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## Electronic structure of the metal-to-insulator transition in VO2

The strongly correlated material  $VO_2$  displays a metal-to-insulator (MIT) transition when going below  $T_{MIT} = 280K$ . Alongside this electronic transition, the material undergoes a structural transition from a rutile structure in the metallic phase to a monoclinic structure in the insulating phase. These simultaneous transitions have created a long-lasting debate within the community: is the electronic transition induced by the structural changes (Peierls transition) or is it happening alongside it (Mott transition) [1]? This question has been nicknamed the chicken-and-egg dilemma [2] of condensed matter. Recent ARPES studies addressed the changes of the electronic structure of  $VO_2$  across the transition [3].

However, a detailed imaging of the evolution of the conduction band spectral function in the transition regime is still lacking. I will present our ongoing ARPES studies on  $VO_2$  where we were able to observe a progressive transfer of spectral weight between two distinct states composing the conduction band.

## References:

[1] Dynamical Singlets and Correlation-Assisted Peierls Transition in VO2, Silke Biermann et al., Physical Review Letters (2005)

[2] Resolving the VO2 controversy: Mott mechanism dominates the insulator-to-metal transition, O. Nájera et al., Physical Review B. (2017)

[3] Photoelectron dispersion in metallic and insulating VO2 thin films, Viktor Jonsson et al., Physical Review Research (2021)

## Affiliation de l'auteur principal

ISMO, Université-Paris-Saclay, Orsay

## Auteur principal: DAVID, Emma (ISMO, Orsay)

**Co-auteurs:** SHIGA, Daisuke (KEK Photon Factory, Japan); JHA-THAKUR, Amitayush (ISMO, Orsay); THEES, Maximilian (ISMO, Orsay); HENRIQUE REZENDE GONÇALVES, Pedro (ISMO, Orsay); ANTEZAK, Alexandre (ISMO, Orsay); CHENG, Xianglin (KEK Photon Factory, Japan); KIM, Taehyun (KEK Photon Factory, Japan); KANDA, Tatsuhiko (KEK Photon Factory, Japan); FRANTZESKAKIS, Emmanouil (ISMO, Orsay); FORTUNA, Franck (ISMO, Orsay); MAYNE, Andrew (ISMO, Orsay); KUMIGASHIRA, Hiroshi (KEK Photon Factory, Japan); SANTANDER, Andres (ISMO, Orsay)

Orateur: DAVID, Emma (ISMO, Orsay)

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