Bulletin of the American Physical Society

APS March Meeting 2019

Monday-Friday, March 4-8, 2019; Boston, Massachusetts

Session C09: Superconductivity: Copper Oxide - Theory 2:30 PM-5:30 PM, Monday, March 4, 2019 BCEC Room: 151A

Sponsoring Unit: DCMP

Chair: Michael Sentef, Max Planck Institute for the Structure and Dynamics of Matter

Abstract: C09.00013 : Continuing search for the origin of HTSC: DFT studies of selected copper oxide proxy structures reviewed and paths forward suggested 4:54 PM-5:06 PM

← Abstract →

Presenter:

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In this presentation, we review our past attempts to uncover the pairing mechanism underlying high temperature superconductivity in copper oxide compounds and suggest possible paths forward.^{1, 2} One such path would be to derive and generalize pairing coupling functions to apply to DFT + U computed eigenstates in order to estimate Cooper pair coupling strengths arising from a combination of both lattice and spin excitations. Interestingly, such interactions were found in transition metal alums some 78 years ago as manifested in linking their respective Debye and Curie temperatures.³ We suggest repeating such experiments today on the copper oxide compounds as a function of hole/electron concentration, along with a possible computational strategy to pursue in the interpretation of the results⁴ to finally resolve the fundamental origin of high temperature superconductivity.

- ¹ Paul Grant, http://meetings.aps.org/link/BAPS.2016.MAR.R25.8
- ² P.M. Grant, Journal of Physics: Conference Series 129 (2008) 012042, doi:10.1088/1742-6596/129/1/012042
- ³ C. Starr, Phys. Rev. 60, 241 (1941)
- ⁴ Paul Grant, https://meetings.aps.org/Meeting/FWS17/Session/B2.10

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