Research Article

Stacking Fault Energy of Si Nanocrystals Embedded in SiO₂

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Si nanocrystals (Si nc) were produced by the implantation of Si⁺ into a SiO₂ film on (100) Si, followed by high-temperature annealing. High-resolution transmission electron microscopy (HRTEM) observation has shown that a perfect dislocation (Burgers vector \( \mathbf{b} = (1/2)\langle 110 \rangle \)) can dissociate into two Shockley partials (Burgers vector \( \mathbf{b} = (1/6)\langle 112 \rangle \)) bounding a strip of stacking faults (SFs). The width of the SFs has been determined from the HRTEM image, and the stacking fault energy for Si nc has been calculated. The stacking fault energy for Si nc is compared with that for bulk Si, and the formation probability of defects in Si nc is also discussed. The results will shed a light on the dissociation of dislocations in nanoparticles.

1. Introduction

Si nanocrystals (Si nc) embedded in a SiO₂ matrix have attracted much attention as a promising candidate for optoelectronics. It has been agreed by most researchers that the defect center at the Si nc/SiO₂ interface plays an important role in the light emission [1, 2]. In order to fully understand light emission mechanism, a comprehensive investigation of the microstructure of the Si nc and their surrounding matrix, especially the defects inside the Si nc or at the interface of Si nc/SiO₂, is necessary.

When analyzing the reasons for the formation of defects in the nanocrystals, the following two factors are usually considered [3]: the size effect and the energy of formation of stacking faults (SFs). The size effect on the formation of defects has been discussed in our previous paper [4]. It has been pointed out that the probability of defects being formed drops sharply when the Si nc are smaller than 5 nm [4]. The probability of defects, for example, SFs, being formed also depends on their energy of formation. However, the influence of stacking fault energy on the formation of defects in Si nc has not been reported yet.

2. Experimental

A 1 μm thick film of amorphous SiO₂ was obtained from universitywafer.com. It was produced by thermal oxidation of (100) Si substrate at high temperature (~1100°C under oxygen flow). The amorphous SiO₂ film was implanted at room temperature with 100 keV Si⁺ (implantation dose of \( 3 \times 10^{17} \text{ cm}^{-2} \)), followed by high-temperature annealing at 1100°C for 1 h under an atmosphere of nitrogen (N₂). After the implantation, hydrogen passivation was carried out at 500°C for 1 h in a forming gas of H₂ (5%) and N₂ (95%). The specimens for transmission electron microscopy examination were prepared in a cross-sectional orientation ([011] zone axis for the Si substrate) using conventional techniques of mechanical polishing, dimpling, and ion thinning. The ion thinning was performed using a Gatan model 691 PIPS. Dark field (DF) examination was carried out on a Philips CM30 microscope operating at 300 kV. High-resolution transmission electron microscopy (HRTEM) observations were performed using a JEOL JEM 2100F transmission electron microscope operating at 200 kV.

3. Results and Discussion

Figure 1(a) shows a typical cross-sectional DF image of this specimen [5]. From Figure 1(a), it can be seen that the Si nc layer extends from 50 ± 5 nm to a maximum depth of 290 ± 10 nm, and the size ranges from 2 to 22 nm in diameter. The Si nc can be divided into three regions, RI, RII, and RIII, according to their sizes. The Si nc in the middle region of
the implanted layer are larger than those near the free surface or the bottom of the implanted layer, which is in agreement with a Gaussian-shaped excess Si concentration depth profile in Figure 1(b). Figure 1(b) is the depth profile of Si nc obtained by means of SRIM code [6], and the sputtering effect has been considered. In Figure 1(b), the average sizes of the Si nc are divided into three major groups: 3, 6, and 12 nm. From Figure 1, it can be seen that the simulated depth profile is consistent with that obtained from TEM observation.

Extensive HRTEM observations of the Si nc show that a perfect dislocation can dissociate into two Shockley partials (Burgers vector $\mathbf{b} = (1/6)(112)$) bounding a strip of SFs. Figure 2(a) shows a typical HRTEM image of a Si nanocrystal with an extended dislocation [6], where the particle is oriented along the [011] direction. The size of the Si nanocrystal is around 15 nm. It can be seen from Figure 2(a) that two Shockley partial dislocations are connected by a strip of SFs. The two Shockley partials are indicated by two arrows and labeled as 30° and 90°, and the dislocation line directions are along [011]. The SFs have a characteristic of an intrinsic stacking fault with a displacement of $\mathbf{R} = -1/3\langle 111 \rangle$ [7]. In order to demonstrate the configuration more clearly, the enlarged HRTEM image of the rectangle-enclosed region in Figure 2(a) is shown in Figure 2(b). The Burgers circuit enclosing the two Shockley partials (indicated by two arrows) is drawn in Figure 2(b). From the Burgers circuit, it can be clearly seen that there is a gap between the starting and ending point, which is indicated by an arrow. The total Burgers vector is determined to be $1/2(110)$. In addition, the SFs region is enclosed by the dashed lines in Figure 2(b). The width of SFs is measured to be $2.97 \pm 0.33$ nm from Figure 2(b).

There are mainly two factors exerting influences on the width of SFs: the interactions between two partial dislocations and the stacking fault energy. On one hand, dislocations lines repel each other with a force $f$ per unit length; on the other hand, stacking fault energy per unit area, $\gamma$, forces the dislocation lines to move closer. At equilibrium, $f = \gamma$, the SF width ($d_0$) can be described as [8, 9]

$$d_0 = \frac{G\mathbf{b}_1 \cdot \mathbf{b}_2}{\gamma},$$

where $G$ is the shear modulus and $\mathbf{b}_1$ and $\mathbf{b}_2$ are the Burgers vectors for two partial dislocations.
In deformed bulk Si, the SFs and the dissociation of a perfect dislocation (Burgers vector \( \mathbf{b} = (1/2)(110) \)) into two Shockley partial dislocations (Burgers vector \( \mathbf{b} = (1/6)(112) \)) were observed using the weak-beam dark-field technique \([10, 11]\) and HRTEM \([12]\). Experimental observations indicated that the size of the stacking faults range from 30 to 50 Å \([11–13]\), and the value for the intrinsic stacking faults energy is calculated to be 55 ± 7 mJ/m² \([10, 11, 14]\) or 69 mJ/m² \([15, 16]\). However, the stacking fault energy for Si nc has not been determined yet.

The equilibrium width for SFs in the Si nc (\( d_{\text{Si nc}} \)) has been measured to be about 2.97 ± 0.33 nm, so the stacking fault energy can be calculated using (1). The value for \( G \) of 6.36 × 10¹⁰ N/m² is used \([17]\), which is the effective value for defects on \{111\} planes. The SF energy for the Si nc (\( \gamma_{\text{Si nc}} \)) is estimated to be 84 ± 9 mJ/m², which is much higher than that for bulk Si reported in the literature \([10, 11, 14–17]\). Therefore, we can deduce that the probability for the formation of SFs in Si nc is less than that in bulk Si. Usually in the nanoparticles, dislocations can slip through the particle and no evident boundary can be observed between the perfect and faulted lattices \([7]\). Here we observed the dissociation of a perfect dislocation into two partial dislocations. The partial dislocations can be pinned at the Si nc/SiO² interface where some impurities are located \([18]\).

4. Conclusions

In conclusion, extended dislocations have been observed in the Si nanocrystals using HRTEM. The width of the SFS is measured to be 2.97 ± 0.33 nm, and the stacking fault energy is calculated to be 84 ± 9 mJ/m². The probability for the formation of SFs in Si nc is less than in bulk Si.

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