

[本期目录](#) | [下期目录](#) | [过刊浏览](#) | [高级检索](#)[\[打印本页\]](#) [\[关闭\]](#)**论文****系列双磷维生素B₁离子盐化合物的合成、结构及分子构象**白凤英^{1,3}, 李晓天¹, 朱广山², 邢永恒³, 曾小庆⁴, 葛茂发⁴

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摘要:

将Ni(ClO₄)₂和NH₄PF₆分别与硫胺素焦磷酸在甲醇体系中反应, 得到了两个新的离子盐型化合物[TPP·ClO₄·H₂O](1)和[TPP·PF₆·CH₃OH](2)(TPP为硫胺素焦磷酸酯). 通过元素分析、红外光谱及X射线衍射等方法对它们进行了表征. 结构分析表明, 它们属于离子型化合物, 而且硫胺素焦磷酸与高氯酸根, 六氟磷酸根形成了大量的氢键网络结构. 结合计算结果进一步分析了化合物的活性及电子结构特征.

关键词: 硫胺素焦磷酸酯; 氢键; 晶体结构; 量子化学计算**Synthesis, Structure and Molecular Configuration of a Series of Ionic Salt Compounds of Pyrophosphate Ester of Thiamine**BAI Feng-Ying^{1,3}, LI Xiao-Tian^{1*}, ZHU Guang-Shan², XING Yong-Heng³, ZENG Xiao-Qing⁴, GE Mao-Fa⁴

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Abstract:

Ionic salt compounds [TPP·ClO₄·H₂O](1) and [TPP·PF₆·CH₃OH](2)(TPP as the pyrophosphate ester of thiamine) were prepared in the methanol system under room temperature, and characterized by elemental analysis, IR spectra, and single X-ray diffraction. Structural analysis shown that they are ionic compounds and there are abundant hydrogen bondings in molecular structures. On the basis of the data of quantum chemical calculation, their potential reaction activity and electronic configuration were also discussed.

Keywords: Pyrophosphate ester of thiamine; Hydrogen bond; Crystal structure; Quantum chemical calculation

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