

# Charge-ordered phases in the extended Hubbard model

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Various charge ordered states are relevant to a broad range of materials, e.g. manganites, cuprates, magnetite, doped transition metal compounds ( $\text{Ti}_{4x}\text{V}_x\text{O}_7$ ,  $\text{WO}_{3-x}$ ), heavy-fermion systems and organic compounds. The extended Hubbard model is one of the simplest models that captures the interplay between strong correlations and charge-ordering effects [1-3]. The model can describe the insulator-metal transition between phases with long-range charge-order [1]. We present studies of the model with both (i) on-site interaction  $U$  and (ii) intersite density-density interactions between nearest-neighbors  $W_1$  and next-nearest-neighbors  $W_2$  beyond the standard two-sublattice assumption [2,3]. In particular, we investigate the effects of next-nearest-neighbor interactions on phase diagrams of the model in different limits and we show that charge-stripes can occur for repulsive  $W_2 > 0$  for both signs of  $W_1$  [2].

## References:

- [1] K. J. Kapcia, S. Robaszkiewicz, M. Capone, A. Amaricci, Phys. Rev. B (2017), *accepted*; arXiv:1611.03455
- [2] K. J. Kapcia, J. Barański, S. Robaszkiewicz, A. Ptok, J. Supercond. Nov. Magn. **30**, 109 (2017).
- [3] K. J. Kapcia, S. Robaszkiewicz, Physica A **461**, 487 (2016).

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