

# Electronic States in Electron-Doped High- $T_c$ cuprate $\text{Nd}_{2-x}\text{Ce}_x\text{CuO}_4$ investigated by Angle-Resolved and Inverse Photoemission Spectroscopies

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The mechanism of superconductivity in conventional superconductors is understood to involve the formation of Cooper pairs via phonons, which is explained by BCS theory. In contrast with “low-temperature” superconductors, cuprate “high-temperature” superconductors were discovered in 1986, which show the highest superconducting transition temperatures ( $T_c$ ), however, as for elucidations of their pairing mechanism, both experimental and theoretical studies have not been achieved yet. One of the most urgent issues in clarifying the mechanism of high- $T_c$  cuprate superconductivity is to understand the origin, namely, boson that mediates Cooper pairing. While electron-boson interactions and strong electron correlations are believed to contribute to the origin of high- $T_c$  cuprate superconductivity [1–3], there is less research into the origin of the boson and the pairing mechanism in electron-doped cuprates than in hole-doped ones. Therefore, to deepen the understanding of the microscopic mechanism of superconductivity through observations of the electronic states in electron-doped cuprates, we have investigated the electronic structure of  $\text{Nd}_{2-x}\text{Ce}_x\text{CuO}_4$  (NCCO), one of the electron-doped high- $T_c$  cuprate superconductors, using angle-resolved photoemission spectroscopy (ARPES) and angle-resolved inverse photoemission spectroscopy (AR-IPES).

In this study, we experimentally performed ARPES and AR-IPES studies to investigate the presence of an emergent band structure arising from the coupling between quantum charge fluctuations and quasiparticles in high- $T_c$  cuprate superconductors. Furthermore, by comparing our experimental results of the electronic structure and calculations obtained from density functional theory (DFT) and dynamical mean-field theory (DMFT) calculations [6].

In addition, to study an insight of the ground state of NCCO, we have studied temperature dependent ARPES study for the spectral weight ( $Z_k$ ) of the quasiparticle band along the nodal direction in ARPES spectra. A previous theoretical study suggested that if resonating-valence-bond-like electron correlations are involved,  $Z_k$  changes markedly across  $T_c$  in the ground state in strongly correlated electron systems such as cuprate superconductors [5]. This prediction has been verified by the temperature dependence of  $Z_k$  reported for hole-doped cuprates [4], and it could offer a new way of understanding the ground state of electron-doped cuprate superconductors.

In the poster presentation, based on experimentally obtained spectra of the electronic structure of NCCO, we discuss the verification of quantum charge fluctuations, including theoretical calculations (DFT+DMFT) and the temperature dependence of the  $Z_k$  spectrum.

## REFERENCES

- 1 A. Lanzara *et al.*, Nature **412**, 510 (2001).
- 2 J. P. Carbotte *et al.*, Nature **401**, 354 (1999).
- 3 H. F. Fong *et al.*, Phys. Rev. Lett. **75**, 316 (1995).
- 4 S. Kudo *et al.*, Phys. Rev. B **92**, 195135 (2015).
- 5 C.-P. Chou *et al.*, J. Phys. Chem. Solids **69**, 2993–2995 (2008).
- 6 H. Yamaguchi *et al.*, arXiv:2505.12639.