BL-28A/2012G075, 2012S2-001 n in the iron based superconductor  $\mathbf{P}_{a}\mathbf{F}_{a}$  (As  $\mathbf{P}_{a}$ )

Superconducting gap in the iron-based superconductor  $BaFe_2(As_{1-x}P_x)_2$ 

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## 1 Introduction

The iron-based superconductor  $BaFe_2(As_{1-x}P_x)_2$  has attracted much attention since the presence of superconducting gap nodes was suggested from penetration depth and thermal conductivity measurements [1]. In order to investigate the superconducting gap, we have studied  $BaFe_2(As_{1-x}P_x)_2$  with x = 0.34 ( $T_c = 24$  K) by angle-resolved photoemission spectroscopy (ARPES). Different doping concentrations (x = 0.3, 0.38 and 0.45) also have been studied by ARPES measurement for systematic research on  $BaFe_2(As_{1-x}P_x)_2$ .

## 2 Experiment condition

High-quality single crystals of BaFe<sub>2</sub>(As<sub>0.66</sub>P<sub>0.34</sub>)<sub>2</sub> were grown using the self-flux method. Experiments were carried out at BL-28A of Photo Factory using Scienta SES-2002 analyzer and linearly polarized light, with the photon energy of hv = 40 eV, and the energy resolution of 9-10 meV. The crystals were cleaved *in situ* at T = 9 K, in an ultra-high vacuum of ~  $1.5 \times 10^{-8}$  Pa.

## 3 <u>Result</u>

Figure 1(a) shows the electron Fermi surfaces (FSs) of BaFe<sub>2</sub>(As<sub>0.66</sub>P<sub>0.34</sub>)<sub>2</sub>. The Fermi-surface angle  $\theta$  is defined as shown in Fig. 1(a), in which  $\theta = 0$  along *X*- $\Gamma$  direction. Symmetrized energy distribution curves blow (T = 10 K, blue solid line) and above (T = 30 K, red dashed line) the  $T_c$  along the inner and outer FSs are shown in panels (b) and (c), respectively. Nearly isotropic gaps both on the inner and outer FSs are opened for this doping concentration.

Optimally doped sample (x = 0.3,  $T_c = 30$  K) measured at BL-28A using circularly polarized light with the photon energy of hv = 40 eV, and at BL5-4 of Stanford Synchrotron Radiation Laboratory (SSRL) using linearly polarized light with hv = 28 eV show that anisotropic gaps open on electron FSs. While anisotropy turn to insignificant in this slightly overdoped composition x =0.34 and previous x = 0.38 one (BL-28A, circularly polarized light, hv = 40 eV).



Fig. 1: (a) Electron Fermi surfaces around the X point  $(h\nu = 40 \text{ eV})$  of BaFe<sub>2</sub>(As<sub>0.66</sub>P<sub>0.34</sub>)<sub>2</sub> ( $T_c = 24 \text{ K}$ ). (b), (c) Symmetrized energy distribution curves at  $k_F$  points on the inner and outer FSs.

Reference:

[1] K. Hashimoto. *et al.*, Phys. Rev. B **81**, 220501 (2010)

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