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 $\rm M \, M \, M \, LaFe_{0..8-x} Cu_x Ni_{0..2} 0_3$ 的合成和烧结性能 Synthesis and Sintering of Cathode Materials $\rm LaFe_{0..8-x} Cu_x Ni_{0..2} 0_3$

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中文关键词: 阴极材料; LaFe $_{0.8-x}$ Cu $_x$ Ni $_{0.2}$ O $_3$; 共沉淀法; 烧结性能

英文关键词: cathode material; LaFe $_{0.8-x}$ Cu $_x$ Ni $_{0.2}$ O $_3$; Co-precipitation method; sintering

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中文摘要:

英文摘要:

LaFe $_{0.8-x}$ Cu $_x$ Ni $_{0.2}$ O $_3$ (x=0.0-0.2) (LFCN), a new cathode material of solid oxide fuel cell (SOFC), was synthesized by Co-precipitation method using sodium bicarbonate. The lattice structures of samples with different x contents were characterized by XRD. Porosity and density of the porous LaFe $_{0.8-x}$ Cu $_x$ Ni $_{0.2}$ O $_3$ (x=0.0-0.1) as a function of sintering temperature were investigated. It was found that the orthorhombic structure could be formed after calcination at 900 °C for 4 h. The particle size of LFCN was about 350 nm. The density of the porous LFCN increased with sintering temperature, but the opposite was true for the porosity. On the other hand, at the same sintering temperature, the porosity of LaFe $_{0.8-x}$ Cu $_x$ Ni $_{0.2}$ O $_3$ (x=0.0-0.1) decreased with increasing x contents. It is indicated that the dopant of Cu on LaFe $_{0.8}$ Ni $_{0.2}$ O $_3$ can facilitate the sintering of the materials. After sintering at 1 100 °C for 4 h, the porous LaFe $_{0.7}$ Cu $_{0.1}$ Ni $_{0.2}$ O $_3$ was still with appropriate structure, and its porosity was 29%.

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