

Structure and intramolecular mobility of N-(thio)phosphoryl(thio)amides: XVI. ^1H , ^{13}C and ^{31}P NMR study of intramolecular dynamics of N,N'-bis(thio)phosphoryl(thio)urea containing an open-chain fragment in CD_2Cl_2 and CD_3CN solutions

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Abstract

The structure and intramolecular transitions of N,N'-bis[N-disopropoxythiophosphorylaminothiocarbonyl]-1,7-diaminoheptane in CD_2Cl_2 and CD_3CN 3-10% solutions were studied by means of ^1H , ^{13}C , and ^{31}P NMR spectroscopy. Combined analysis of the NMR data confirmed a high lability of the molecules with the realization of two conformational forms of the macromolecule, the amide-amide proton exchange, and existence of various tautomeric forms. © 2011 Pleiades Publishing, Ltd.

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