

Addendum

To the following article:

The γ - and the δ -effects in ^{13}C NMR spectroscopy in terms of nuclear chemical shielding (NCS) analysis

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In the paper, ^{13}C NMR chemical shifts were calculated at the *ab initio* quantum mechanical level of theory and discussed with respect to possible sources for the observed substituent effects. Previous work in this context at this level of theory was appropriately cited. It is important, however, to note that empirical, molecular mechanics, and semiempirical calculations have also been previously performed^{1–17} and examined for this same purpose.

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