

## Angle-resolved photoemission studies of Cu(104)-O surface

Kazutoshi YAGI-WATANABE\*<sup>1</sup>, Yasushi WAKIMOTO<sup>2</sup>, Daisuke OGARANE<sup>2</sup>, Takanobu GOTO<sup>2</sup>, Daiichiro SEKIBA<sup>2</sup>

<sup>1</sup>National Institute of Advanced Industrial Science and Technology (AIST), Tsukuba Central-2, Umezono 1-1-1, Tsukuba, Ibaraki 305-85681, Japan

<sup>2</sup>Inst. of Phys., Univ. of Tsukuba, Ten-nodai 1-1-1, Tsukuba, Ibaraki 305-8571, Japan

### Introduction

On metal crystal surfaces with atomically rough structures, faceting can be observed due to adsorption-induced reconstruction. The vicinal (11 $n$ ) copper surface is known to facet to (104) in the presence of adsorbed oxygen [1]. The Cu(104)-O surface is characterized by the formation of Cu-O-Cu-O chains which are aligned along the [010] direction. The structure model of the Cu(104)-O surface is shown in Fig. 1. This surface is very similar to that proposed for the Cu(001)(2 $\sqrt{2}$ ×2 $\sqrt{2}$ )R45-O structure. In a recent paper [2], we report the electronic structure of Cu(001)(2 $\sqrt{2}$ ×2 $\sqrt{2}$ )R45-O surface studied by angle-resolved photoemission spectroscopy (ARUPS) and tight-binding calculation. We examine the electronic structure of the Cu(104)-O surface comparing with that of the Cu(001)(2 $\sqrt{2}$ ×2 $\sqrt{2}$ )R45-O surface.

### Experimental

The angle-resolved photoemission measurements were made on BL-18A with an ADES-500 spectrometer. The clean Cu(104) surface was prepared by repeated Ar<sup>+</sup> sputtering and annealing cycles. We observed LEED patterns with sharp double spots, consistent with a regularly stepped surface structure. The Cu(104)-O surface was prepared by 50 Langmuir oxygen exposure at room temperature. The Cu(104)-O structure was verified by LEED.

### Results and Discussion

Fig.2 shows angle-resolved energy distribution curves (AREDCs) of Cu(104)-O surface taken along [010] with  $h\nu=20$  eV, p-polarization. O2p-derived features are indicated by tick marks in the figure. There are some oxygen-induced features in a region of -7 to -4 eV binding energy. This agrees well with the previous results of ARUPS for Cu(001)(2 $\sqrt{2}$ ×2 $\sqrt{2}$ )R45-O surface. We also observe an oxygen-induced feature at about -1.8 eV. We estimate this to be anti-bonding states. These results will be examined in terms of tight-binding model and compared with the electronic structures for Cu(001)(2 $\sqrt{2}$ ×2 $\sqrt{2}$ )R45-O and Cu(110)p(2×2)-O surfaces [3].

### References

- [1] N.Reinecke et al., Surf. Sci. 454-456, 94 (2000).  
 [2] D.Sekiba et al., Surf. Sci., 470, 43 (2000).

[3] R. Ozawa et al., Surf. Sci., 346, 237 (1996).

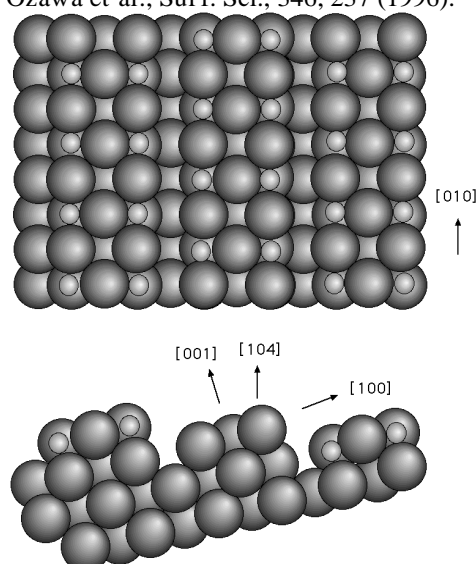


Fig.1. Structure model of the Cu(104)-O surface. Top: top view, bottom: side view. Small balls indicate oxygen atoms.

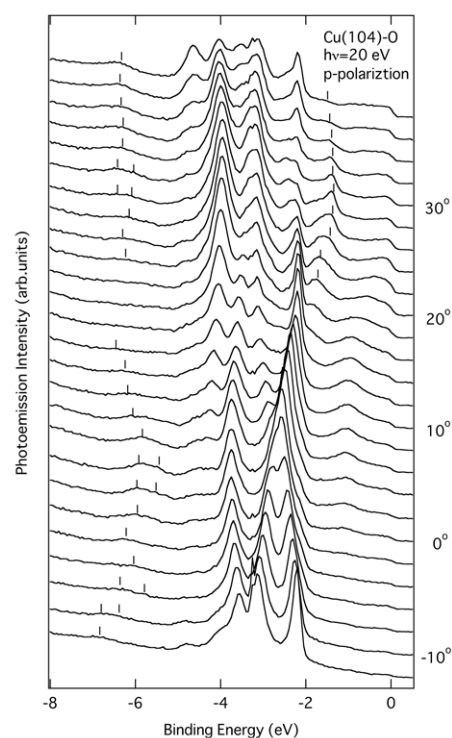


Fig.2. AREDCs of Cu(104)-O surface.

\* yagi-watanabe@aist.go.jp