

# Single Crystal Studies of Charge Density Wave Physics in Quasi-1D Metals $RNiC_2$

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Ternary carbides  $RNiC_2$  ( $R$  = rare earths and Y) crystallize in the non-centrosymmetric orthorhombic  $CeNiC_2$ -type structure, space group  $Amm2$ . These compounds attracted attention, because of various interesting properties such as superconductivity, magnetism, multiple charge density wave (CDW) transitions related to quasi-one-dimensional electronic features of  $RNiC_2$ , and finally, a complex interplay of CDW order and rare earth magnetism in  $RNiC_2$  (see Ref. [1] for a review).

Here, we report on crystal growth, crystallographic characteristics explored via single crystal XRD, as well as magnetic, thermodynamic, and electronic properties studied by a variety of techniques revealing orientation dependent, anisotropic features of selected  $RNiC_2$  single crystals. Based on heat capacity, thermal expansion and electrical resistivity studies, we analyse the nature of multiple CDW transitions occurring in  $RNiC_2$  crystals (e.g., orthorhombic to incommensurate CDW and incommensurate to commensurate CDW transitions). We discuss their relations to electronic and crystal structure changes and, employing Clausius-Clapeyron and generalized Ehrenfest relations, we discuss resulting pressure and strain dependencies of the observed CDW phases.

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[1] V. Babizhetskyy, B. Kotur, V. Levytskyy, H. Michor, Chapter 298: Alloy systems and compounds containing rare earth metals and carbon, in: J.-C.G. Bünzli, V.K. Pecharsky (Eds.), Handbook on the Physics and Chemistry of Rare Earths, Vol. 52, North-Holland, Amsterdam, pp. 1-263. (2017).