
Electronic transport and crystallization kinetics of melt-spun Ni_{33.3}Zr_{66.7} ribbon studied by electrical resistivity measurements

Smili Billel^{*1}, Ivan Kaban², and Jean-Georges Gasser³

¹SMILI – Laboratory of Inorganic Materials Chemistry, University Badji Mokhtar of Annaba BP 12, Annaba 23000, Algeria, Algeria

²KABAN – IFW Dresden, Institute for Complex Materials, Helmholtzstr. 20, 01069 Dresden, Germany, Germany

³GASSER – Laboratoire de Chimie et Physique - Approche Multi-échelle des Milieux Complexes Chimie, Physique et Matériaux, Université de Lorraine, 1 Boulevard Arago - 57078 – France

Abstract

In this paper, rapidly quenched Ni_{33.3}Zr_{66.7} ribbons were successfully prepared by melt spinning technique. The electrical and thermal transport properties of as-spun Ni_{33.3}Zr_{66.7} alloys were studied in detail by a combination of electrical resistivity and absolute thermoelectric power measurements over a temperature range from 25 °C up to 400 °C. The non-crystalline structure of the samples was fully confirmed by X-ray diffraction (XRD) and scanning electron microscope (SEM). Moreover, the crystallization kinetics of Ni_{33.3}Zr_{66.7} glassy alloy has been investigated during isochronal and isothermal annealing treatments based on the electrical resistivity measurements. The crystallization activation energy, E_x , for a series of electrical resistivity measurements at various heating rates, was calculated in the order of 371.4 kJ/mol and 382.2 kJ/mol by means of Kissinger and Ozawa methods, respectively. The Johnson-Mehl-Avrami (JMA) analysis was applied to the isothermal crystallization kinetics, and the local Avrami exponent has been determined in the range from 2.97 to 3.23 with an average value of $n = 3.1$, revealing that isothermal crystallization mechanism is diffusion-controlled three-dimensional growth crystallization mechanism, as well as an increasing nucleation rate. In addition, the local activation energy for crystallization, E_α , calculated from the Arrhenius equation decreases at the crystallized volume fraction $0.2 \leq \alpha \leq 0.8$ and giving an average value of 376.2 kJ/mol.

Keywords: Metallic glasses, Electronic transport properties, Thermal stability, Crystallization kinetics, Activation energy

*Speaker