

Control of orientation of C₆₀ monolayers using steps

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To improve the performance of devices using organic thin films, we have to be able to grow high quality films. Understanding the growth mechanism of the organic thin films would greatly contribute to this ability. For this purpose, we have been investigating the initial stages of epitaxial growth of C₆₀ molecules on Si(111). In this work, we found that the orientation of the C₆₀ monolayer can be controlled using steps on Si(111).

We grew C₆₀ molecules on nominally flat and vicinal Si(111) surfaces at the substrate temperature of about 400 K. Before the C₆₀ growth, we deposited Au on the surface to form the 6x6 structure. Single-layer steps and multilayer steps existed on the nominally flat and vicinal surfaces, respectively. The structure of the C₆₀ layers was investigated using low-energy electron microscopy (LEEM) and low-energy electron diffraction (LEED).

Figure 1 shows a LEEM image after the C₆₀ deposition on the vicinal Si(111) surface misoriented 2 degrees to the <11-2> direction. Straight lines running almost vertically are multilayer steps, and two-dimensional C₆₀ islands seen as dark patches nucleate along the steps. Figure 2 shows a contrast-reversed LEED pattern after further deposition of C₆₀ on the same substrate. The diffraction spots from the substrate are indicated by dotted circles. The other spots originating from the C₆₀ monolayer constitute a triangular lattice, indicating that all the C₆₀ islands have the same orientation. Several domains with different orientations nucleate on the nominally flat and <1-10>-misoriented vicinal Si(111) surfaces. The height and direction of the steps are the keys to controlling the orientation of the C₆₀ monolayer.

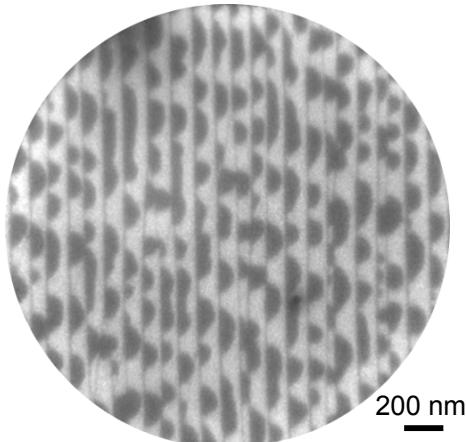


Fig. 1. LEEM image after C₆₀ growth on vicinal Si(111) misoriented 2 degrees to <11-2>. The electron energy is 10 eV.

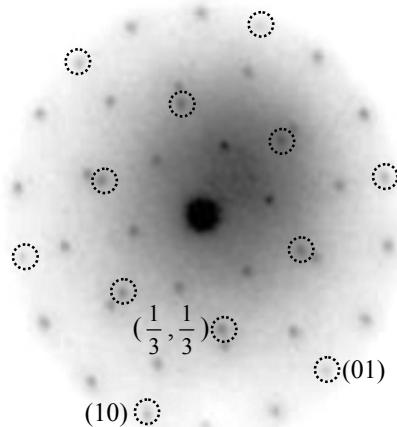


Fig. 2. LEED pattern after C₆₀ growth on vicinal Si(111) misoriented 2 degrees to <11-2>. The electron energy is 15.2 eV.