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Surface x-ray scattering study of a thallium-induced Si(111)-1×1 surface

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We report the atomic arrangement of a thallium (TI)-induced 1x1 phase on Si(111) surface determined by utilizing synchrotron x-ray scattering. The adsorption of TI on the Si(111)-7×7 surface has attracted attention due to its peculiar behavior of the "inert pair effect" distinctly different from that of other group III elements such as AI, Ga, and In. Early studies thus claimed that TI atoms occupied T1 sites directly on top of the unreconstructed Si layer (a 1×1 phase at 1.0 monolayer) due to the monovalent character of TI.⁽¹⁾ This feature is in sharp contrast with the trivalent nature found in the $\sqrt{3} \times \sqrt{3}$ phase observed commonly to all other group III elements. However, this has been challenged by observing a stable $\sqrt{3} \times \sqrt{3}$ phase and by finding a T4 site as the most stable binding site.⁽²⁾ The T4 sites for TI have been further confirmed by LEED I(V) analysis.⁽³⁾ From analysis of the crystal truncation rods, we also confirm that TI occupy the T4 binding sites as proposed in previous study and the bondlength of TI-Si agrees quite well within 4%.

[참고문헌]

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