

# Unusual magnetic properties of layered $M\text{CrS}_2$ compounds

A. N. Yaresko,<sup>1</sup> D. A. Kukusta,<sup>1,2</sup> A. V. Ushakov,<sup>1,3,4</sup> and D. I. Khomskii<sup>3</sup>

<sup>1</sup>*Max-Planck-Institut für Festkörperforschung,  
Heisenbergstraße 1, D-70569 Stuttgart, Germany*

<sup>2</sup>*Institute for Metal Physics, 36 Vernadskyi Bld., UA-03680 Kiev, Ukraine*

<sup>3</sup>*II. Physikalisches Institut, Universität zu Köln,  
Zùlpicherstraße 77, D-50937 Köln, Germany*

<sup>4</sup>*Institute for Theoretical Physics, Clausthal University of Technology,  
Leibnizstraße 10, D-38678 Clausthal Zellerfeld, Germany*

## Abstract

A chromium sulfide  $\text{AgCrS}_2$  has recently attracted considerable attention due to its multiferroic properties and unusual double-stripe magnetic order which appear below the Néel temperature of 50 K.  $\text{AgCrS}_2$  belongs to a family of layered  $M\text{CrS}_2$  ( $M^+=\text{Li, Na, K, Cu, Ag, Au}$ ) compounds in which magnetic  $\text{Cr}^{3+}$  ( $3d^3$ ) ions form a triangular lattice. Although charge and orbital degrees of freedom in these compounds are frozen, geometrical frustrations of magnetic interactions, inherent in the triangular lattice, lead to fascinating magnetic properties: depending on the size of an  $M^+$  ion magnetic order in Cr planes changes from non-collinear  $120^\circ$  (Li) antiferromagnetic, to double stripes (Ag, Au), and, finally, to ferromagnetic (K). In order to understand this strong variation of the magnetic properties we calculated band structures and total energies  $E(\mathbf{q})$  as a function of a wave vector  $\mathbf{q}$  for various spin spiral structures. Effective exchange coupling constants  $j$  between Cr spins are then estimated by fitting  $E(\mathbf{q})$  to an appropriate classical Heisenberg model. We found that depending on the  $M$  size the nearest neighbor coupling  $j_1$  changes from strongly antiferromagnetic in  $\text{LiCrS}_2$  to ferromagnetic in  $\text{KCrS}_2$ , whereas the coupling  $j_3$  between the 3-rd Cr neighbors is strong and remains nearly constant in all the compounds. In  $\text{AgCrS}_2$  with  $j_1 \ll j_3$  the double stripe magnetic order is stabilized by monoclinic distortions of the crystal structure. We discuss the microscopic origin of various  $j$  and show that similar considerations help to understand the magnetic properties of other frustrated Cr compounds  $A\text{Cr}_2\text{S}_4$  with a spinel structure.