

## REAL TIME UHV-HRTEM OBSERVATION OF Si(111) $\sqrt{3}\times\sqrt{3}$ -Pd SURFACE AND DYNAMIC MOTION OF Pd CLUSTERS

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Recently, the development of advanced quantum devices requires the understanding of self-organized growth of metal or semiconductor nano-islands on semiconductor surfaces. For this purpose, the characterization and visualization tools for the surface growth with atomic resolution have been increasingly desired. Ultrahigh vacuum high-resolution transmission electron microscopy (UHV-HRTEM) has been a powerful tool to visualize atomic structure not only in a bulk but also on a surface. Moreover, well-established electron-diffraction and imaging theory can support the modeling of atomic structure. Individual atoms and clusters on a surface have often been imaged by HRTEM so far [1-2]. It is reported that Si(111) $\sqrt{3}\times\sqrt{3}$ -Bi was observed with HRTEM but image of surface atoms were overlaid with that of a substrate atoms [3]. A combination of HRTEM with computational process has provided clear atomic images of Si(111) $5\times 2$ -Au and Si(111) $7\times 7$  surface without substrate image [4-5]. However, the atomic motions of atoms and/or clusters on the surface cannot be analyzed with such off-line processes.

This paper reports an atomic structure of a Si(111) $\sqrt{3}\times\sqrt{3}$ -Pd surface and the motion of Pd clusters on the surface with UHV-TEM in real time. Figure 1 shows a schematic drawing of Pd deposition. In-situ TEM observation was performed using a 200 kV UHV-TEM equipped with a field emission gun. Pd less than 1 ML was deposited onto a Si (111) $7\times 7$  surface using an electron beam evaporator attached to the microscope column at about 550 K. Figure 2A shows a conventional HRTEM image of the Si(111) surface, on which Pd exhibits the  $\sqrt{3}\times\sqrt{3}$  structure on the thin specimen (10 nm or less). Since the image was taken at near the Scherzer defocus, Si(111) $1\times 1$  lattice fringes (0.19nm) are dominantly observed and the surface structure is unclearly overlaid. From the image simulation, we found that the Si  $1\times 1$  lattice image disappears and the  $\sqrt{3}\times\sqrt{3}$  structure is enhanced when the convergent angle is larger than 1 mrad at the condition of over-focus of 20-50 nm and the thickness of 7.5-10.3 nm, and that the contrast enhancement of Pd clusters against the  $\sqrt{3}\times\sqrt{3}$  structure is maximized at the over-focus of around 34nm and the convergent angle of around 3 mrad. Furthermore, we introduced an objective aperture so that electron diffraction of 0.20 nm or less are filtered off. Figure 2B shows a HRTEM image of the same area, indicating that atomic arrangement of the  $\sqrt{3}\times\sqrt{3}$  structure is clearly seen without any Si(111) $1\times 1$  lattice fringes. At this condition, Pd clusters are visualized as a set of three bright dots as shown in Figure 3. We proposed two types of structure models for the cluster, e.g., Pd-trimer and silicide models, and the simulated image calculated using these models were compared with the experimental one. Figure 4 shows the each structure model and simulated images at 2 and 3 mrad convergent angles, indicating that the Pd-trimer model is consistent with our results rather than the silicide one.

### References

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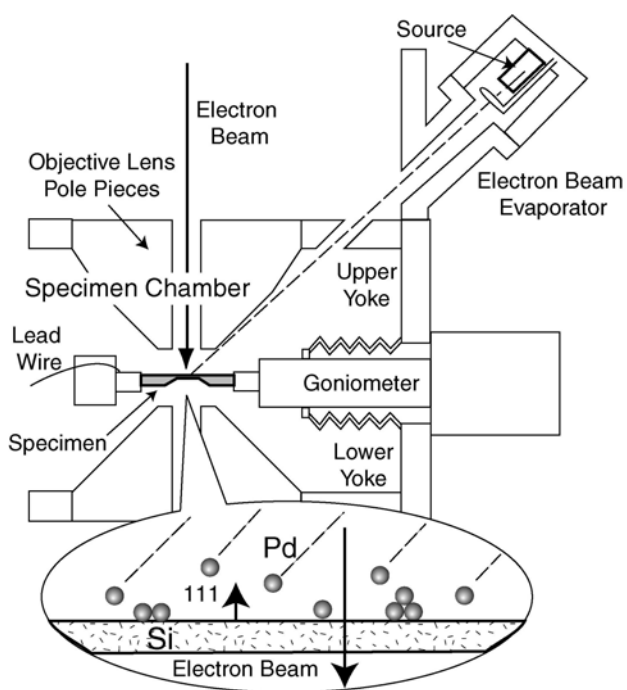


FIG 1. A schematic drawing of Pd deposition on Si (111) in the UHV-TEM

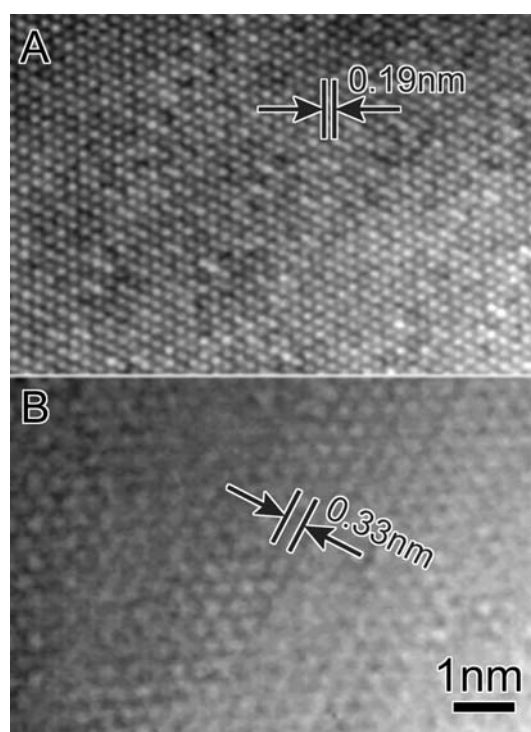


FIG 2. HRTEM images of  $\text{Si}(111)\sqrt{3}\times\sqrt{3}\text{-Pd}$  taken by a conventional manner(A) and taken on the present experimental condition(B).

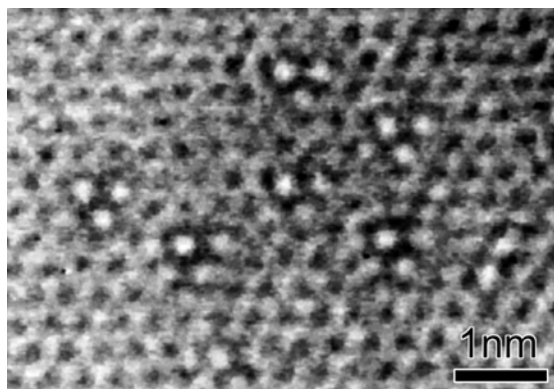


FIG 3. HRTEM image of Pd clusters moving around on the  $\sqrt{3}\times\sqrt{3}$  layer.

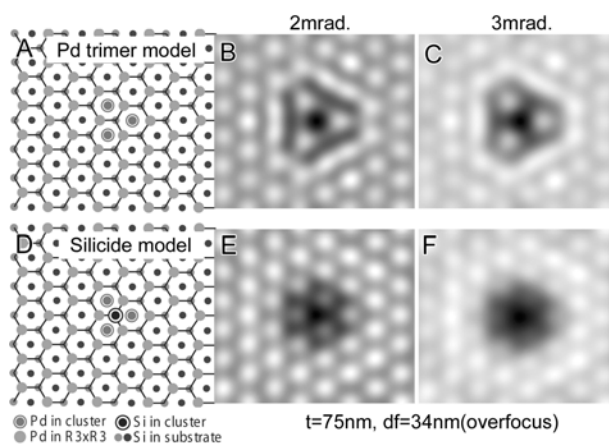


FIG 4. Structure models for a Pd cluster and their simulation images.