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水与正丁醇二元体系最低共沸混合物的热容

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摘要 用精密绝热量热计测定了由水和正丁醇组成的最低共沸混合物在78至320 K温区的热容, 并建立了该体系热力学函数和温度的定量函数关系。根据测定的热容曲线表明, 最低共沸混合物在111.9±1.2

K发生玻璃态转变, 179.26±0.77 K和269.69±0.14 K发生固-液相变。还计算了其相变焓和相变熵。考虑了物理贡献和化学贡献两部分的作用, 拟合了超额热容与温度的函数关系。讨论了物理贡献和化学贡献两部分的作用对超额焓和超额熵的影响。

关键词 [绝热量热学](#), [热容](#), [超额热容](#), [最低共沸混合物](#), [水](#), [正丁醇](#)

分类号

Measurements of the Heat Capacity of an Azeotropic Mixture

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Abstract Molar heat capacities of *n*-butanol and the azeotropic mixture in the binary system [water ($x=0.716$) plus *n*-butanol ($x=0.284$)] were measured with an adiabatic calorimeter in a temperature range from 78 to 320 K. The functions of the heat capacity with respect to thermodynamic temperature were established for the azeotropic mixture. A glass transition was observed at (111.9±1.2) K. The phase transitions took place at (179.26±0.77) and (269.69±0.14) K corresponding to the solid-liquid phase transitions of *n*-butanol and water, respectively. The phase-transition enthalpy and entropy of water were calculated. A thermodynamic function of excess molar heat capacity with respect to temperature was established, which took account of physical mixing, destructions of self-association and cross-association for *n*-butanol and water, respectively. The thermodynamic functions and the excess thermodynamic ones of the binary systems relative to 298.15 K were derived based on the relationships of the thermodynamic functions and the function of the measured heat capacity and the calculated excess heat capacity with respect to temperature.

Key words [adiabatic calorimetry](#) [heat capacity](#) [excess heat capacity](#) [azeotropic mixture](#) [water](#) [n-butanol](#)

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