Riccardo Piombo



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ABOUT ME

I am a condensed matter physicist proficient in several programming languages and expert in numerical simulation of strongly correlated system, and phase change memories.

• WORK EXPERIENCE

Rome, Italy
POST DOCTORAL RESEARCHER CREF

Development of optimal transport methods in economic complexity from maximum entropy probability distributions.

Rome, Italy

POST DOCTORAL RESEARCHER UNIVERSITÀ DI ROMA - LA SAPIENZA

Neuromorphic computing - Ab-Initio and classical molecular dynamics simulations of chalcogenide heterostructures using neural network interatomic potentials.

EDUCATION AND TRAINING

Rome, Italy

PHD IN PHYSICS Università di Roma - La Sapeinza

I focused on the chemistry of silver fluoride (AgF2) as a potential analog to cuprates, materials exhibiting hightemperature superconductivity. I modelized several AgF2 experimental data, including valence band photoemission and Meitner-Auger spectroscopy. I strongly relied on DFT and exact diagonalization computations to estimate AgF2 Coulomb interaction and its charge-transfer energy. My investigations revealed that AgF2 is a charge-transfer correlated insulator within the Zaanen-Sawatzky-Allen scheme. Moreover, AgF2 possesses a more covalent nature compared to cuprates despite sharing similarities in electronic excitations: we provided detailed insights into AgF2 charge-transfer gap, d-d excitations and a potential proximity to a charge-transfer instability. Finally, I explored with exact diagonalization techniques the spin spectra of several cuprates: La2CuO4, Sr2CuO2Cl2 and CaCuO2.

Final grade Excellent | Thesis Correlation and Cuprate-like Physics in Silver Fluorides

Rome, Italy M.SC. IN THEORETICAL PHYSICS Università di Roma - La Sapeinza

I studied a simplified model of AgF or the parent compound of this hypothetical new class of high-\$T_c\$ superconductors: one-dimensional linear chains formed by Ag and F sites with strong inter-site and on-site Coulomb interactions. Here, I only considered the charge degrees of freedom to study the valence instability of the compound. The hypothesis was the following: when holes localized on the silver ions tend to move collectively on the fluorine ions, excitonic soft modes appear during the phase transition. I analyzed the phase diagram and the dynamical charge structure factor of the system using a Quantum Lattice Model Simulator and Green's function theory to determine the characteristics of the transitions close to the valence instability.

Final grade 109/110 | **Thesis** Valence instabilities and electronic soft modes

PUBLICATIONS

2024

Ab Initio Study of Novel Phase-Change Heterostructures

Phys. Rev. B, 106, 035142

2022

Multiple-Magnon Excitations Shape the Spin Spectrum of Cuprate Parent Compounds

Phys. Rev. B 103, L140409

2022

Charge Transfer and d – d Excitations in AgF2

Phys. Rev. Research 4, 023108

HONOURS AND AWARDS

Contributi Premiali regione lazio - Regione lazio

Selected researchers received a €2000 grant to encourage more and higher-quality scientific publications, awarded after a thorough peer-review of their work.

Avvio alla Ricerca Tipo II Starting Grant

NANO PCM - New Advanced NeurOmorphic Phase-Change Materials Project AR2221816834DCFB funded with €4000.

Various HPC projects

More than 160000 cpu/hours in CINECA ISCRA projects (1 KNL CPU hour \approx \in 0.10)

CONFERENCES AND SEMINARS

Rome EPCOS

Contributed talk: Structural and Dynamical Properties of the Liquid Phase of New Phase-Change Heterostructures

LANGUAGE SKILLS

Mother tongue(s): ITALIAN

Other language(s):

	UNDERST	UNDERSTANDING		SPEAKING	
	Listening	Reading	Spoken production S	ı	
ENGLISH	C1	C2	C1	C1	C1

Levels: A1 and A2: Basic user; B1 and B2: Independent user; C1 and C2: Proficient user

"Autorizzo il trattamento dei miei dati personali ai sensi del Decreto Legislativo 196/2003, coordinato con il Decreto Legislativo 101/2018, e dell'art. 13 del GDPR (Regolamento UE 2016/679) ai fini della pubblicazione in Amministrazione Trasparente - Sapienza come da normativa vigente."