### **JAEJUN YU**

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### Profile

Currently Dean of the College of Natural Sciences, Seoul National University, and a professor in physics at Seoul National University. Expert in theoretical and computational condensed matter physics. Research focuses on electronic, magnetic, and other physical properties of novel transition metal oxides and related materials. Excellence in teaching science to non-science majors. Well-versed in university administration and management in promoting research and education.

# Experience

- Dean, College of Natural Sciences, Seoul National University 2022 ~ present
- Director, Center for Theoretical Physics, Seoul National University 2019 ~ 2022
- Dean, Faculty of Liberal Education, Seoul National University 2019 ~ 2022
- Visiting Scientist, MPIPKS, Dresden, Germany 2018.01 ~ 07
- Associate Dean, Faculty of Liberal Education, Seoul National University 2014 ~ 2016
- Treasurer, the Korean Physical Society 2014 ~ 2016
- General Council Member, APCTP (Asia Pacific Center for Theoretical Physics) 2014 ~ 2016
- Associate Dean for Academic Affairs, College of Natural Sciences, Seoul National University
   2012 ~ 2014
- Associate Editor, Philosophical Magazine and Philosophical Magazine Letters, UK 2012 ~ present
- Visiting Professor, The Institute of Scientific and Industrial Research, Osaka University, Japan — 2012 Summer
- Associate Dean for Student Affairs, College of Liberal Studies, Seoul National University 2011~2012
- Visiting (Sabbatical) Scientist, National Renewable Energy Laboratory; Colorado, USA 2010 ~ 2011
- Principal Investigator, Center for Strongly Correlated Materials Research (CSCMR), Seoul National University — 2008 ~ 2012
- Visiting Scholar, Department of Physics and Astronomy, Rutgers University; New Jersey, USA — 2006 Summer
- Executive Editor, Journal of the Korean Physical Society (JKPS) 2007 ~ 2007
- Assistant and Associate Professor, Physics Department, Sogang University 1992 ~ 1999
- Research Associate, Physics Department, the University of Illinois at Urbana-Champaign;
   Illinois, USA 1991 ~ 1992

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 Research Associate, Physics Department, Northwestern University; Illinois, USA — 1988 ~ 1990

## Education

- Seoul National University, Seoul, Korea B.S in Physics, 1984
- Northwestern University, Illinois, USA M.S. in Physics, 1986
- Northwestern University, Illinois, USA Ph.D. in Physics, 1988

### **Awards**

- SNU Awards: Excellence in Teaching, Seoul National University; Seoul, Korea 2012
- Graduate Student Awards, 1988 MRS Spring Meeting; Reno, Nevada, USA 1988

## Research Interests

Jaejun Yu's research interests lie primarily in condensed matter theory of strongly correlated electron materials and computational materials physics. Of particular interest is understanding and discovering new physical properties of novel materials, including transition metal compounds and low-dimensional systems. He is also involved in developing theoretical and computational methods and their application to study real materials' electronic and magnetic properties. Current research focuses on understanding topological band structures of novel transition metal compounds, e.g., Ir-oxides or chalcogenides, exploring new two-dimensional materials manifesting interesting topological characteristics of topological insulators, Chern insulators, or fractional Chern insulators, and manipulating defects in metal oxides for *p*-type oxide semiconductors, catalysts, and possible device applications.

# Research Highlights

### **CITATION METRICS**

According to Google Scholar (<a href="https://goo.gl/hT33xf">https://goo.gl/hT33xf</a>), the number of citations received by his publications is more than 13,271 as of June 2024. And, the impact measures are the h-index of 53 and the i10-index of 137 out of 170 published papers.

The list of publications is available at <a href="https://goo.gl/hT33xf">https://goo.gl/hT33xf</a>.

# SELECTED LIST OF PUBLICATIONS

- B. J. Kim, Hosub Jin, S. J. Moon, J-Y. Kim, B-G. Park, C. S. Leem, <u>Jaejun Yu, et al.</u> "Novel J<sub>eff</sub>= 1/2 Mott state induced by relativistic spin-orbit coupling in Sr<sub>2</sub>IrO<sub>4</sub>." Physical Review Letters 101, 076402 (2008). [cited by 1819]
- <u>Jaejun Yu</u>, Arthur J. Freeman, and J-H. Xu. "Electronically driven instabilities and superconductivity in the layered La<sub>2-x</sub>Ba<sub>x</sub>CuO<sub>4</sub> perovskites." Physical Review Letters 58, 1035 (1987). [cited by 712]
- Hosik Lee, Young-Woo Son, Noejung Park, Seungwu Han, and <u>Jaejun Yu</u>. "Magnetic ordering at the edges of graphitic fragments: Magnetic tail interactions between the edge-localized states." Physical Review B 72, 174431 (2005). [cited by 554]
- S. J. Moon, Hosub Jin, Kyung Wan Kim, W. S. Choi, Y. S. Lee, <u>Jaejun Yu</u>, G. Cao, *et al.* "Dimensionality-Controlled Insulator-Metal Transition and Correlated Metallic State in 5d

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Transition Metal Oxides  $Sr_{n+1}Ir_nO_{3n+1}$  (n=1,2, and  $\infty$ )." Physical Review Letters 101, 226402 (2008). [cited by 568]

• S. Massidda, <u>Jaejun Yu</u>, A. J. Freeman, and D. D. Koelling. "Electronic structure and properties of YBa<sub>2</sub>Cu<sub>3</sub>O<sub>7- $\delta$ </sub>, a low dimensional, low density of states superconductor." Physics Letters A 122, no. 3-4 (1987): 198-202. [cited by 608]

#### **RESEARCH ACHIEVEMENTS**

Transition metal oxides. He started his research career by publishing papers on the electronic band structures of famous high  $T_c$  Cu-oxide superconductors. The first one of the series was published in Phys. Rev. Lett. (1987) (cited more than 720 times) with Prof. A. J. Freeman at Northwestern University, which is one of the first reports on the electronic structure theory of high T<sub>c</sub> Cu-oxides, uncovering the 2D nature of the band structure arising from the strong hybridization of Cu d and O p orbitals within the 2D square lattice of the Cu-O layer, and became a solid standard of following electronic structure calculations. Later his research evolved by building on DFT (density functional theory) calculations and elaborating model analyses for the physical properties of real materials. He has made numerous contributions to understanding novel transition metal oxides and compounds based on the DFT works. After high T<sub>c</sub> cuprates, he continued his research on related transition metal oxide materials, including CMR (colossal magnetoresistance) manganese oxide and associated compounds, heterostructures, and superlattices of transition metal oxides, and extended his scope to the system with orbital degrees of freedom competing with effective spin interactions, and the understanding of spin and orbital orderings and results of optical spectroscopy in collaboration with experimental groups. He recently explored several oxide superlattices and clarified the role of oxygen vacancies in generating charge carriers at the interface of LaAlO<sub>3</sub> overlays on the SrTiO<sub>3</sub> (001) surface. Further, he searched for a new type of oxide semiconductors with ptype carriers, e.g., BaSnO<sub>3</sub>, and investigated the role of defects in various oxide materials with applications to electronic devices and catalysts.

 $J_{\text{eff}}$ =1/2 physics. Gaining experience in 3d transition metal oxides, he looked into the physics of 4d and 5d transition metal oxides, where the spin-orbit-coupling (SOC) interactions become significant and can play a crucial role in determining their electronic and magnetic properties. While moving from 3d to 4d and 5d metal oxides, the bandwidth gets broadened, and the strength of on-site Coulomb interactions becomes relatively weaker. Thus, in contrast to the insulating ground state of its 3d oxides, it is natural to expect a metallic ground state for the 4d and 5d compounds. To one's surprise, Sr<sub>2</sub>IrO<sub>4</sub> was reported to be insulating with an optical gap of about 0.5 eV, contrary to the metallic Sr<sub>2</sub>RhO<sub>4</sub> of its 4d counterpart. Indeed, it remained a puzzle until he came up with a solution that the SOC combined with a moderate on-site Coulomb interaction can give rise to a new kind of entangled quantum state manifesting the spin-orbit coupled J-state with an effective angular momentum of  $J_{\rm eff}$ =1/2. The idea of  $J_{\rm eff}$ =1/2 was first suggested in his seminal paper of Phys. Rev. Lett. (2008) (cited more than 1,550 times) has significantly influenced various aspects of correlated electron physics with both spin and orbital degrees of freedom, such as superconductivity, quantum magnetism, topological insulators, metal-insulator transition, and so on. Recently, by expanding the idea of the spinorbit coupled state, he has proposed that a (111)-oriented BaBiO<sub>3</sub> bilayer sandwiched between large-gap perovskite oxides is a promising candidate for topological electronics by realizing a gap-tunable and topology-tunable system, enabling the switch between quantum spin Hall and quantum valley Hall phases.

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**2D** materials. Magnetism in 2D materials has attracted much interest due to its potential applications to spintronics and nano-electronic devices. He has made interesting observations on atomically thin materials' charge and magnetic states during the last decade, primarily in graphene and single-layer boron nitride (BN) systems. One of his early works on graphene nanoribbon revealed the formation of local moments and their interactions at the edge of graphitic fragments—this work was published in Phys. Rev. B (2005) (cited more than 520 times) has provided a clear and lucid picture of the magnetic edge states, consequently leading to enormous activities for possible manipulation of ferromagnetic edge states for their potential applications. He has recently proposed that the charge and magnetic conditions of the transition-metal adatom on graphene or single-layer BN can be effectively controlled by either defects or external electric fields.

**DFT code (OPENMX) project.** In collaboration with Prof. Taisuke Ozaki at ISSP, the University of Tokyo, he has been actively involved in the development of the DFT code called OpenMX (Open source package for Material eXplorer; http://www.openmx-square.org), which is a software package for nano-scale material simulations based on DFT, aiming for the realization of large-scale *ab initio* electronic structure calculations. Indeed, the early development of non-collinear spin DFT code within OpenMX, incorporating both spin-orbit coupling and LDA+U, made it possible to discover the  $J_{\rm eff}$  = 1/2 state in Sr<sub>2</sub>IrO<sub>4</sub> Ir oxides. Taking advantage of its local orbital basis formulation, he developed an approach to derive effective Hamiltonians based on Wannier functions derived from DFT calculations. He utilizes this approach to analyze various materials' topological band structure and magnetism.

## Selected List of Conference Activities

- Asian Workshop on First-Principles Electronic Structure Calculations, 1998 present, serving for 19 years as a core member of the International Organizing Committee, contributing to the DFT community in east Asian countries for the computational condensed matter physics and materials sciences.
- Korea-Japan-Taiwan Symposium on Strongly Correlated Electron Systems, 2000 present, serving for 17 years as a core member of the International Organizing Committee, contributing to the promotion of strongly correlated electron physics in the physics community in Korea, Japan, and Taiwan.
- APCTP-IACS (Korea-India) Joint Symposium on Emergent Phenomena in Novel Oxide Materials and Low Dimensional Systems, 2008 present, serving for 14 years as a convener of the bilateral symposium for the promotion of scientific activities and exchange program among the physics communities of Korea and India.
- APCTP-QMS (Quantum Materials Symposium), 2014 present serving as a core member of the Organizing Committee and now as an organizing chair contributing to the condensed matter physics community, especially of strongly correlated electron materials, in Korea.
- International Conference on Strongly Correlated Electron Systems (SCES) 2023, 2022 2023, serving as a co-chair of the Conference, which will be held in Incheon, Korea, from 2nd to 7th July 2023.