

The Serbian Society for Ceramic Materials
Institute for Multidisciplinary Research (IMSI), University of Belgrade
Institute of Physics, University of Belgrade
Center of Excellence for the Synthesis, Processing and Characterization of
Materials for use in Extreme Conditions "CEXTREME LAB" - Institute of
Nuclear Sciences "Vinča", University of Belgrade
Faculty of Mechanical Engineering, University of Belgrade
Center of Excellence for Green Technologies, Institute for Multidisciplinary
Research, University of Belgrade
Faculty of Technology and Metallurgy, University of Belgrade

PROGRAMME and the BOOK of ABSTRACTS

6CSCS-2022

6th Conference of
the Serbian Society for Ceramic Materials
June 28-29. 2022. Belgrade Serbia

Edited by:
Branko Matović
Aleksandra Dapčević
Vladimir V. Srdić

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P-42

PHASE TRANSFORMATIONS DURING CYCLIC ANNEALING OF Ti₃Al-BASED INTERMETALLIC

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Alterations that emerged in the intermetallic Ti₃Al-based alloy microstructure during the cyclic annealing at 600 °C and 900 °C in the air atmosphere were studied by combining experimental and theoretical methods. The phase transformations in the Ti-24Al-11Nb (at.%) alloy microstructure were monitored during the alloy high-temperature processing up to 120 h. The conducted cyclic annealing resulted in the external oxide scale formation and the microstructural features alteration. Obtained experimental results indicated that the annealing temperature and duration significantly affected the external scale composition and structure, as well as the alloy phase composition. Namely, the alloy microstructural characterization revealed that a starting two-phase $\alpha_2+\beta$ microstructure changed to a great extent during the cyclic annealing. The increase of the cyclic annealing temperature induced the intensive $\alpha_2\rightarrow\beta$ phase transformation and as a result, new α_2'' and O-Ti₂AlNb phases appeared in the investigated alloy microstructure. The experimentally observed appearance of the O-Ti₂AlNb phase in the Ti₃Al-based alloy microstructure was confirmed by the results of theoretical investigations. *Ab initio* modeling also enabled the prediction and study of the additional structures in the O-Ti₂AlNb compound using the first-principles calculations.