

Structure-property relationships in AlFe_2B_2 : A magnetofunctional layered system

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Introduction

As magnetism becomes developed into prospective and expanding technologies (encompassing electronics, data storage, and energy conversion) the need for novel, inexpensive magnetic materials produced from abundant elements becomes increasingly evident. Previous reports on the ferromagnetic intermetallic compound AlFe_2B_2 show it fits these criteria and has promise for magnetic refrigeration based on its near-room-temperature magnetic phase transition, which dictates the operating temperature range for applications [1,2]. In this work the ferromagnetic properties of AlFe_2B_2 single crystals are clarified along crystallographic directions of the layered crystal structure (Figure 1). Single crystals possess an unbroken periodic arrangement of atoms through the entire sample, allowing correlations between structure and magnetic response to be studied. In particular, understanding the energetic preference of the magnetic moments to align along one crystallographic direction, (known as the magnetocrystalline anisotropy energy) can be achieved. Understanding the magnetocrystalline anisotropy can shed light on the intrinsic and extrinsic parameters responsible for the magnetic response of AlFe_2B_2 . This information is a blueprint for discovering pathways to tailor this system for advanced applications

Materials and Methods

The objective of isolating single crystals for study was approached by first synthesizing bulk AlFe_2B_2 by arc-melting the constituent elements (99.9% purity) in a 3 Al: 2 Fe: 2 B molar ratio that produced a multiphase alloy of AlFe_2B_2 in an $\text{Al}_{13}\text{Fe}_4$ matrix. X-ray diffraction (XRD) was performed to verify the crystal structures and lattice parameters of phases present. Optical and scanning electron microscopy (SEM) were used to confirm your microstructure and phase distribution within the sample. The arc-melted charges were subsequently etched in 50 % v/v HCl for 30 hours to dissolve the surrounding Al-rich $\text{Al}_{13}\text{Fe}_4$ phase, releasing single crystals. The structure of these extracted single crystals was examined using transmission electron microscopy (TEM). Magnetic characterization was carried out using superconducting quantum interference device (SQUID) magnetometry in magnetic fields up to $\mu_0 H_{app} = 5$ T and temperatures in the range $50 \text{ K} \leq T \leq 390 \text{ K}$. The magnetic transition temperature was determined as the inflection point in a curve generated by plotting the derivative of magnetization as a function of temperature and temperature (dM/dT vs. T).

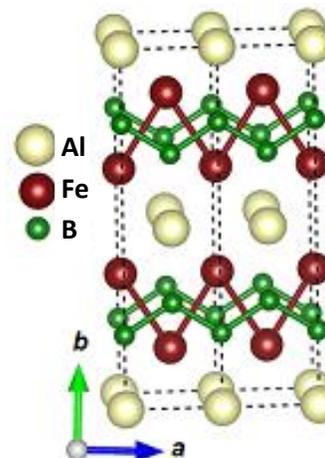


Figure 1. AlFe_2B_2
orthorhombic unit cell

Results and Discussion

SEM and XRD confirm that arc-melted AlFe_2B_2 is multiphase, consisting of primarily AlFe_2B_2 and $\text{Al}_{13}\text{Fe}_4$. After 10 minutes of etching a layered morphology was observed along the thickness of AlFe_2B_2 crystallites. The microstructure of the extracted AlFe_2B_2 crystals after etching for 30 hours was found to consist of layers ~ 100 nm in thickness (Figure 2(a)). Based on this observed layering, it is concluded that extracted crystallites are not complete single crystals. The crystallographic orientation of individual layers of extracted crystallites, and of the magnetic easy axis, was confirmed (Figure 2(a)) via TEM, which also allowed determination of the lattice parameters ($a = 2.930(5)$ Å $c = 2.870(5)$ Å). An appreciable magnetocrystalline anisotropy at $T = 50$ K is observed in this system by measuring the magnetization along the in-plane crystallographic directions. It is found that the magnetic field necessary to saturate the magnetization of the sample along the c axis is approximately ten times that required along the a -axis (Figure 2(b)).

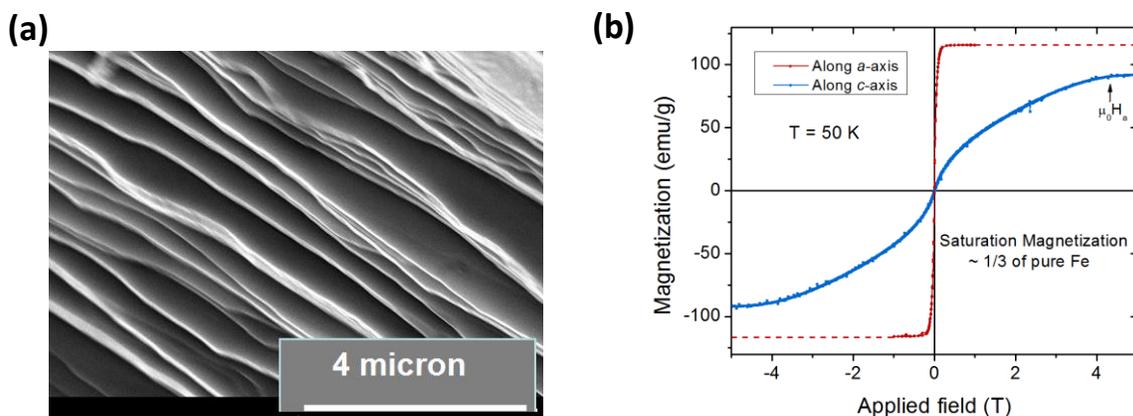


Figure 2. (a) Scanning electron micrograph of AlFe_2B_2 single crystal showing hierarchical layering ~ 100 nm in thickness. (b) Magnetization versus applied field along a and c in-plane directions used to determine the anisotropy field ($\mu_0 H_a$) ~ 4 T at 50 K.

Conclusions

Correlations between structure and magnetism help to provide a better understanding of the bonding in the crystal structure to develop pathways to enhance the functional behavior. This order of magnitude difference in energy required to align the magnetic moments along the two in-plane crystallographic directions confirms that the structure has a dramatic effect on the magnetism in AlFe_2B_2 . New findings provide insight into the evolution of magnetism in AlFe_2B_2 which is critical for the engineering of a magnetic material with applications in mind. Results suggest that intrinsic property variation such as elemental substitution may be an effective method of altering the functional temperature range of this system as well as the strength of the magnetization. AlFe_2B_2 has potential as a new cost effective magnetofunctional material and may provide solutions to current limitations in sensor and solid state cooling applications.

References

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