

Giant Rashba spin splitting on monolayer-Pb-covered Ge(111) surface

Koichiro Yaji^{1,2}, Akito Kakizaki¹, and Tetsuya Aruga^{2,3}

¹*Institute for Solid State Physics, The University of Tokyo, Kashiwa, Chiba 277-8581, Japan*

²*JST CREST, Saitama 332-0012, Japan.*

³*Graduate School of Science, Kyoto University, Kyoto 606-8502, Japan.*

1. Introduction

Rashba spin splitting of a surface state band on a semiconductor surface allows us to open a novel physics such as a spin transport at the surface. The Rashba spin-split bands on the semiconductor surfaces have been identified on Si(111) and Ge(111) surfaces so far [1-4]. Among them, monolayer lead (Pb) covered Ge(111) surface (Pb/Ge(111)- β) is remarkable because one of the surface state bands crosses Fermi level with the Rashba spin splitting [4]. We have pointed out that the Pb/Ge(111)- β surface can become an ideal prototype for the spin transport study. On the other hand, the spin structures of the other surface-state bands are unclear. In the present article, we report on a giant Rashba spin-split band of Pb/Ge(111)- β

near the \bar{M} point measured with angle-resolved photoelectron spectroscopy (ARPES) and spin- and angle-resolved photoelectron spectroscopy (SARPES).

2. Experimental methods

A Ge(111) substrate was prepared by several cycles of 0.5-keV Ar^+ bombardment and annealing up to 900 K for a minute. Clean surface of Ge(111) was checked by observing a sharp (2×8) LEED pattern. Pb was deposited onto the surface at room temperature from an alumina crucible heated with a tungsten filament. The surface was then annealed at 570 K for three minutes to prepare a well-ordered wide terrace of Pb/Ge(111) with $(\sqrt{3} \times \sqrt{3})R30^\circ$ periodicity. The Pb coverage of Pb/Ge(111)- β was checked by an intensity ratio between Ge $3d$ and Pb $5d_{5/2}$ core-level spectra, a full-width at half-maximum of the Pb $5d_{5/2}$ core level spectra, and the observation of the sharp $(\sqrt{3} \times \sqrt{3})R30^\circ$ LEED pattern. ARPES and SARPES were performed at BL19A in Photon Factory. The sample temperature was set to room temperature during the measurements.

3. Results and discussion

Figure 1(b) shows the ARPES image taken along $\bar{\Gamma} \bar{M} \bar{\Gamma}$ of $(\sqrt{3} \times \sqrt{3})R30^\circ$ surface Brillouin zone (SBZ) shown in Fig. 1(a). The photon energy ($h\nu$) was set to 21 eV. We

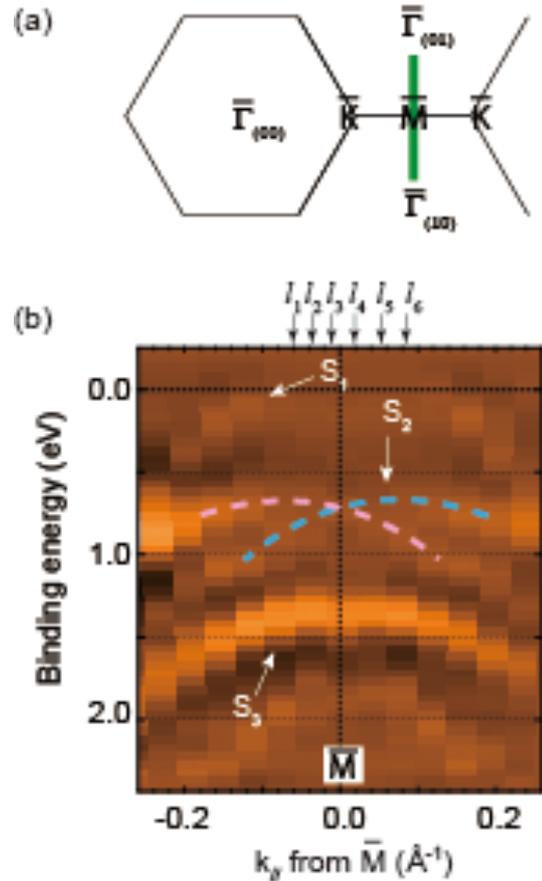


Fig. 1 (a) SBZ of Pb/Ge(111)- β . (b) ARPES image taken along green line in the SBZ shown in (a). Dashed curves represent two branches of S_2 .

found three Pb-induced bands named S_1 , S_2 and S_3 . These bands were also observed with $h\nu = 25, 30$ eV. This indicates that S_1 , S_2 and S_3 are surely surface states or surface resonances. The S_1 band, which shows *free-electron-like* metallic feature, splits into two due to the Rashba effect [4]. The S_2 and S_3 bands are surface resonances because the states appear within the projection of Ge bulk bands. In Fig. 2, the SARPES spectra taken at k points shown by the arrows in Fig. 1(b) are shown. The red and blue spectra represent spin-up and spin-down states, respectively. We found that the spin-up and spin-down branches of S_2 are degenerate at the \bar{M} point. The peaks of S_2 show the spin splitting at k points away from \bar{M} . The peak positions of the spin-up and spin-down branches are inverted around \bar{M} . We therefore conclude that the S_2 band splits into two due to the surface Rashba effect. The Rashba parameter is roughly evaluated to be 1.9 eV \AA , which is comparable with the other giant Rashba systems [1-3]. On the other hand, we could not observe the Rashba spin splitting of S_3 because the intensity of the spin-down state of S_3 is always larger than that of the spin-up state even at \bar{M} .

We found that Pb/Ge(111)- β exhibits the giant Rashba spin-split band around \bar{M} . Similar surface-state bands with the giant Rashba spin splitting around \bar{M} have been identified on Bi adsorbed Si(111) and Ge(111) surfaces [1-3]. Here, the Rashba spin-split bands of Bi/Si(111) and Bi/Ge(111) are ascribed to the bonding orbitals between the Bi atoms near the T_1 sites and the Si (Ge) atoms in the topmost layer of the substrates. Concerning Pb/Ge(111)- β , there are two positions for the Pb adsorption in the unit cell, where one Pb atom is located at an H_3 site and three Pb atoms are located near the T_1 sites. The Pb atoms near the T_1 sites and the Ge atoms in the topmost layer make the Pb-Ge bonding orbitals, which is the origin of the S_2 band [5]. This hybridization can produce the strong perpendicular asymmetry of the wave function, which explains the giant Rashba effect around \bar{M} of Pb/Ge(111)- β . The resemblance of the local adsorption geometries at T_1 and the hybridization between the adsorbates and the substrates may explain the similar giant Rashba spin splitting in Bi/Si(111), Bi/Ge(111) and Pb/Ge(111)- β .

4. Summary

We have studied the electronic structure of Pb/Ge(111)- β by ARPES and SARPES. We found the surface-state band (S_2) with the giant Rashba spin splitting around \bar{M} . The Rashba parameter is evaluated to be 1.9 eV \AA , which is comparable with the other giant Rashba systems such as Bi/Si(111) and Bi/Ge(111). Also, the feature of the giant Rashba band for Pb/Ge(111)- β is similar to those for Bi/Si(111) and Bi/Ge(111). The resemblance comes from the similar local adsorption geometries at T_1 sites and the characters of the bands.

[References]

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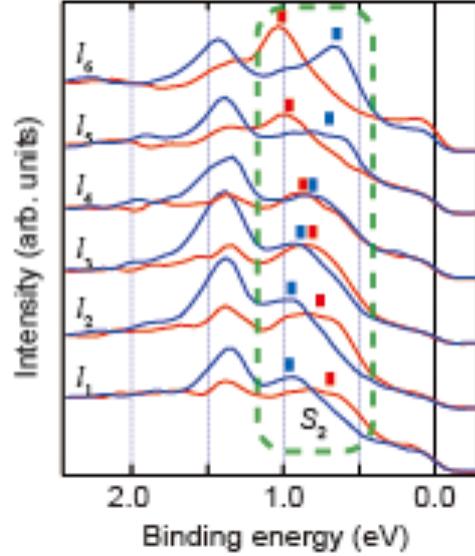


Fig. 2 SARPES spectra taken with the arrows shown in (b). Red and blue lines represent the up and down spin states, respectively.