

Structure and Intramolecular Lability of N-(Thio)phosphoryl(thio)amides. IV. ^1H , ^{13}C and ^{31}P NMR Study of Dynamic Processes in Solutions of N,N'-Bis(diisopropoxythiophosphorylaminothiocarbonyl)-1,10-diaza-18-crown- 6 Ethers

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Abstract

^1H , ^{13}C , and ^{31}P NMR spectroscopy was used to study the structure of N,N'-bis(diisopropoxythiophosphorylaminothiocarbonyl)-1,10-diaza-18-crown- 6 ethers in CD_3CN , CD_2Cl_2 , and $(\text{CD}_3)_2\text{CO}$ solutions. A tautomeric equilibrium was detected, involving the amide (with $\text{C}=\text{S}$ trans to $\text{P}=\text{S}$), two prototropic, and one phosphorylotropic forms. It is found that the macroheteroring has two conformations: with trans,cis, and trans N-substituents. The conformational equilibrium is solvent-dependent.
