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Correlation effects in KFe_2As_2, $RbFe_2As_2$ and $CsFe_2As_2$[1] STEFFEN BACKES, HARALD O. JESCHKE, ROSER VALENTI, University Frankfurt — We perform a systematic LDA+DMFT study of the iron-pnictide series KFe_2As_2[2], $RbFe_2As_2$ and $CsFe_2As_2$ and compare with available experiments. We find not only strong orbital-dependent renormalizations and Fermi surface topology changes compared to the local-density approximation but also interesting features at higher binding energies. We discuss the observation of a possible orbitally-selective instability as a function of isoelectronic doping[3].

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[2]Backes *et al.*, New J. Phys. **16**, 083025 (2014)
[3]Backes *et al.*, in preparation

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