

Structure-property orientation relationship of a $g/a_2/Ti_5Si_3$ in as-cast-45Al-2Nb-0.7Cr-0.3Si intermetallic alloy

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ABSTRACT:

Structure-property relationship of the as-cast $g-Ti-45Al-2Nb-0.7Cr-0.3Si$ based intermetallic alloy was examined. The phases stable at room temperature in the alloy were $g/a_2/Ti_5Si_3$, respectively; while their crystal structures were described by means of chemical formulas, Pearson's number, space groups, lattice parameters, atomic positions and occupancy numbers using a Vienna ab initio simulation package (VASP) software for materials design. Moreover, high resolution electron backscatter diffraction (HSEBSD) was used to analyse the orientation relationship of both the as-cast (b-solidifying) and heat-treated (a-solidifying) microstructural phases. The results showed that the formation of $g/a_2/Ti_5Si_3$ follows the Blackburn orientation relationship (BOR): (b-solidifying) $\{110\}_b // \{0001\}_a$ and $\langle 111 \rangle_b // \langle 2^{-1}10 \rangle_a$, and for (a-solidifying) $\frac{1}{4} (111)_g // (0001)_a$ and $\langle 1^{-1}10 \rangle_g // \langle 11^{-2}0 \rangle_a$.