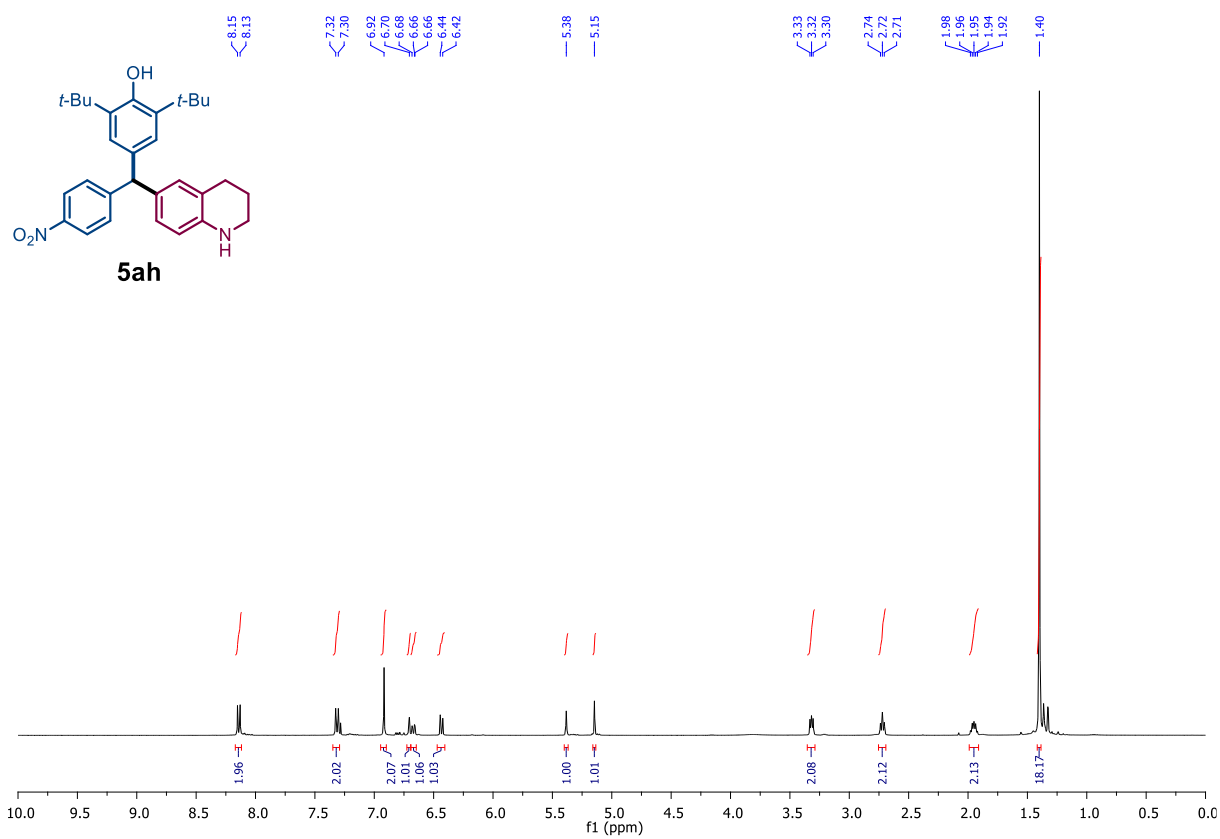


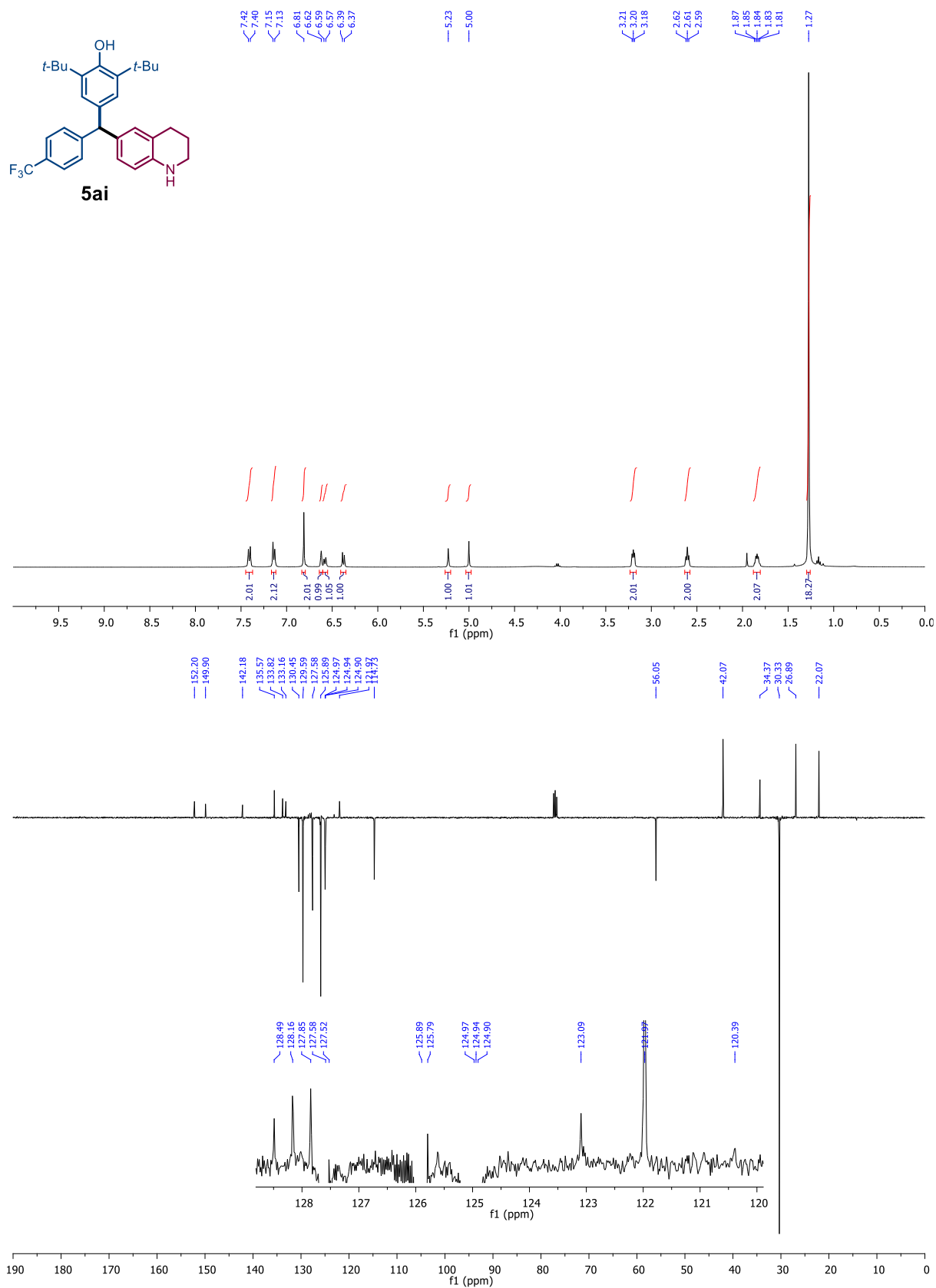


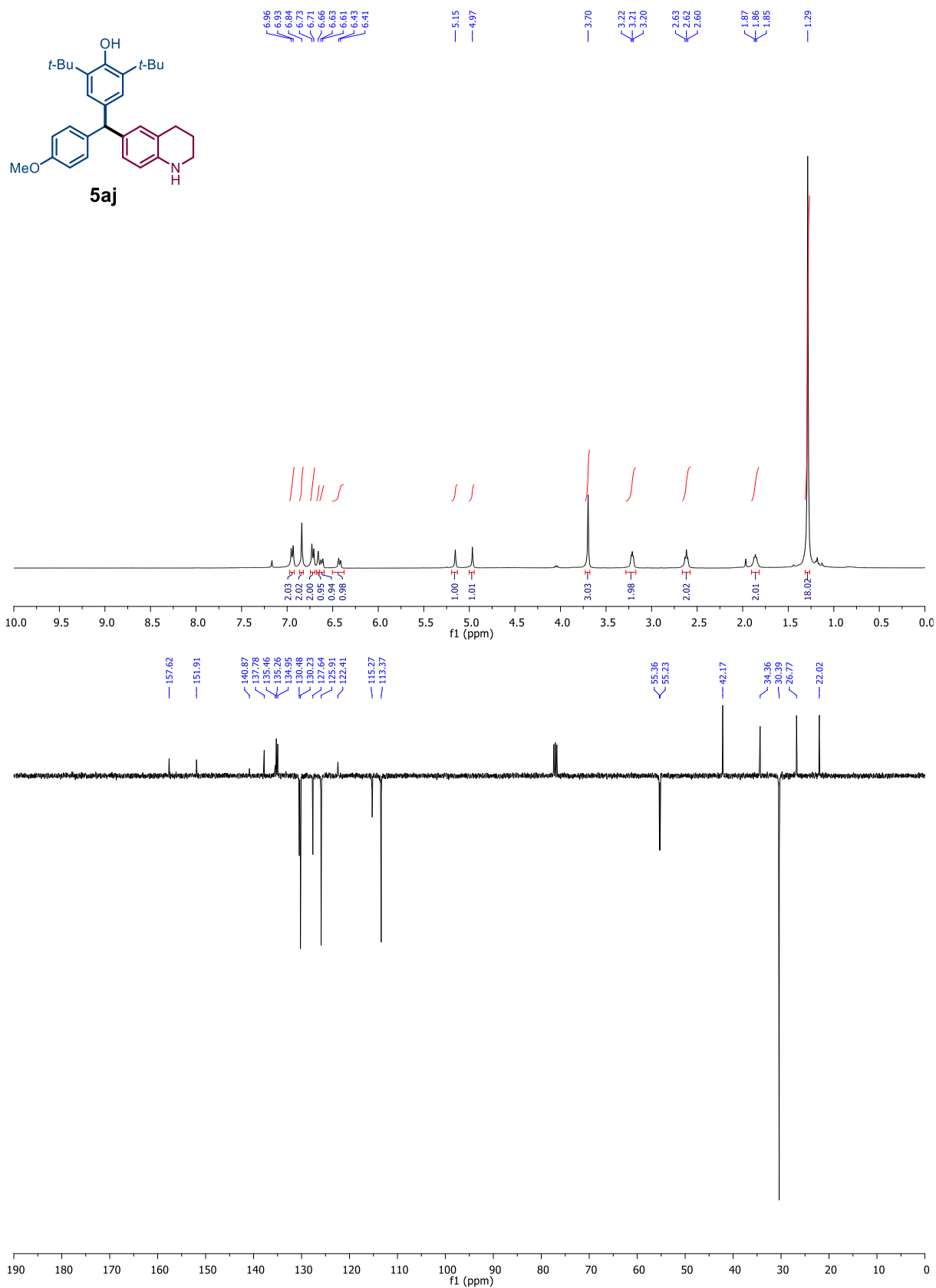


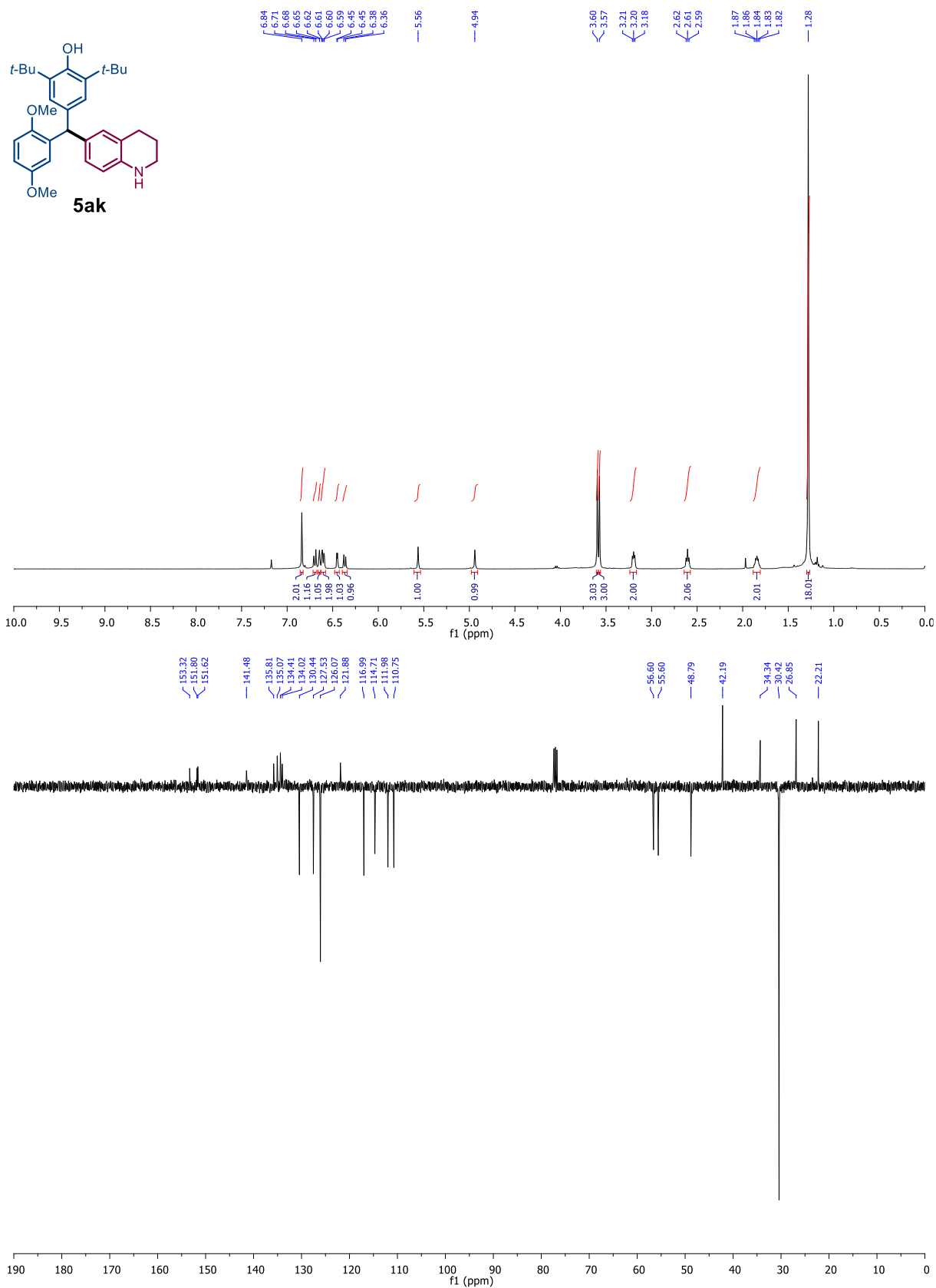


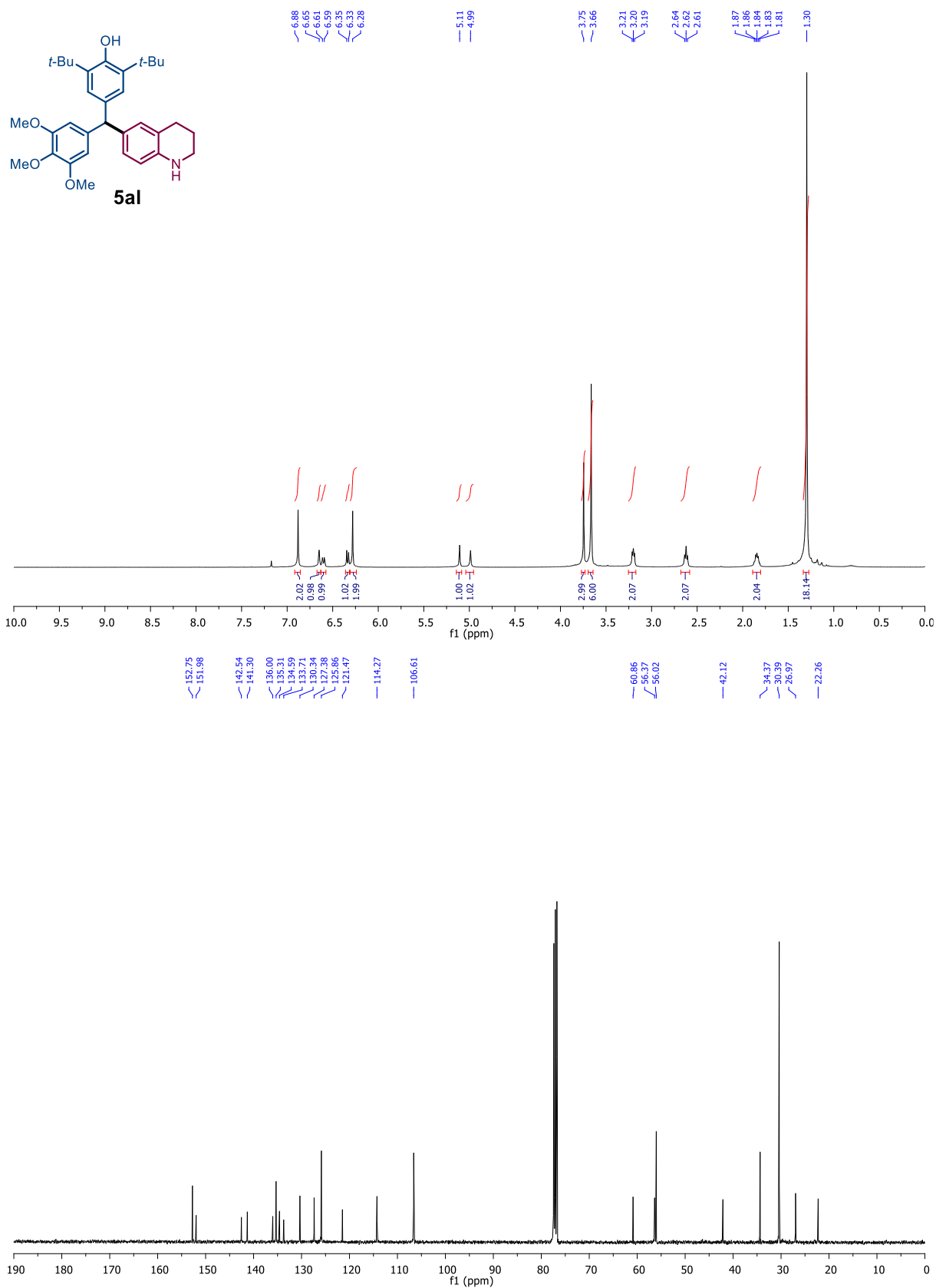


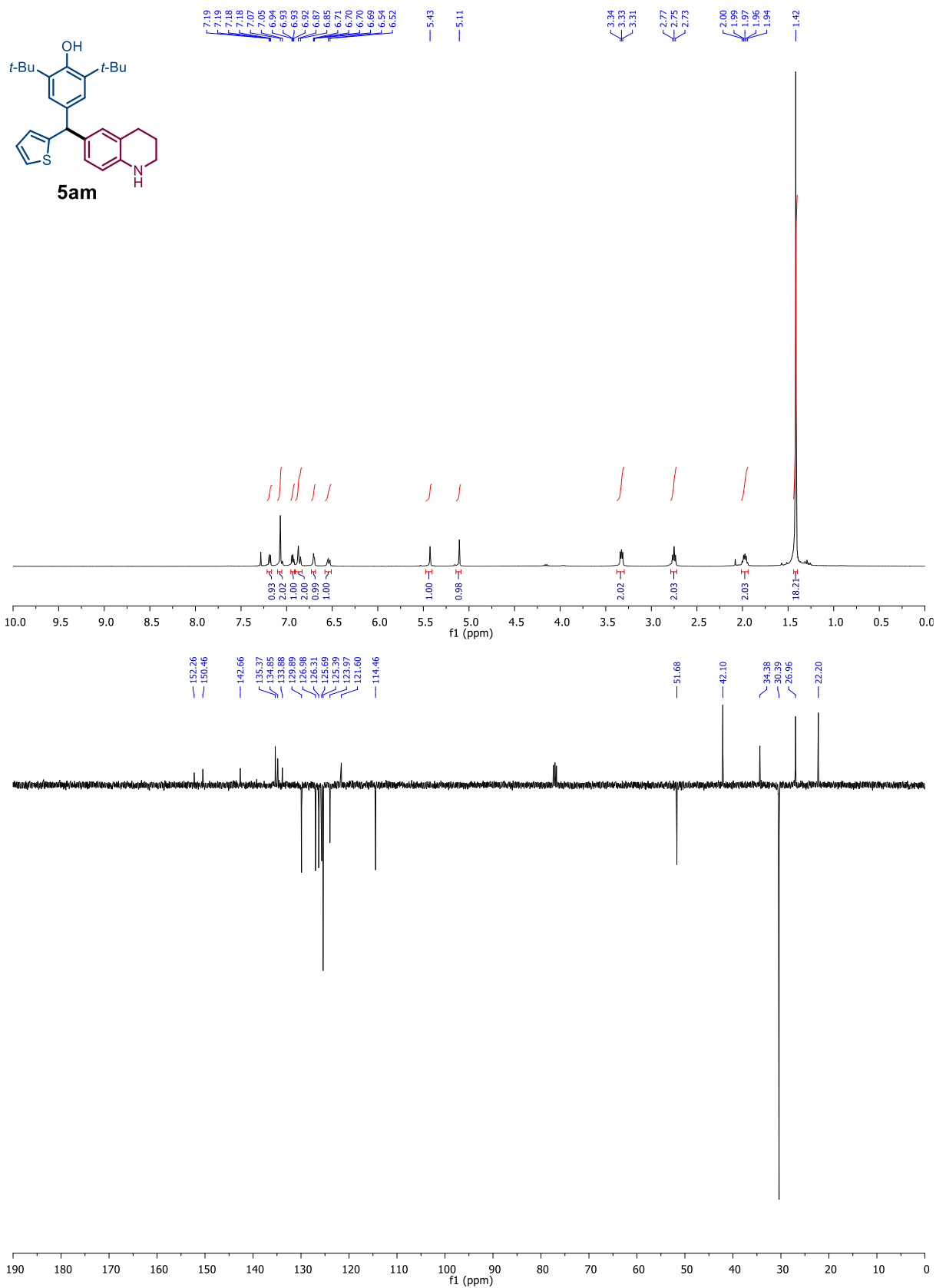


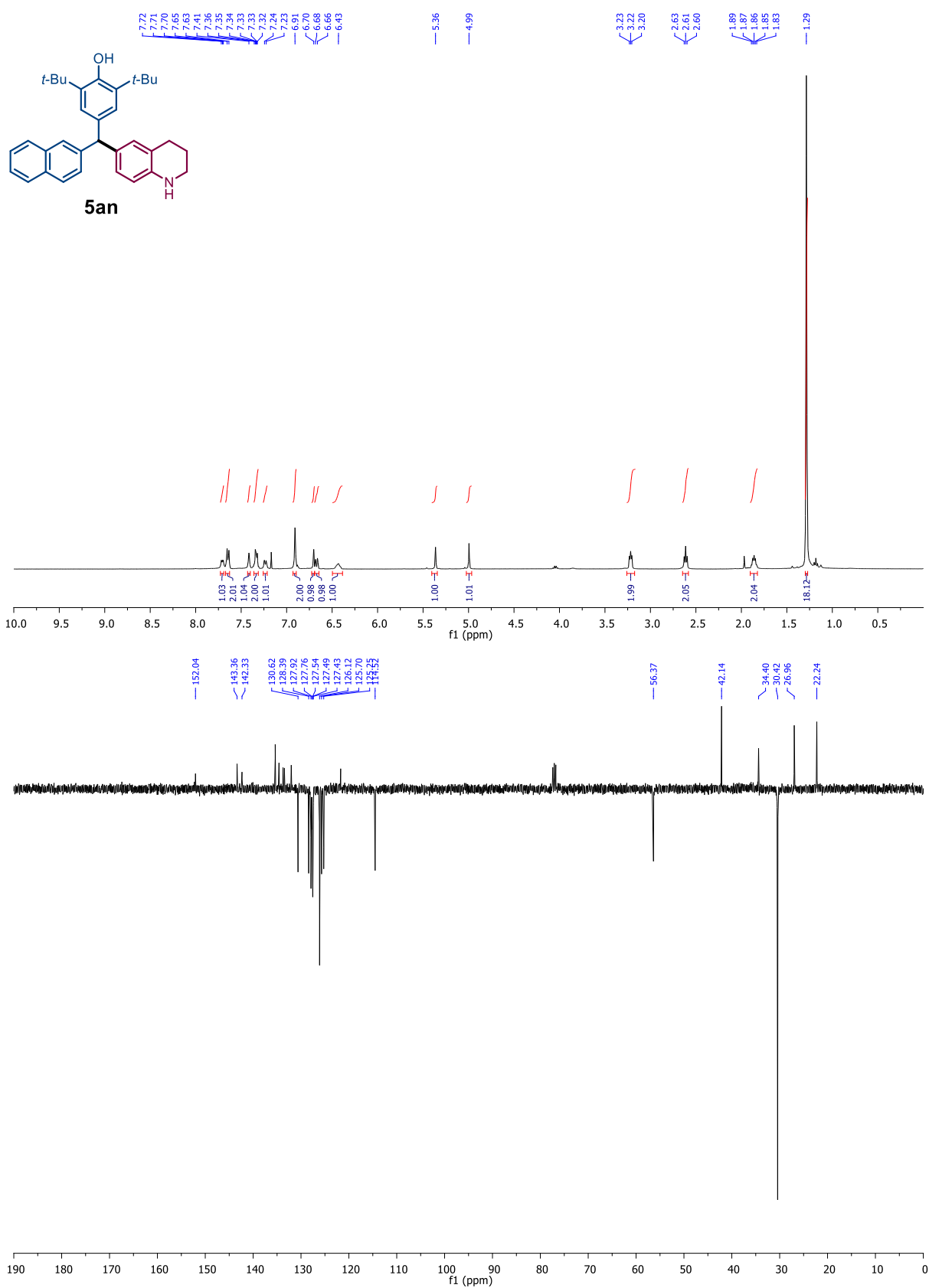


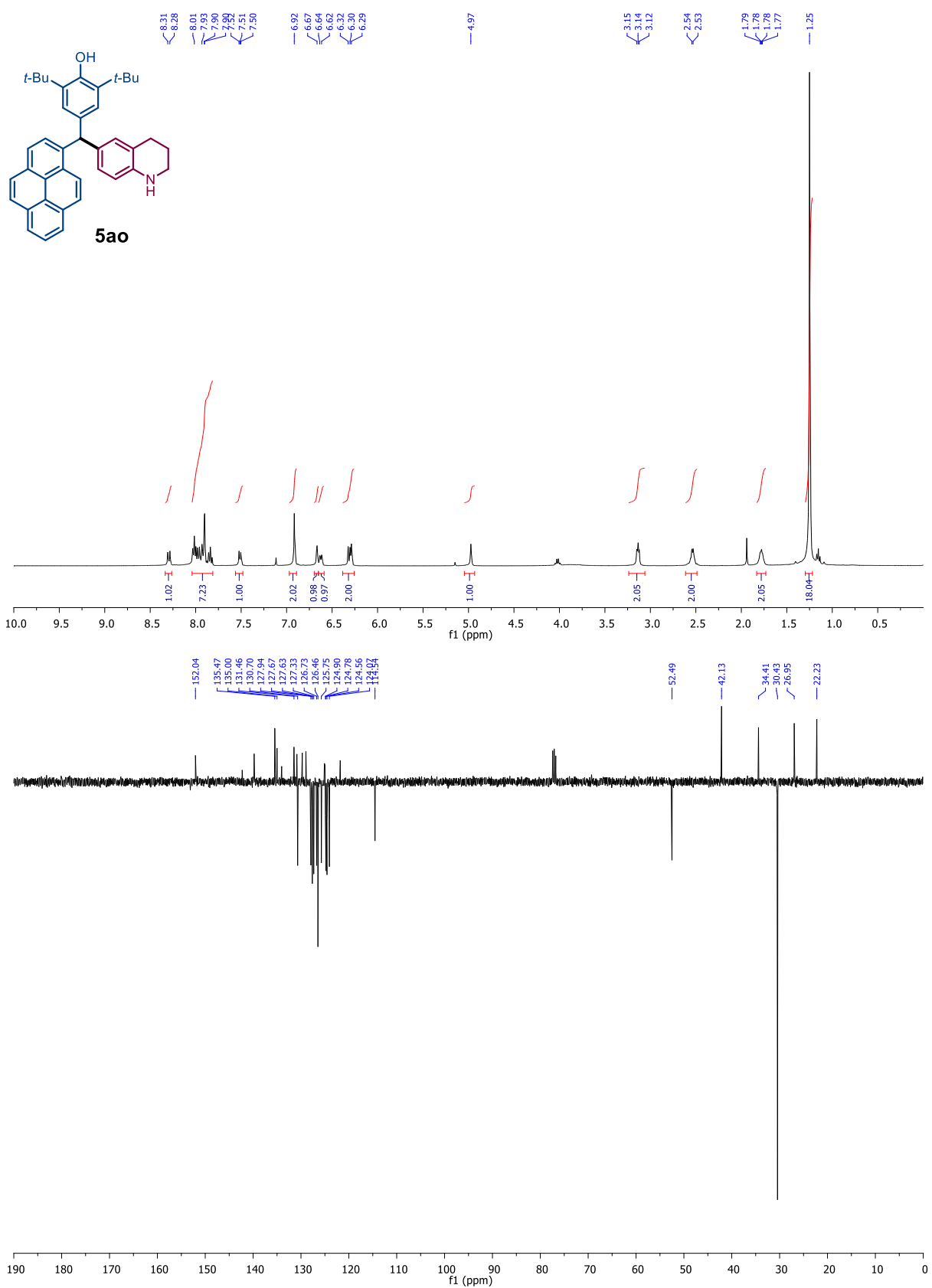


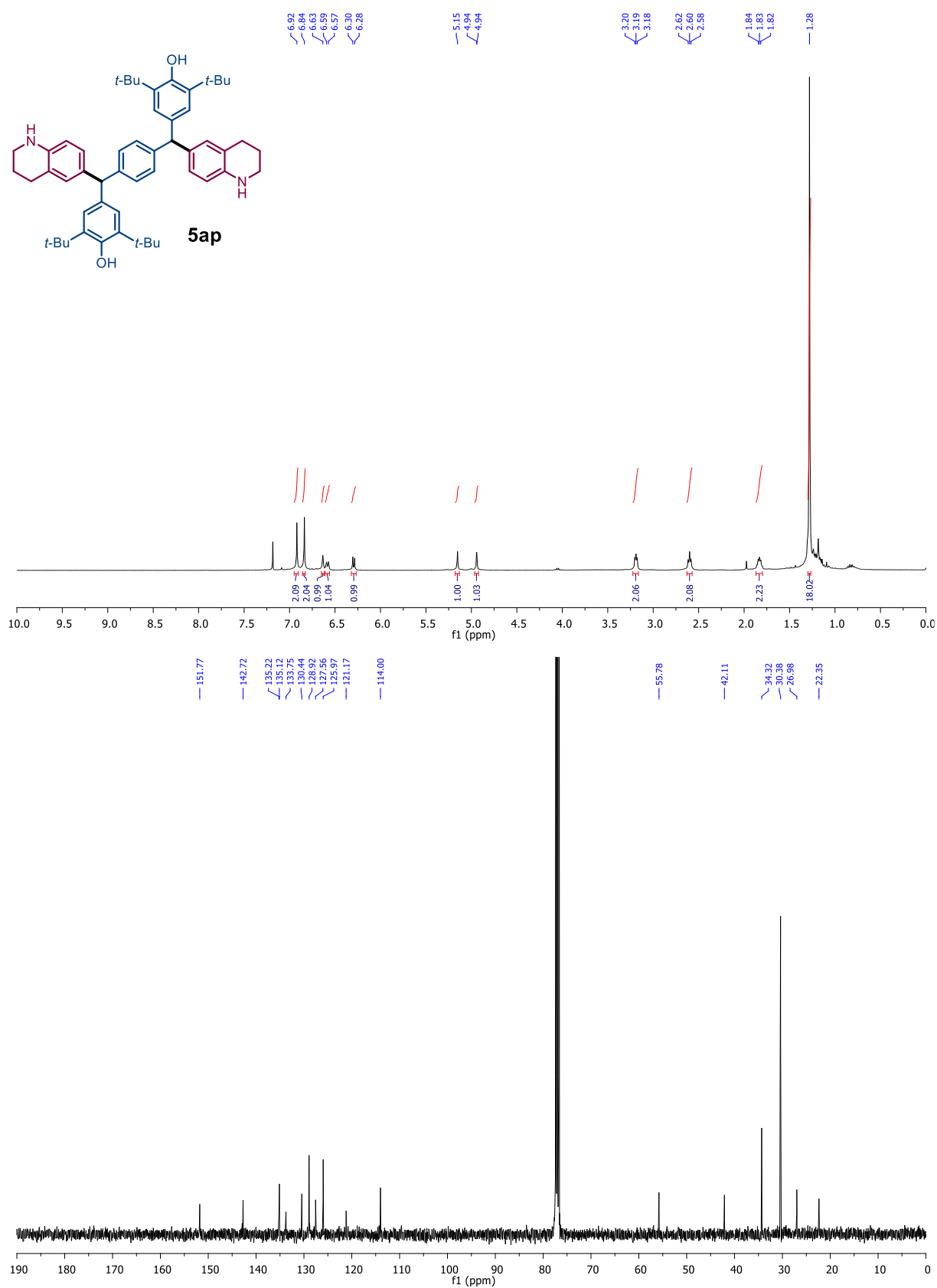


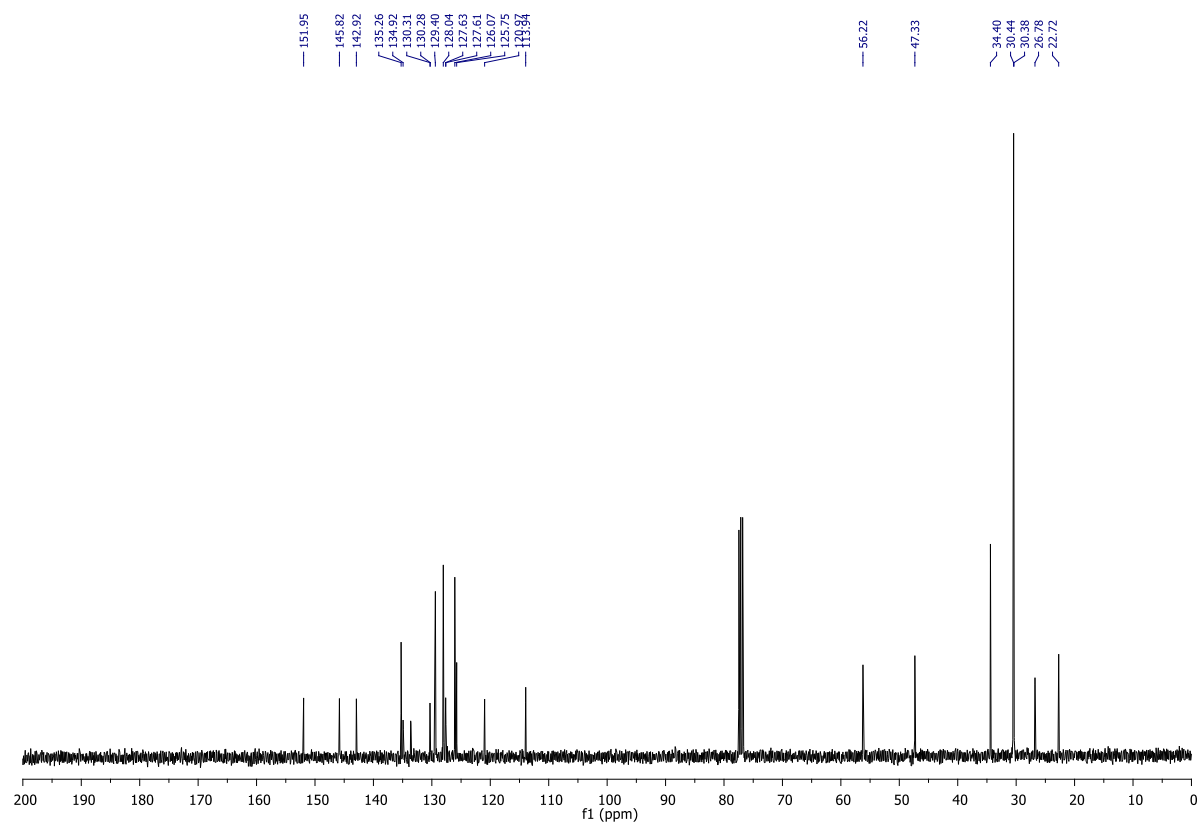
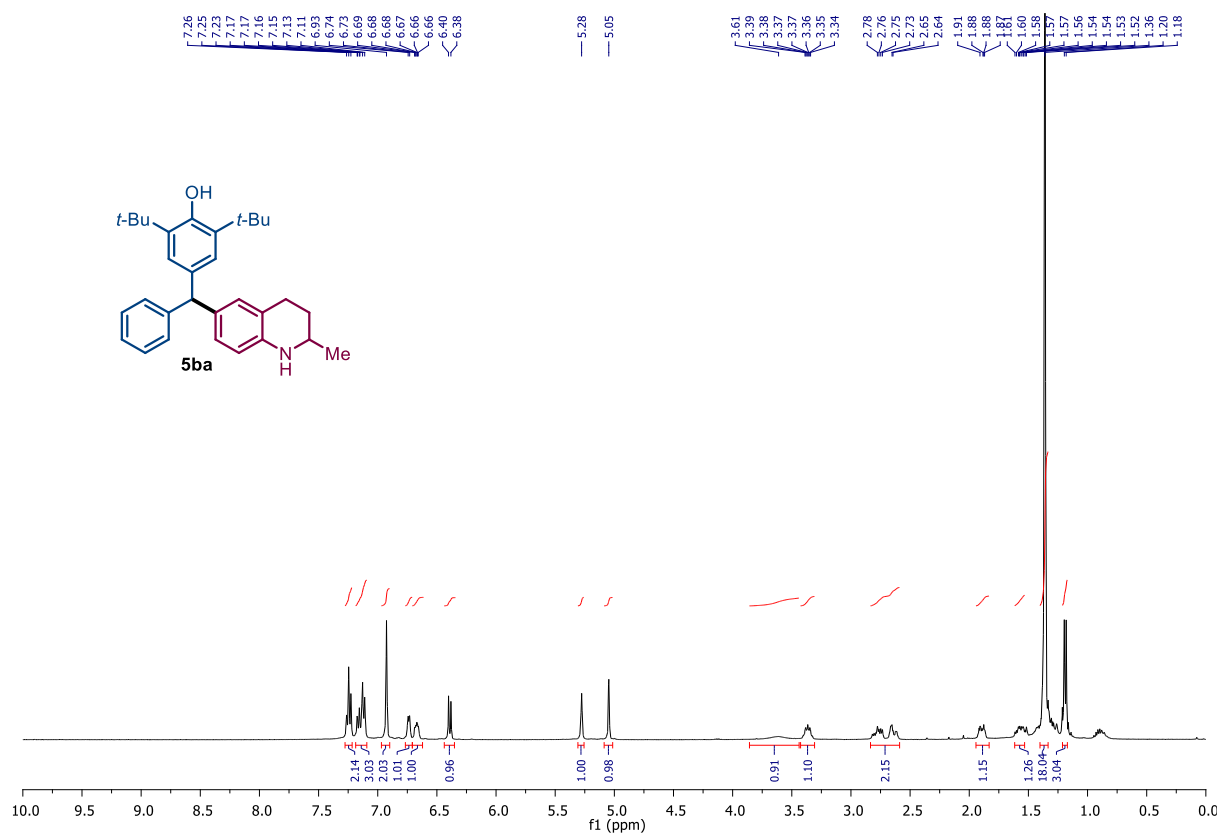


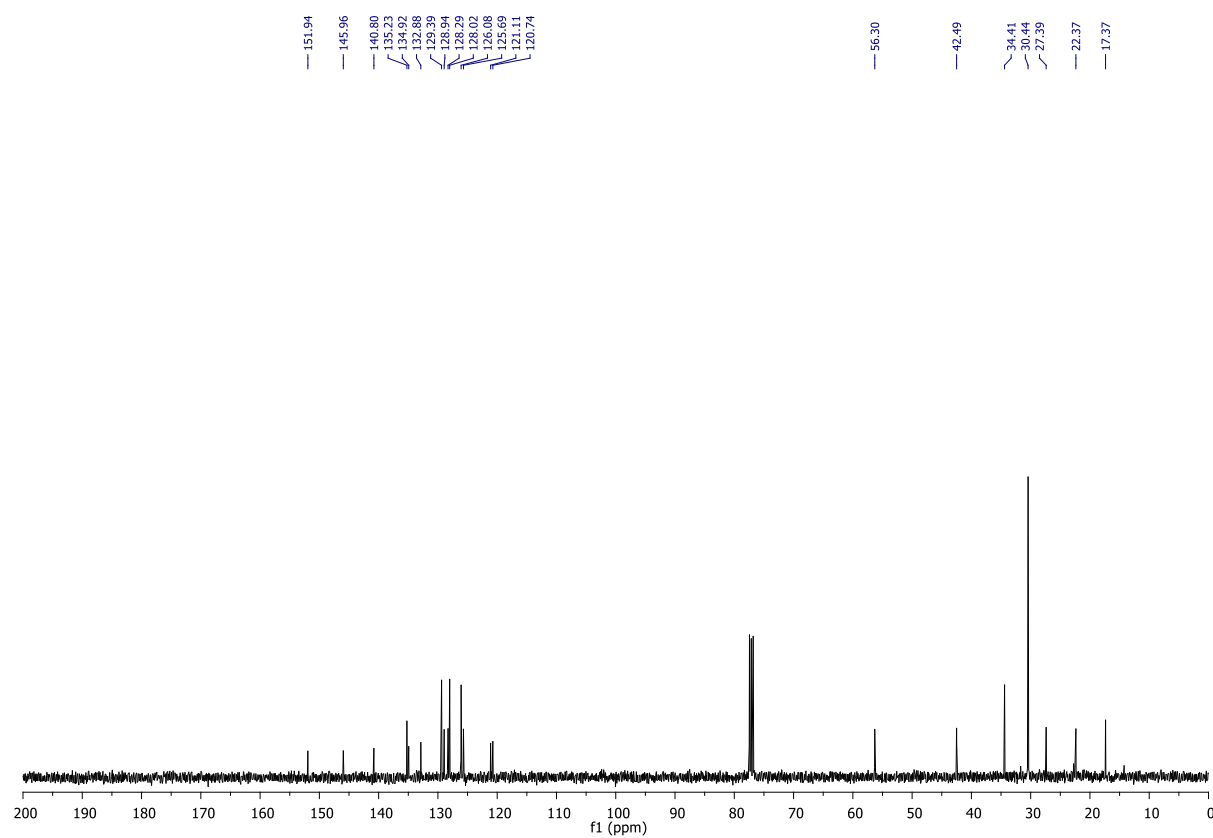
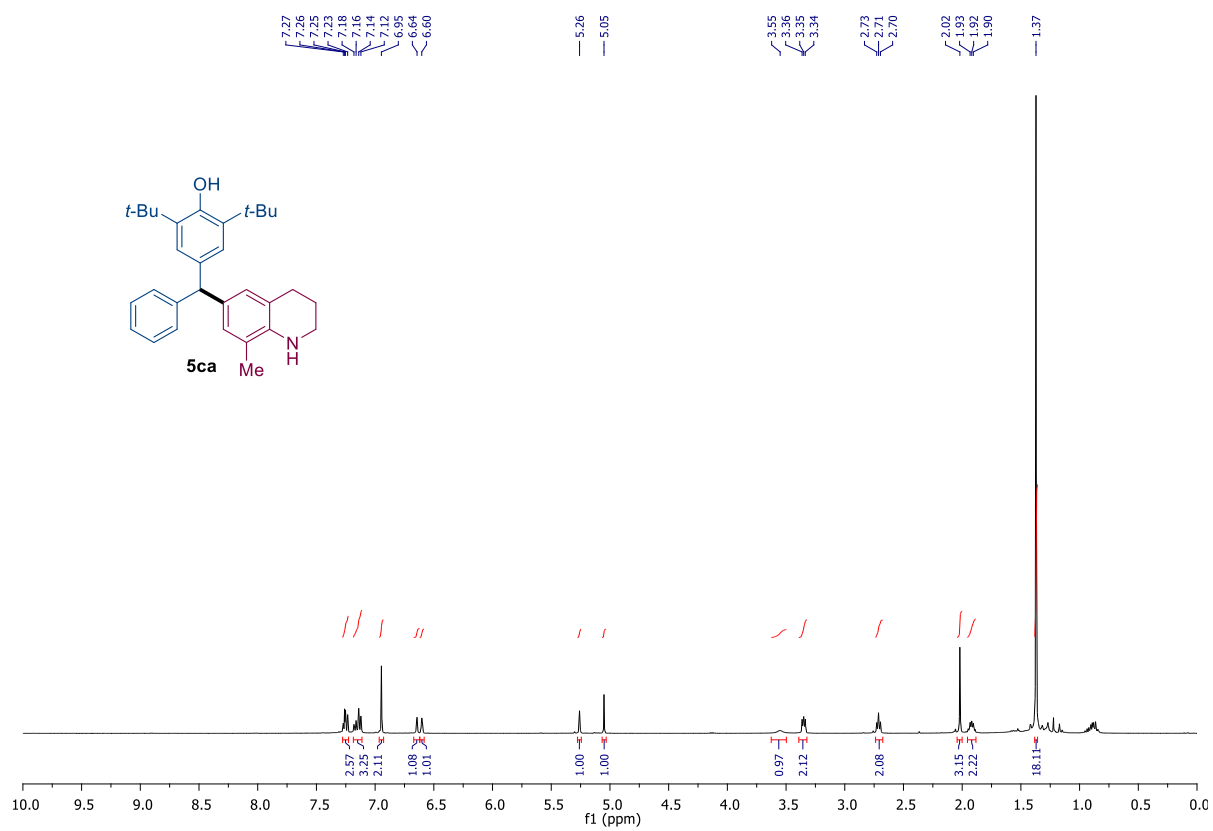


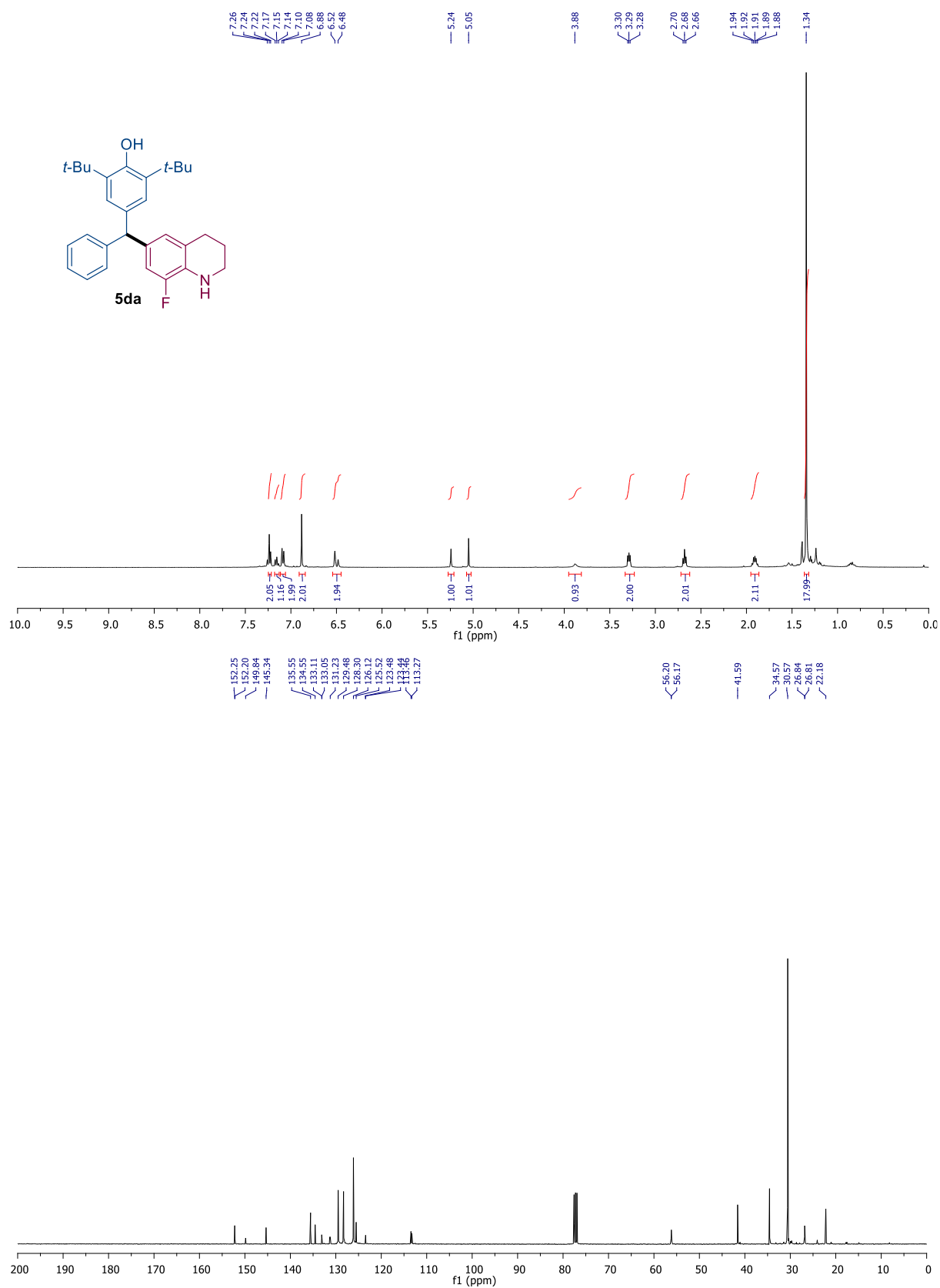


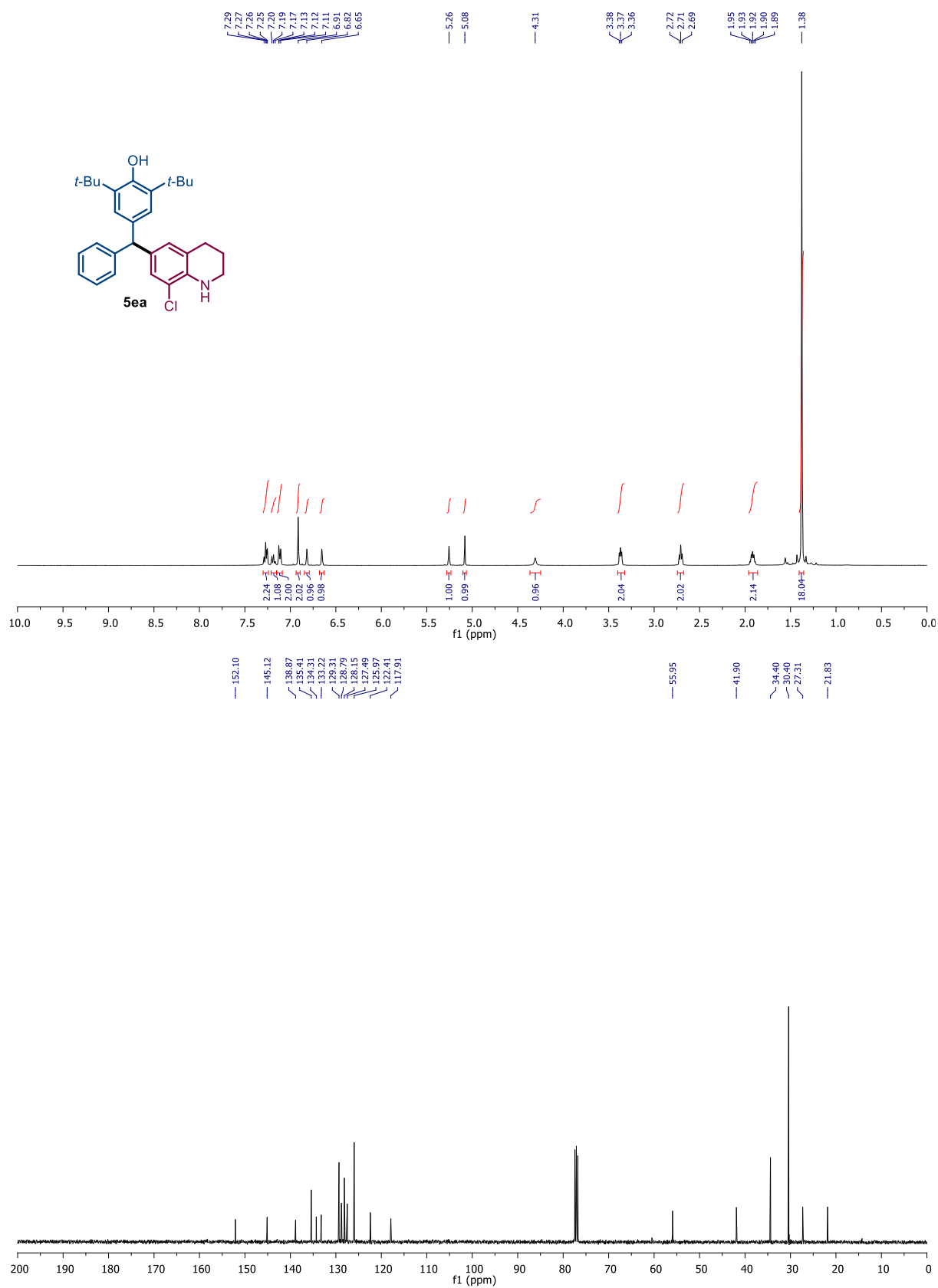


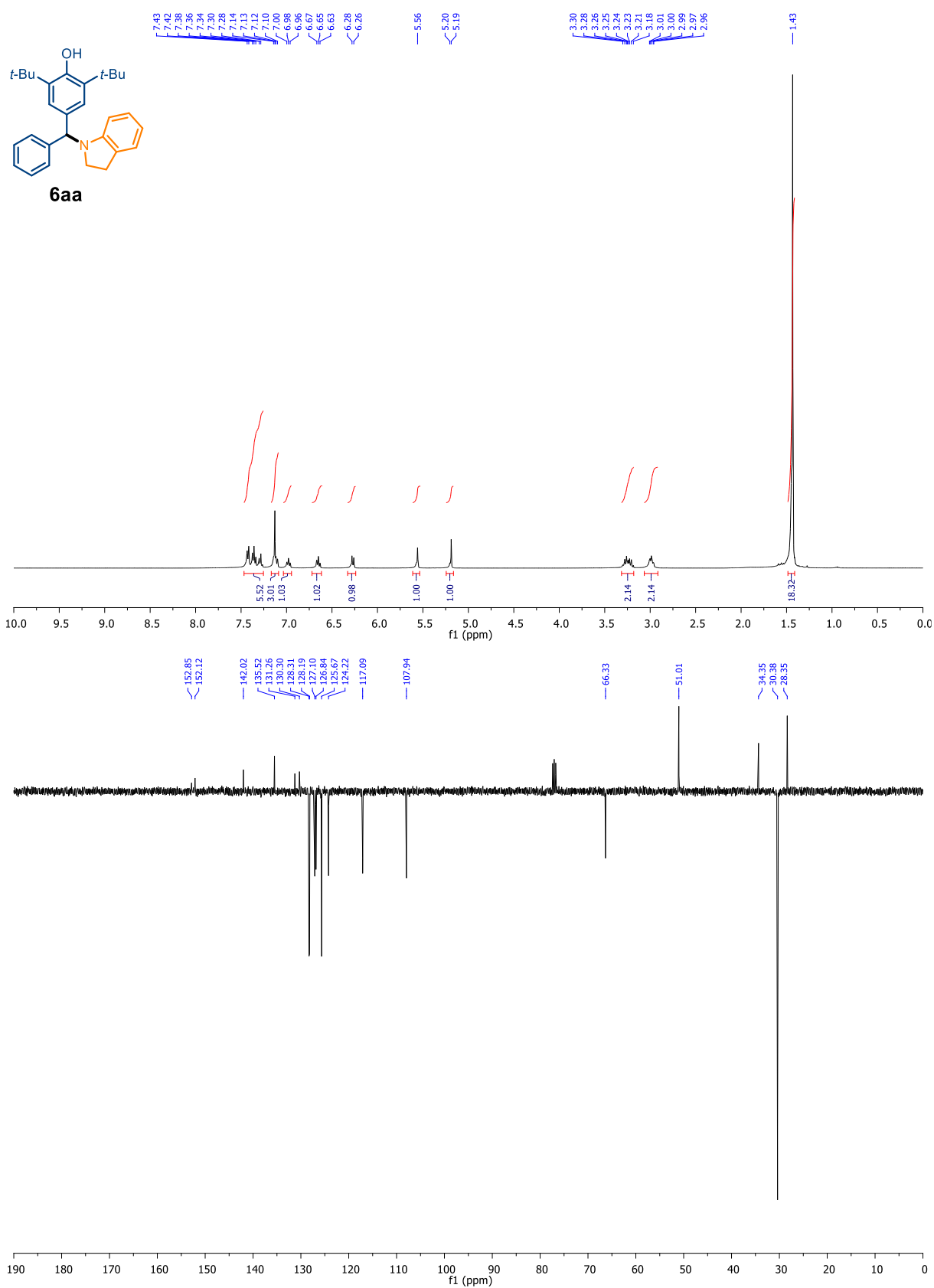


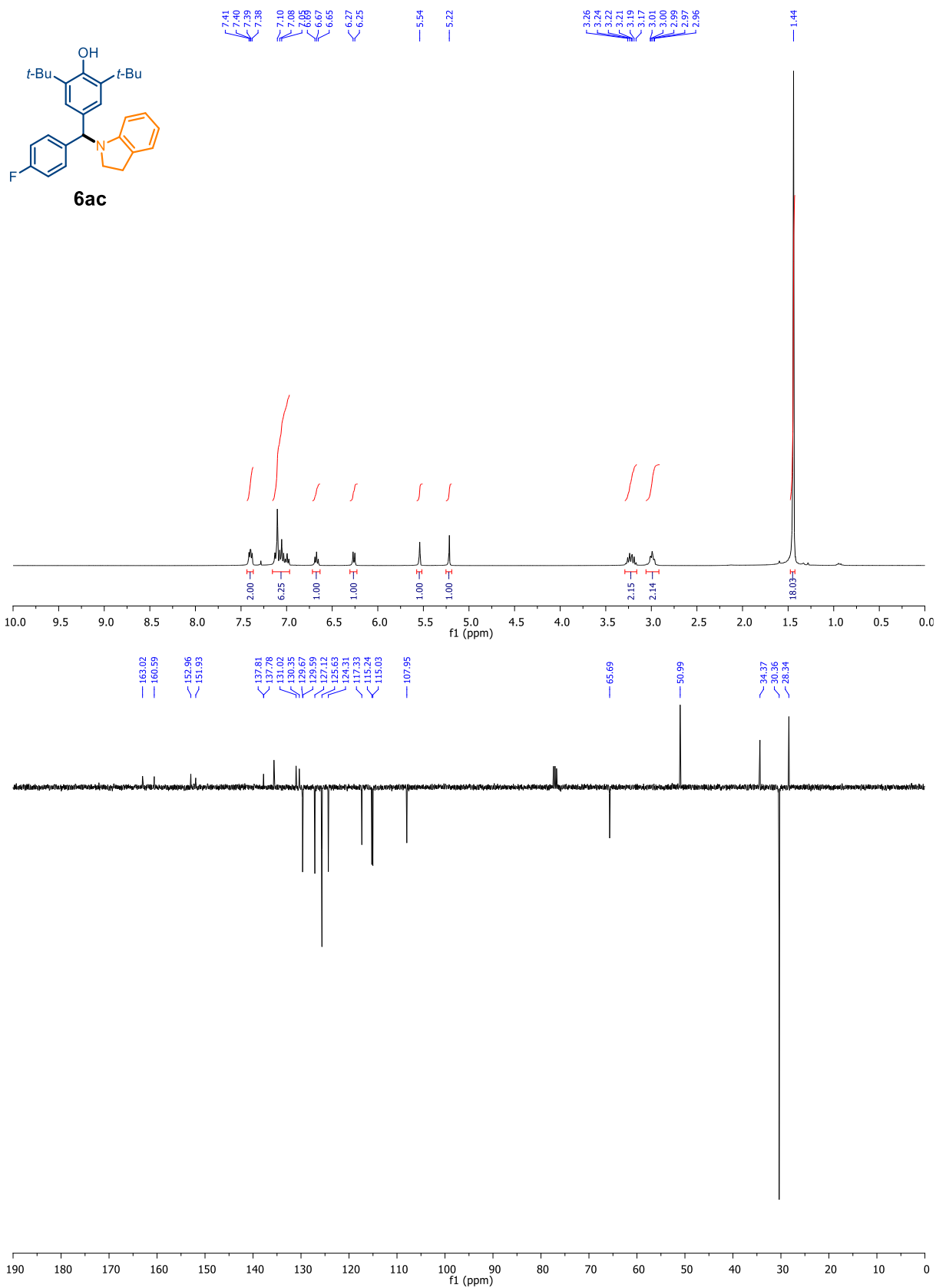


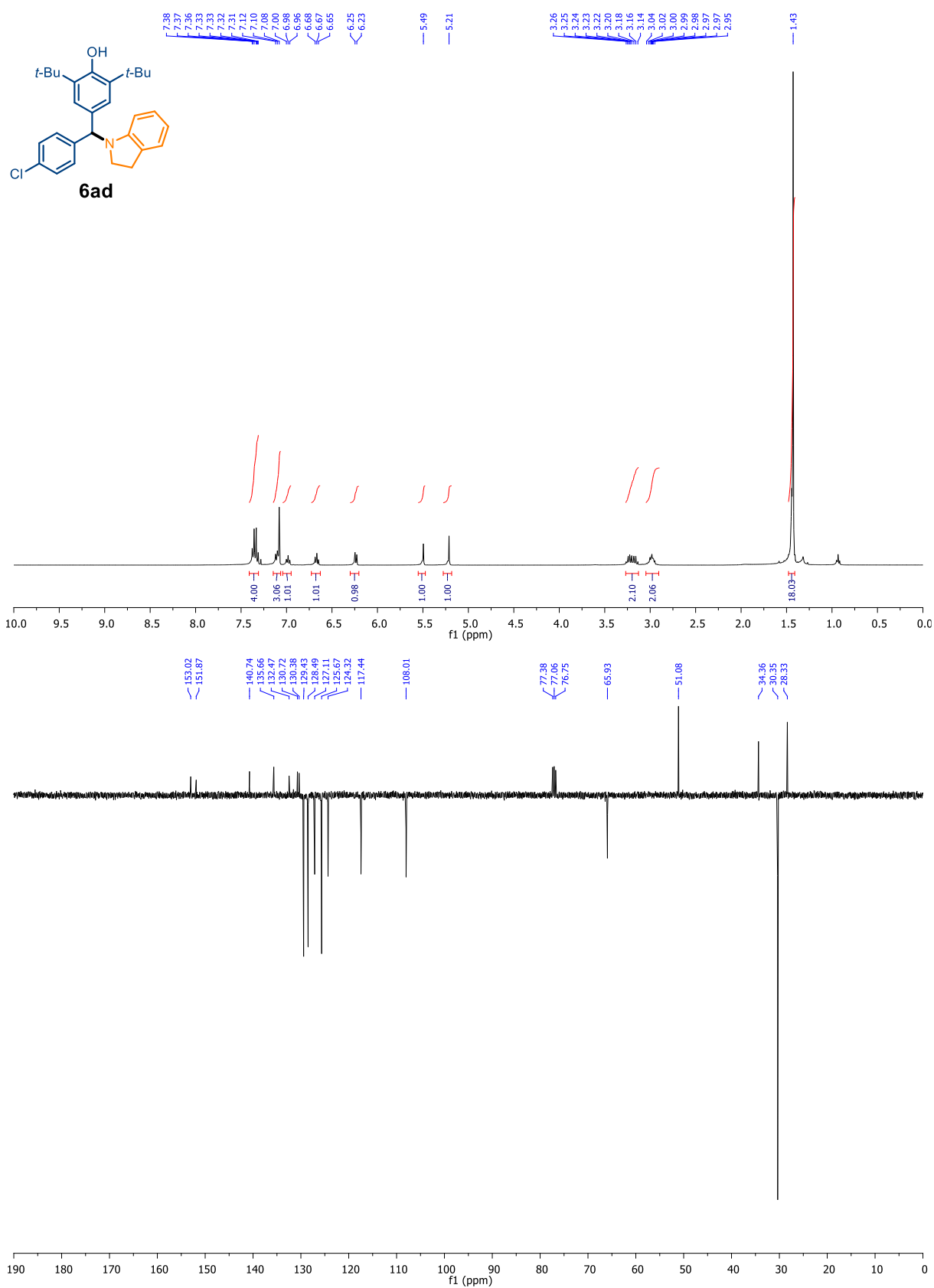


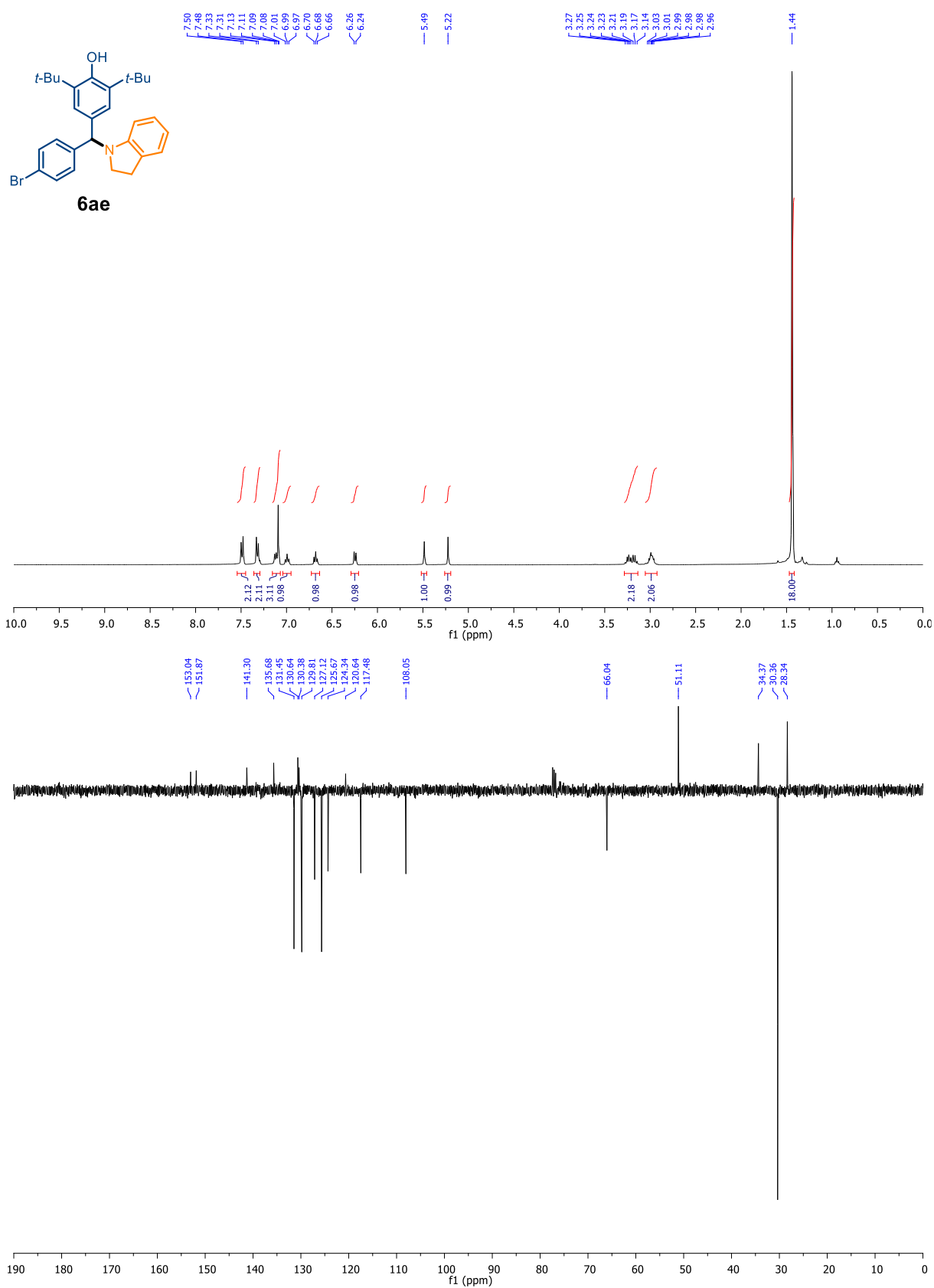


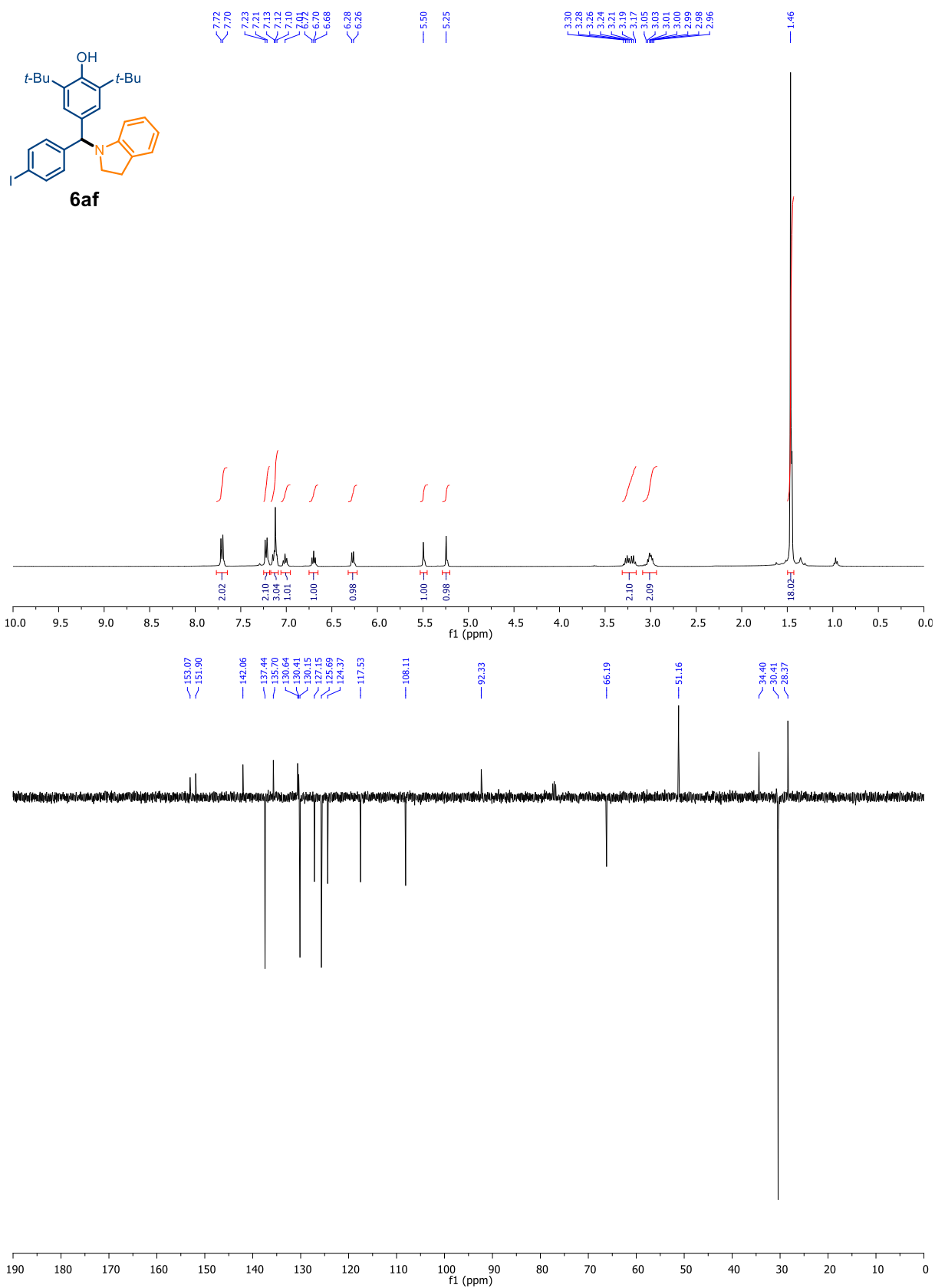


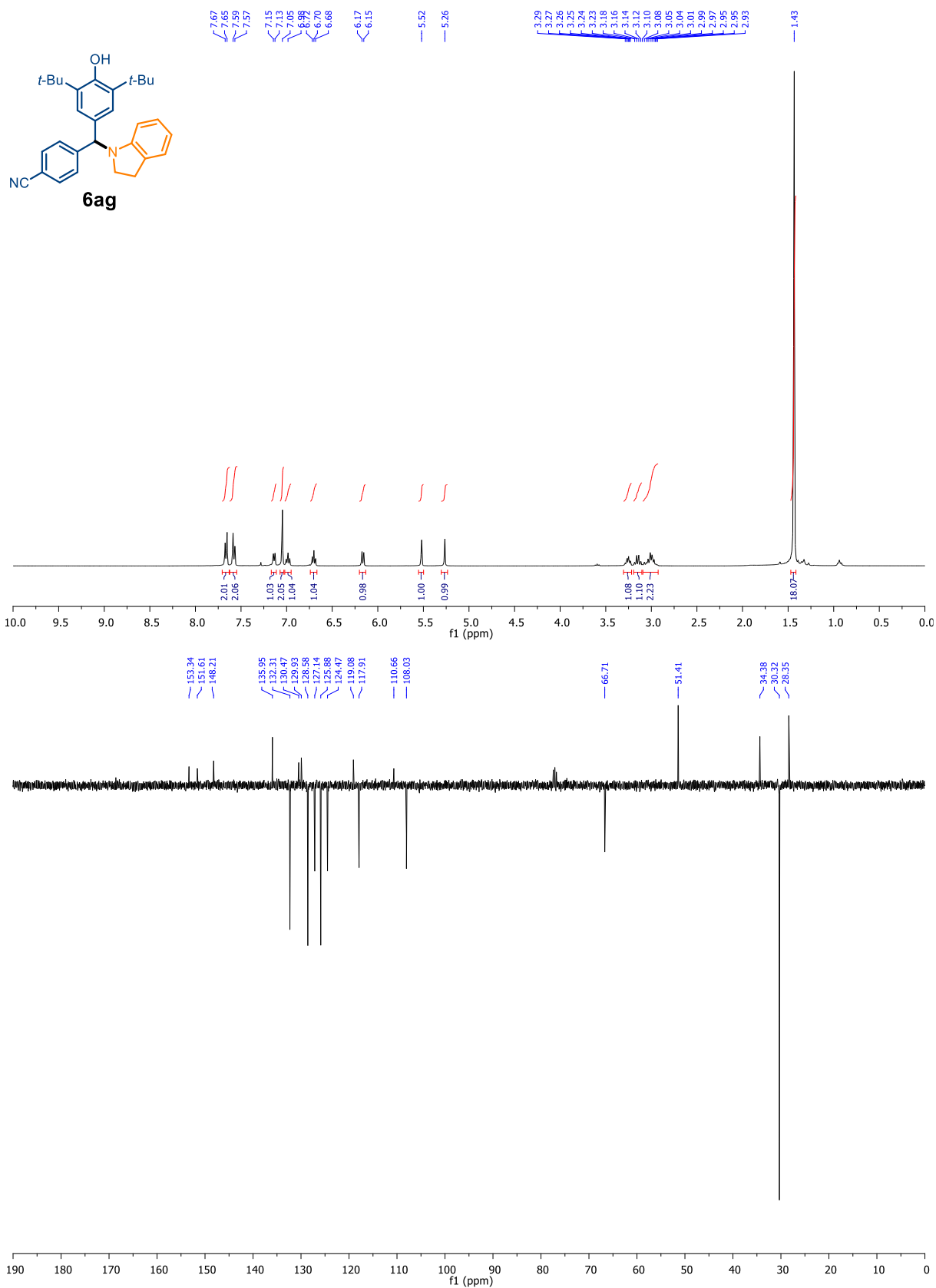


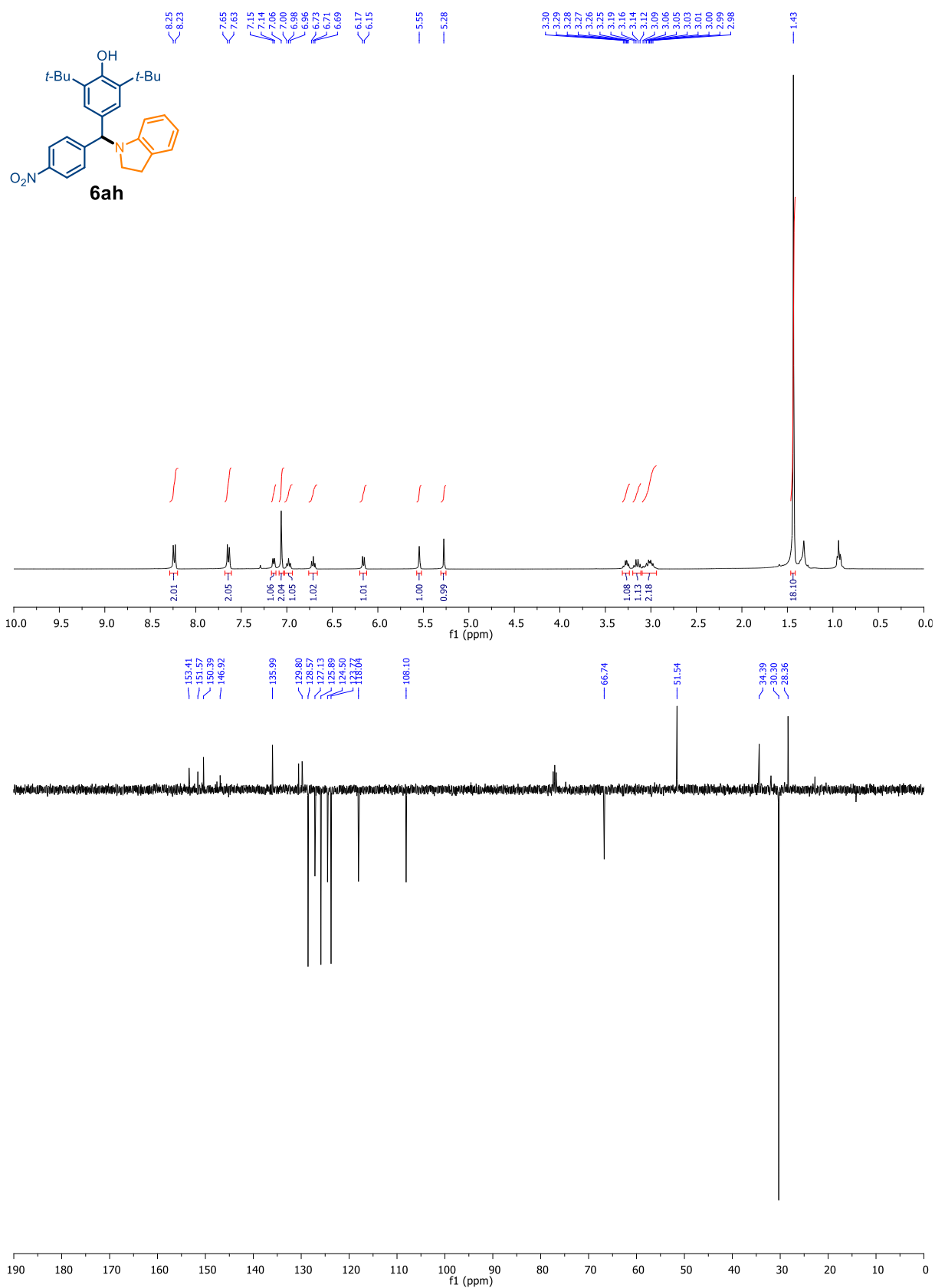


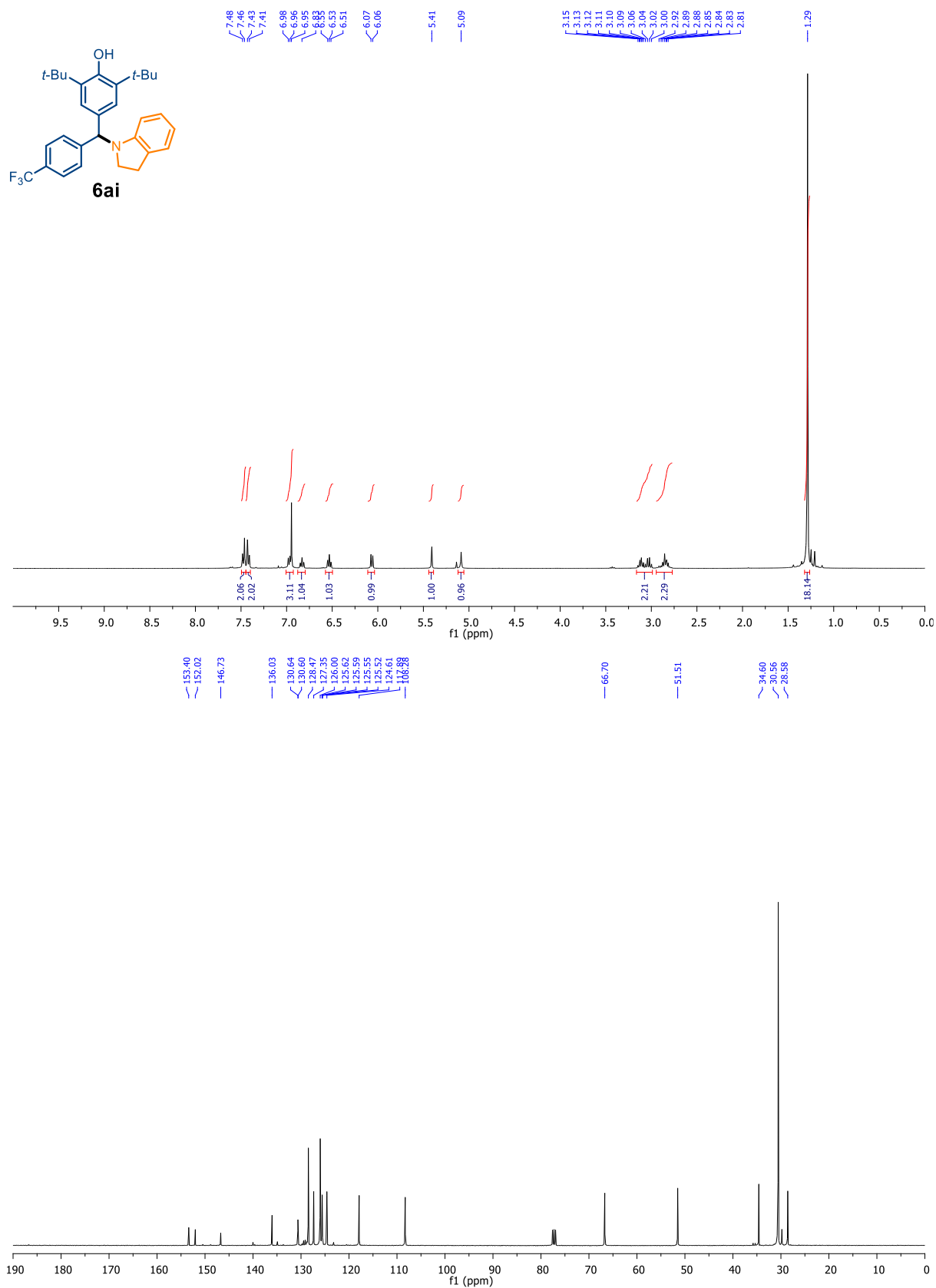


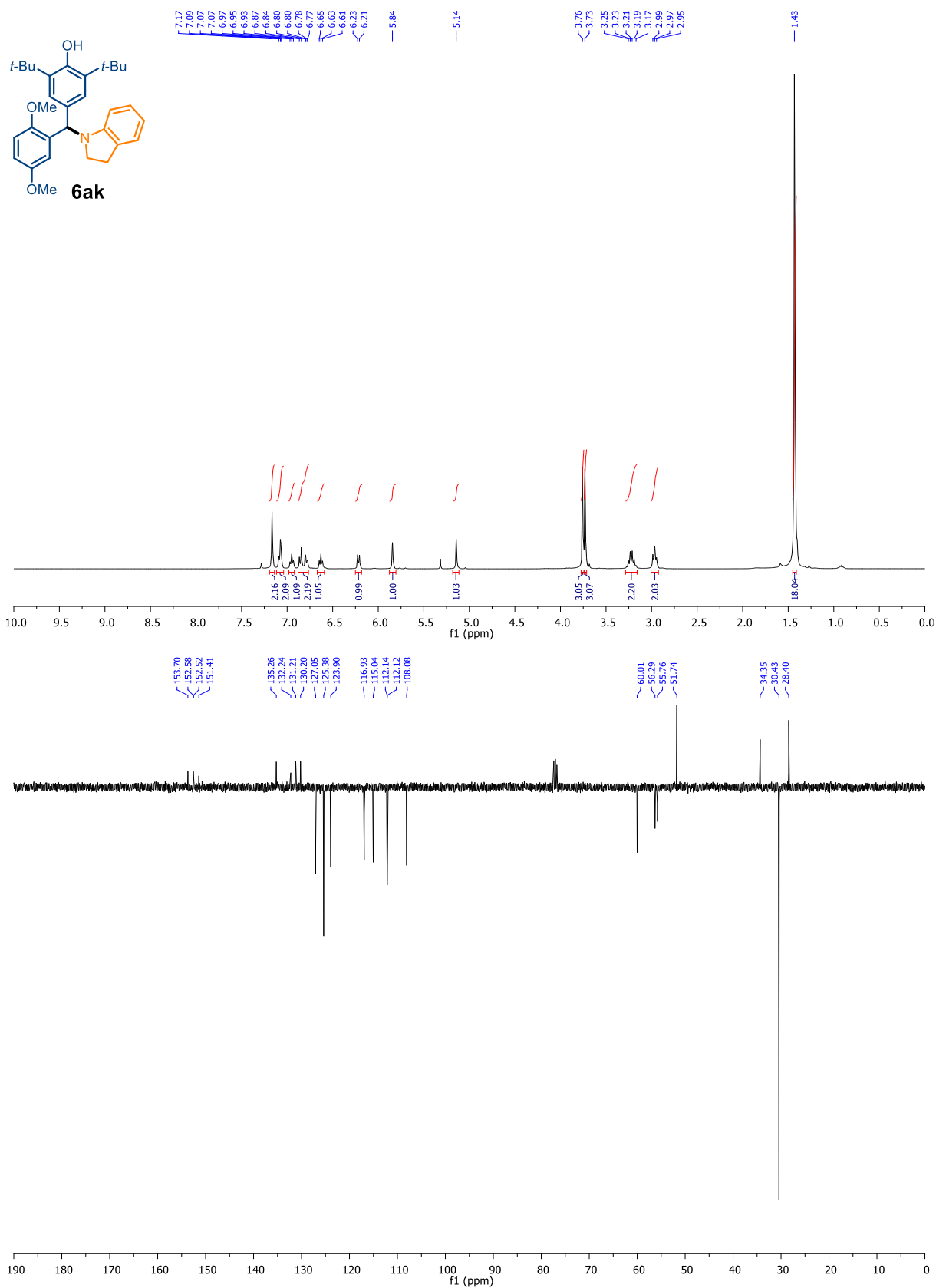


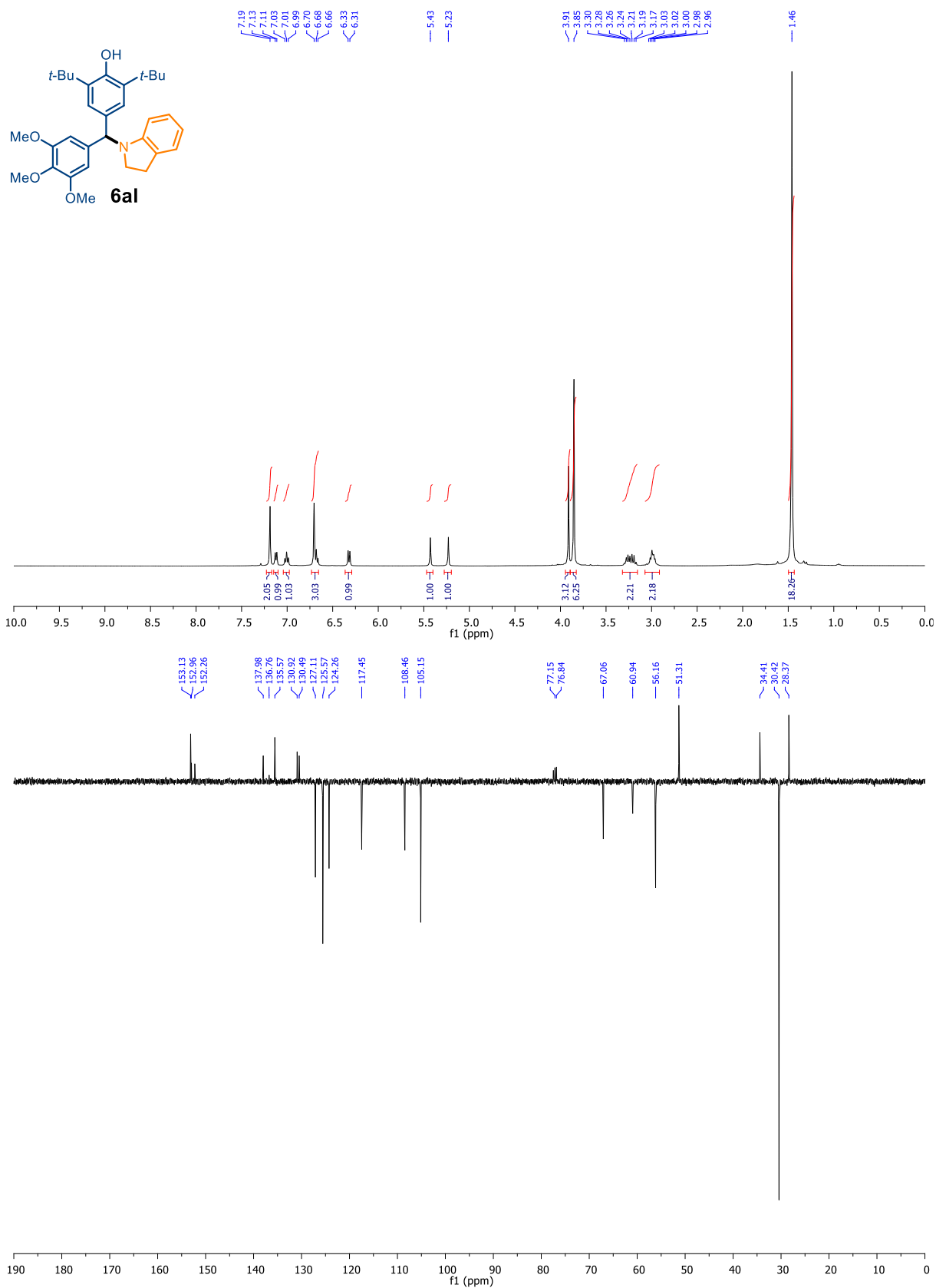


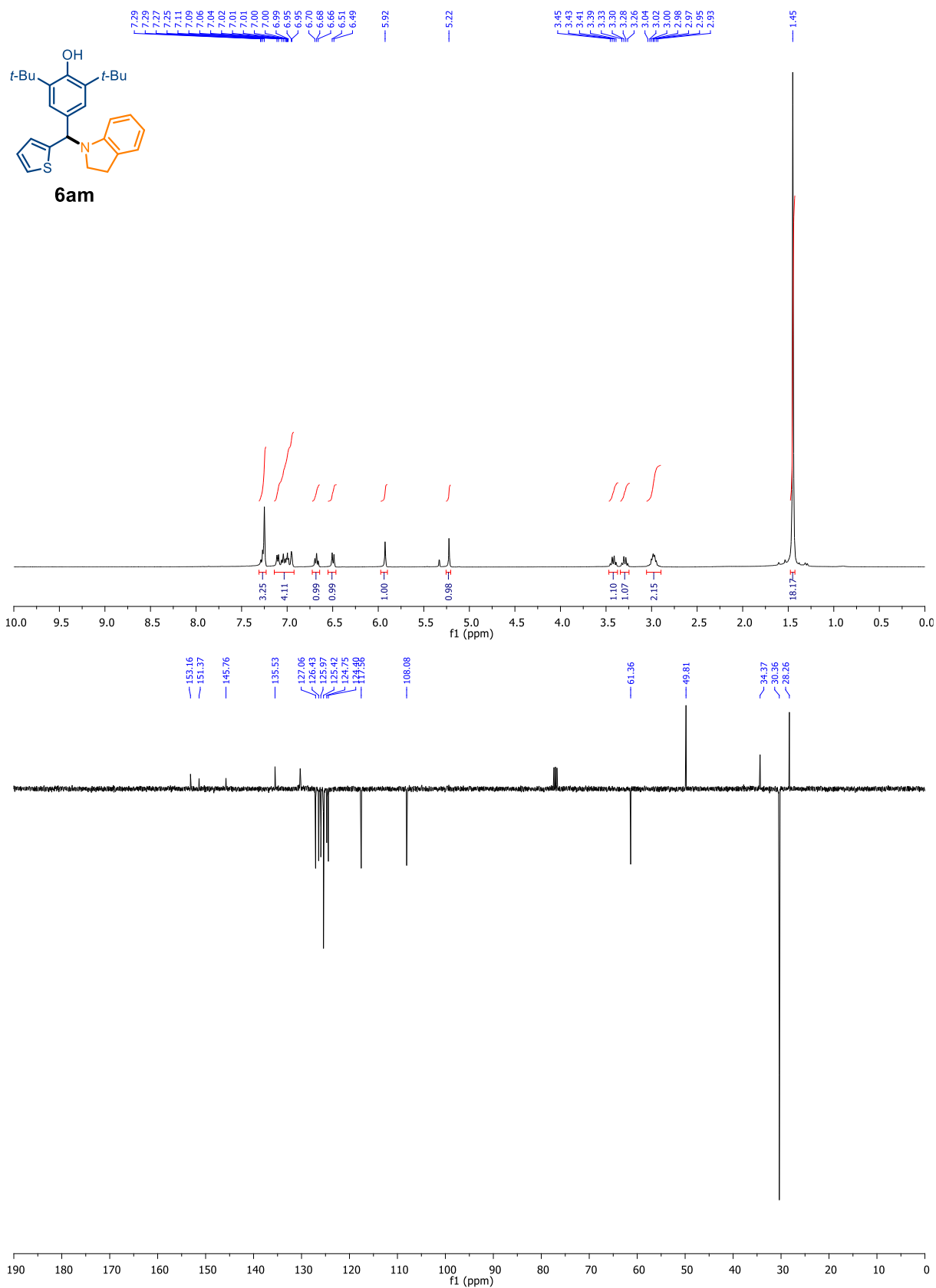


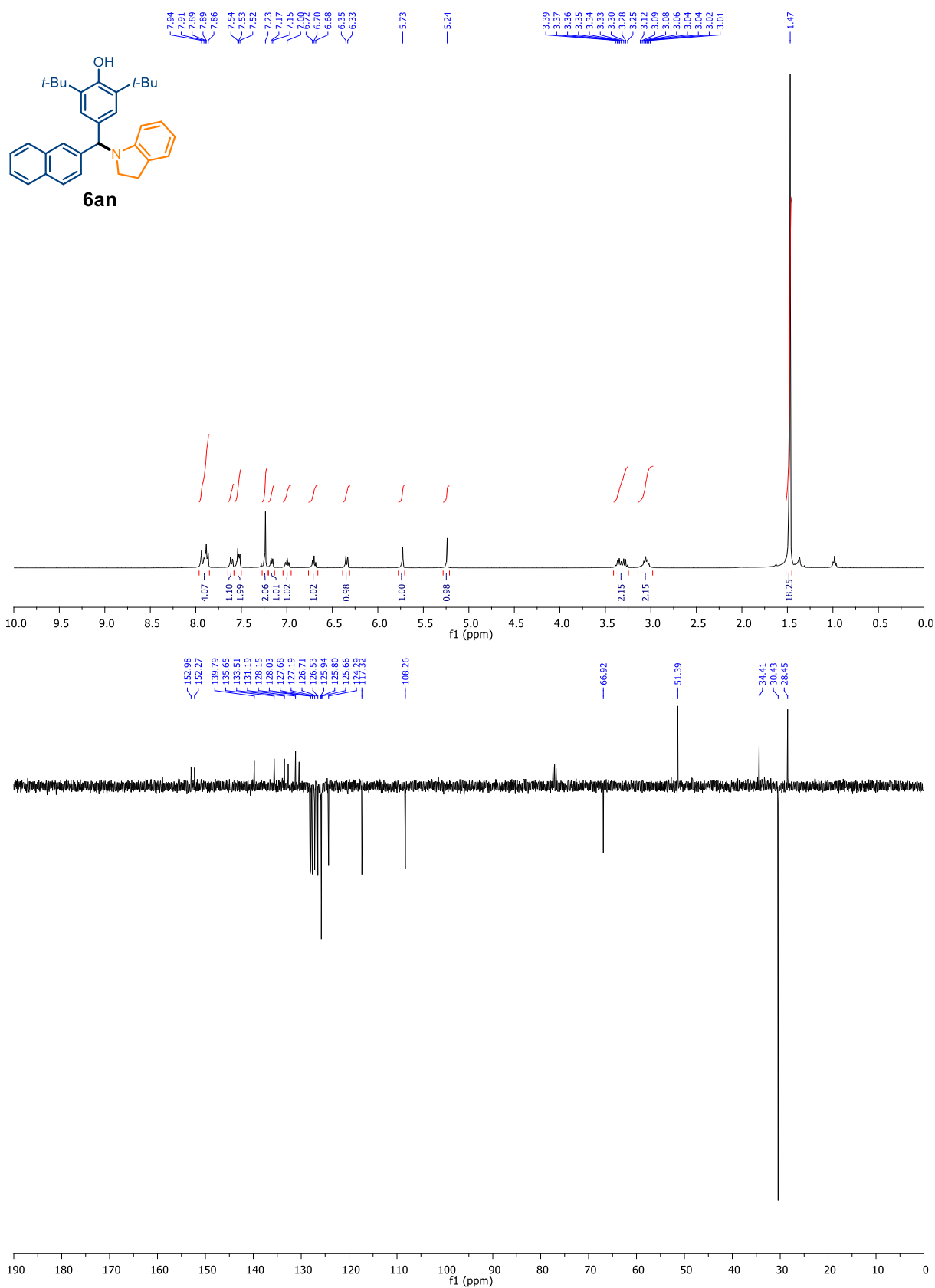


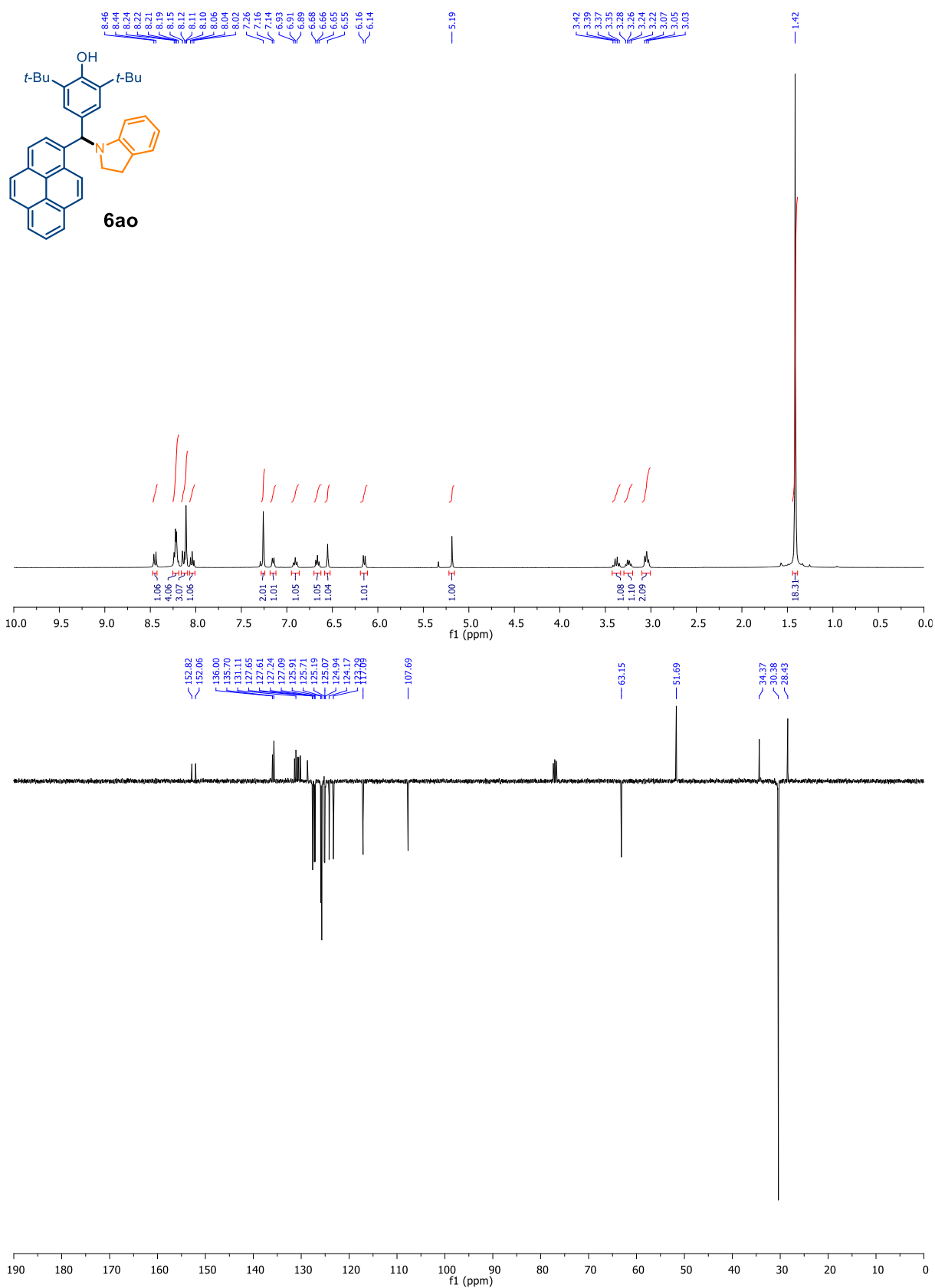


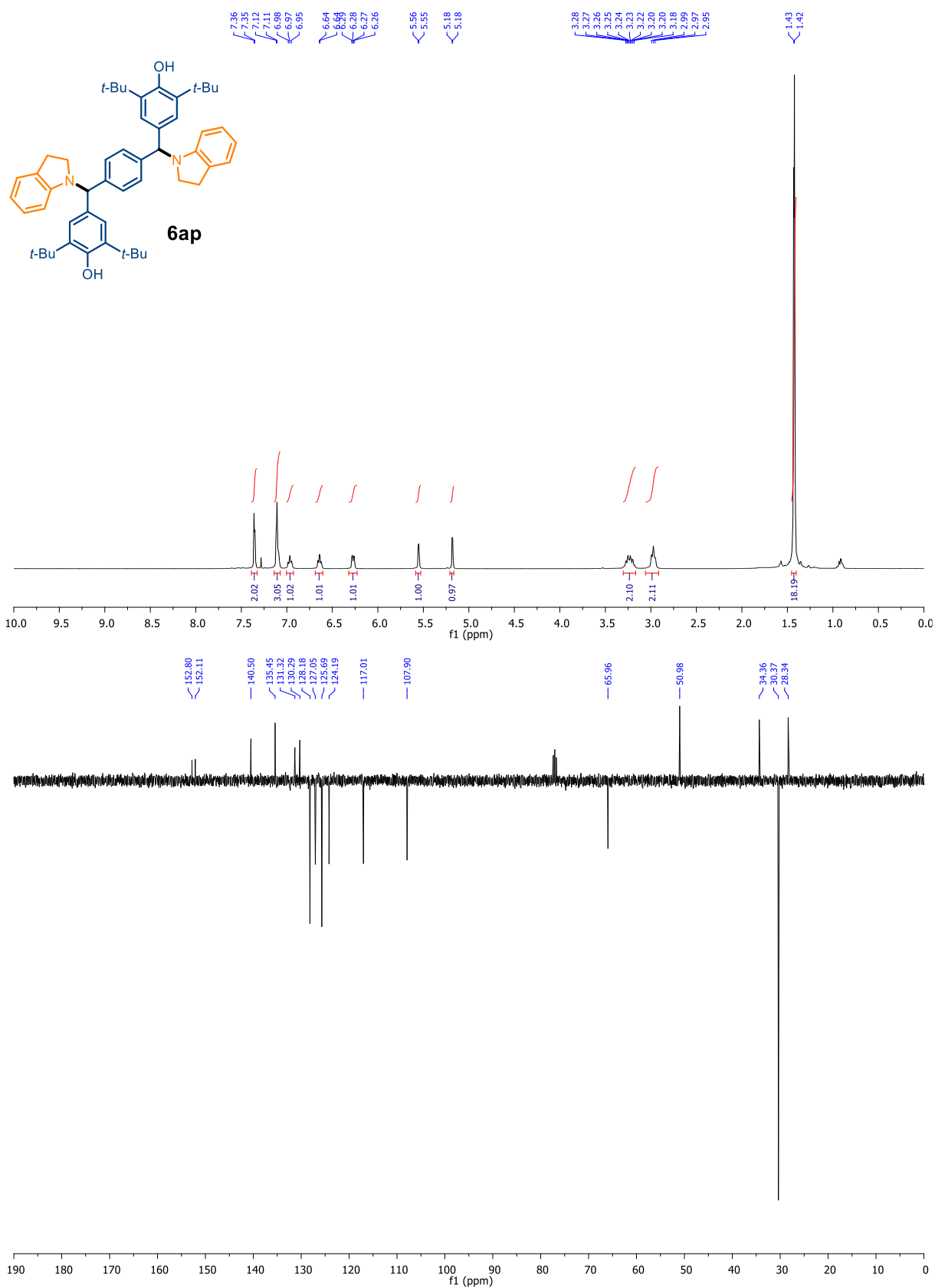


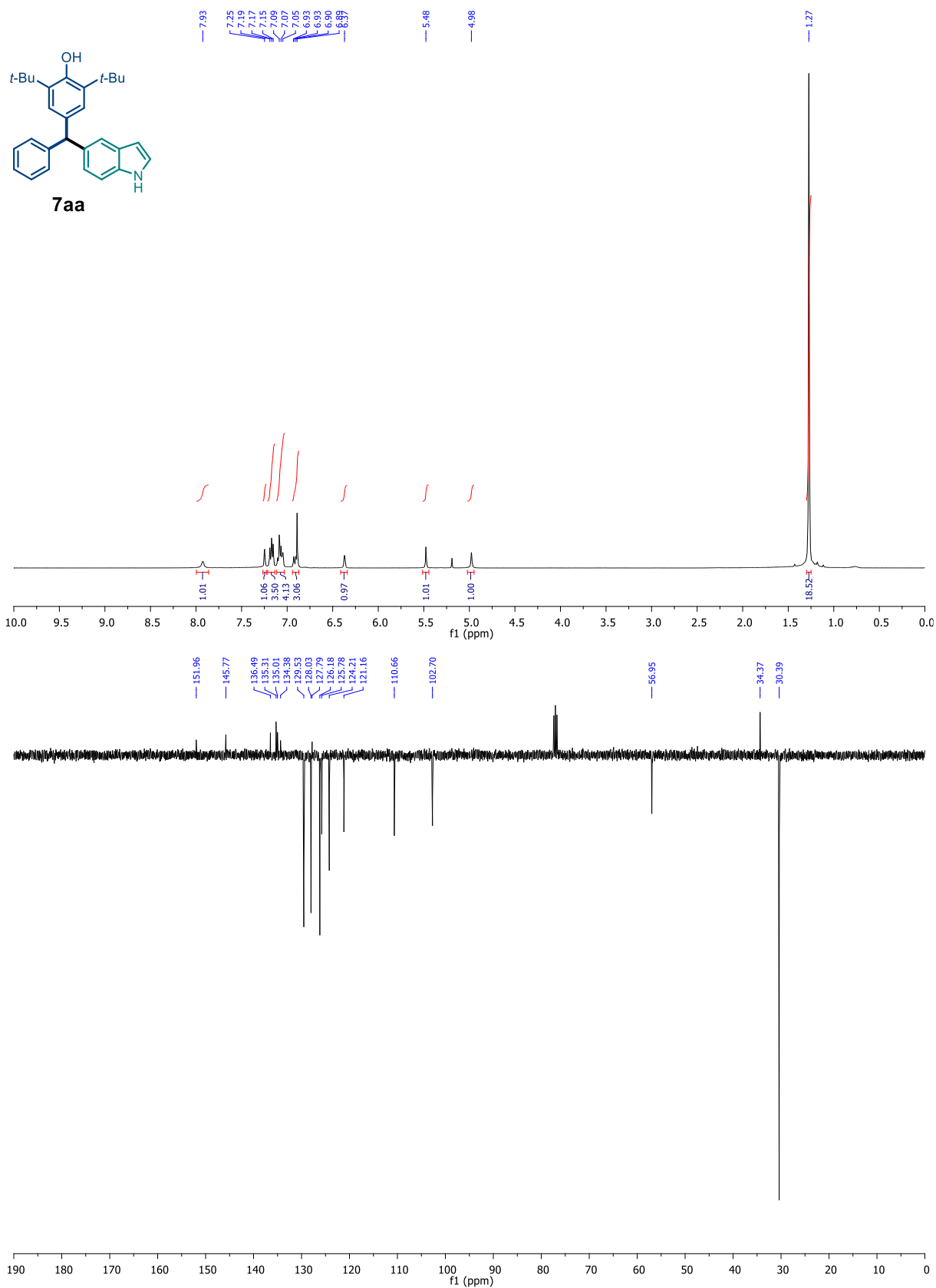


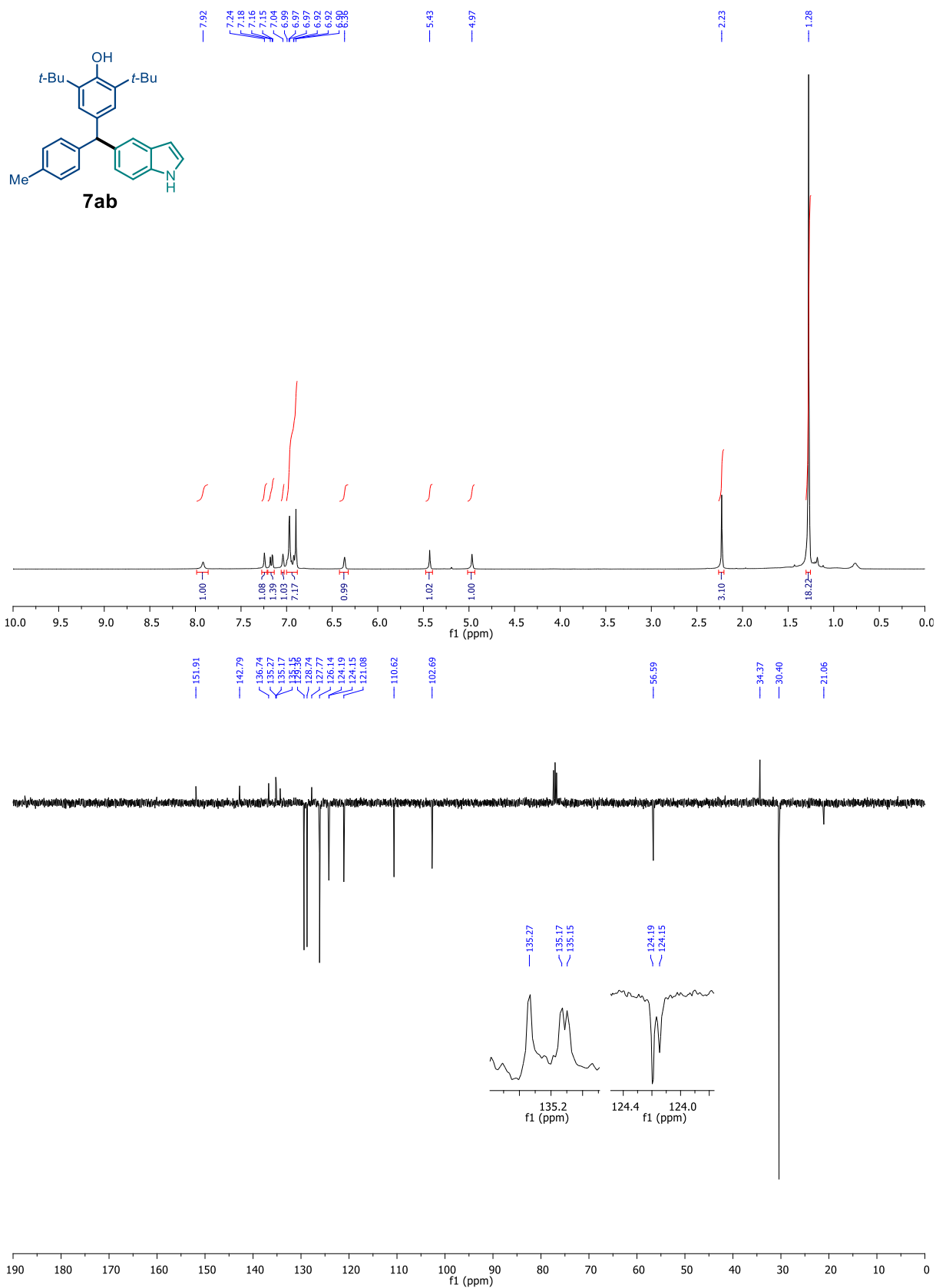


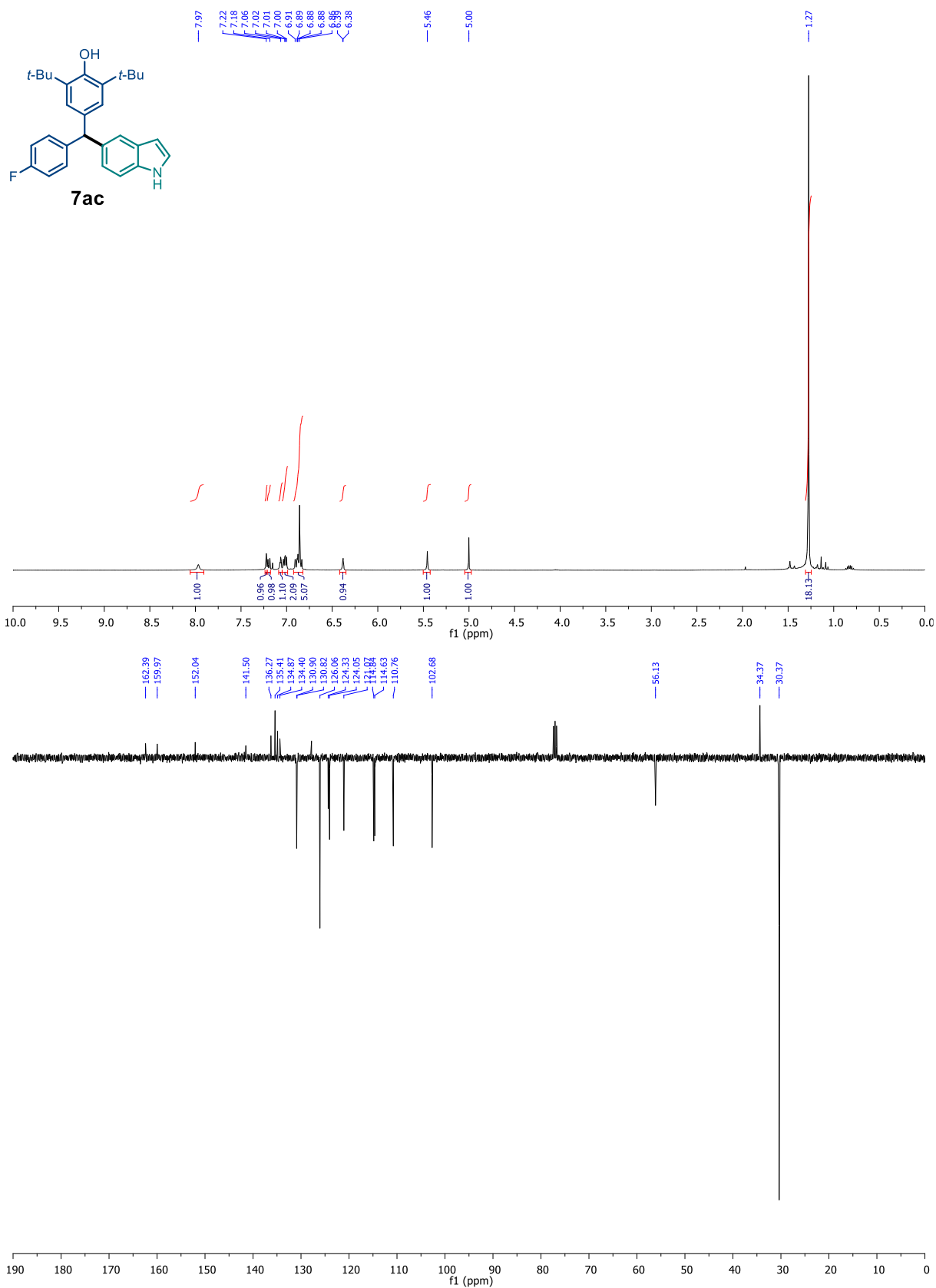


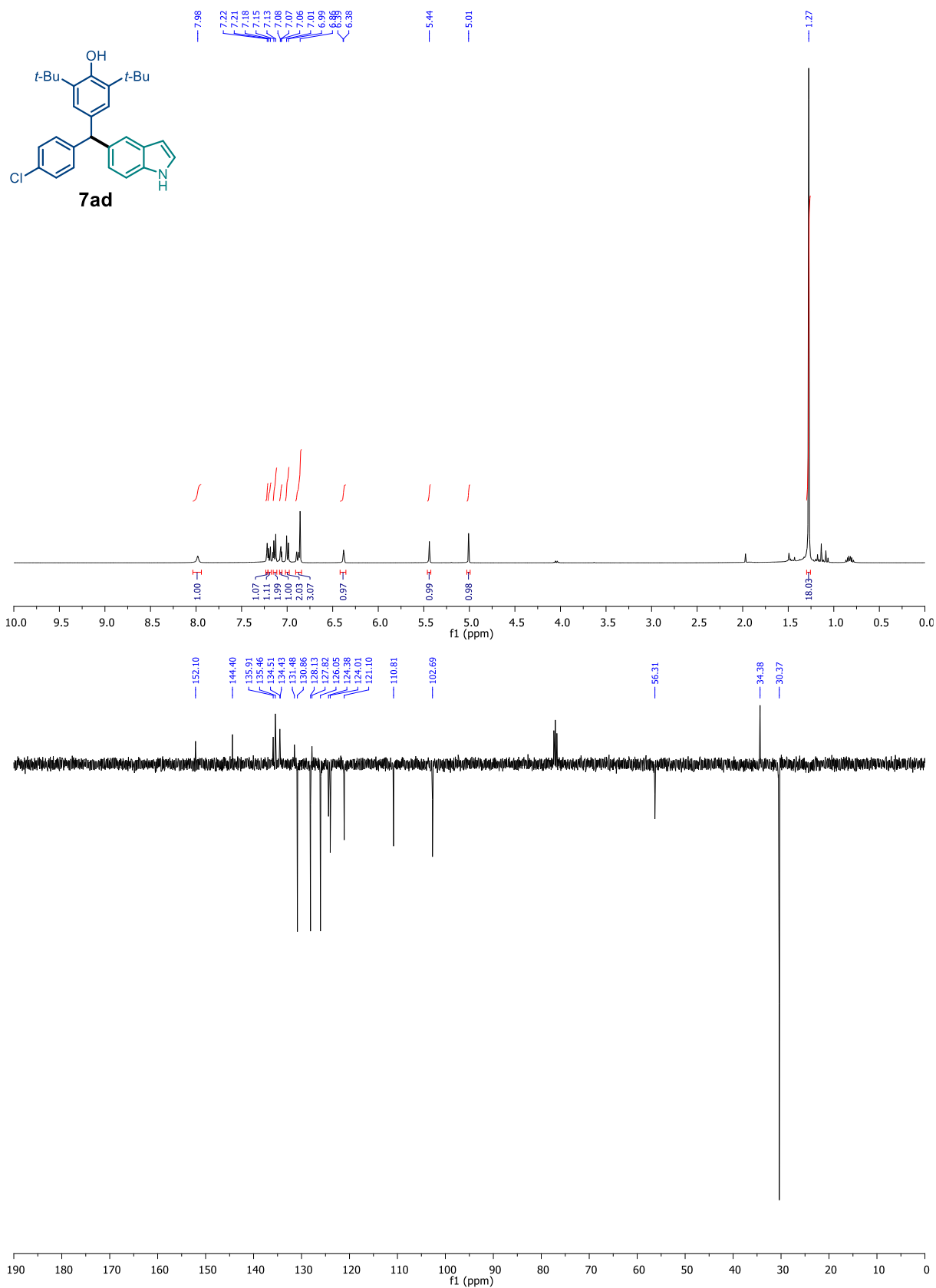


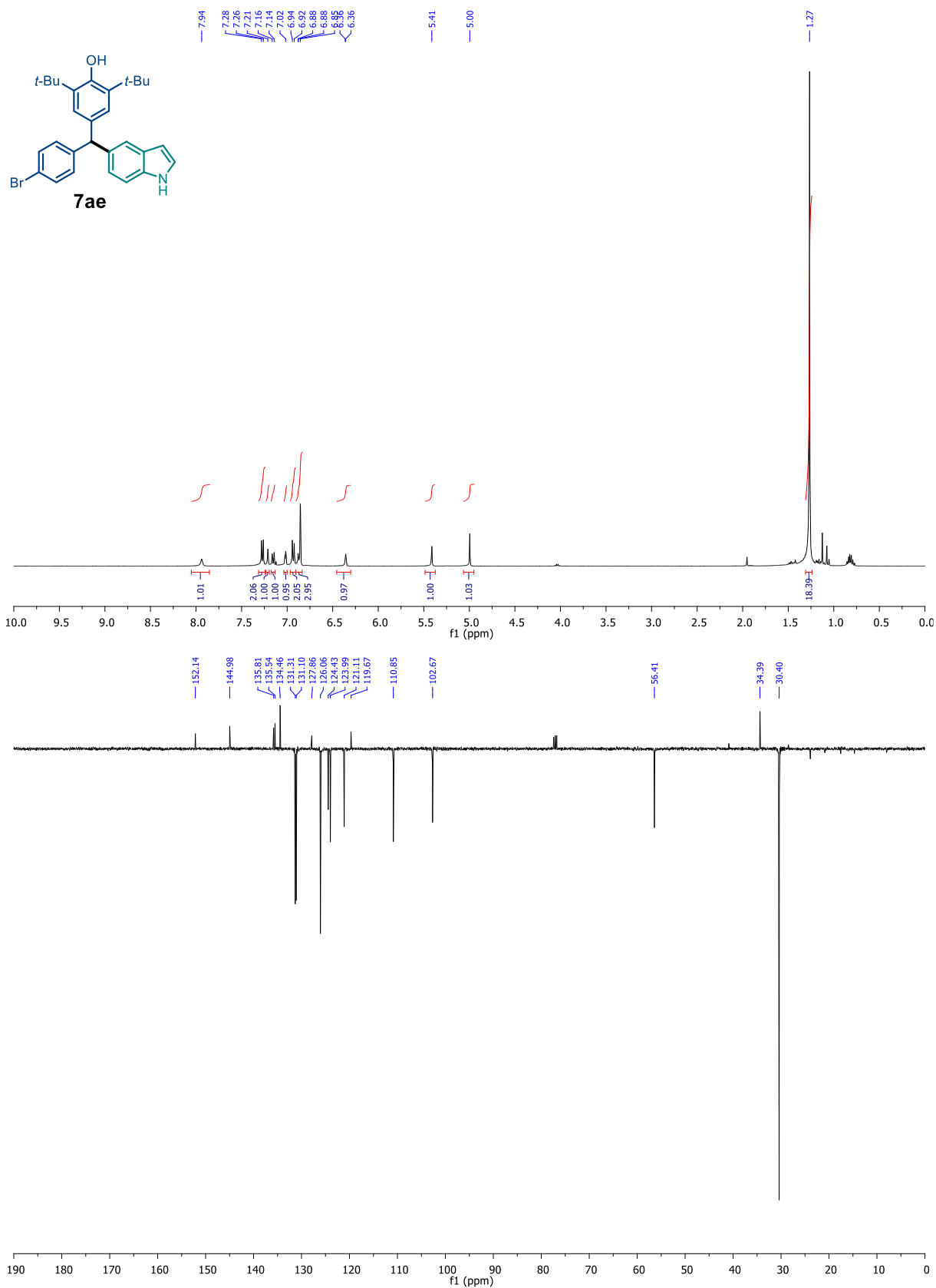


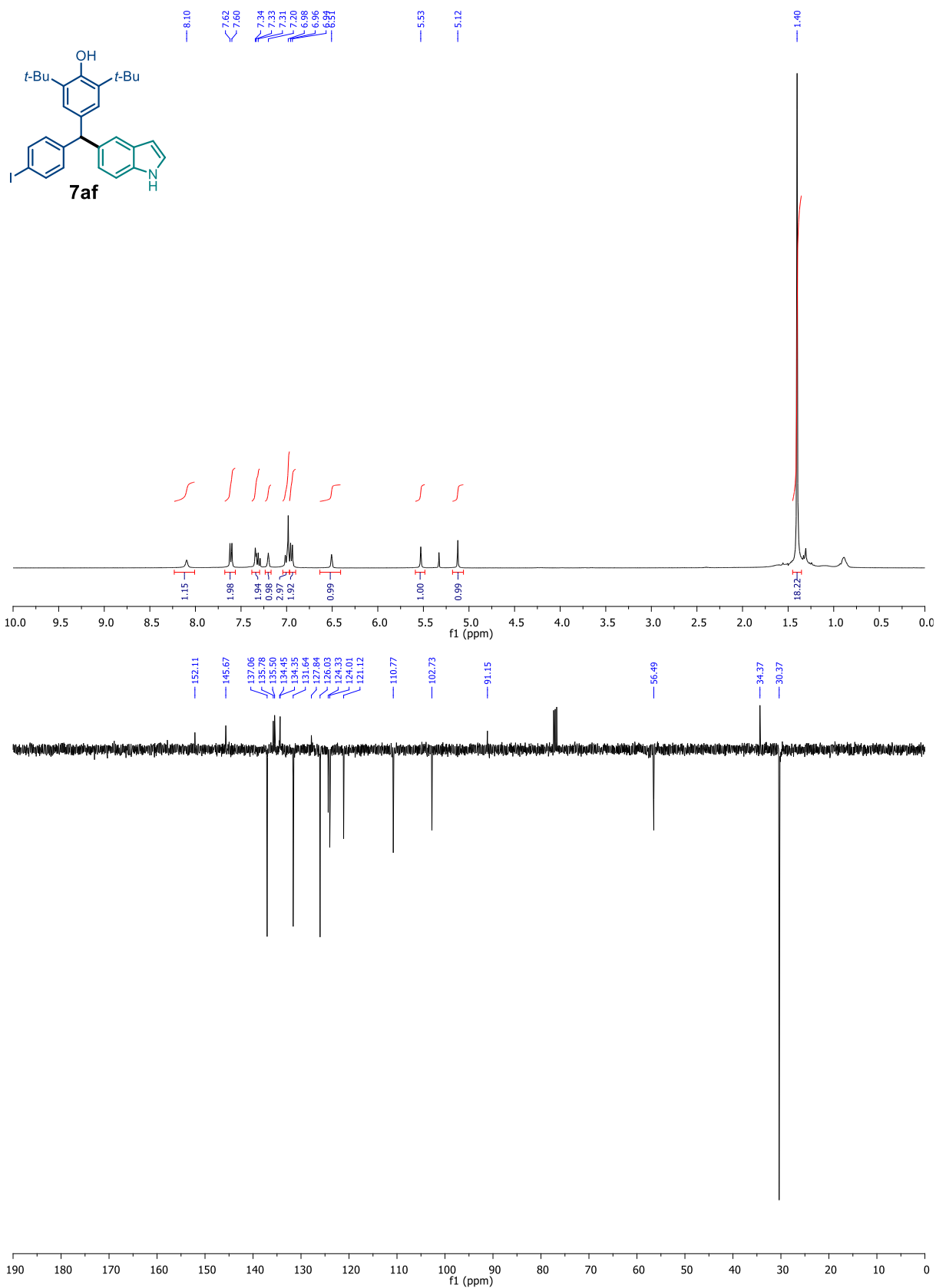


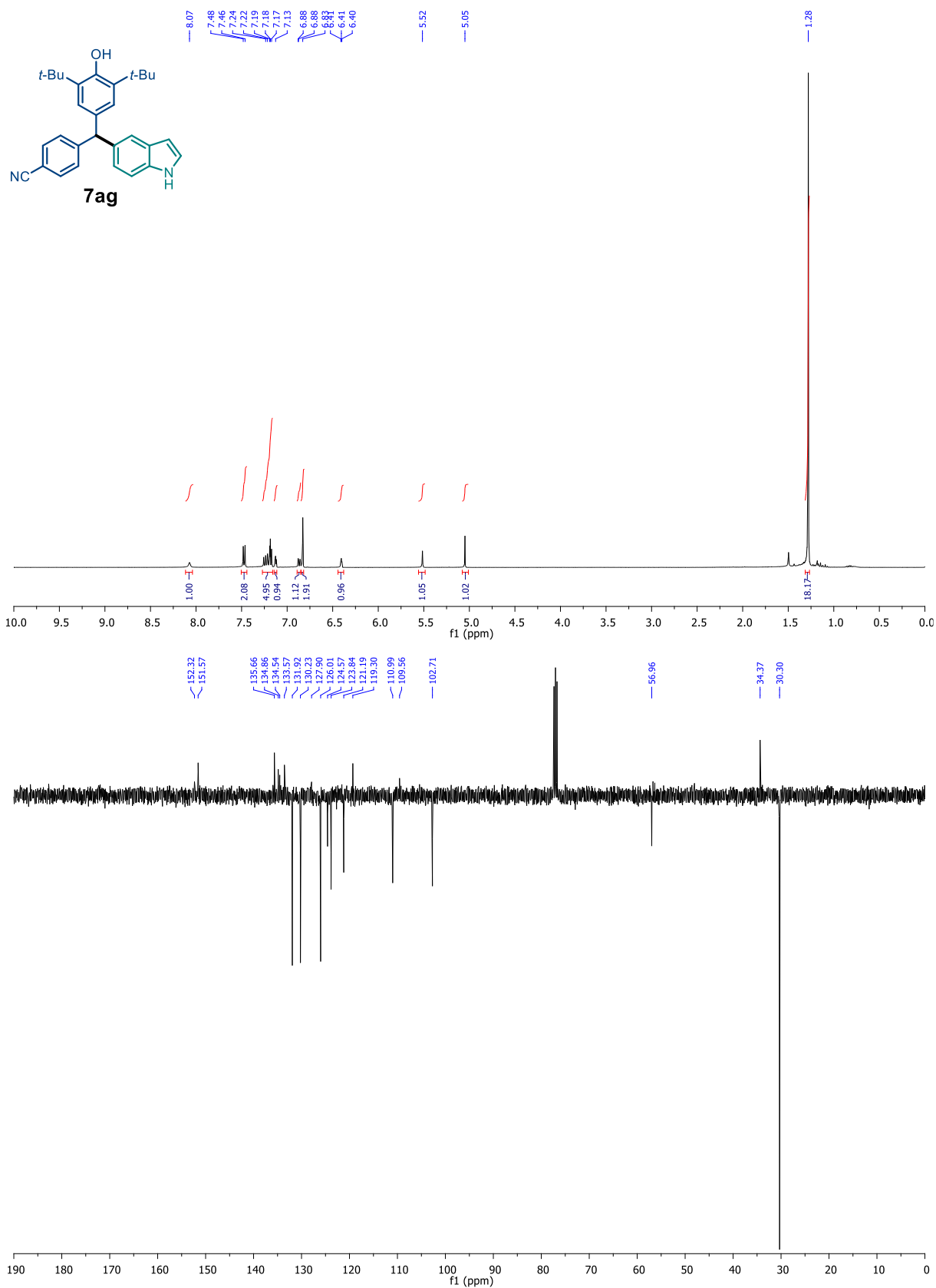


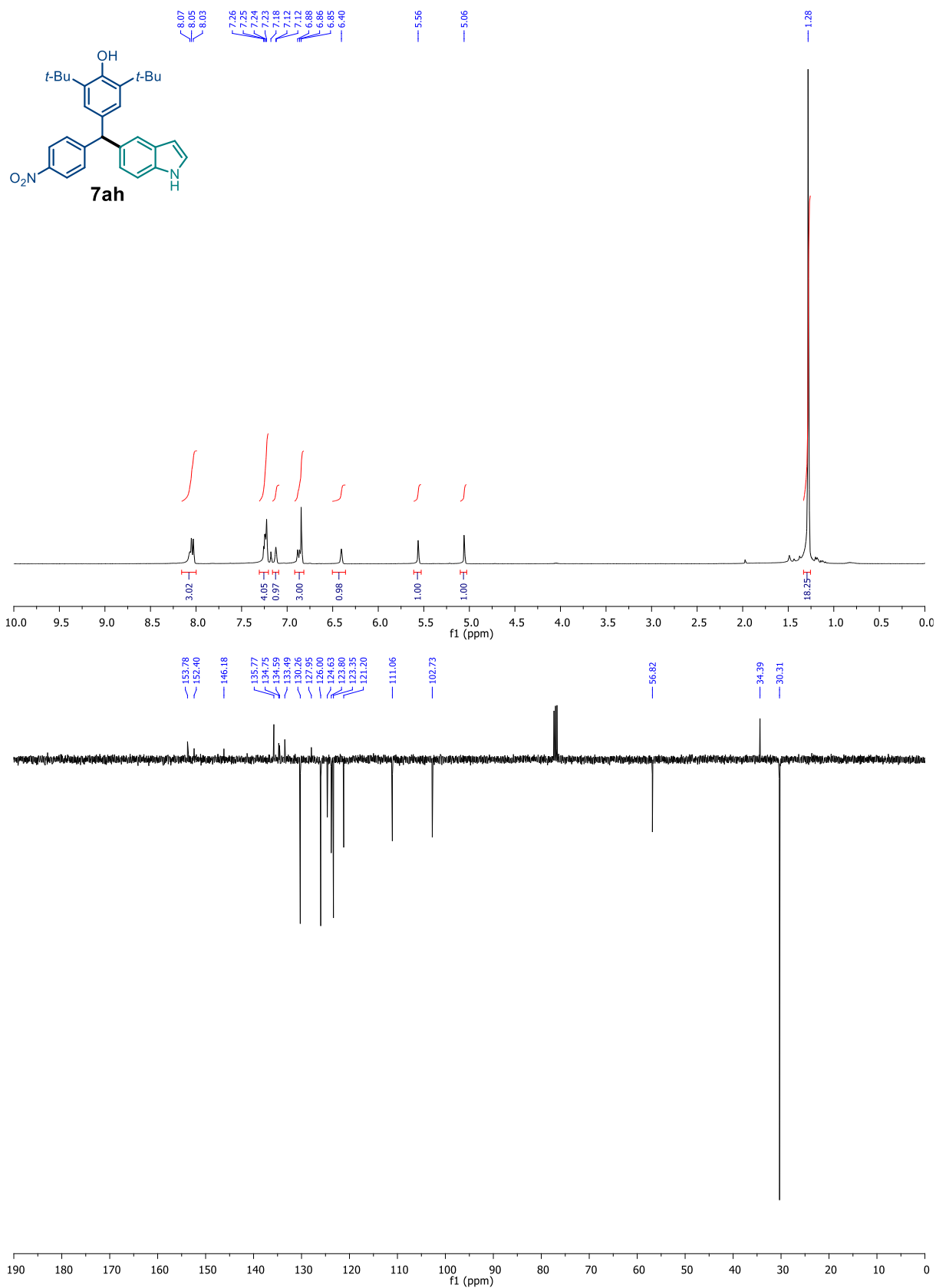


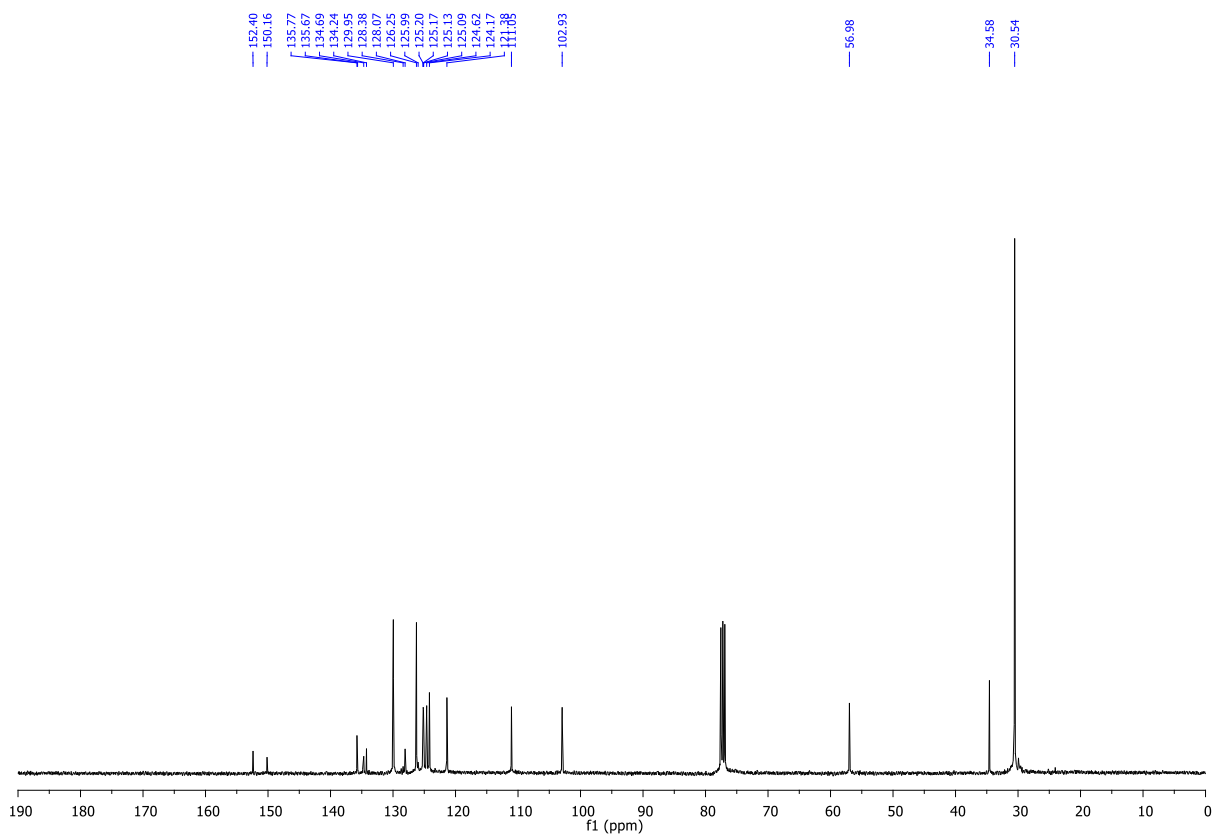
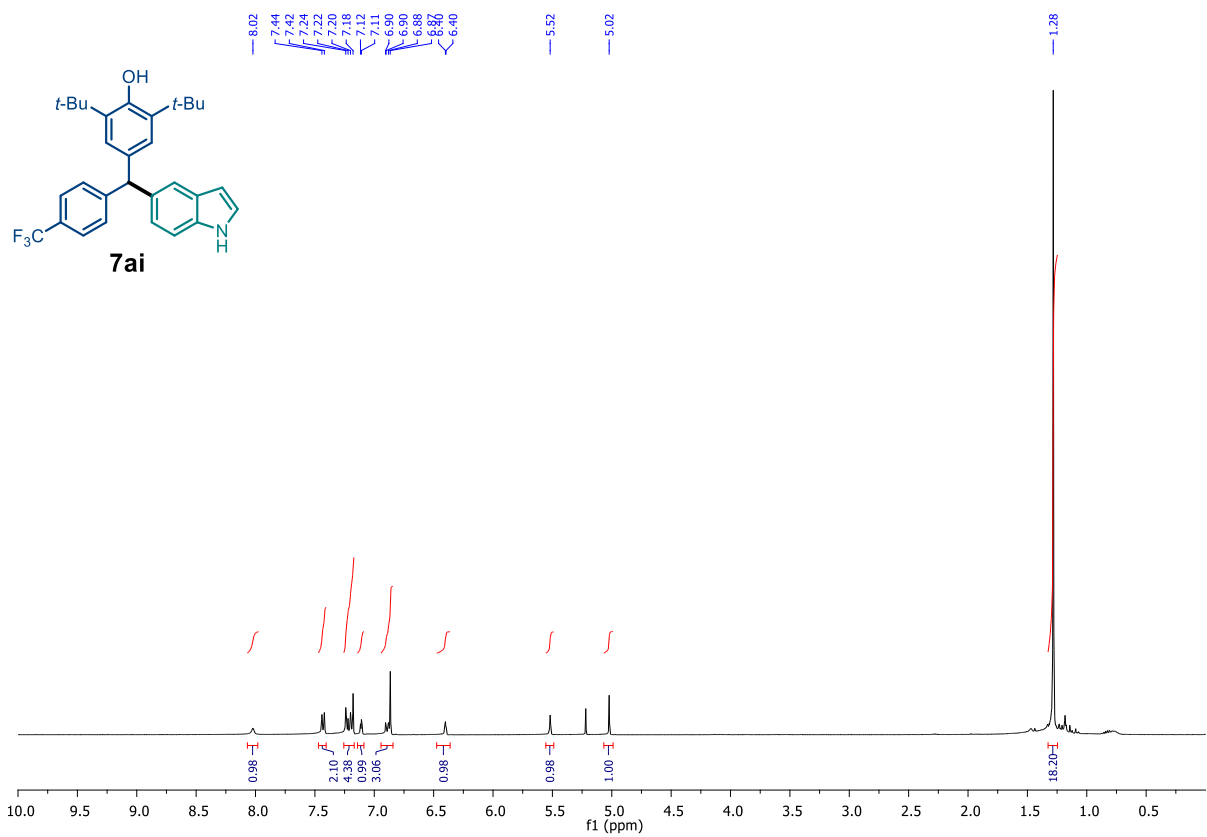


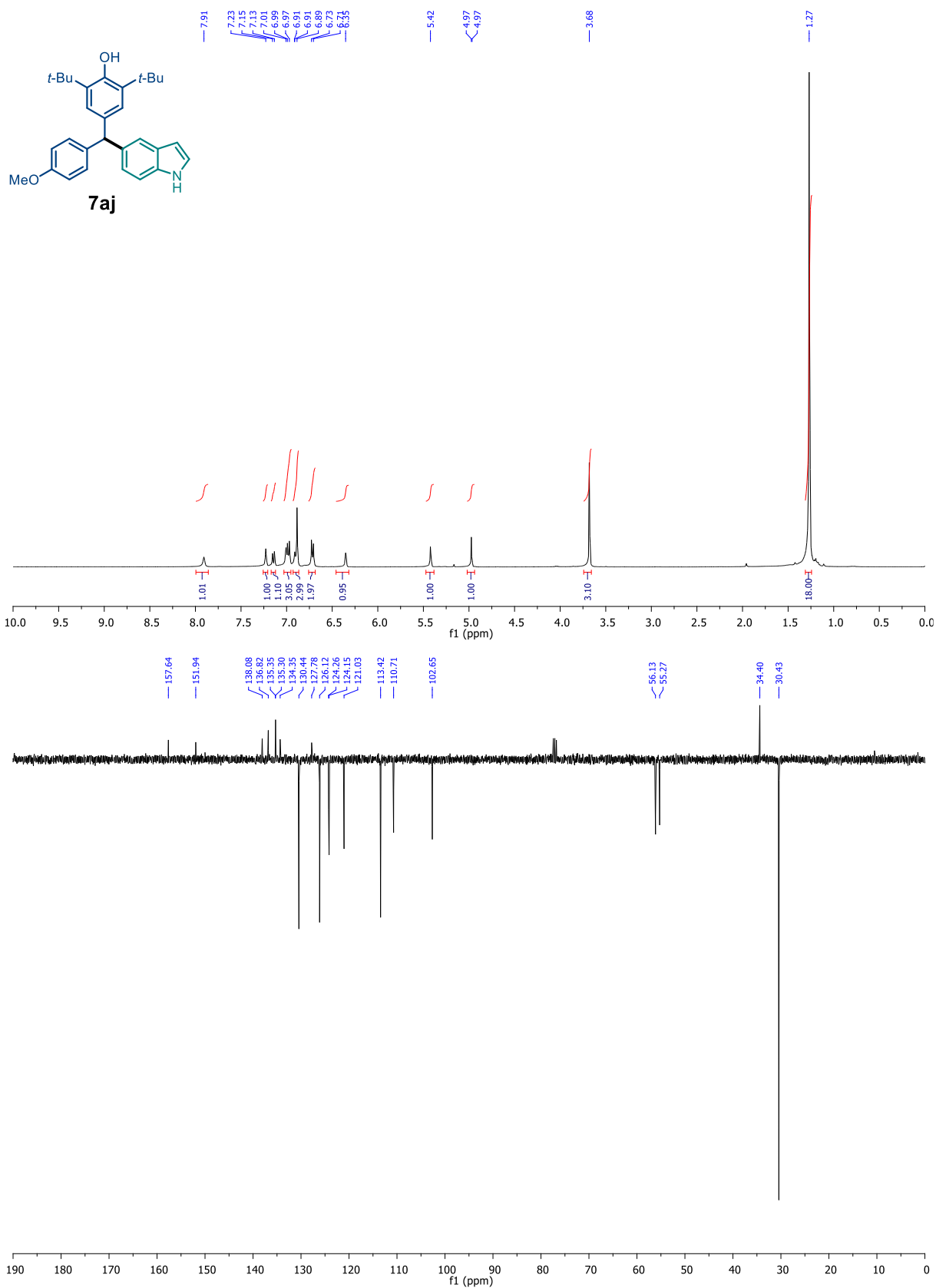


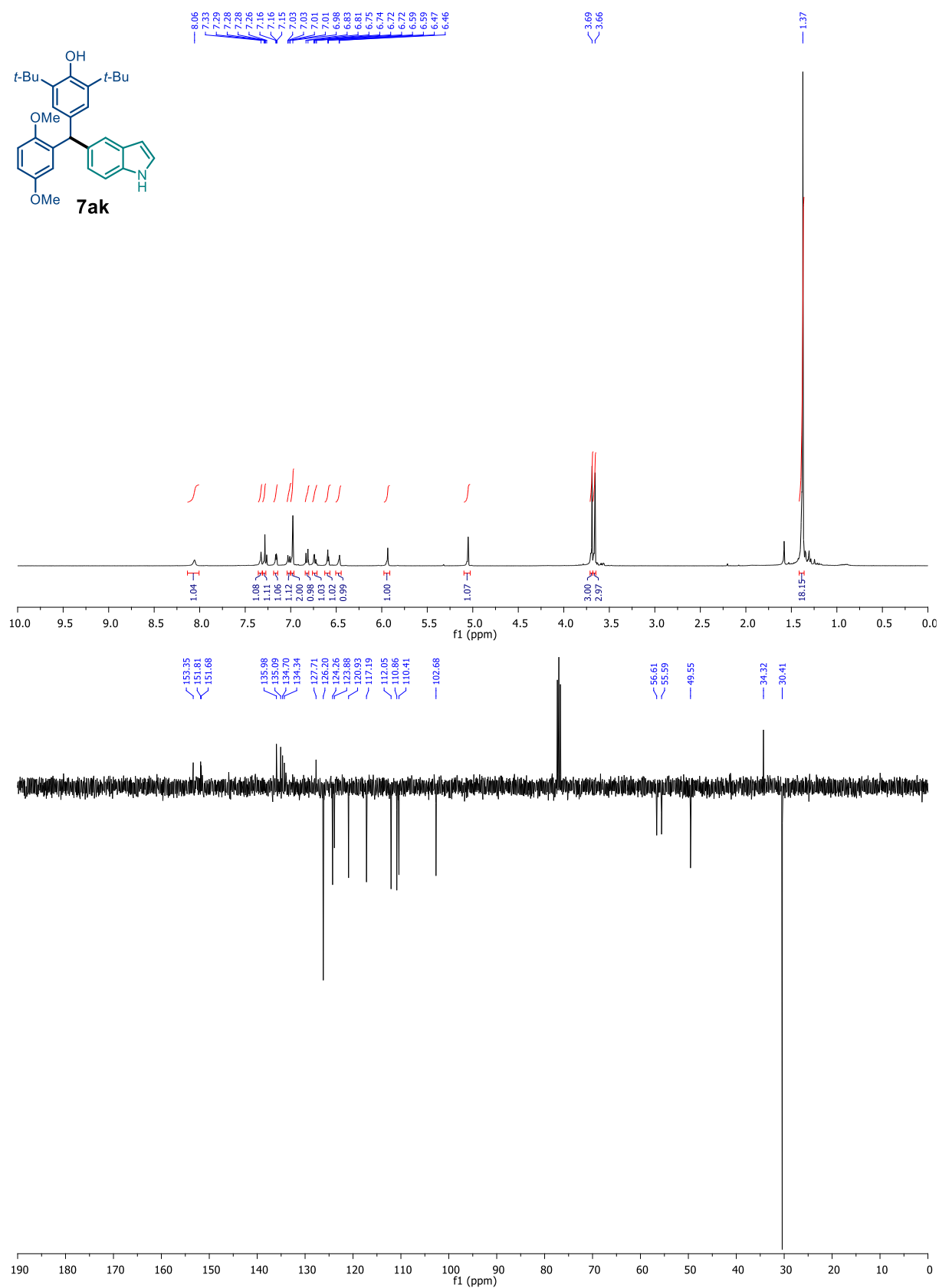
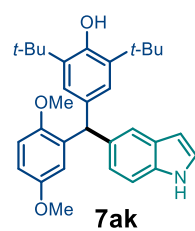


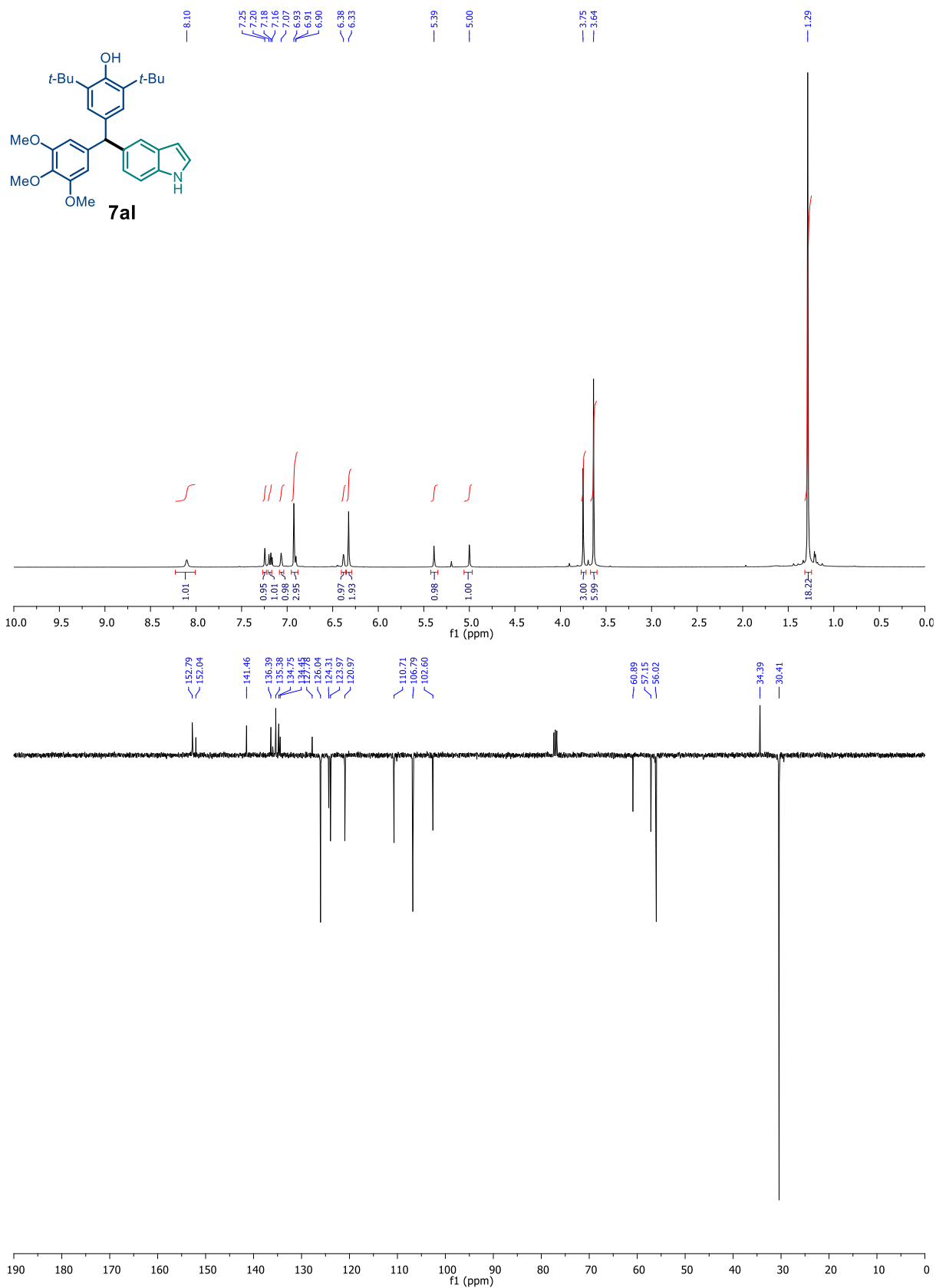


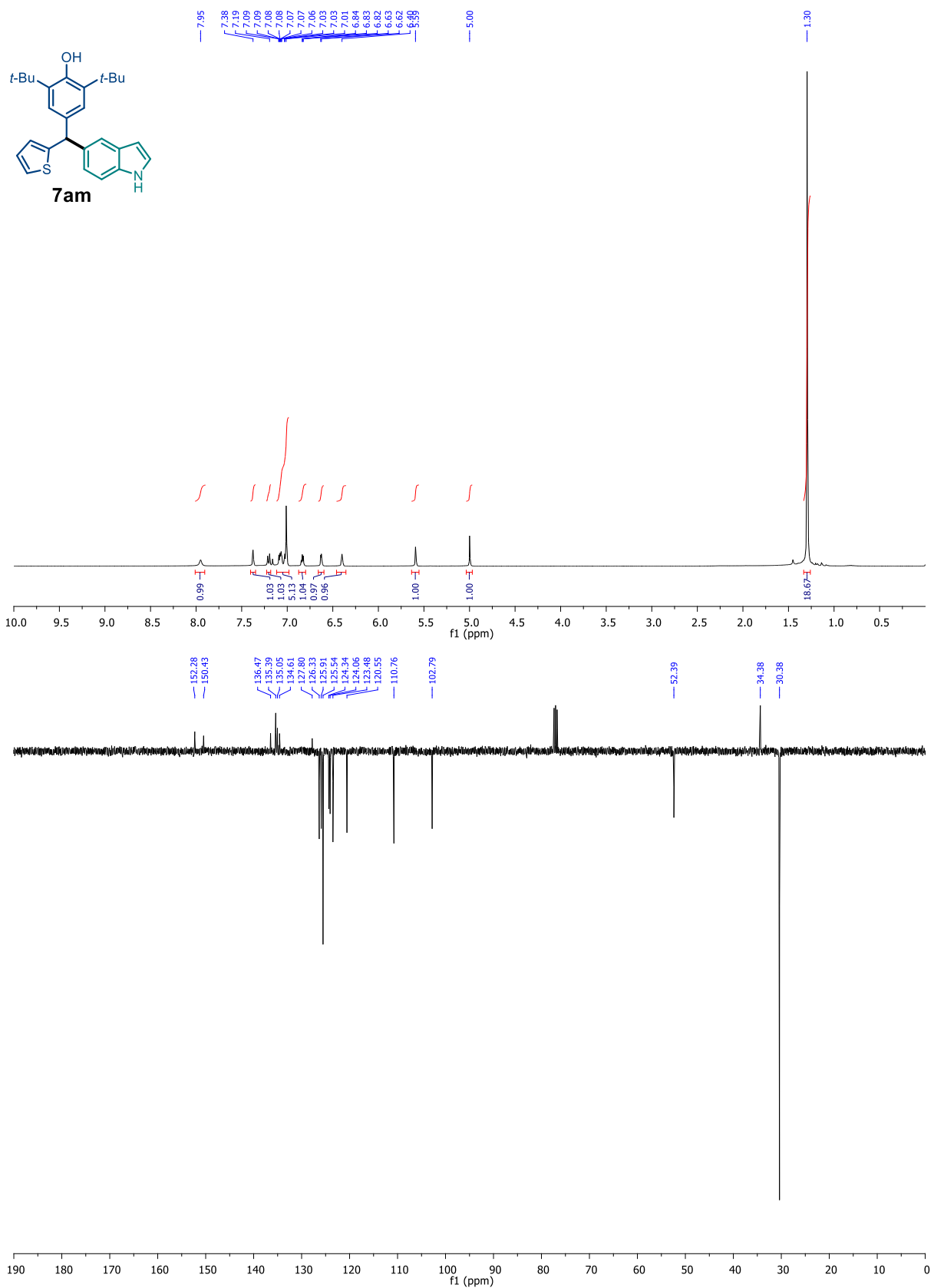


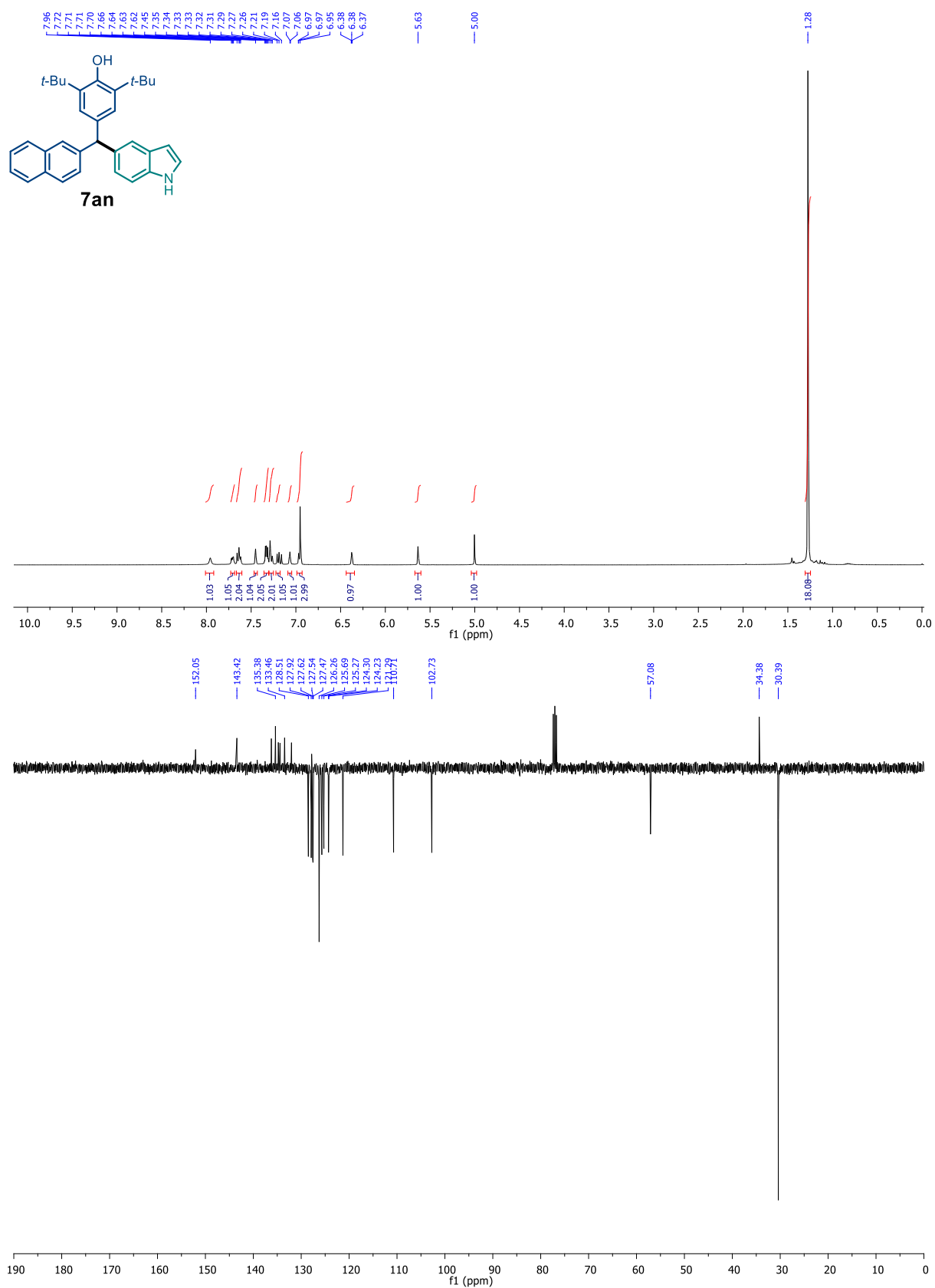


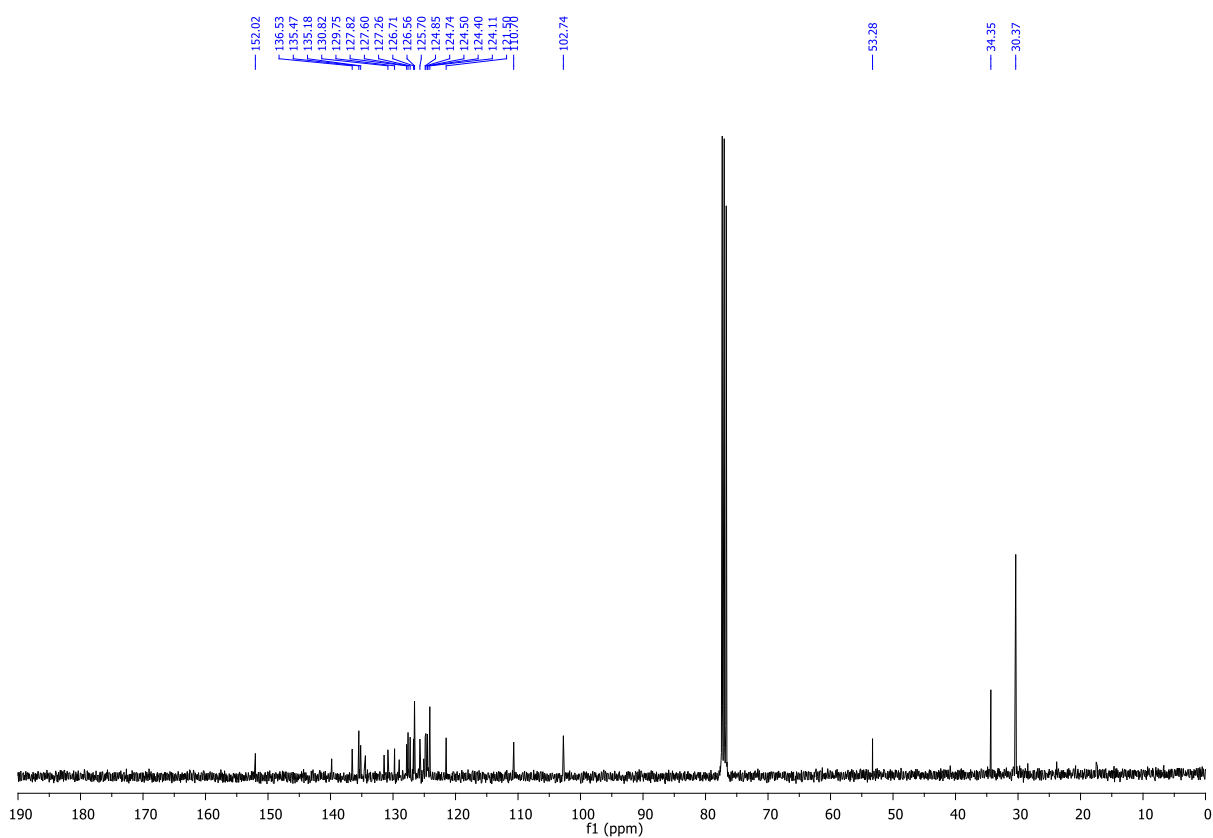
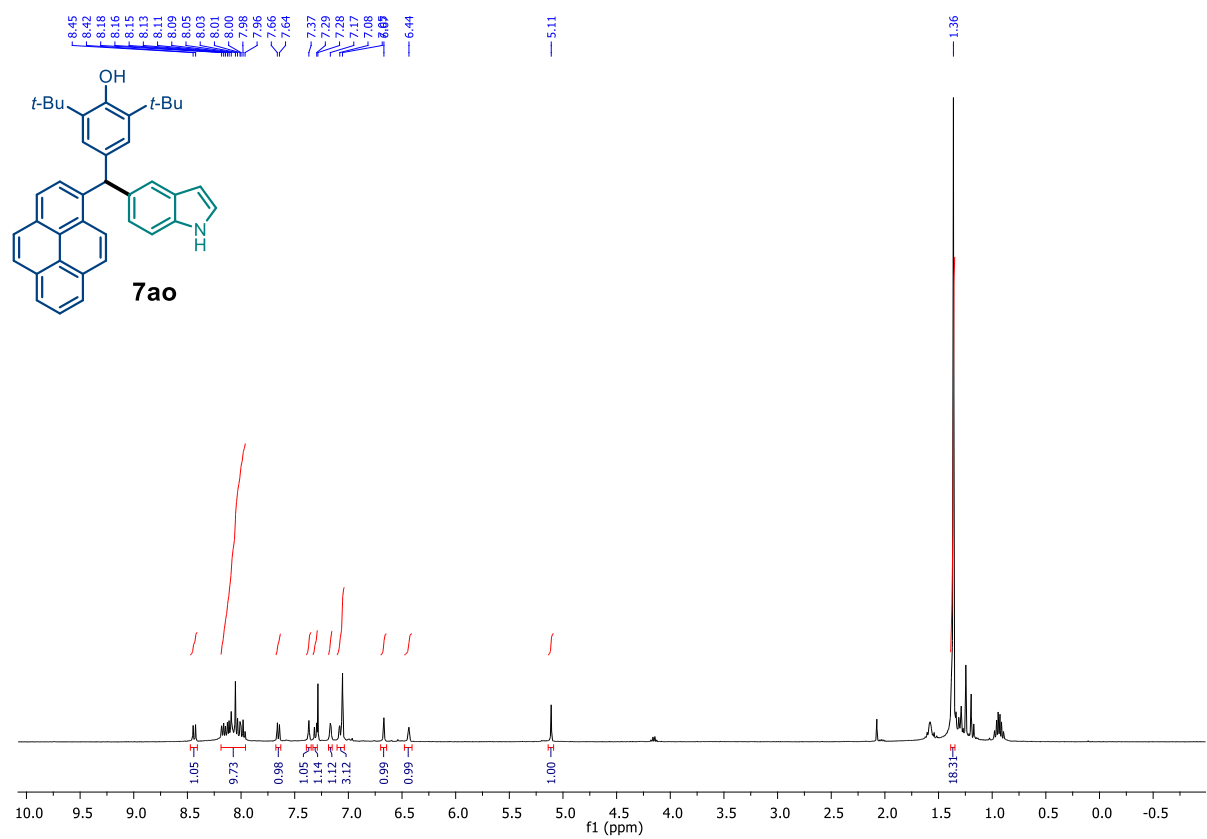


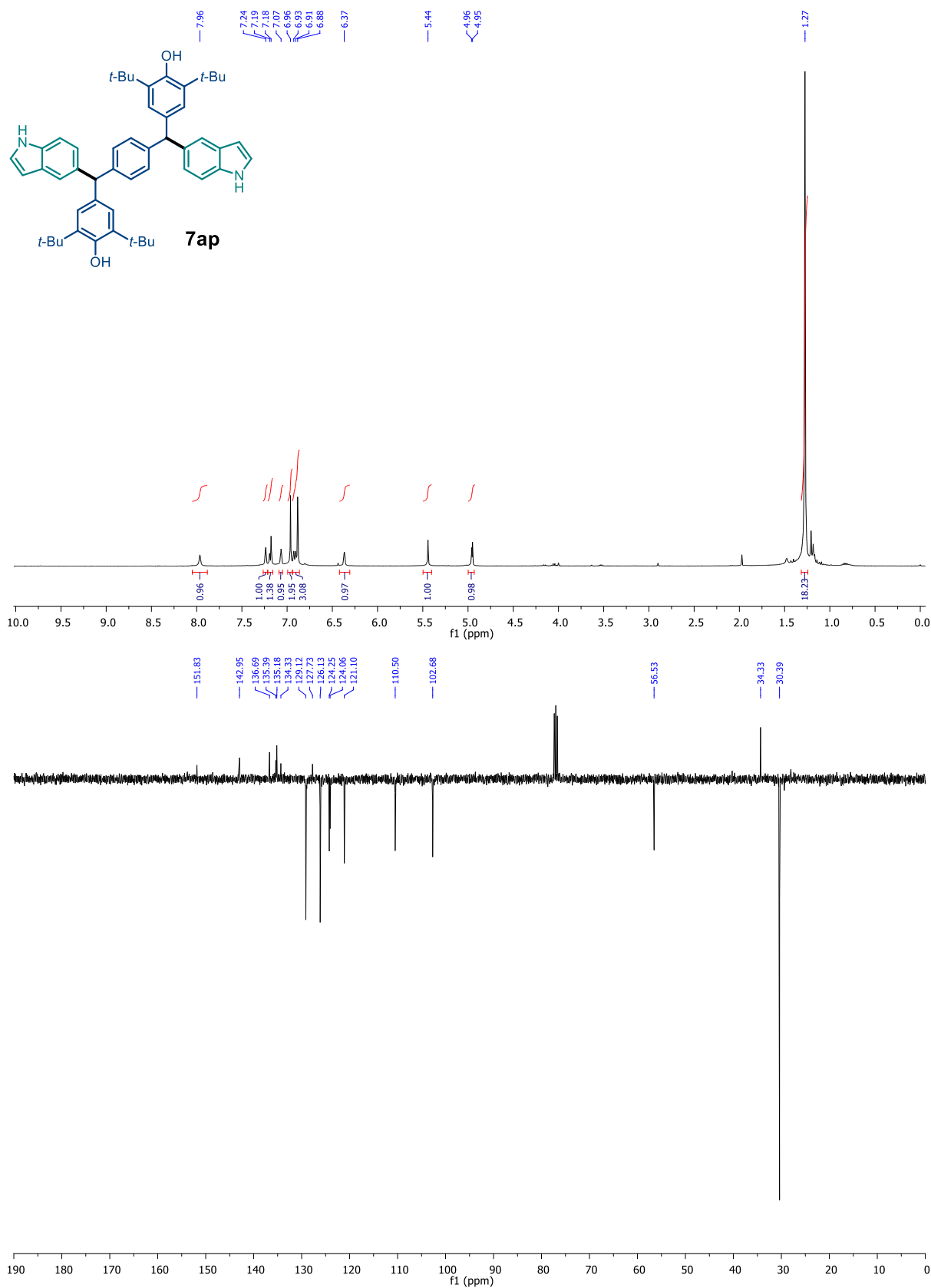


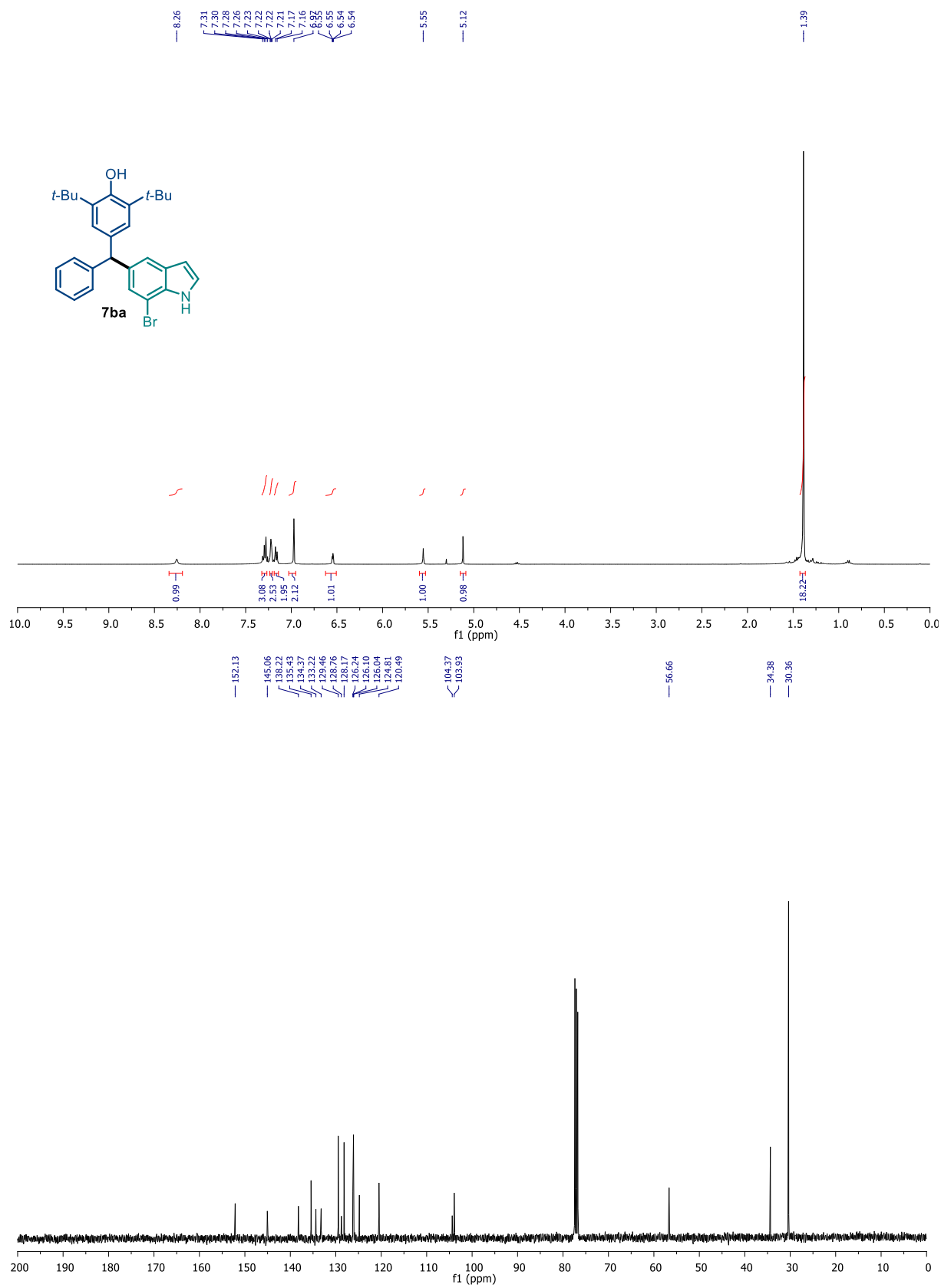


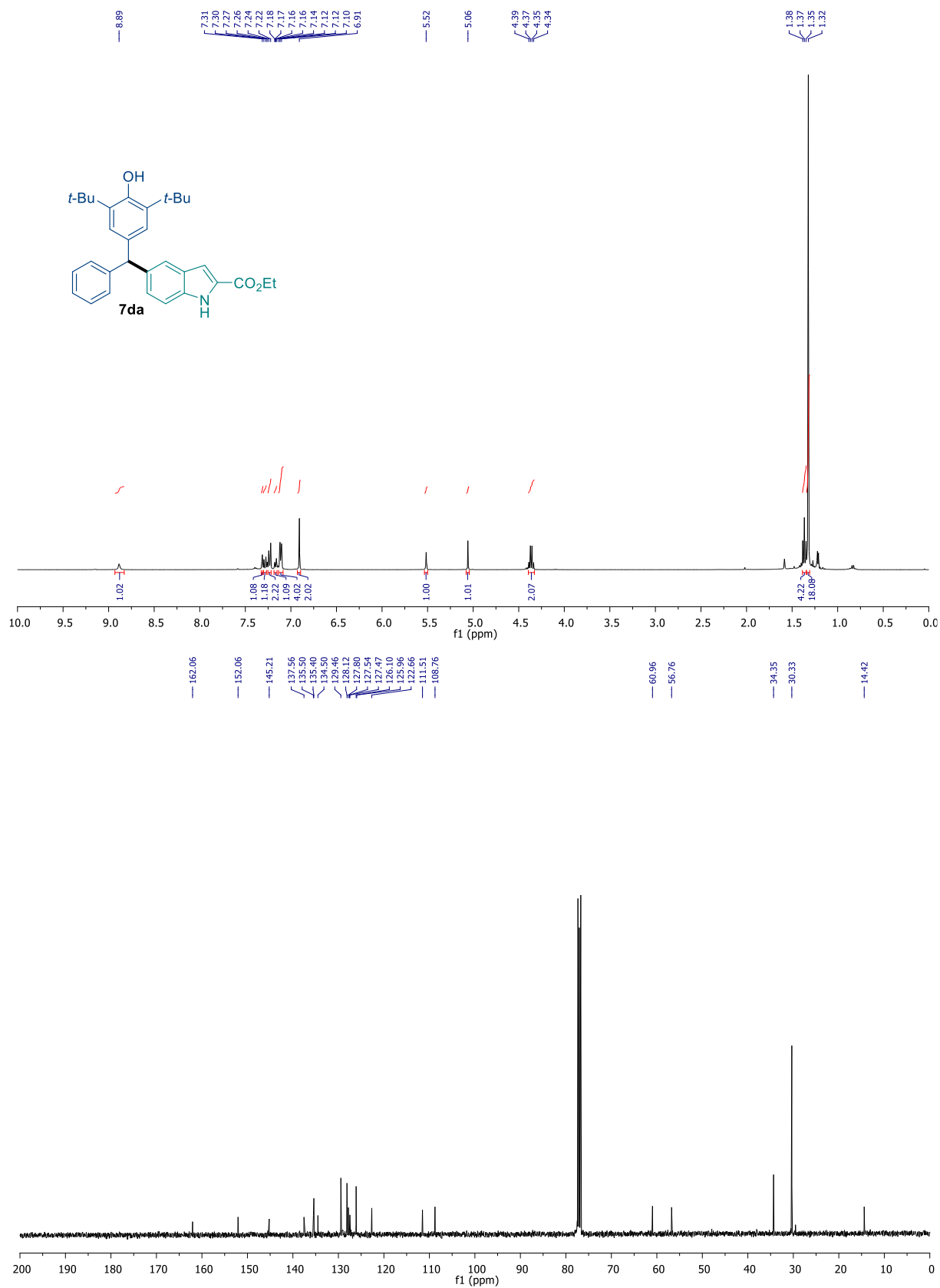


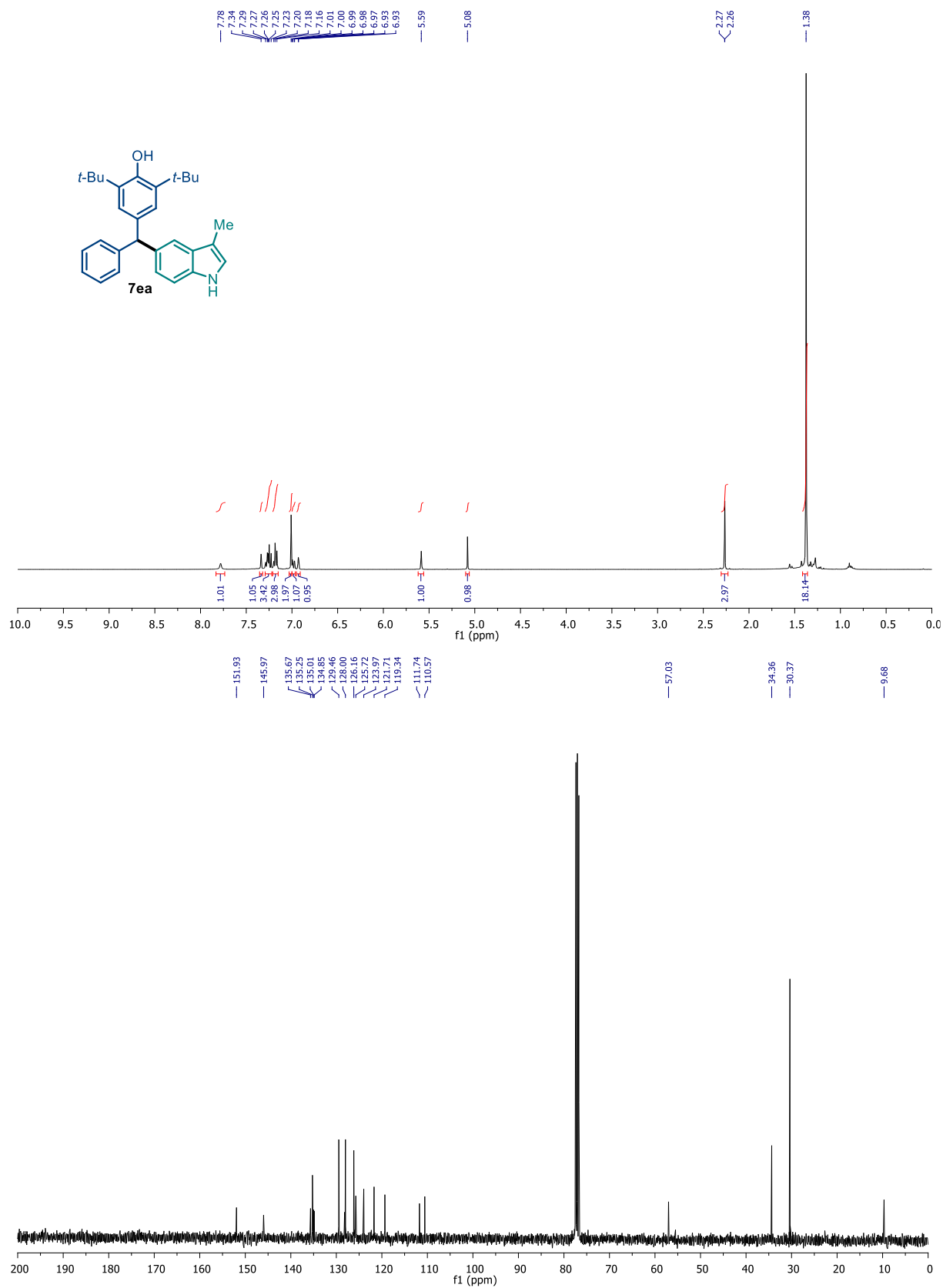


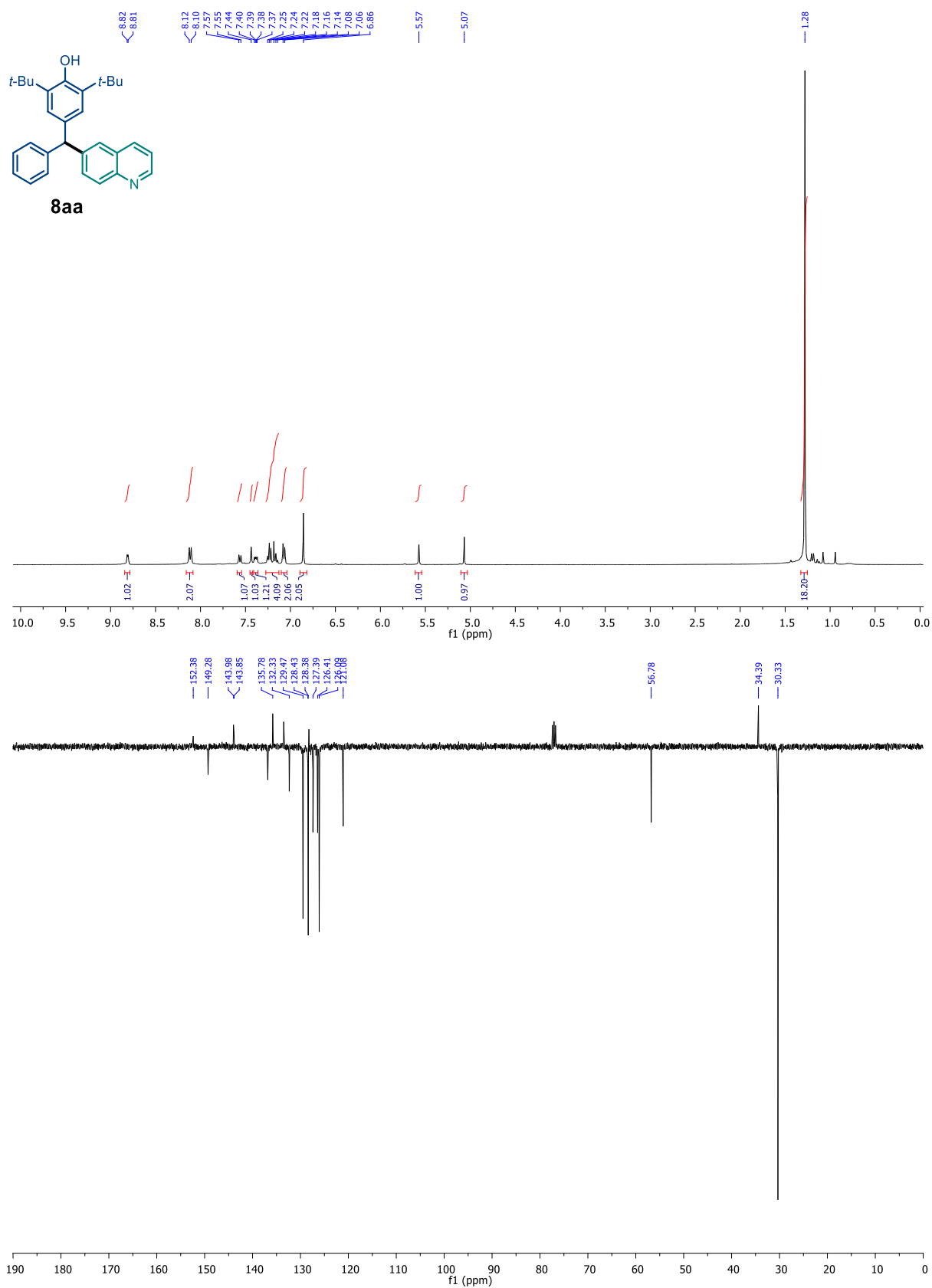


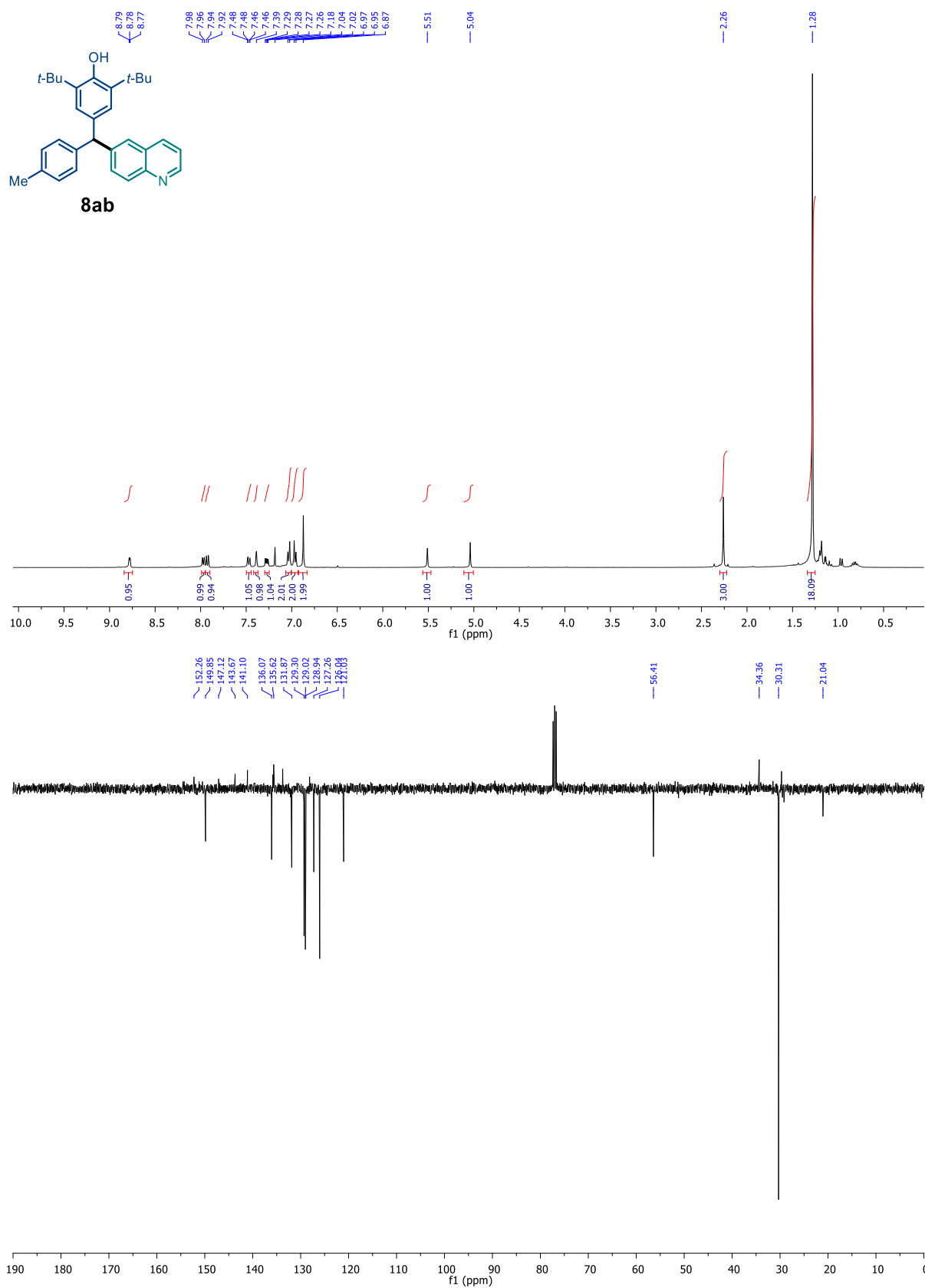


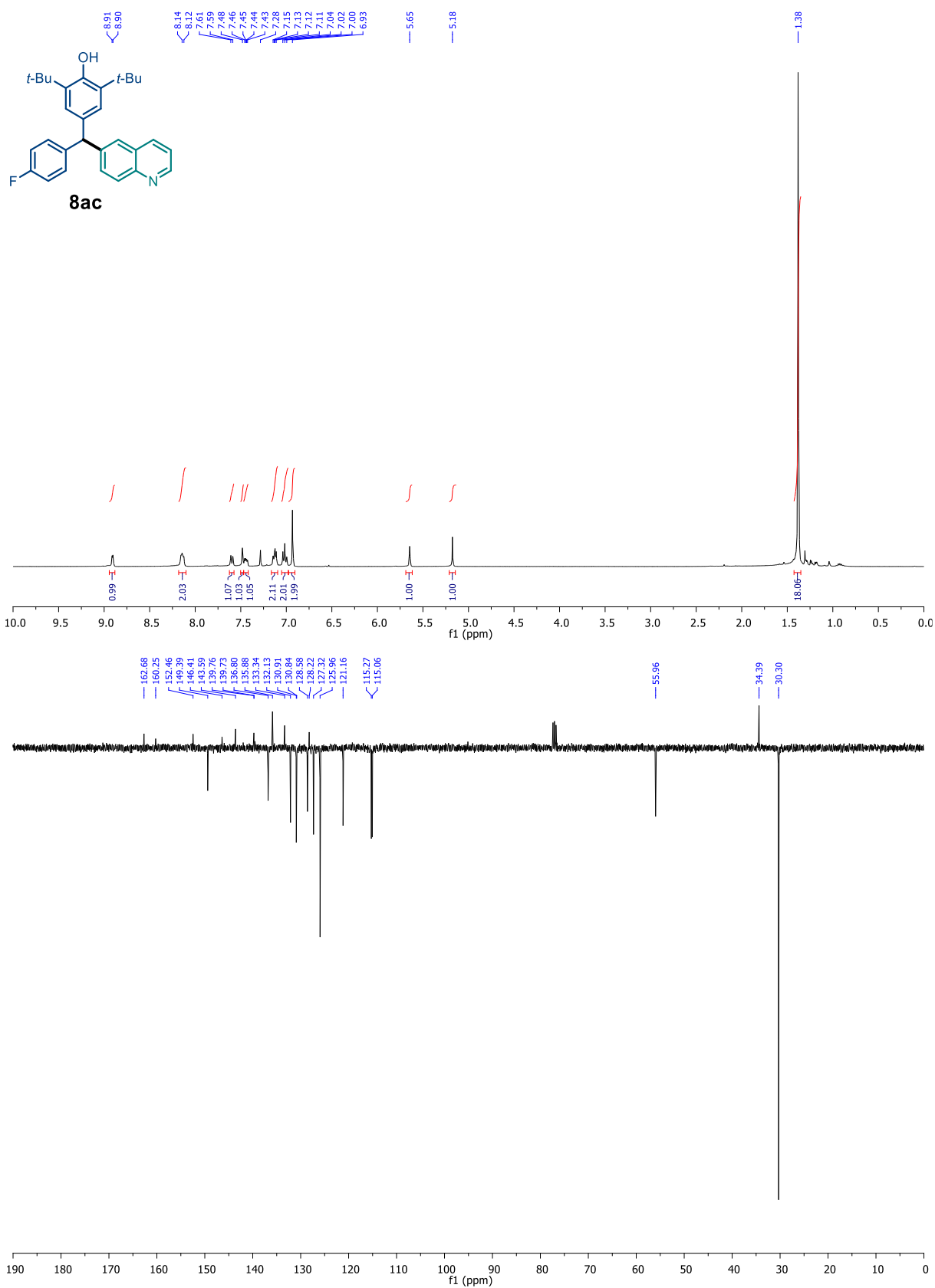


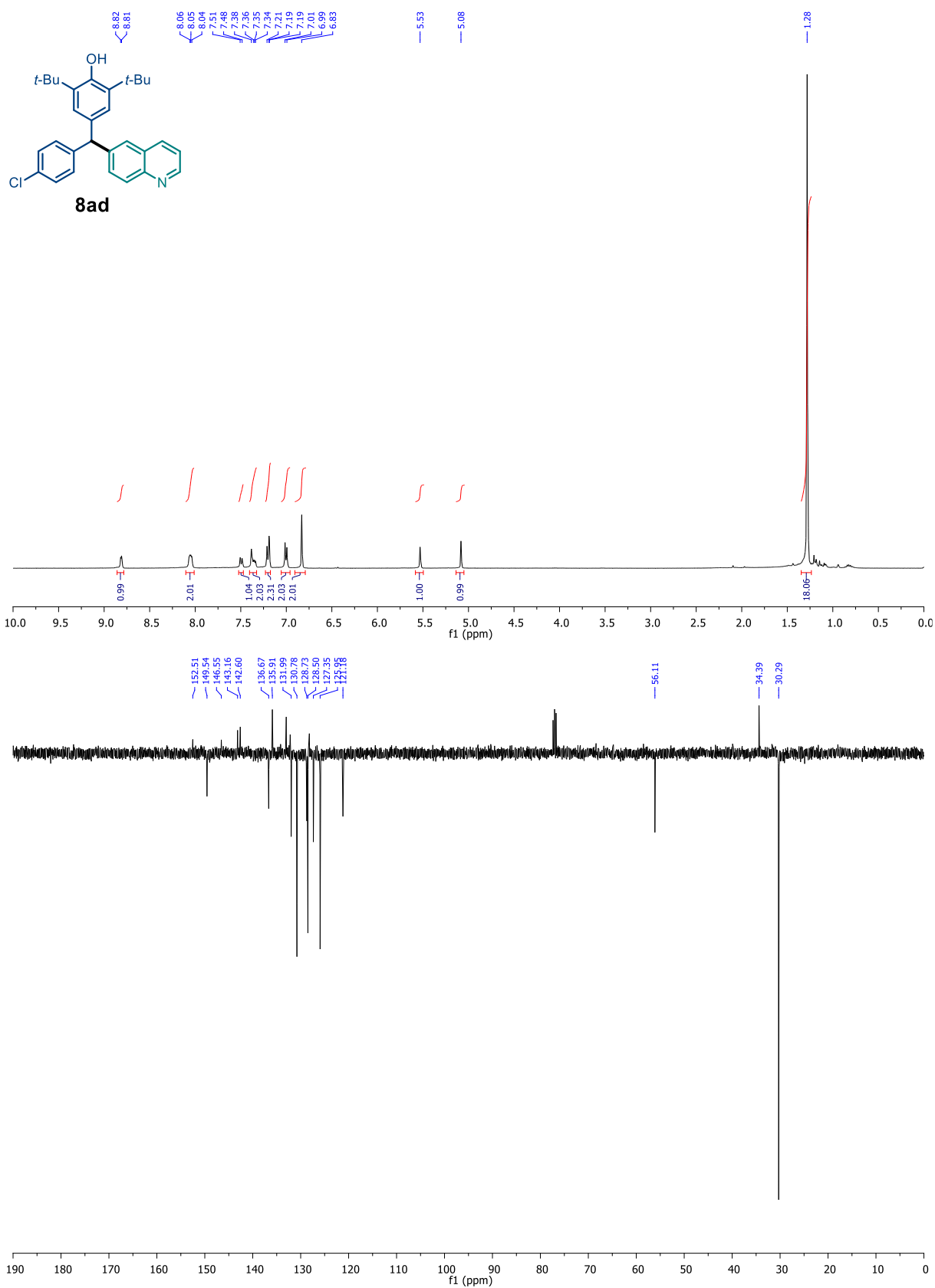


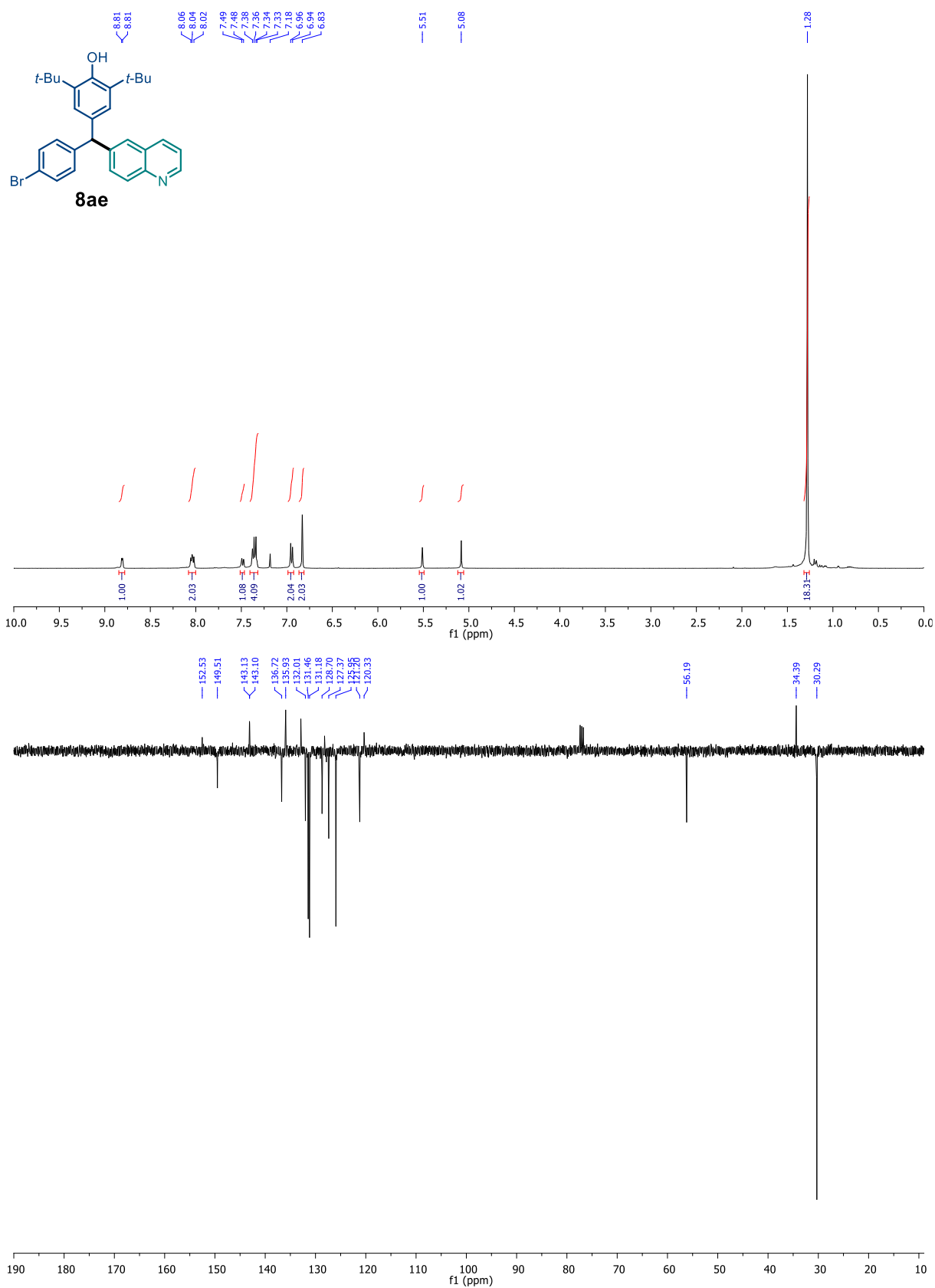


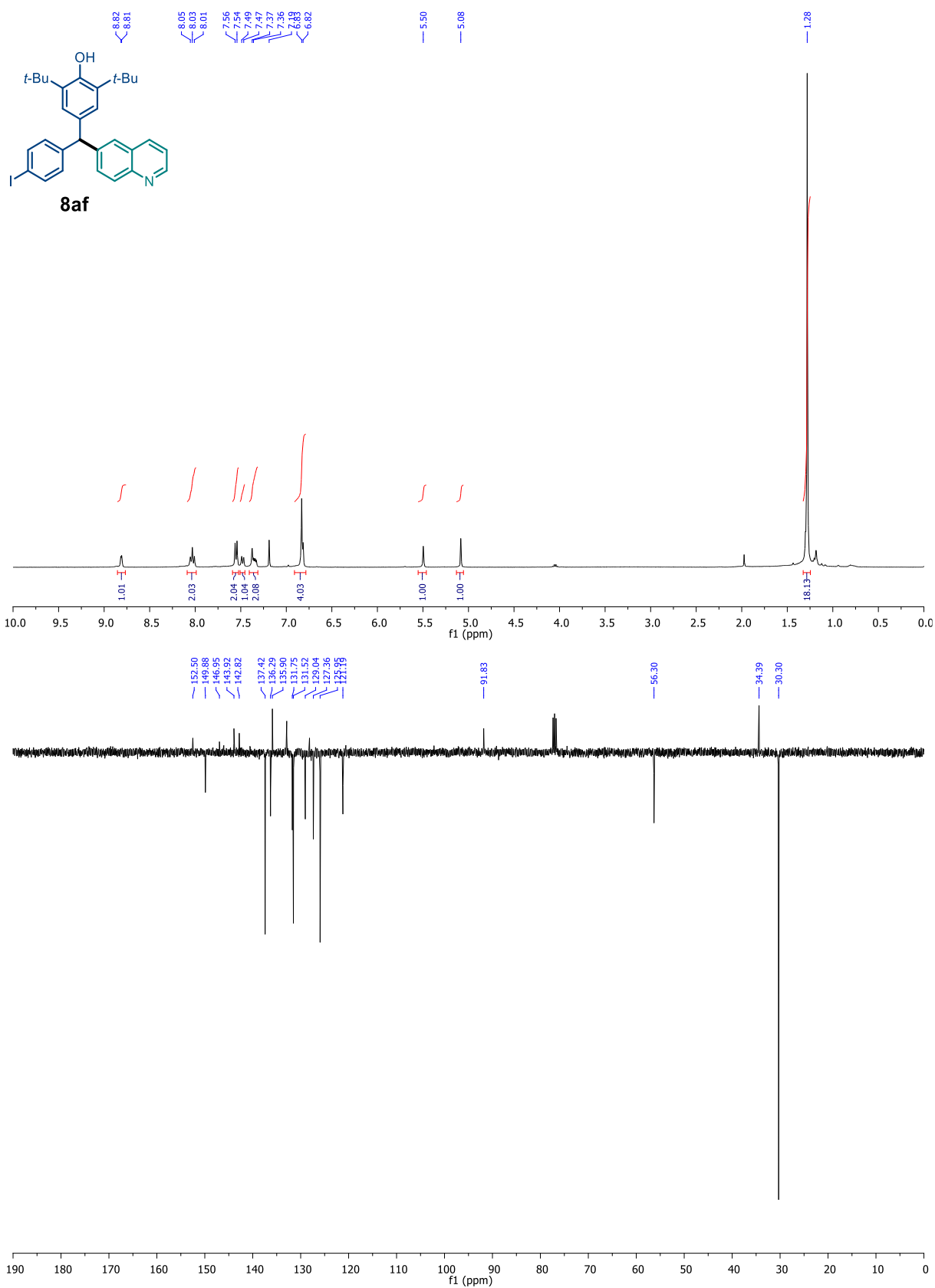


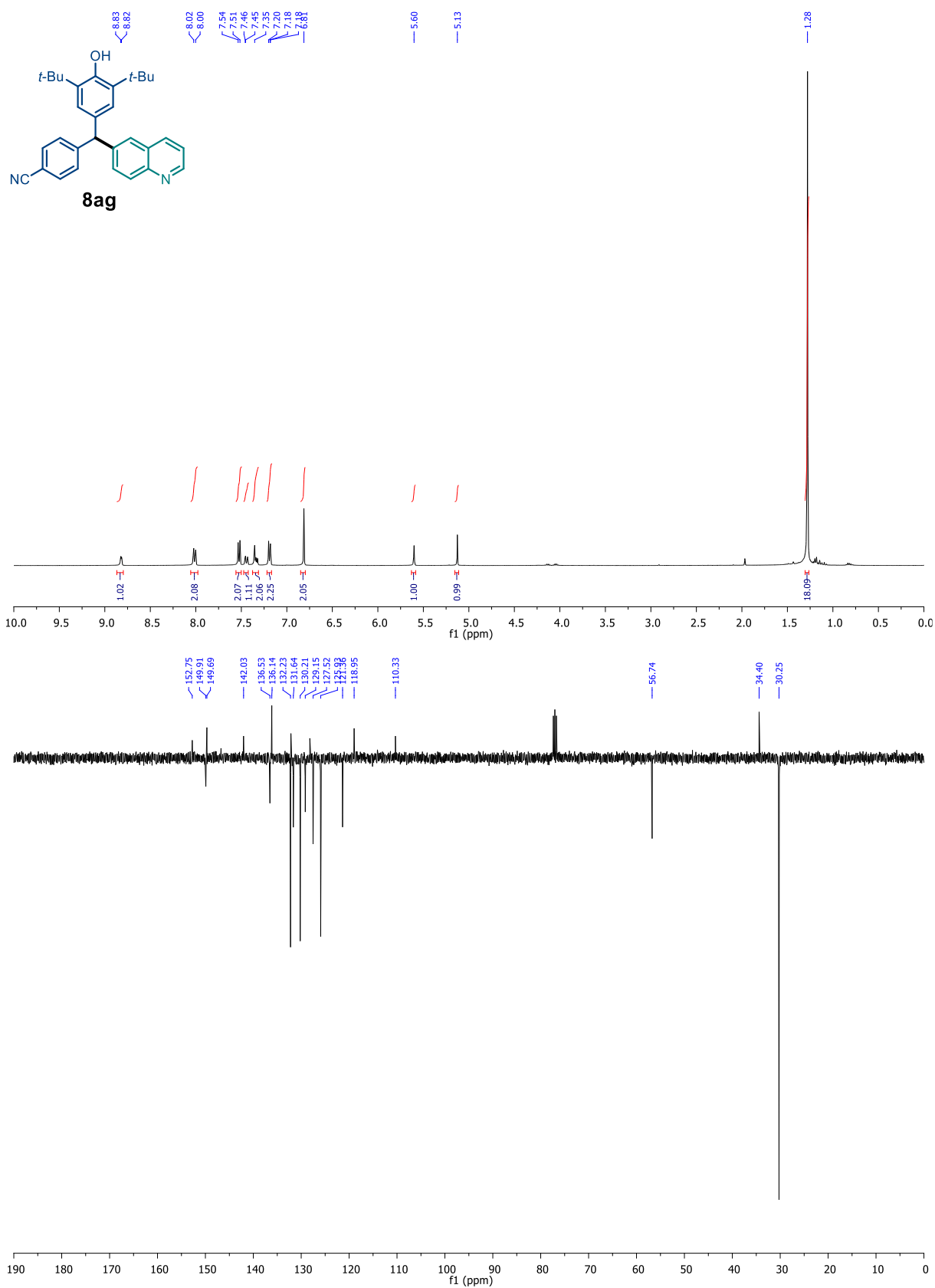


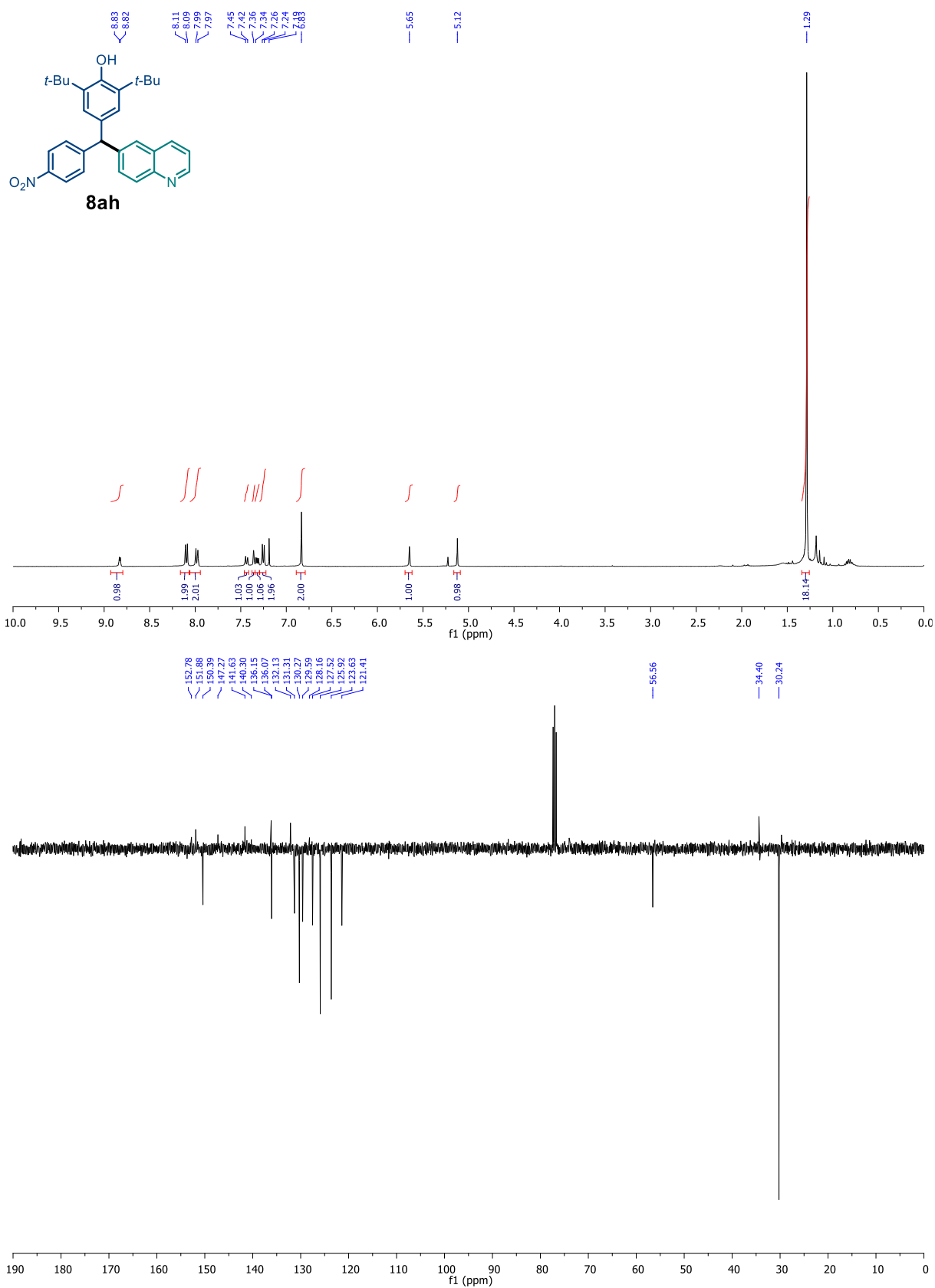


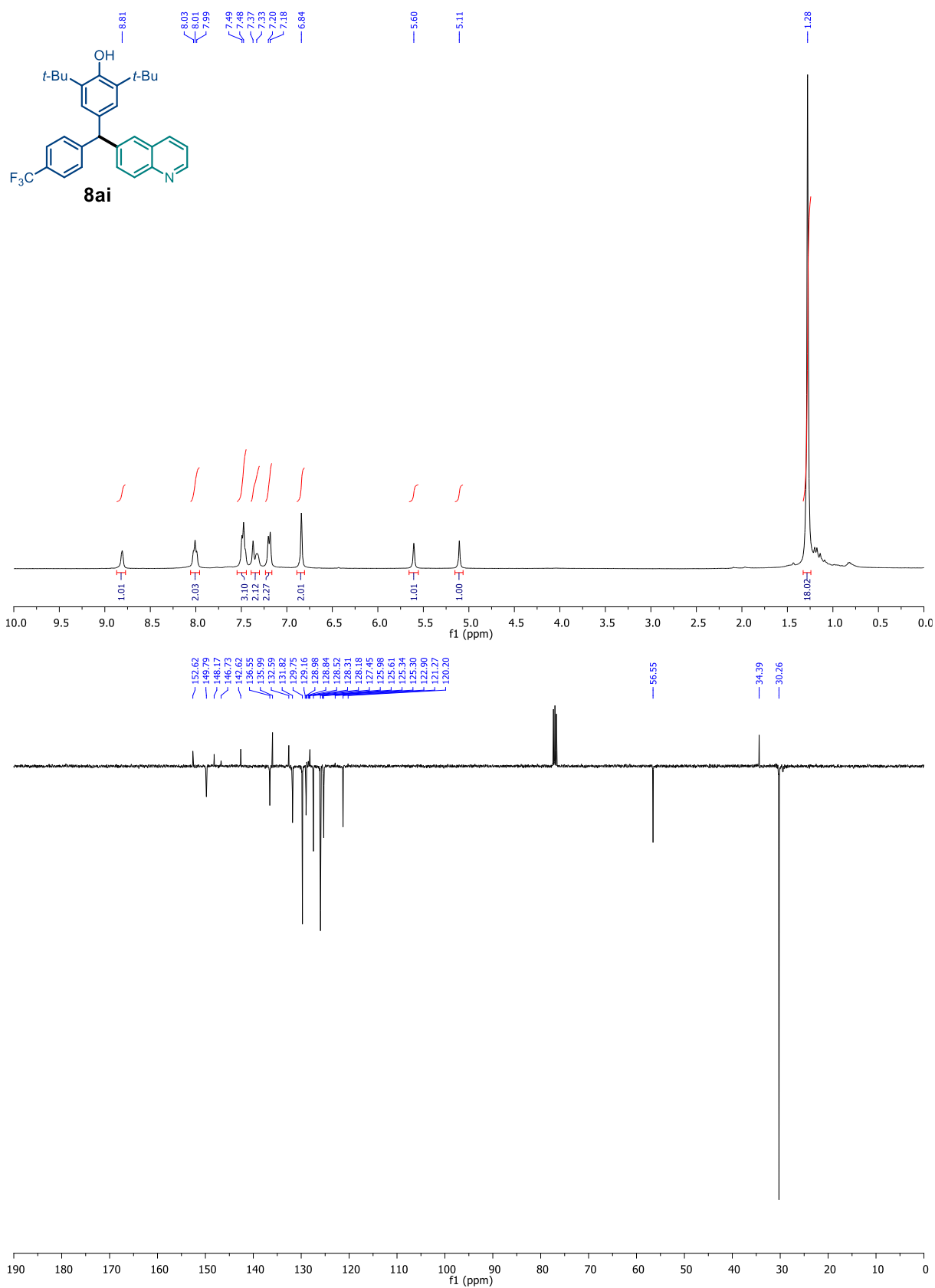


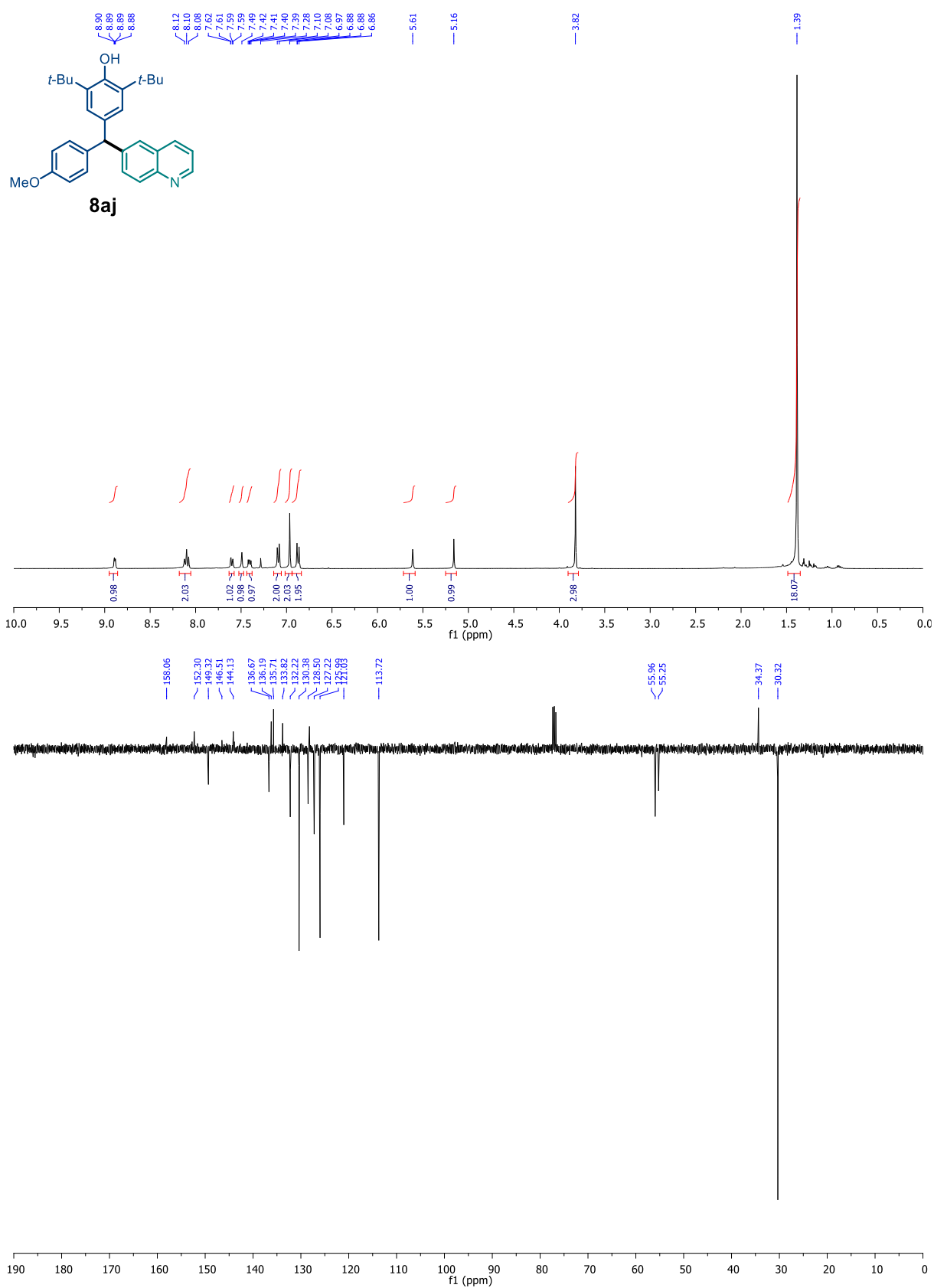


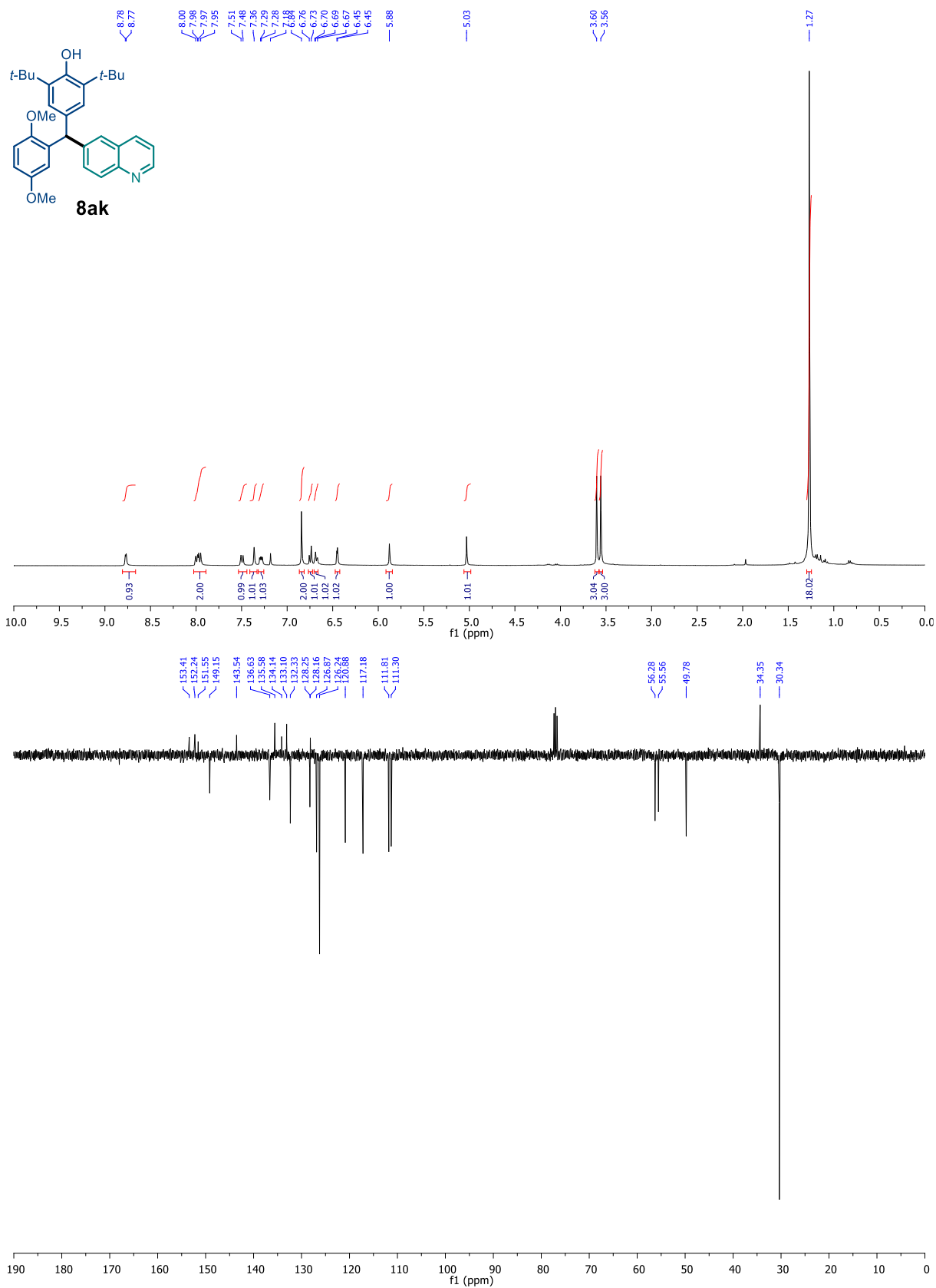


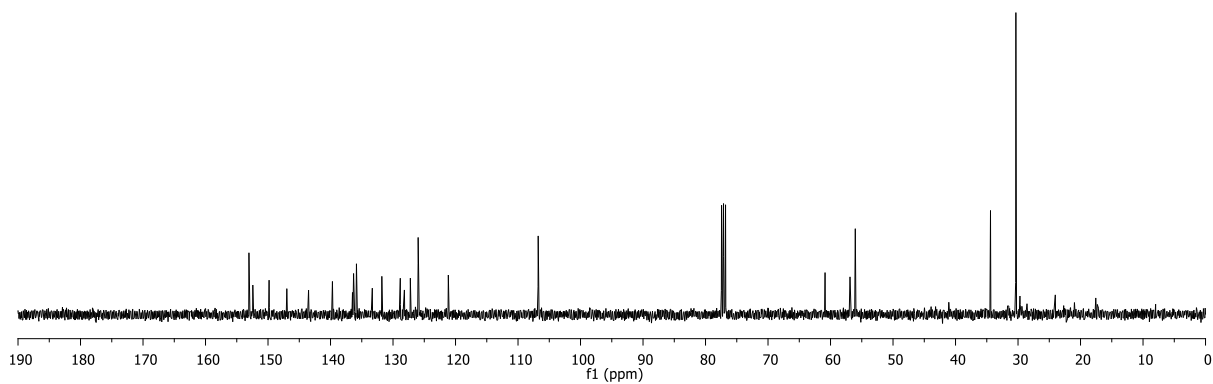
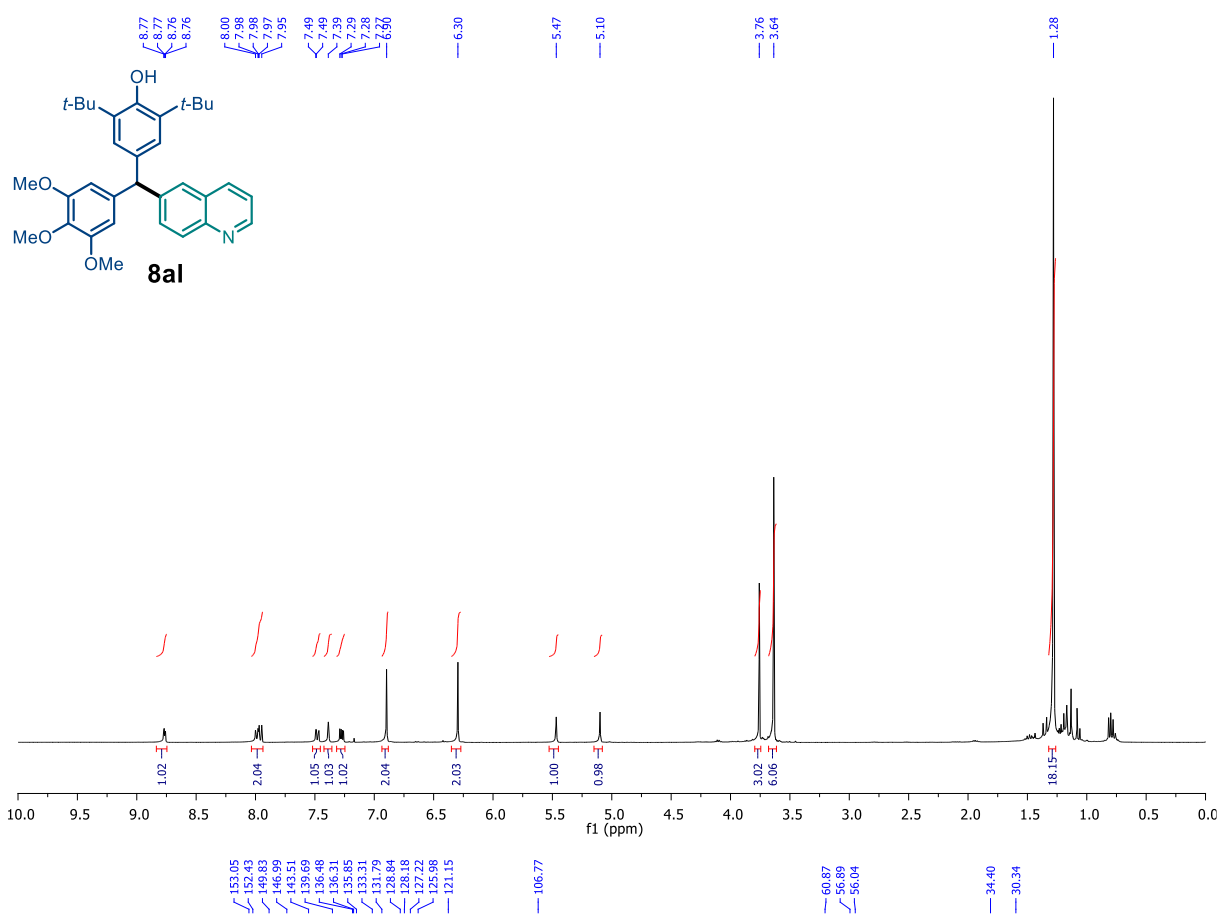


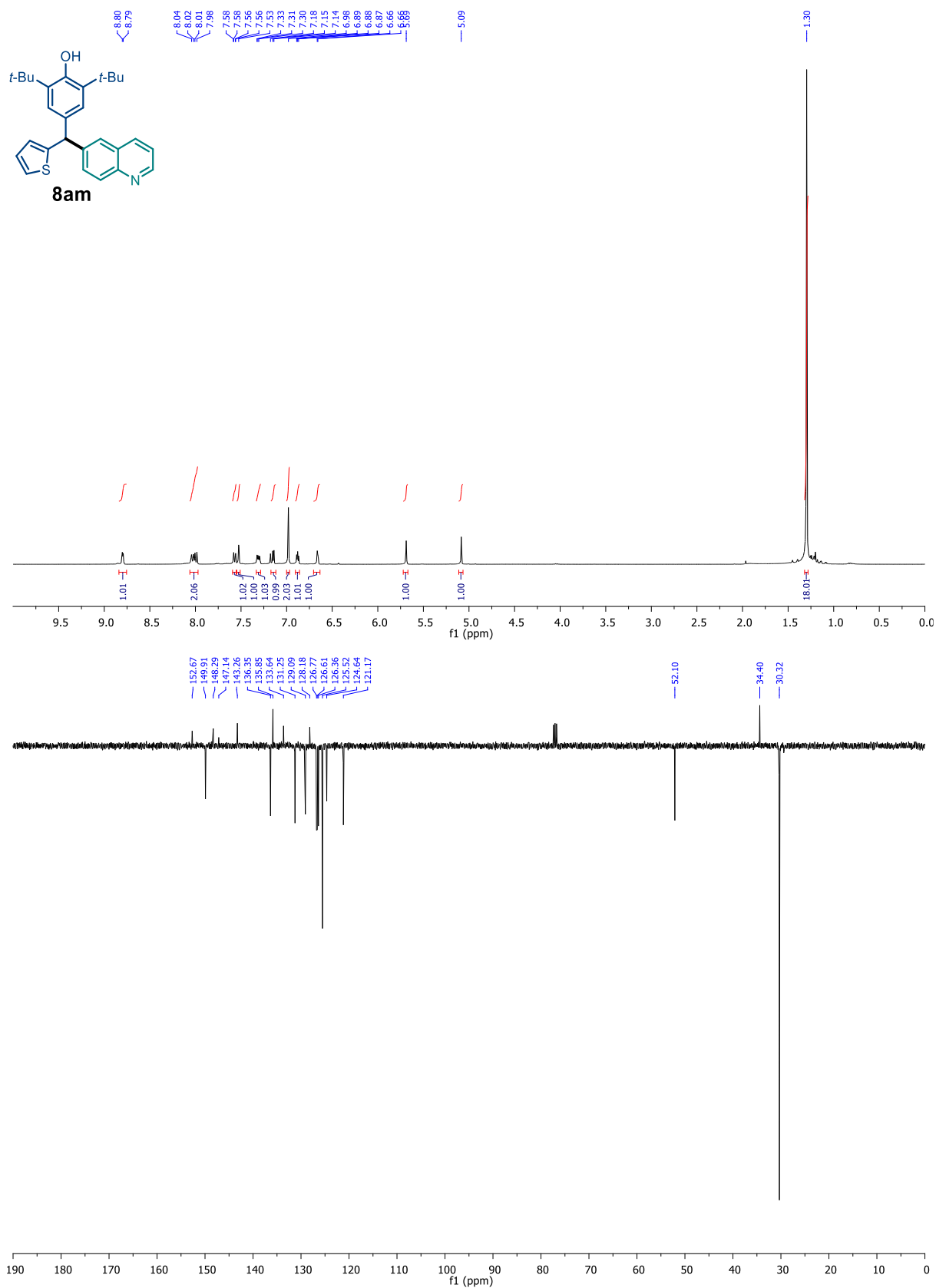


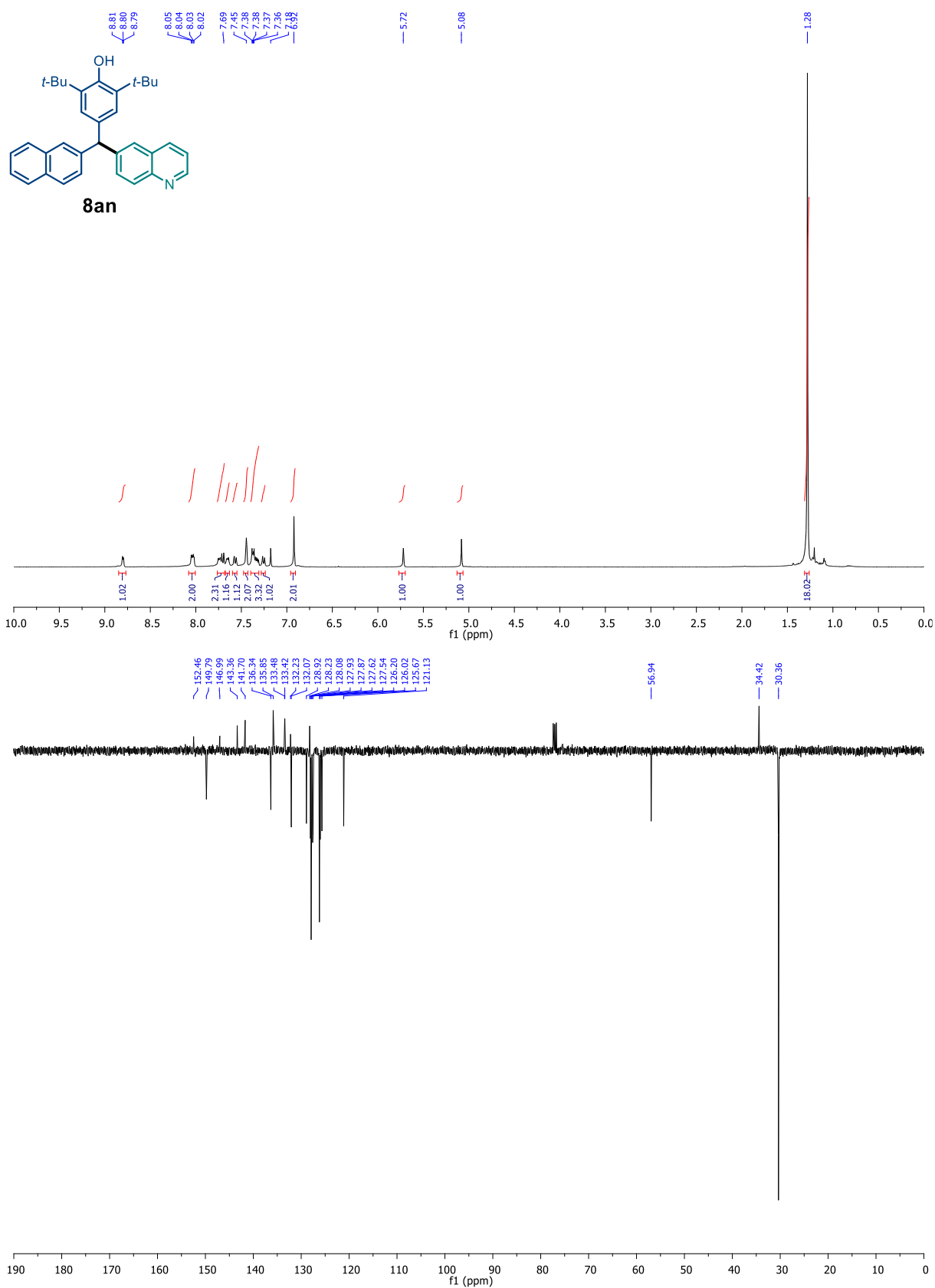


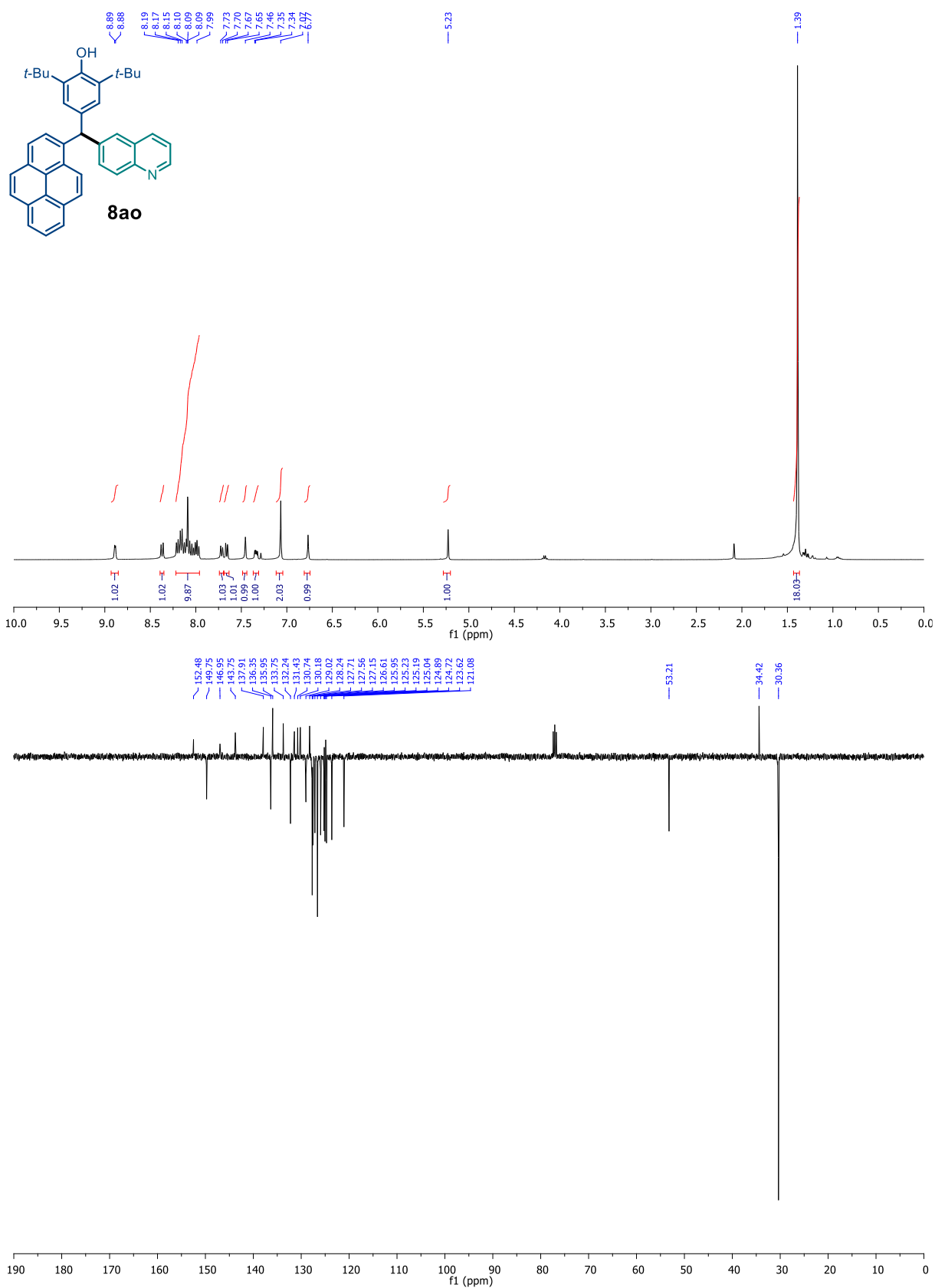


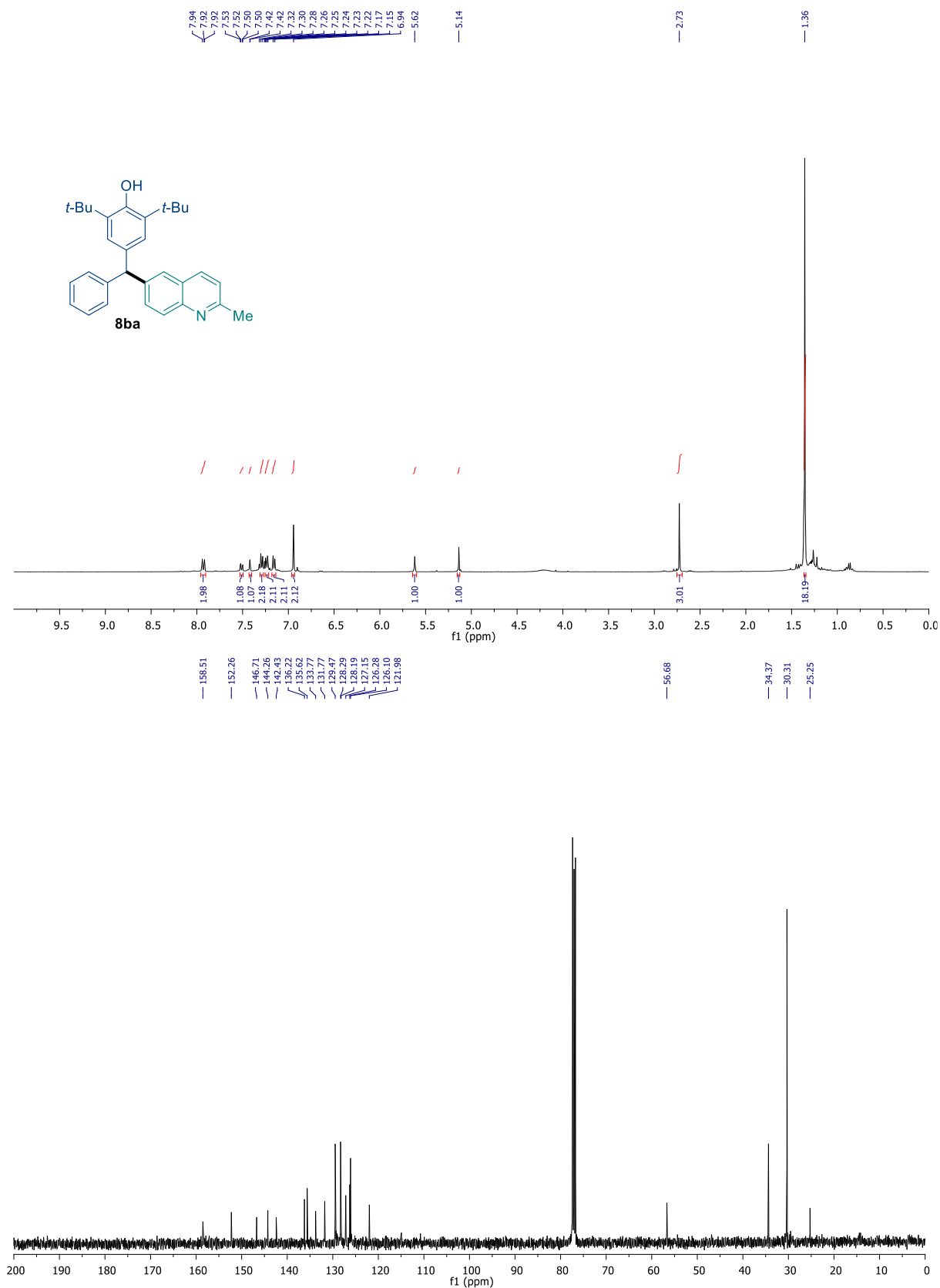


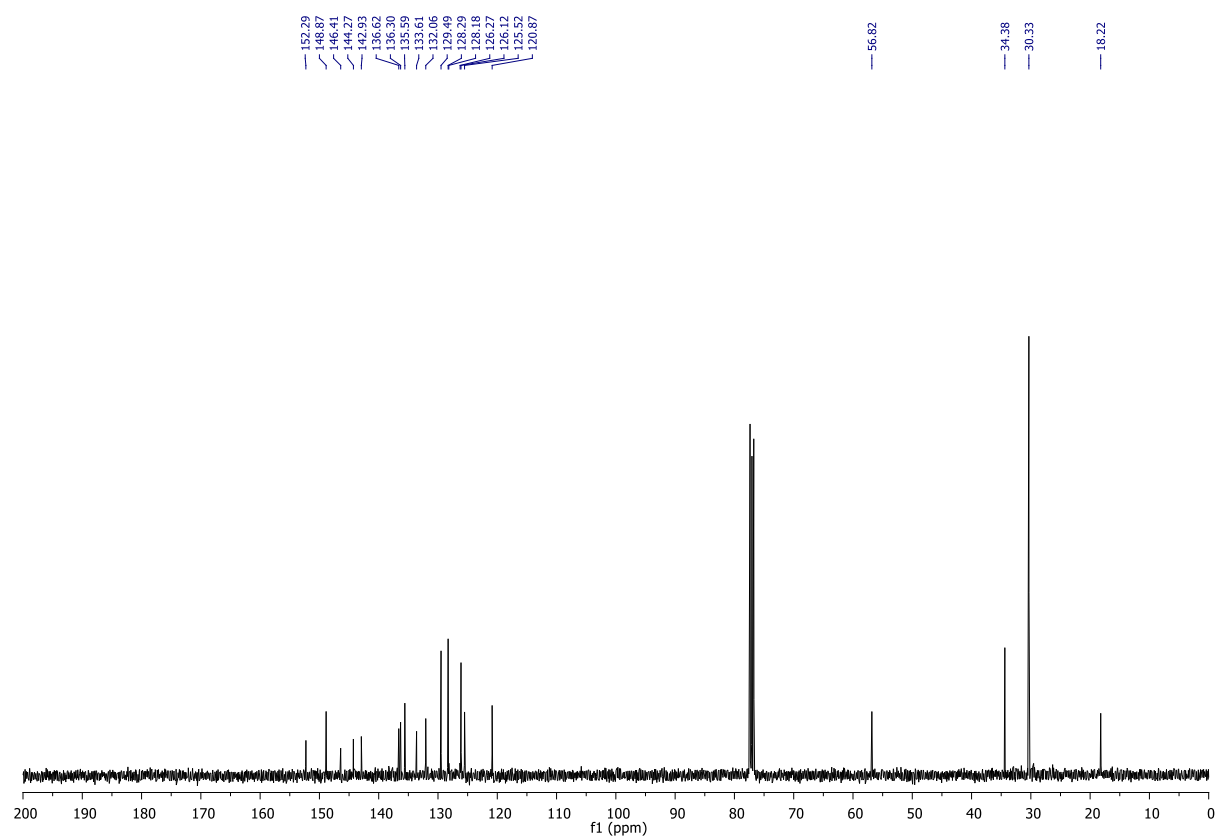
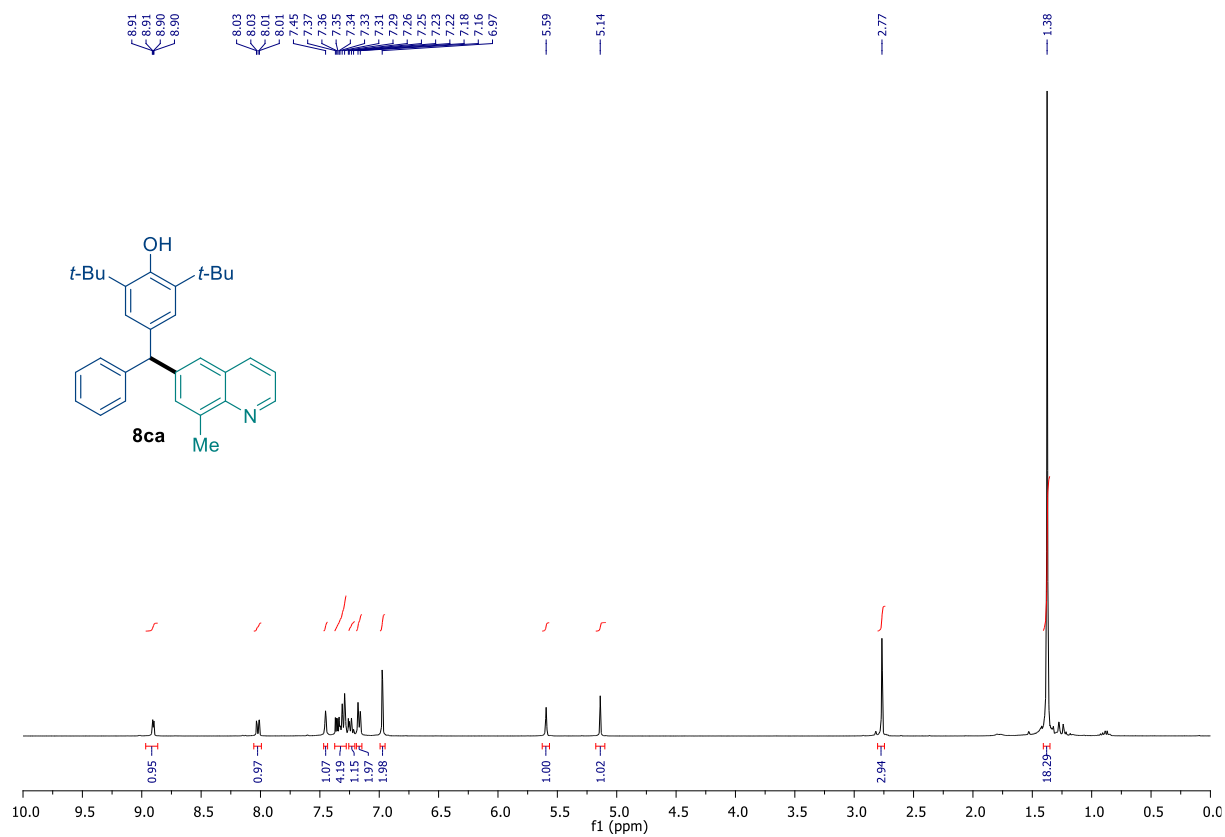


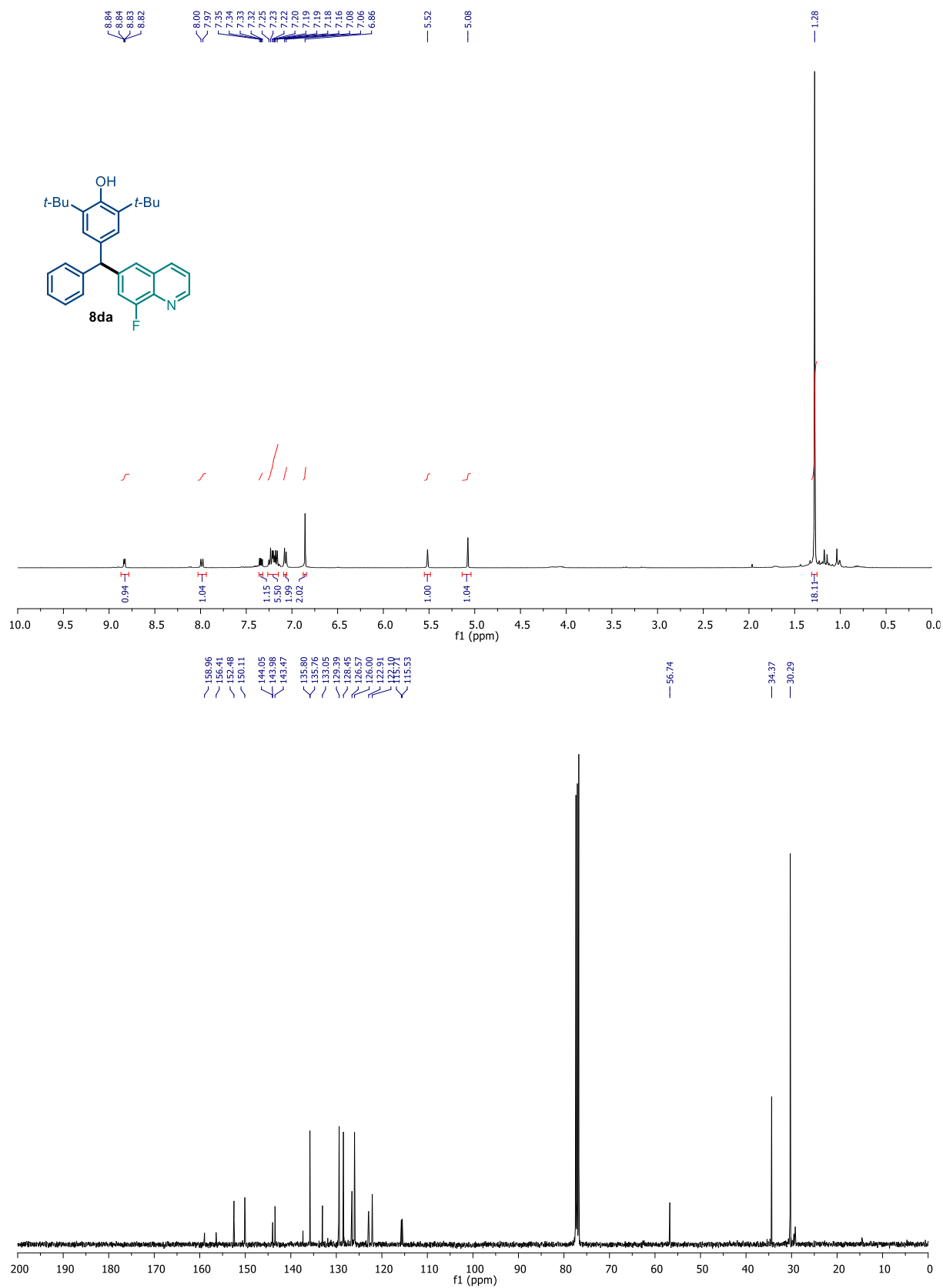


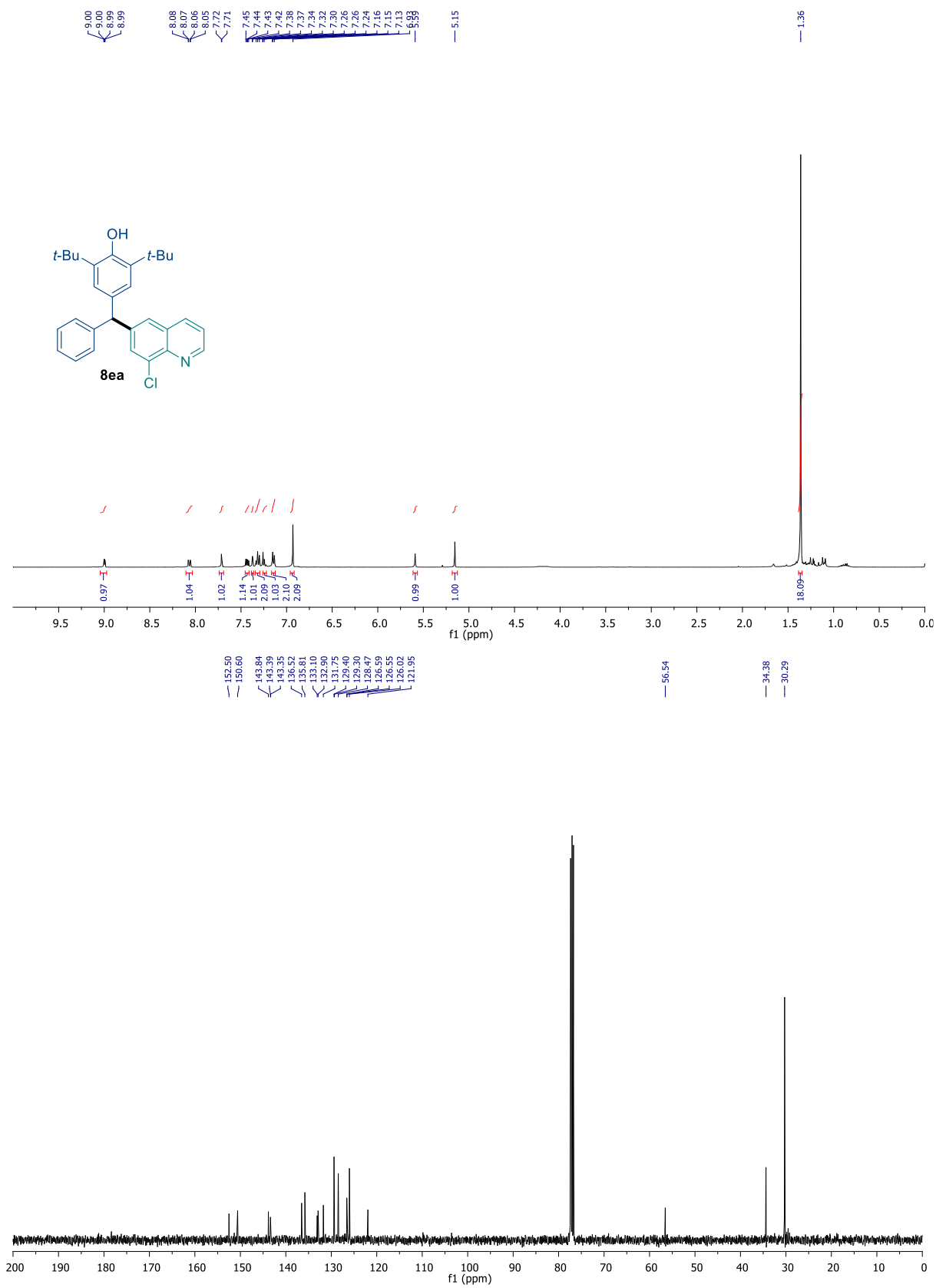


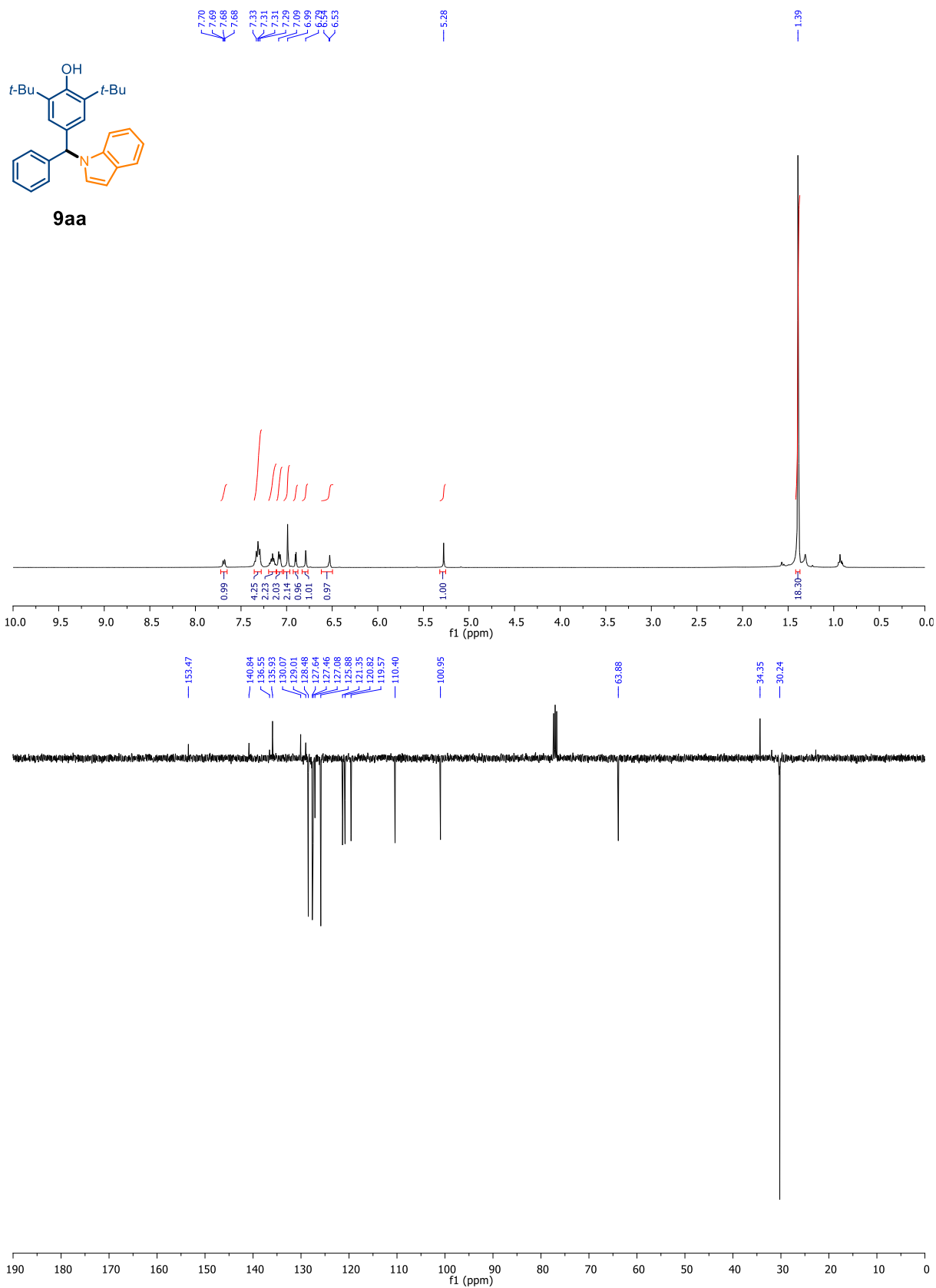




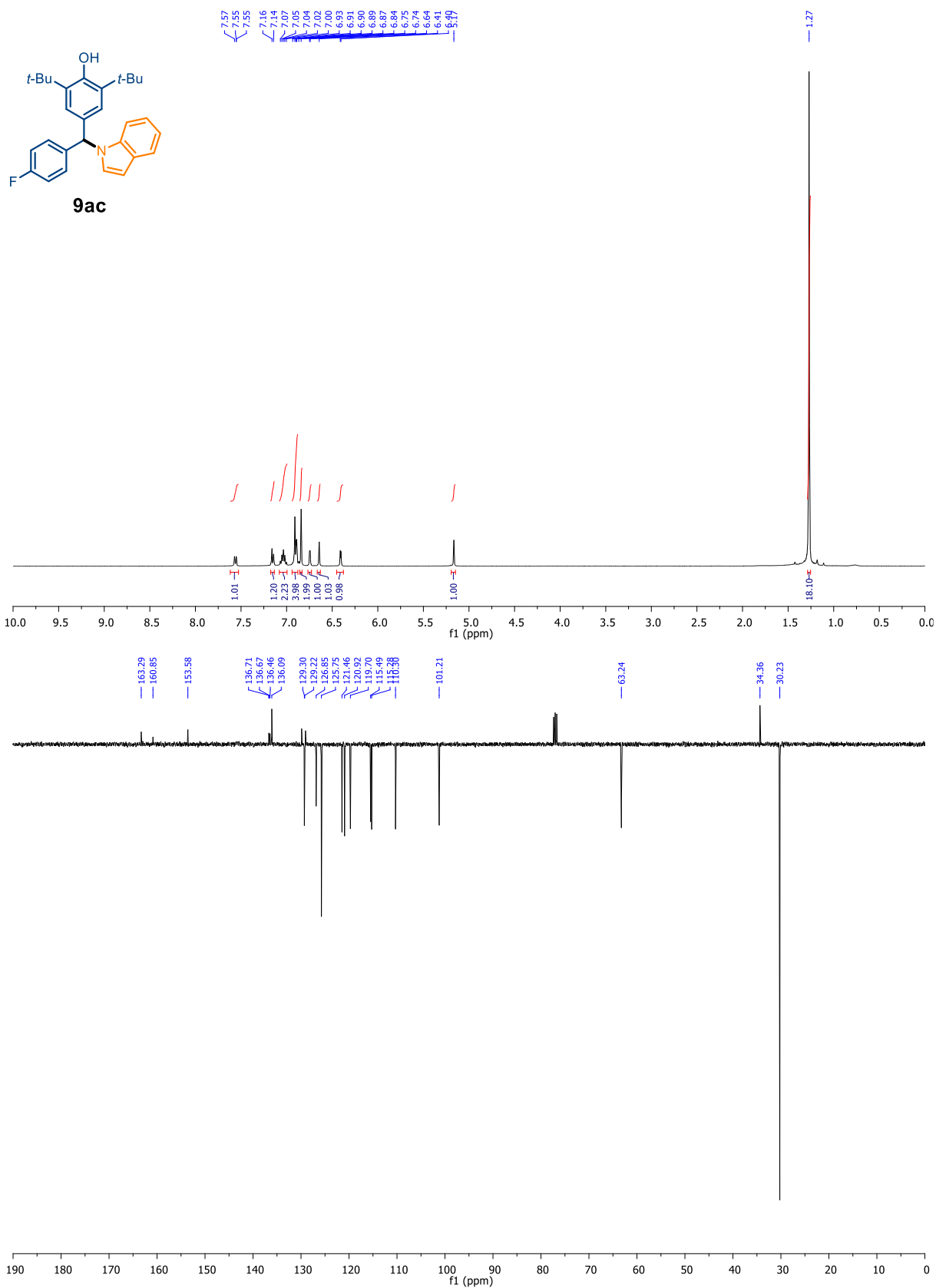


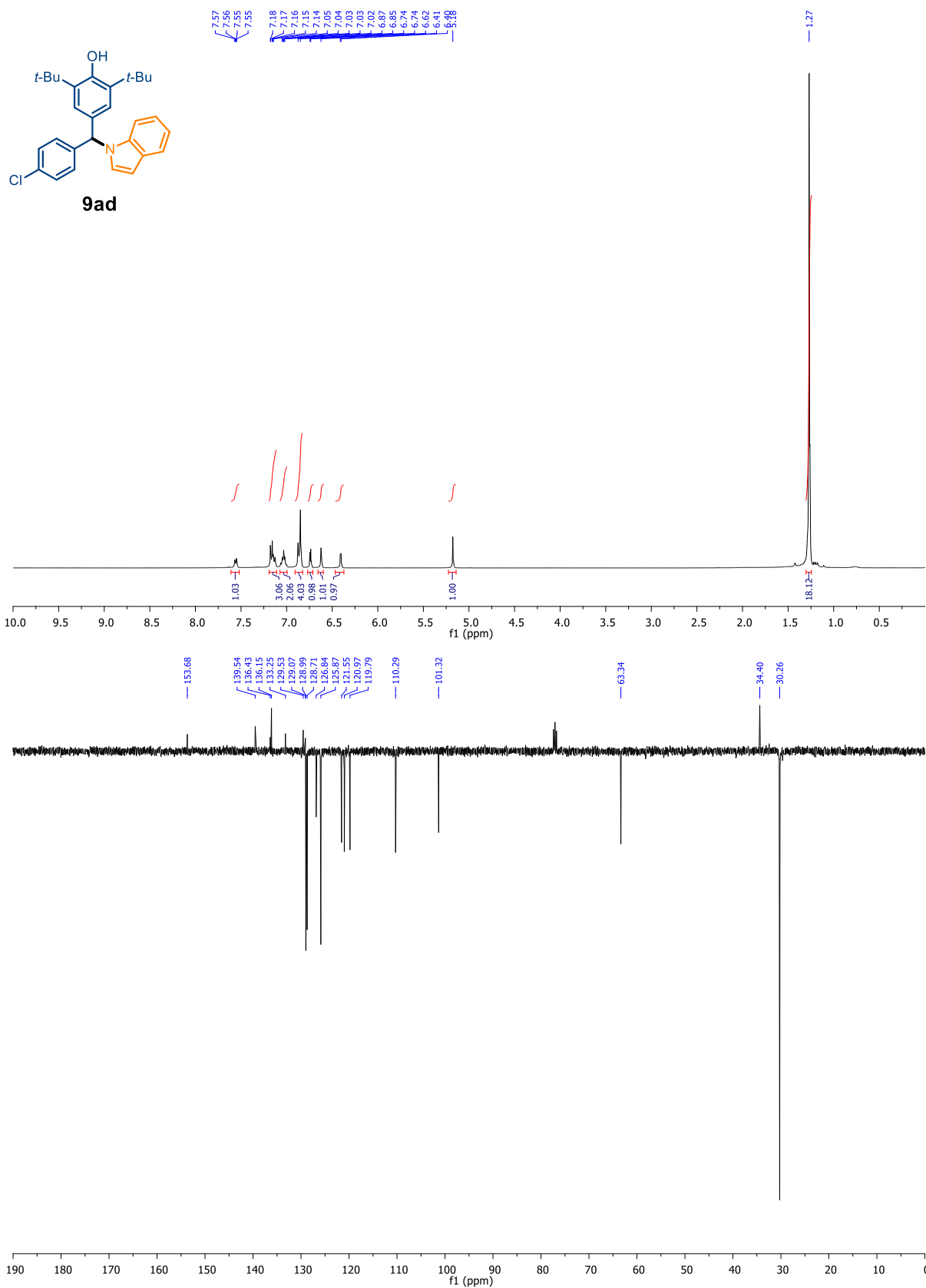


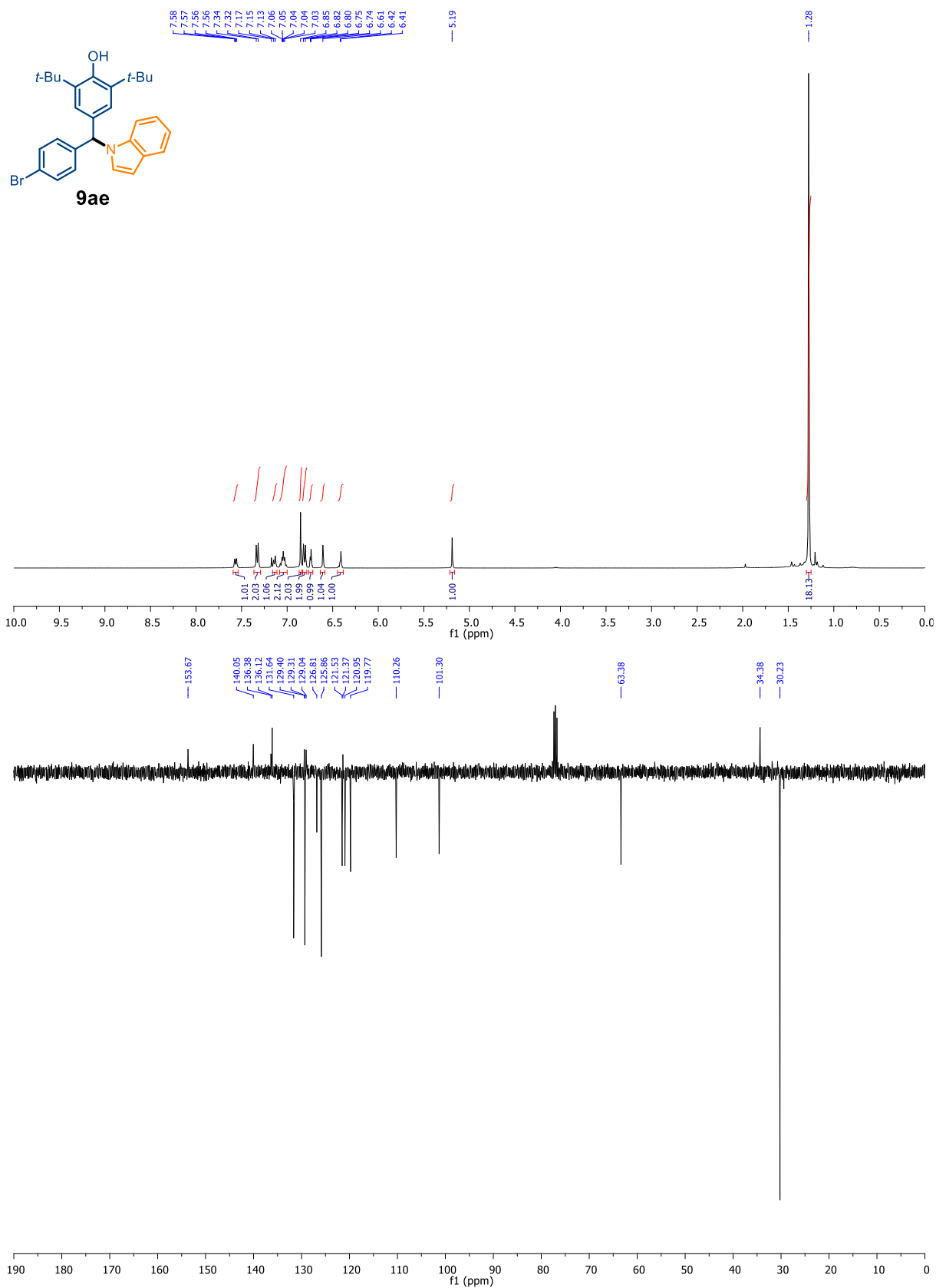


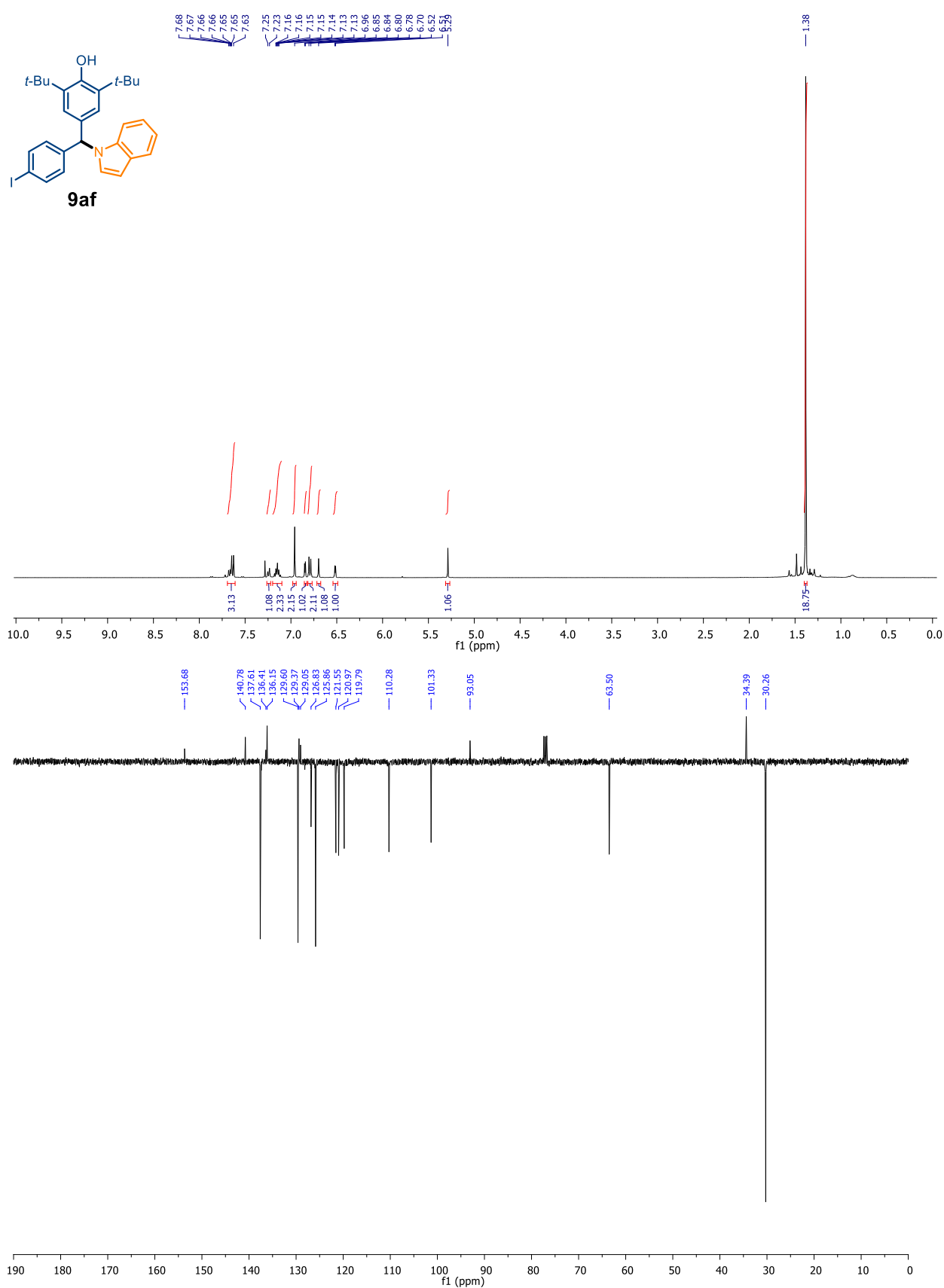


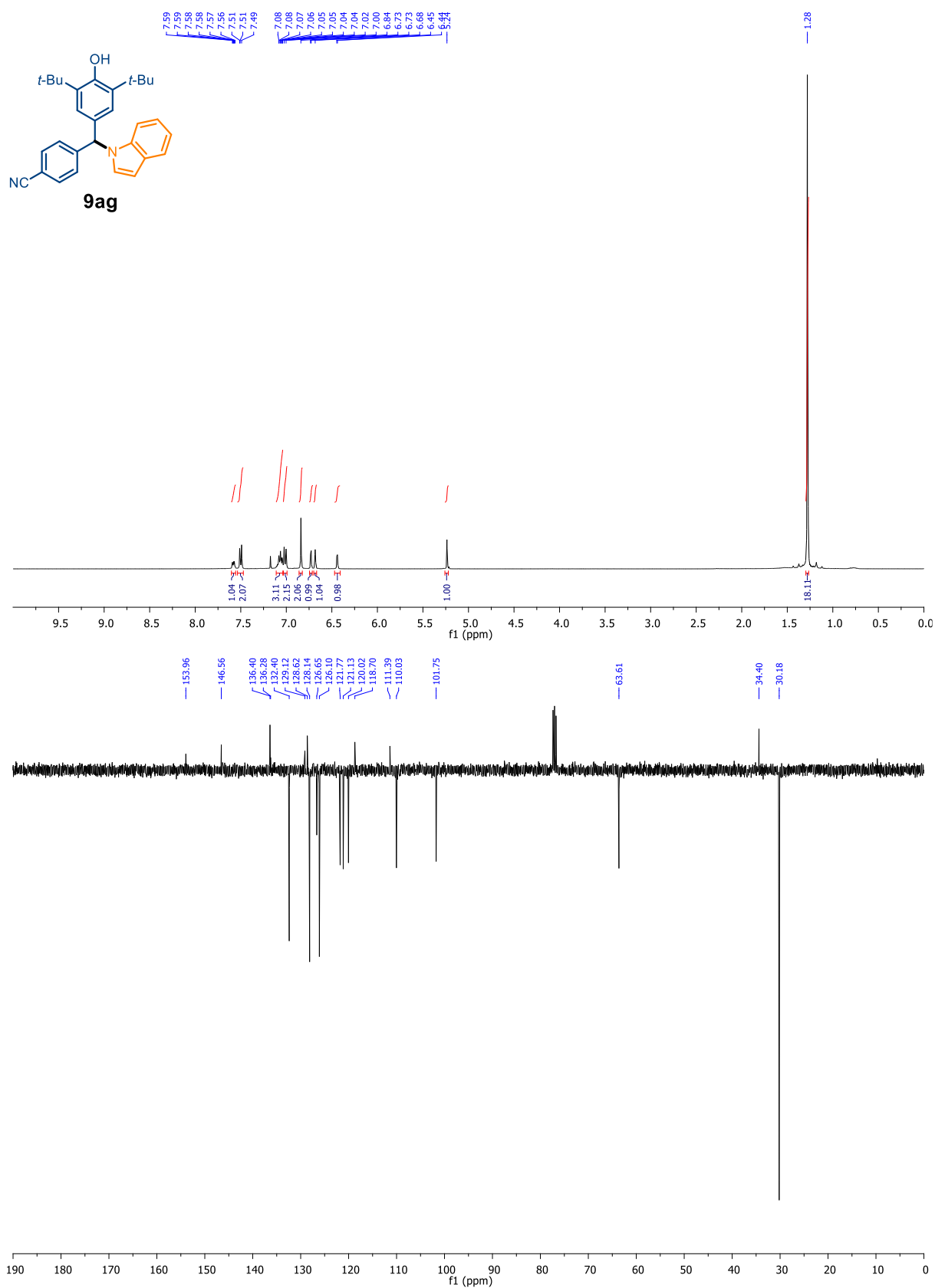


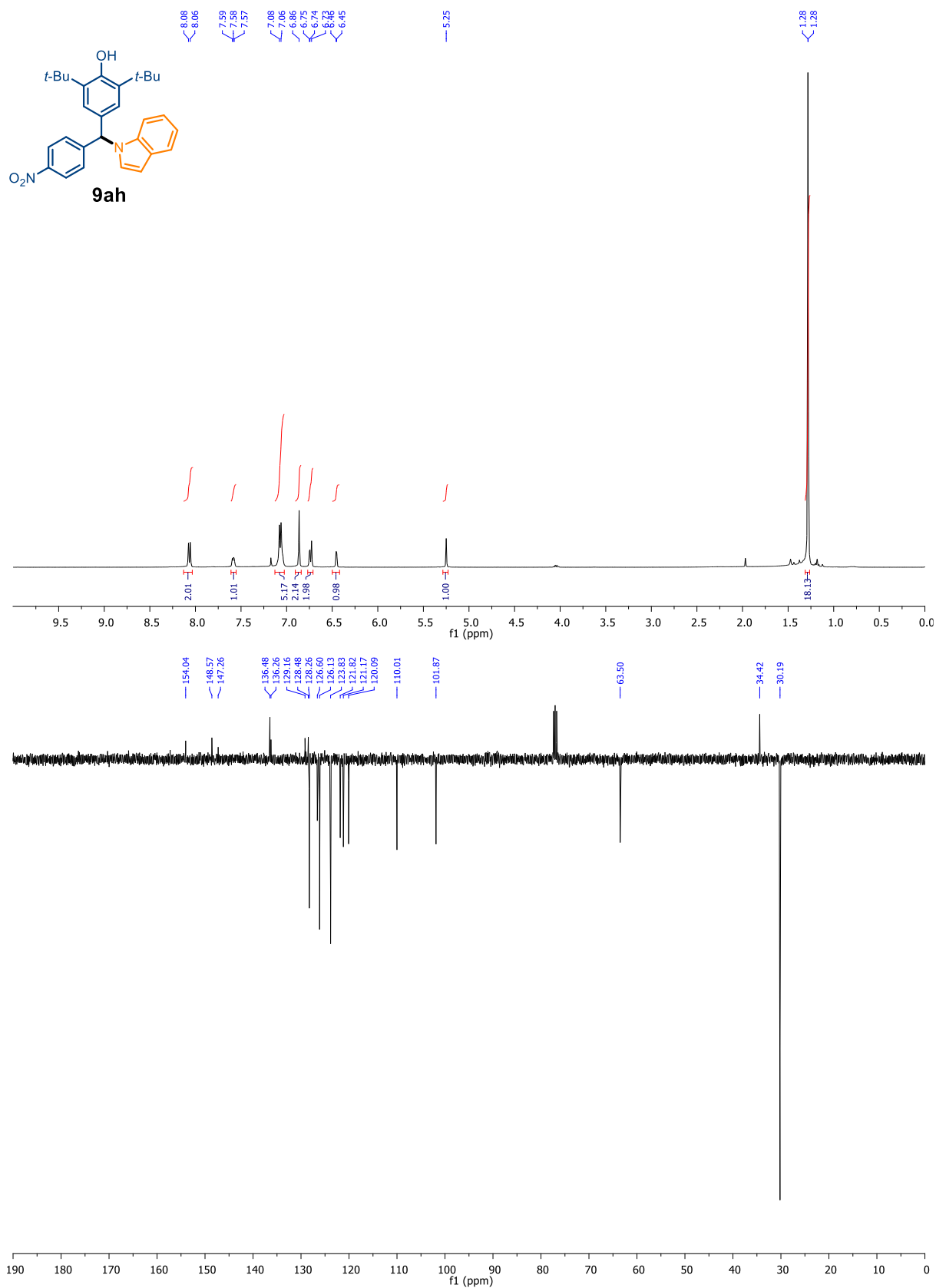


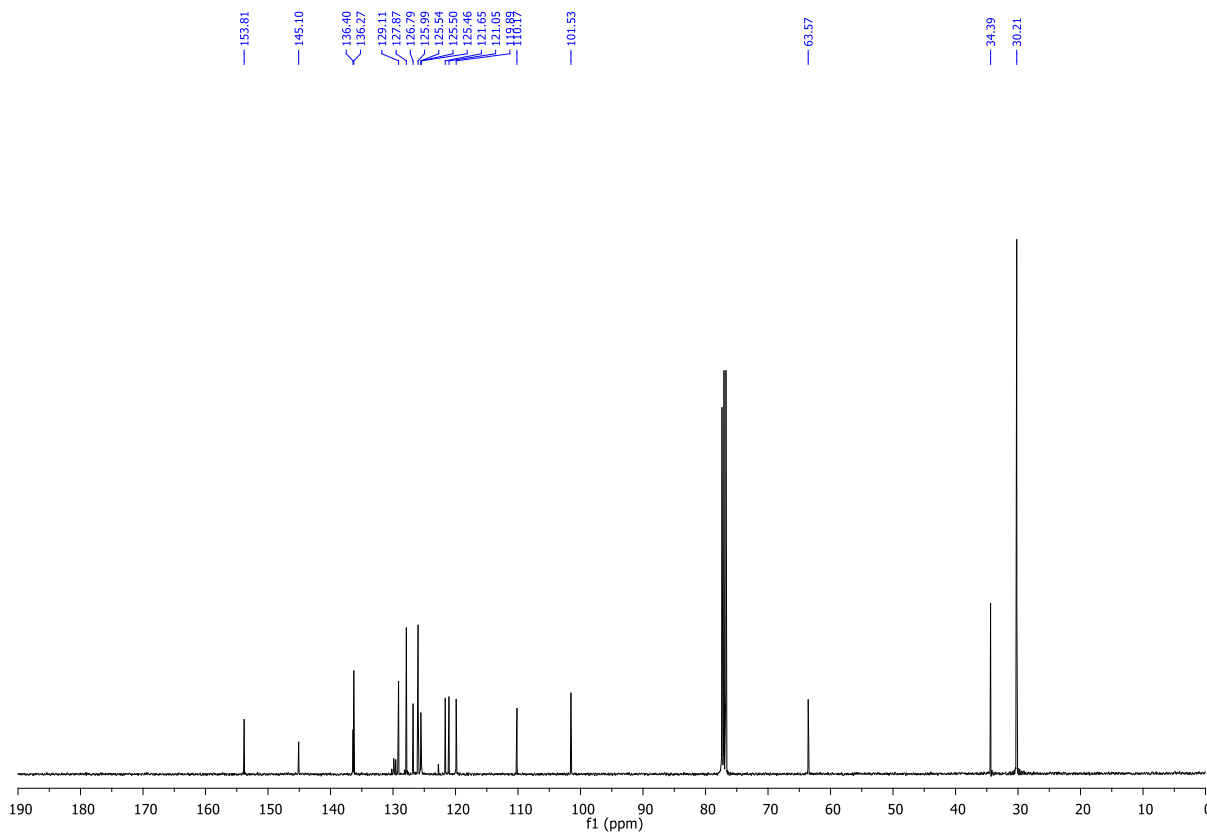
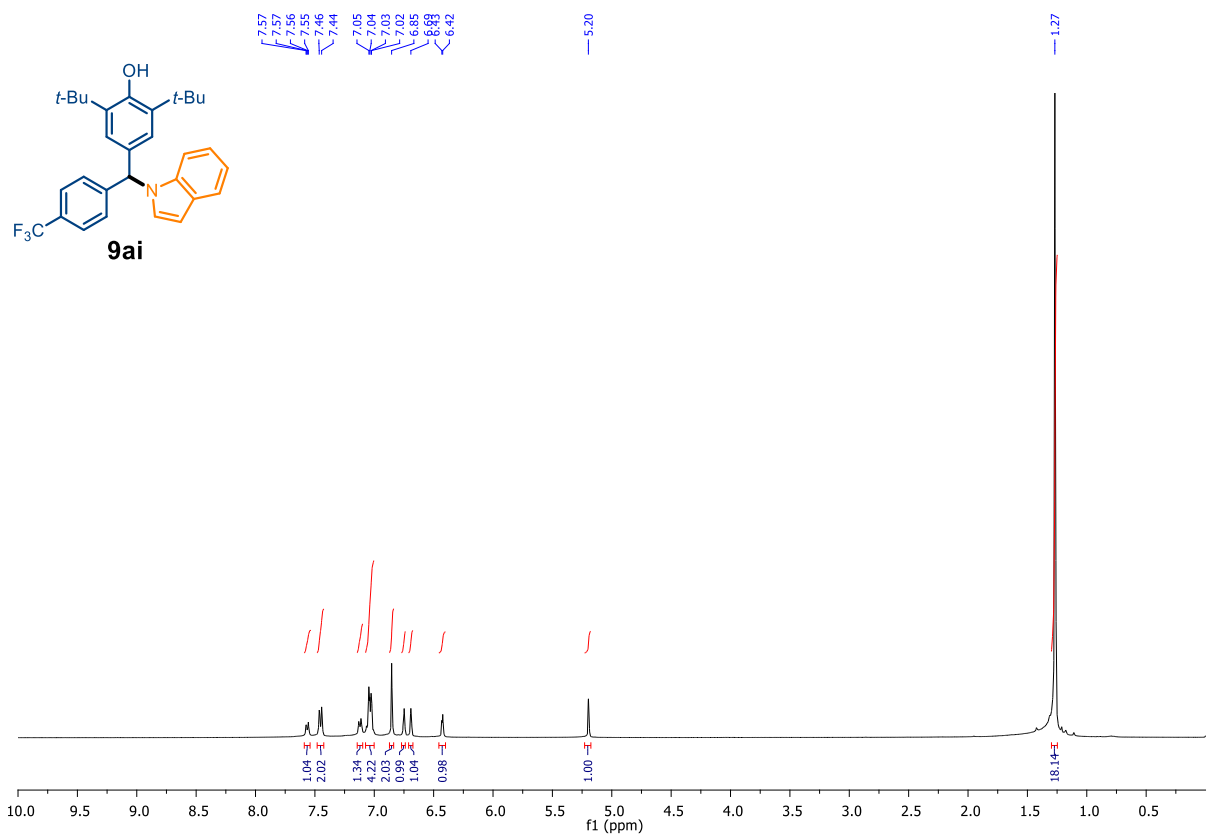


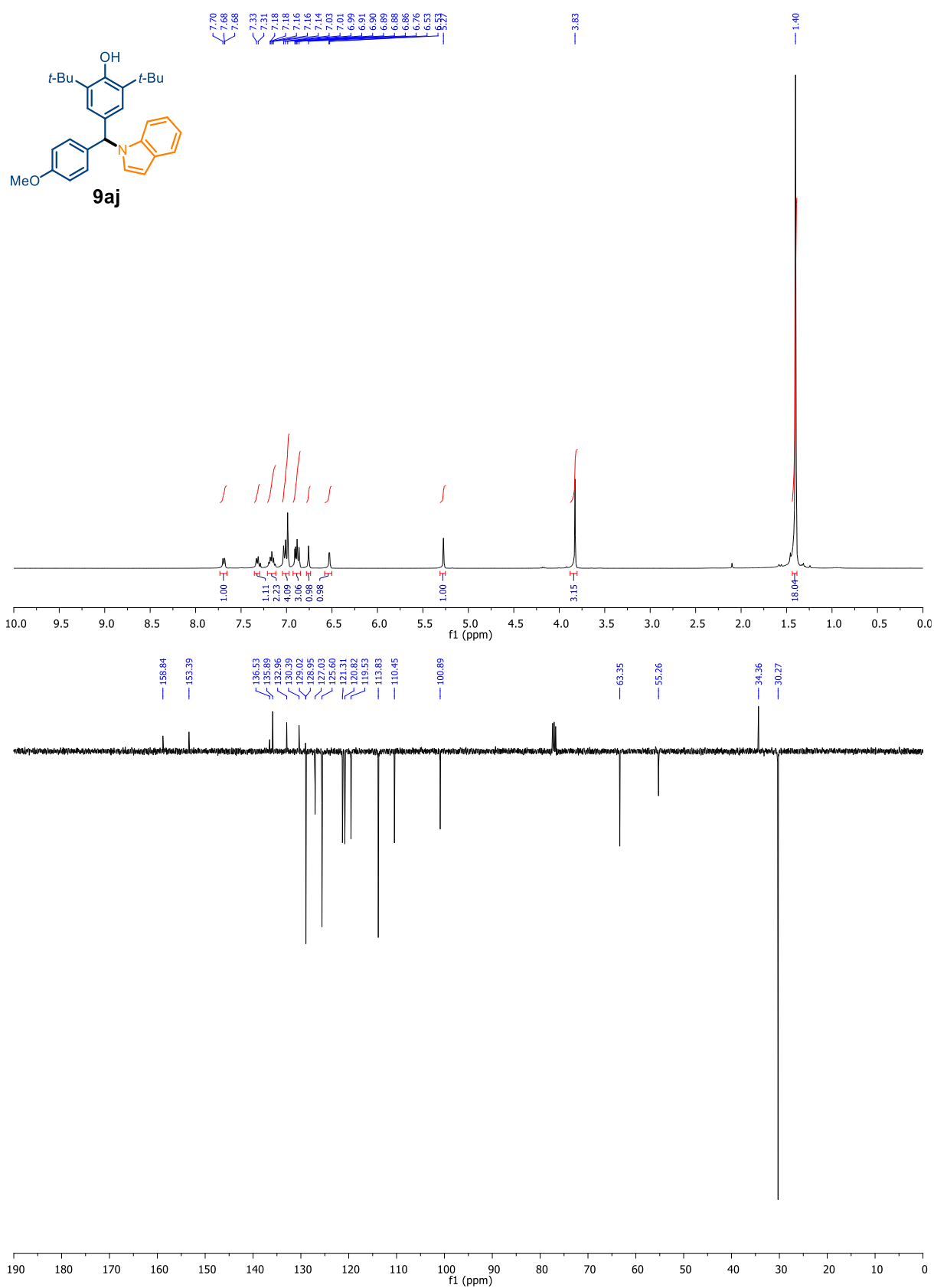


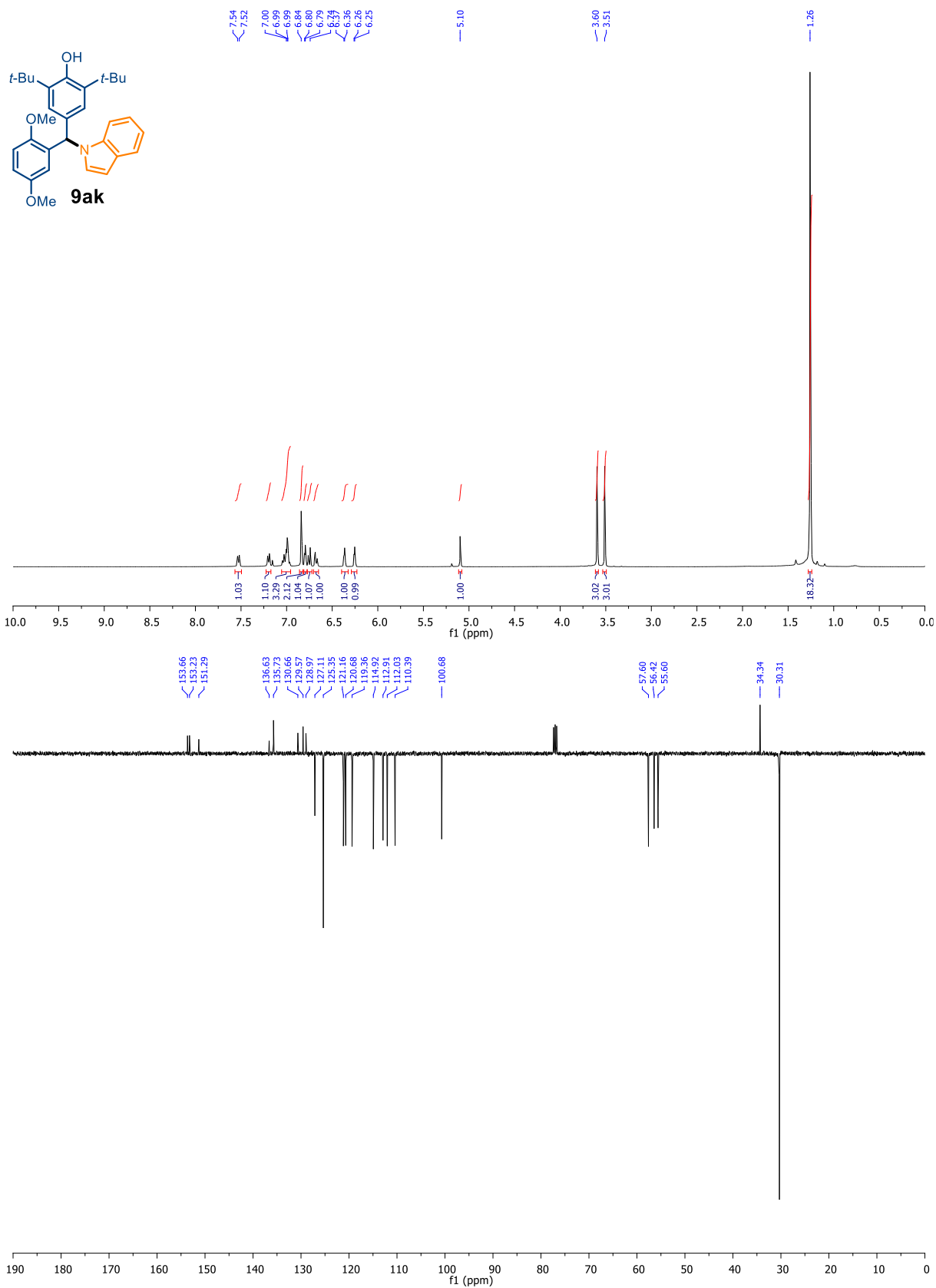


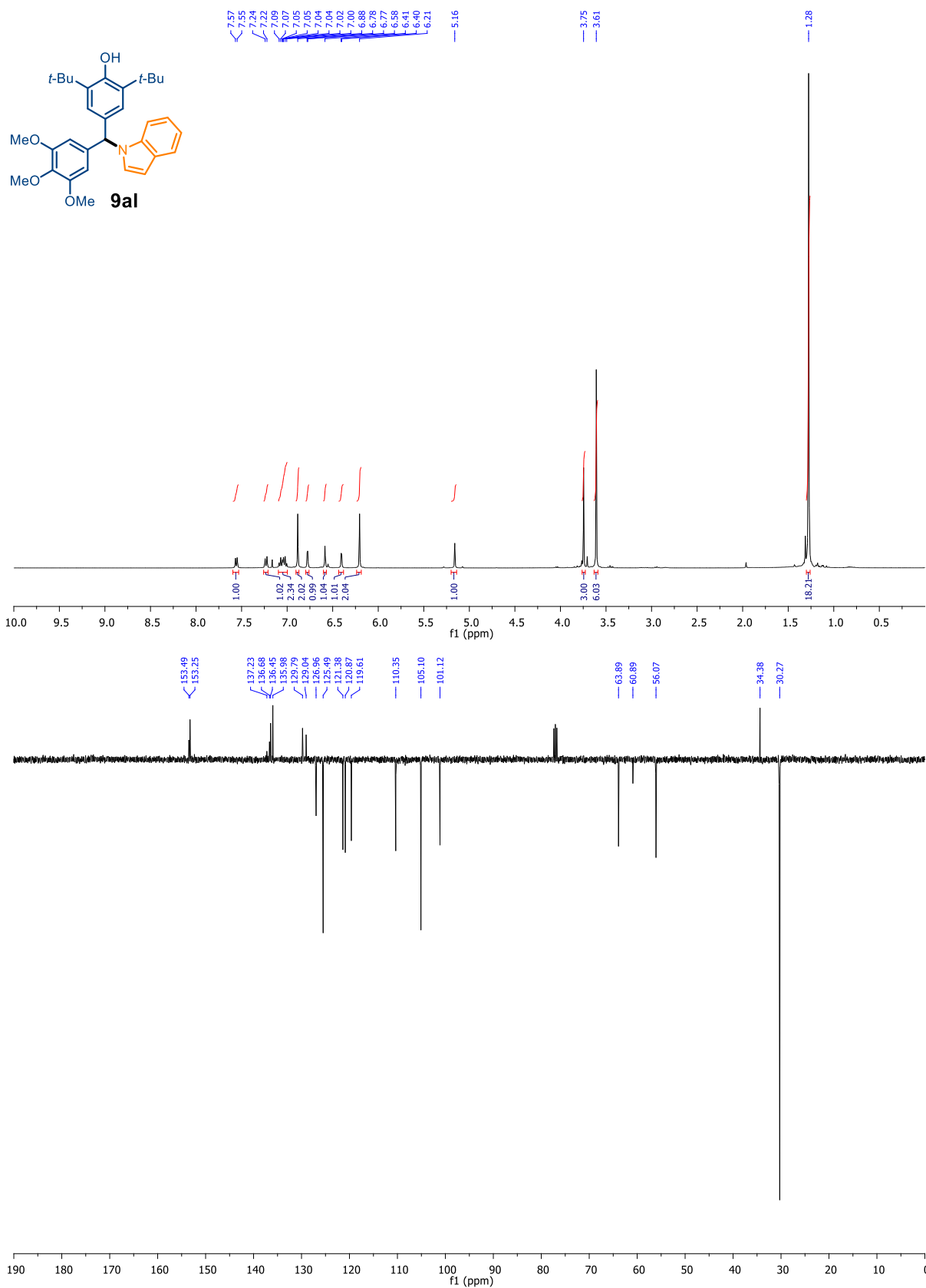


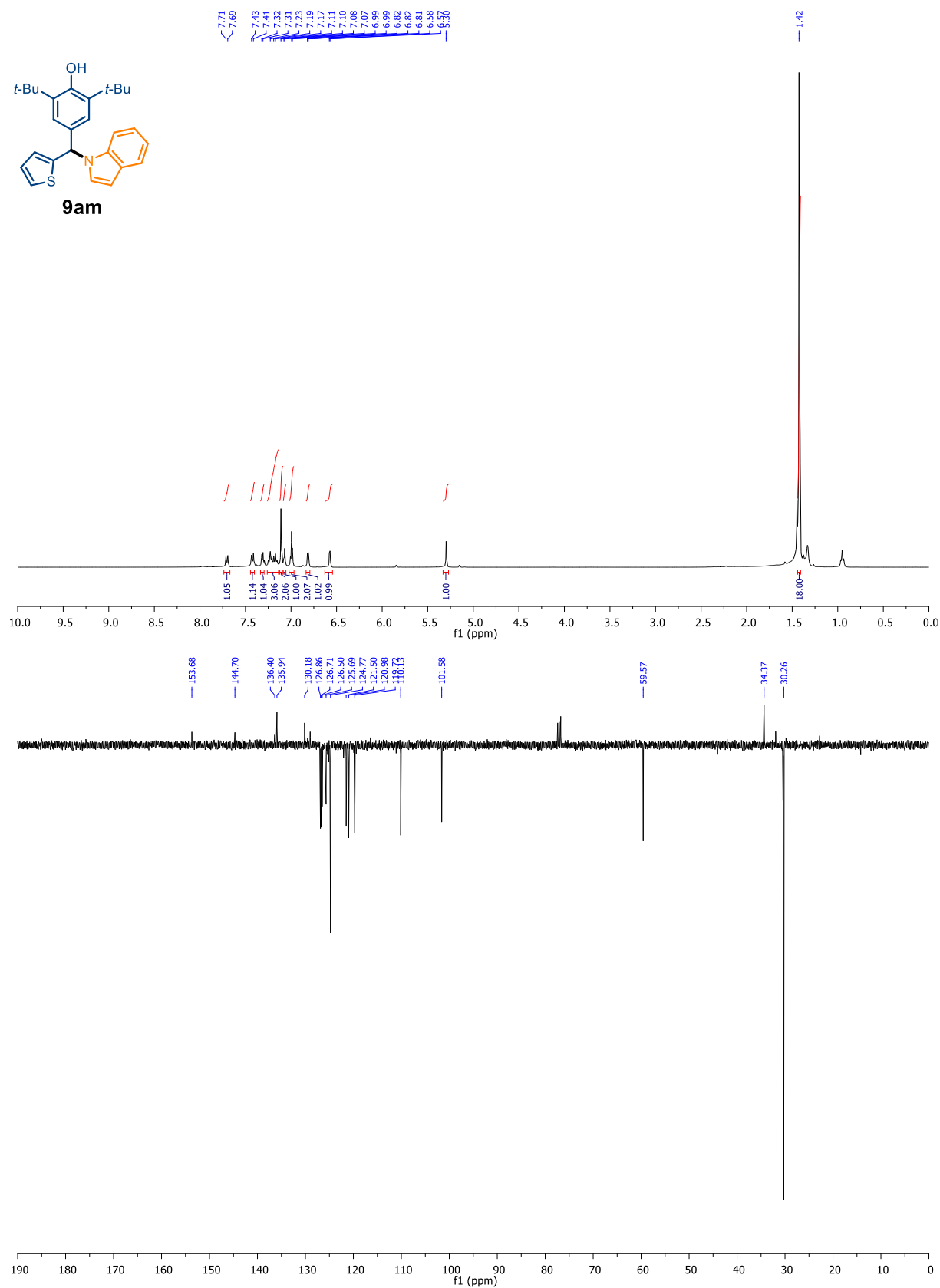
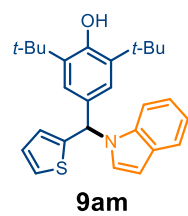


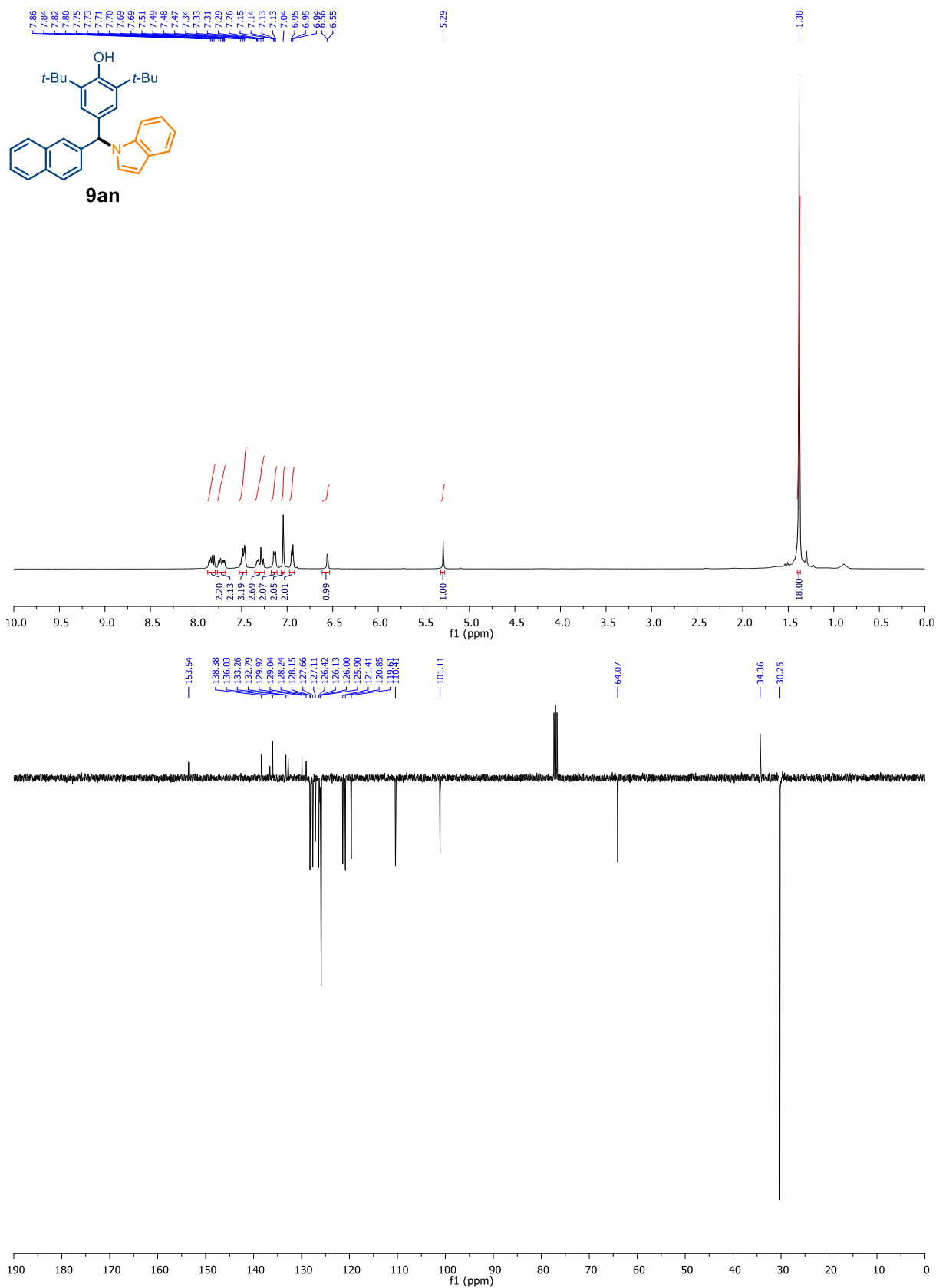


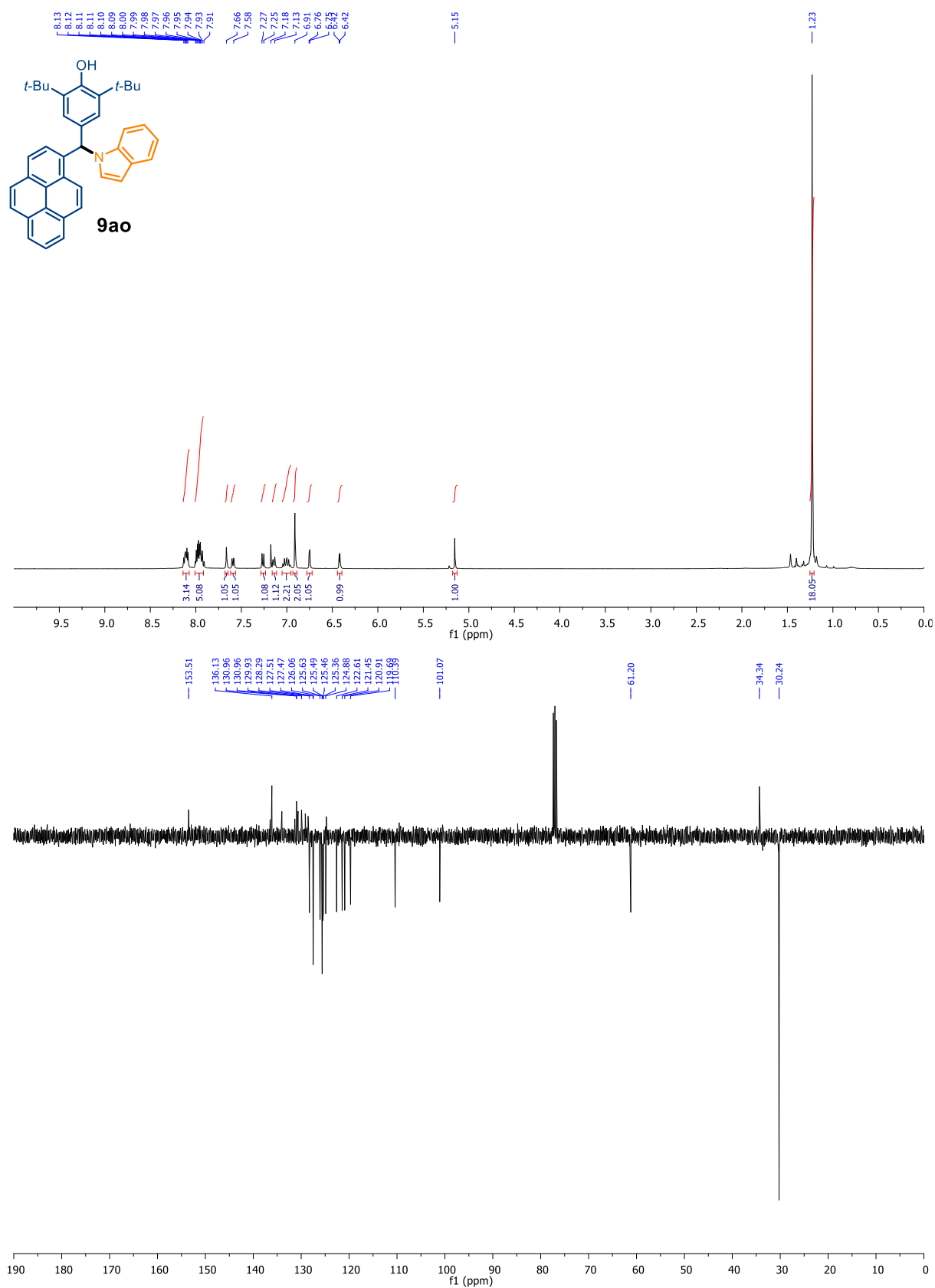


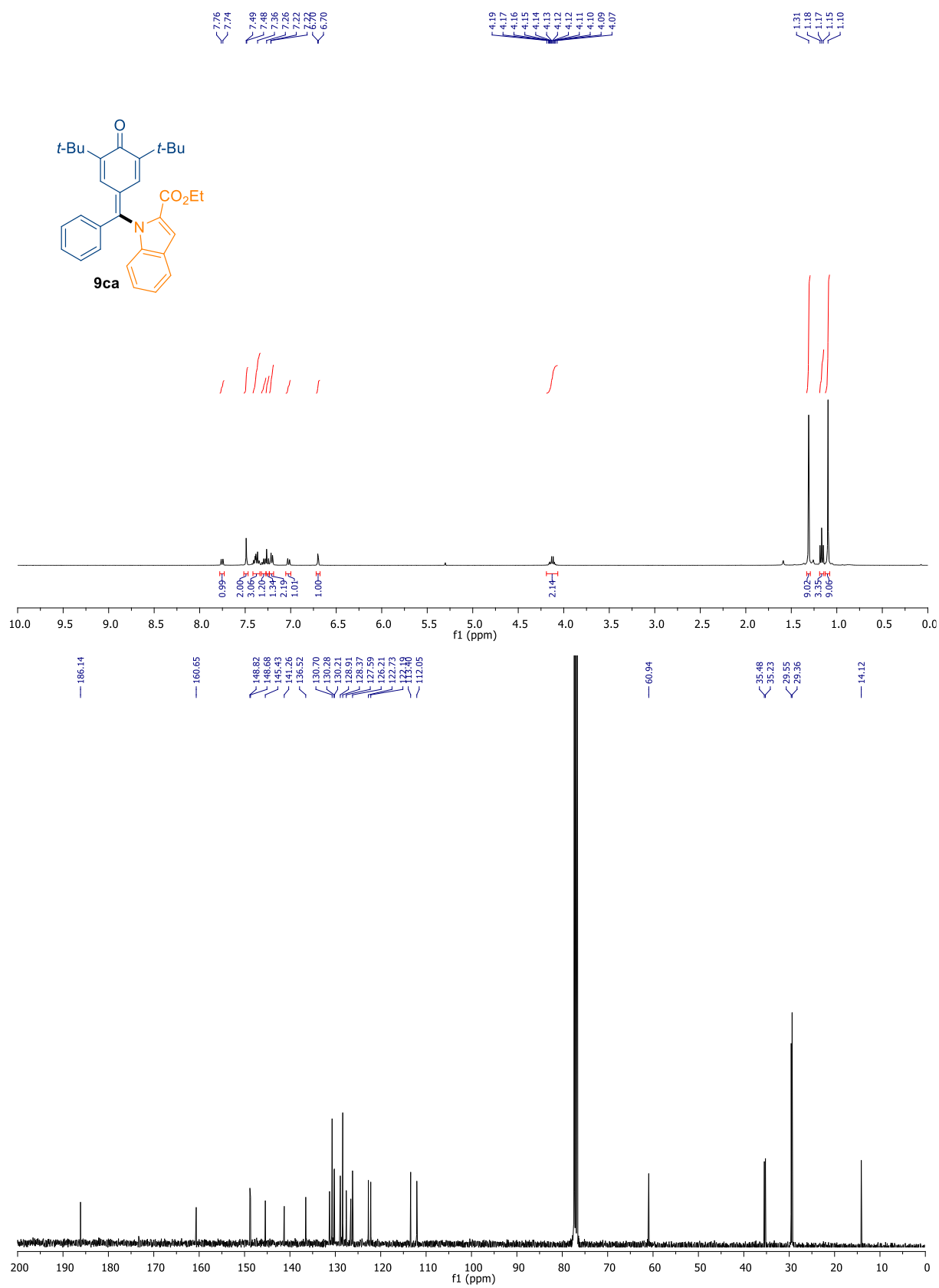


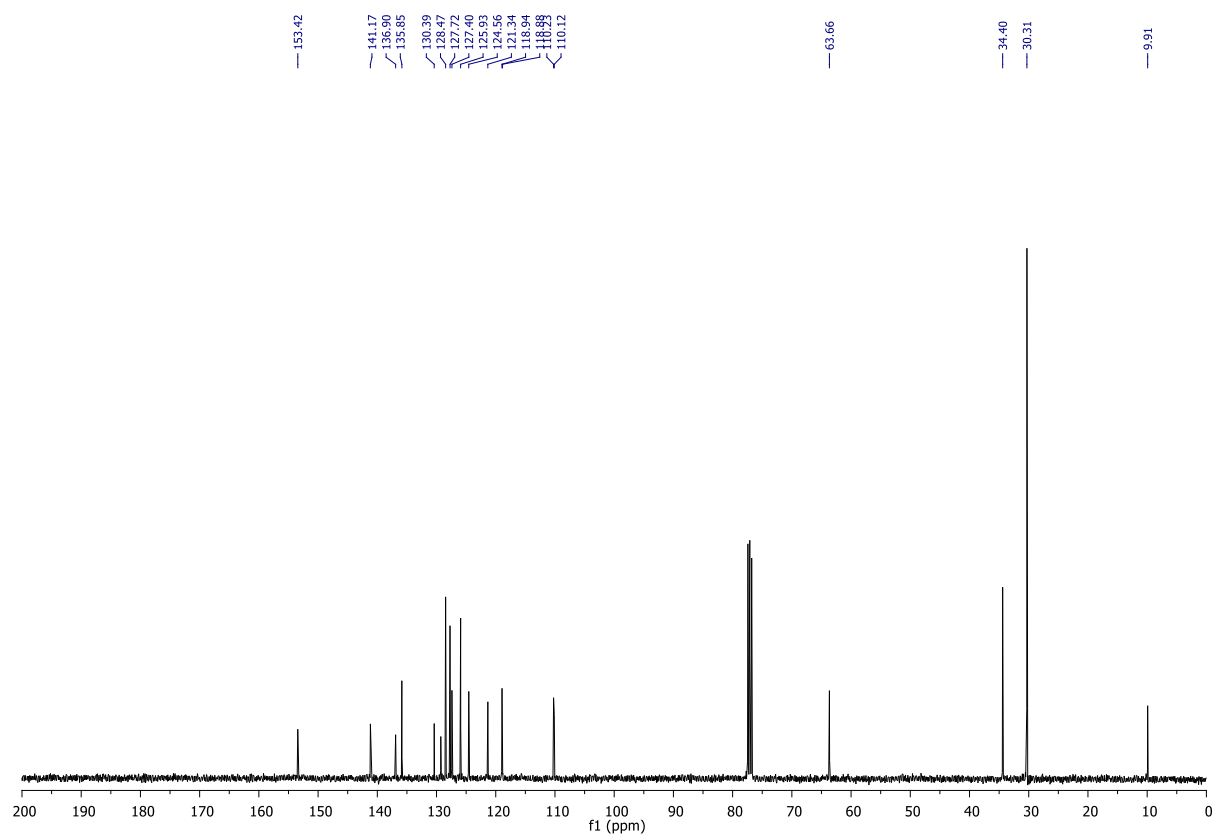
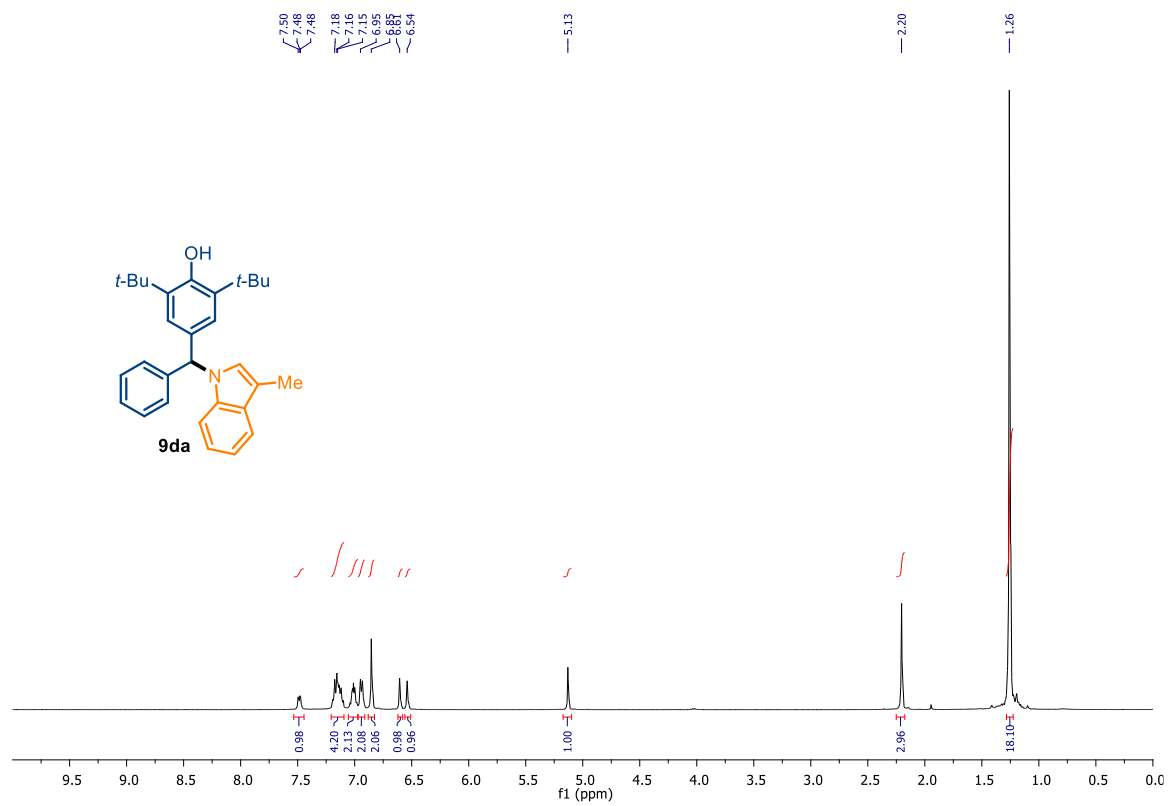


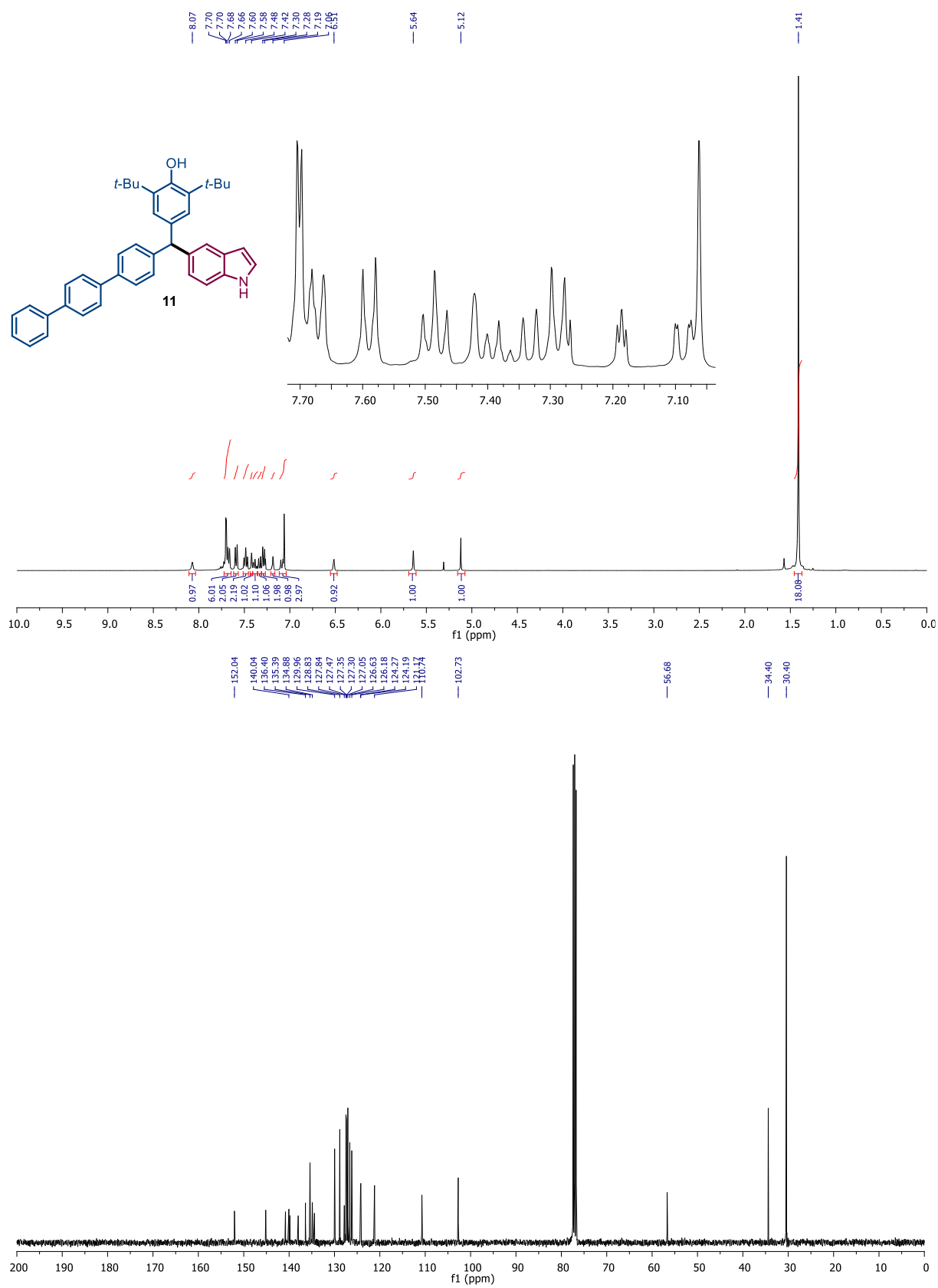


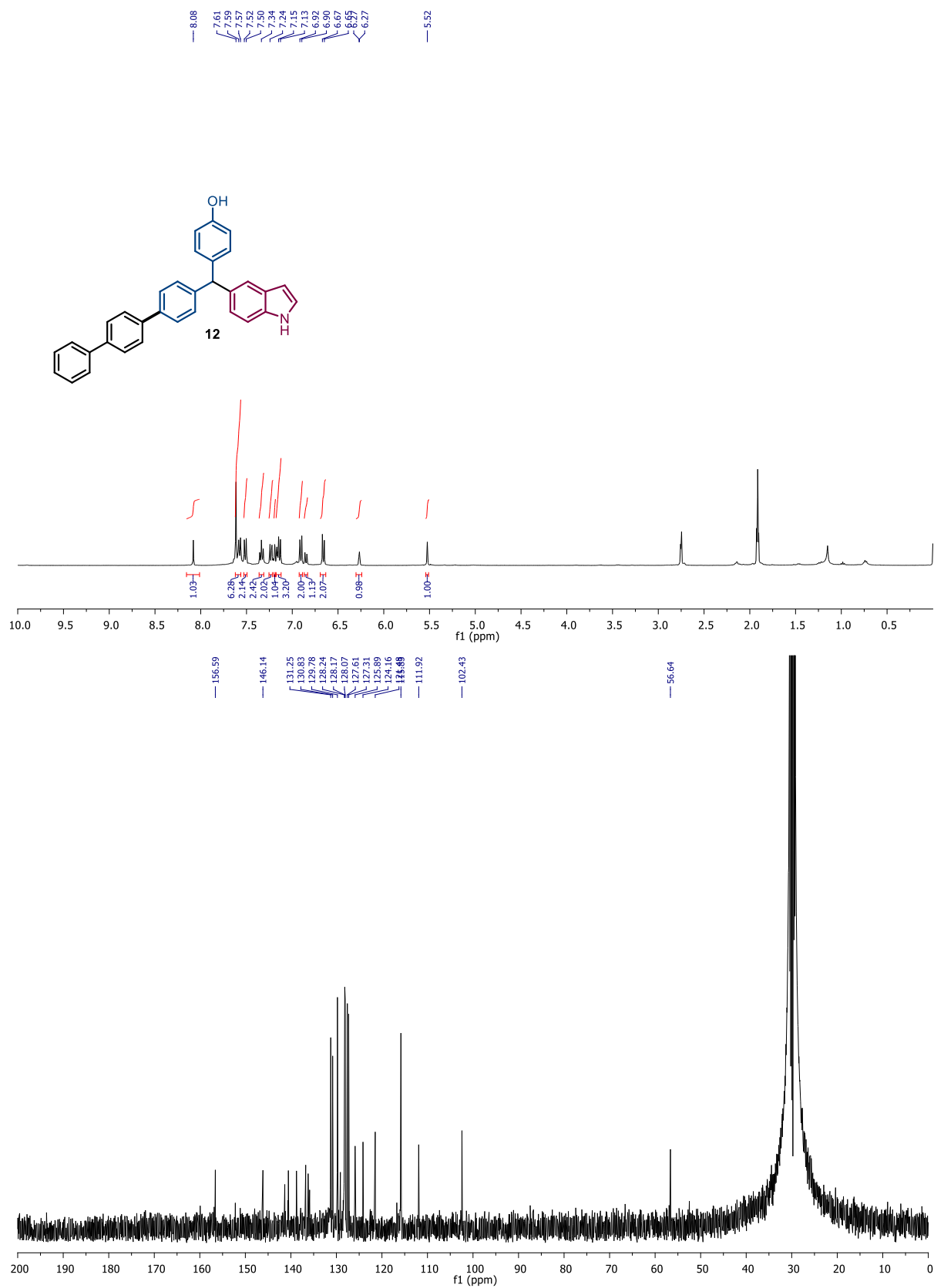


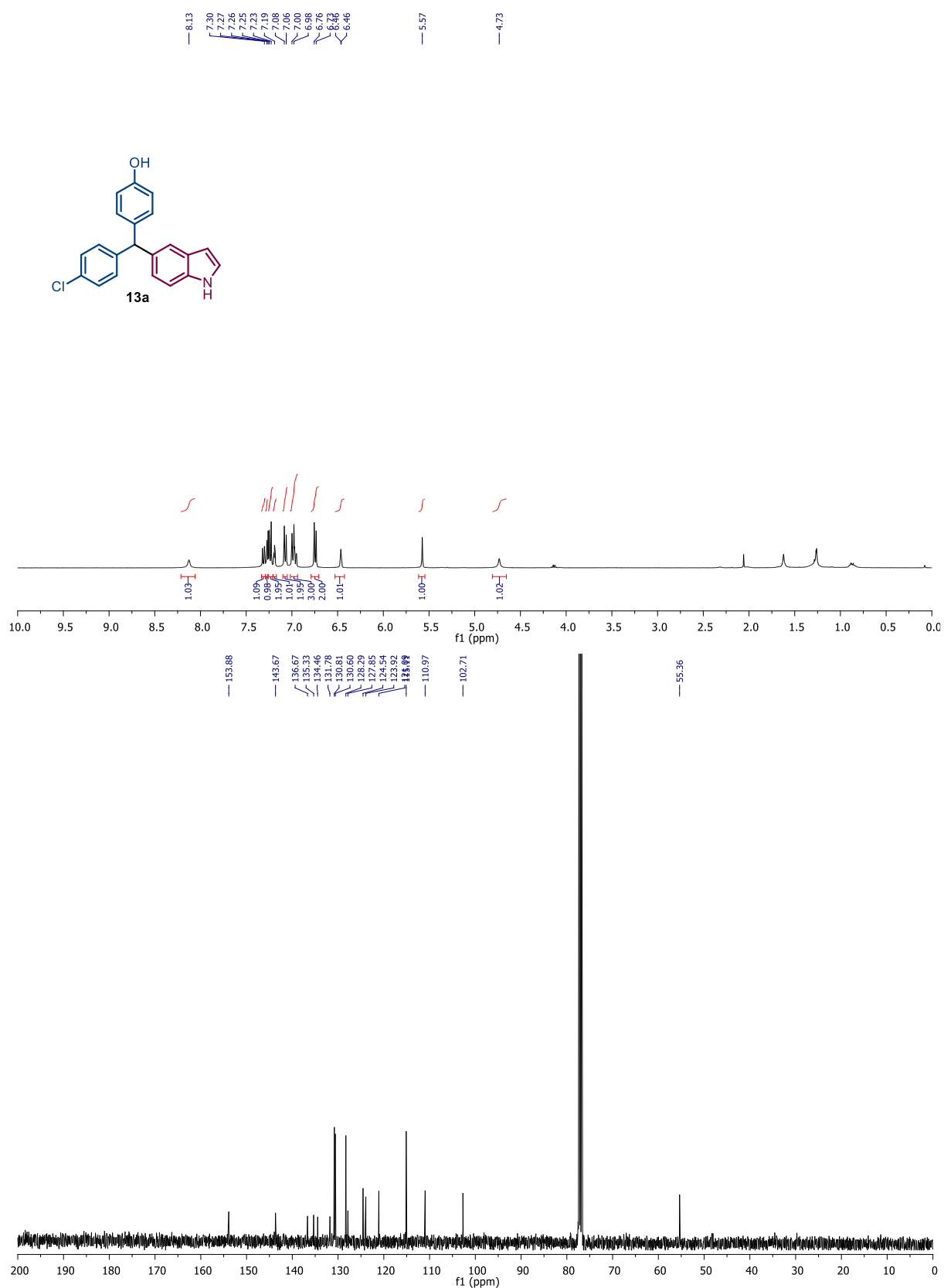


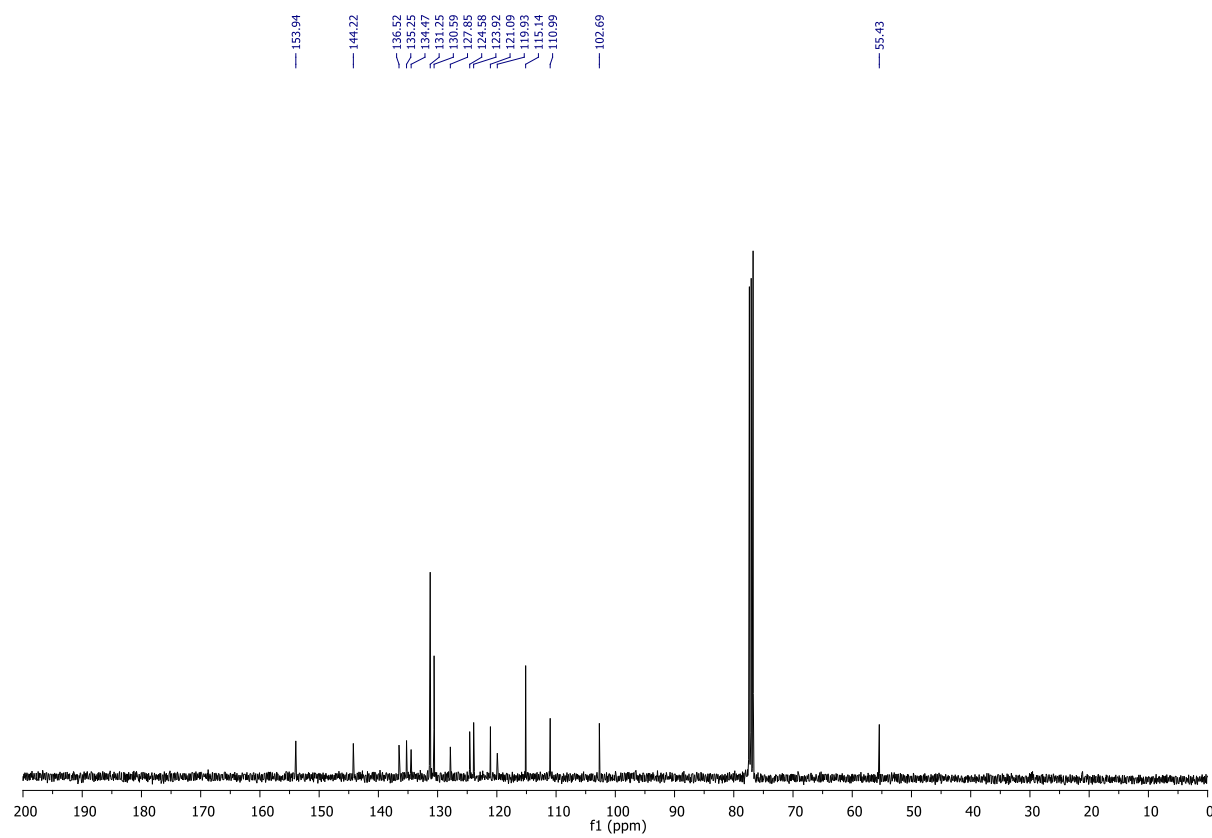
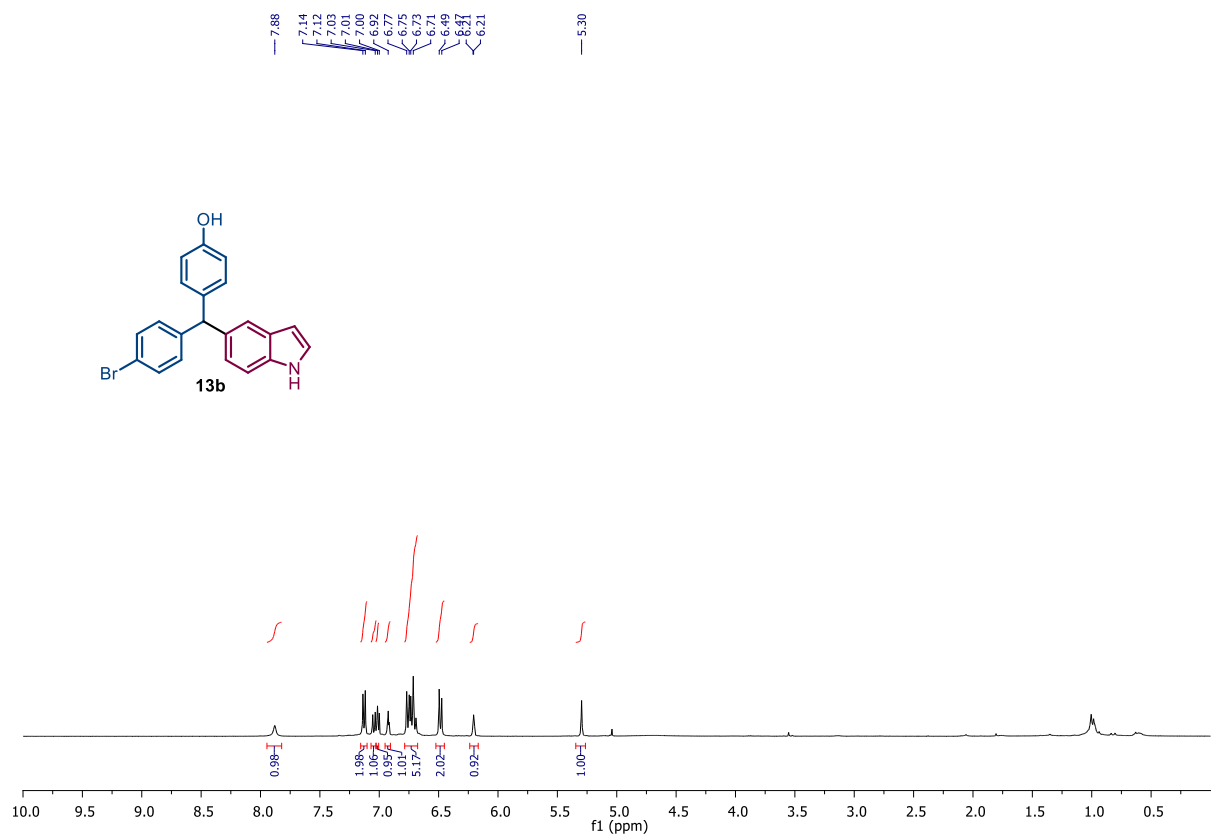




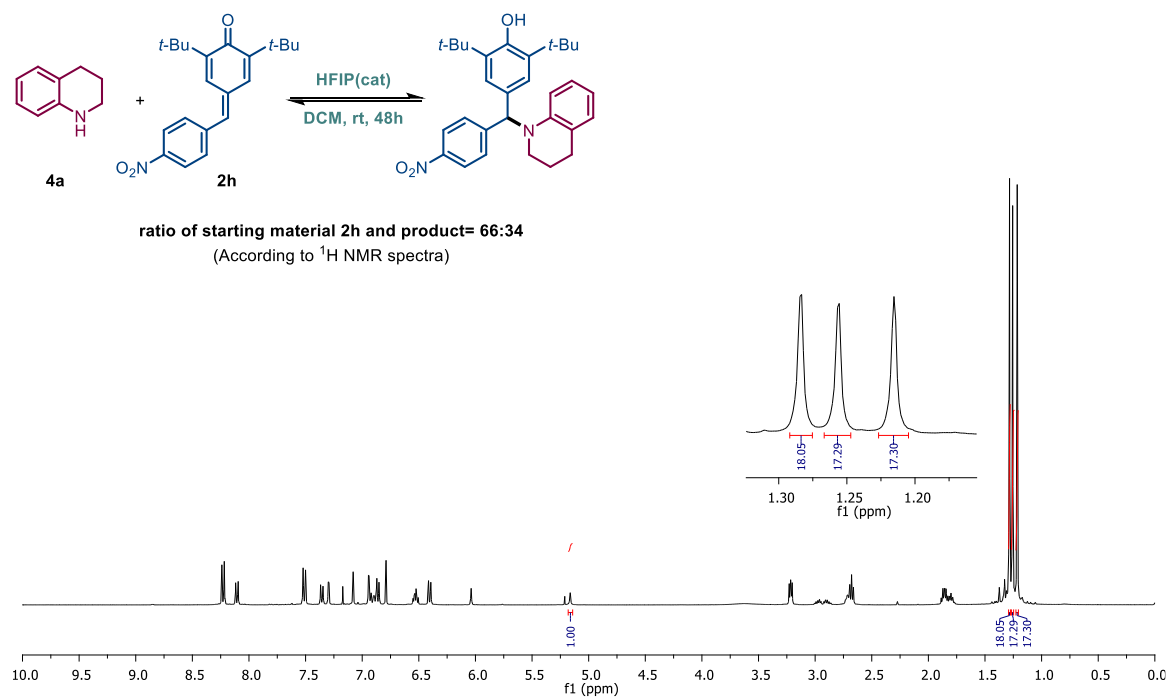
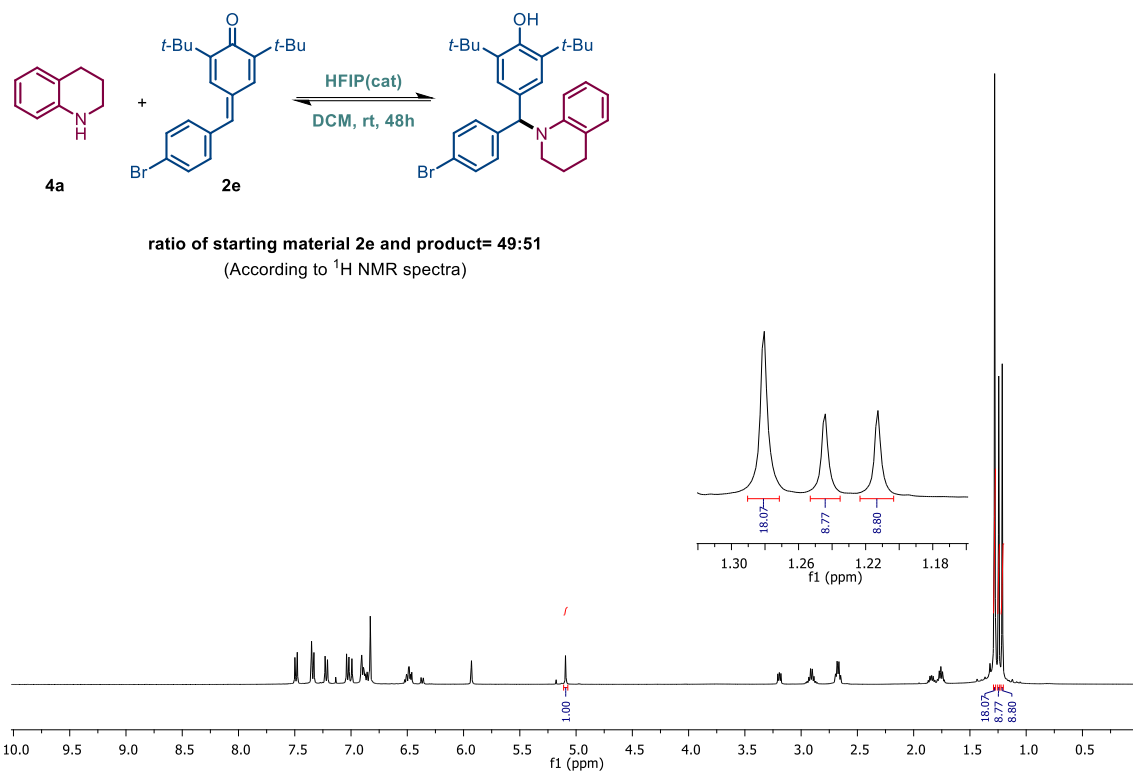


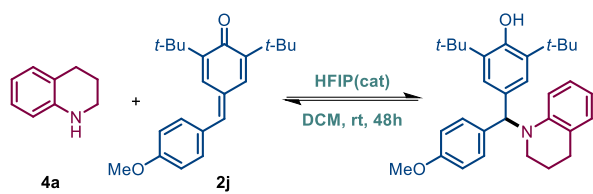






2. ^1H NMR monitoring of reactions between 4a and *p*-QMs (2e, 2h, and 2j)





Ratio of starting material 2j and product= 51:49

(According to ^1H NMR spectra)

