Cu(II)与 α -氨基肟配体PnAO配位反应的热力学及热化学研究

林华宽,张渊明,古宗信,陈荣悌

南开大学化学系

收稿日期 修回日期 网络版发布日期 接受日期

摘要 用pH法测定了PnAO的质子化常数及Cu(II)-PnAO,Cu(II)-5-R-Phen-PnAO配合物的稳定常数. 用直接量热法测定了PnAO的质子化热及Cu(II)二元,三元配合物的生成热.Cu(II)三元配合物体系与5-R-Phen的质子化体系之间存在线性热力学函数关系.并对取代基效应作了解释 关键词 稳定常数 铜络合物 生成热 量热法 取代基效应 配位反应 质子化常数 氨基肟

大键词 <u>穩定吊数</u> <u>铜络台物</u> <u>生成热</u> <u>重热法</u> <u>取代基效应</u> <u>配位反应</u> <u>质于化吊数</u> <u>氨基肟</u> 分类号 0642 0611.662

Thermodynamic and calorimetric study on the coordination reaction of copper(II) with α -aminoxime ligand PnAO

LIN HUAKUAN, ZHANG YUANMING, GU ZONGXIN, CHEN RONGTI

Abstract The pH titration method was used to measure the protonation constants of PnAO [2,2'-(1,3-propyldiamine)-bis(2-methyl-3-butanone oxime)], and the stability constants of binary and ternary complexes. The heats of protonation and of complex formation were measured by calorimetry. The results showed that some linear free energy relationships (LEER) exists between the protonation constants of 5-R-Phen (R = Me, Cl, NO2) and the stability constants of Cu(II)-PnAO-RPhen ternary system. Also found was the existence of linear enthalpy relanationships (LER) in ternary system. The substituent effect is discussed.

Key wordsSTABILITY CONSTANTCOPPER COMPLEXFORMATION HEATCALORIMETRYSUBSTITUENT EFFECTCOORDINATION REACTIONPROTONATION CONSTANT

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