

Simulation of oriented NMR spectra : combining molecular dynamics and chemical shift tensor calculations

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Theoretical prediction and assignment of vicinal ^1H – ^1H coupling constants of diastereomeric 3-alkoxy-6,7-epoxy-2-oxabicyclo[3.3.0]octanes

Aav, Riina; Pehk, Tõnis; **Tamp, Sven**; **Tamm, Toomas**; Kudrjašova, Marina; Parve, Omar; **Lopp, Margus** Magnetic resonance in chemistry 2011 / p. 76-82 : ill <https://analyticalsciencejournals.onlinelibrary.wiley.com/doi/full/10.1002/mrc.2712>