

# Kristallstruktur und Eigenschaften von Chrom(II)-Tetrachloroaluminat

Crystal Structure and Properties of Chromium(II)-Tetrachloroaluminate

Rainhard Müller<sup>a</sup>, Grigori V. Vajenine<sup>b</sup> und Ulrich Keßler<sup>a</sup>

<sup>a</sup> Institut für Anorganische Chemie, Rheinische Friedrich-Wilhelms-Universität Bonn,  
Gerhard-Domagk-Straße 1, D-53121 Bonn, Germany

<sup>b</sup> Max-Planck-Institut für Festkörperforschung, Heisenbergstraße 1, D-70569 Stuttgart, Germany;  
Institut für Anorganische Chemie, Universität Stuttgart, Pfaffenwaldring 55, D-70569 Stuttgart,  
Germany

Reprint requests to Dr. Ulrich Keßler. Fax: (+49)(0)228/73-5660. E-mail: ukessler@uni-bonn.de

*Z. Naturforsch.* **2007**, 62b, 625 – 632; received February 14, 2007

Chromium(II)tetrachloroaluminate can be prepared by reaction of chromium, chromium trichloride and aluminum trichloride in the molar ratio 1 : 1.94 : 5.79 or by reaction of stoichiometric amounts of chromium dichloride and aluminum trichloride. Sublimation of the compound yields phase-pure samples of colourless crystals.  $\text{Cr}(\text{AlCl}_4)_2$  crystallizes in the non-centrosymmetric orthorhombic space group  $Pca2_1$  (No. 29),  $a = 1511.38(3)$ ,  $b = 604.71(1)$ ,  $c = 1301.76(3)$  pm. The aggregation of  $\text{Cr}(\text{AlCl}_4)_2$  molecules leads to a square-bipyramidal coordination of the chromium atom typical of high-spin  $d^4$  ions due to the Jahn-Teller effect. The relationship to centrosymmetric  $\text{Pd}(\text{AlCl}_4)_2$  ( $P2_1/c$ ) which is built up of similar molecules was established by means of a group-subgroup symmetry tree. The temperature dependence of the magnetic susceptibility agrees with Curie-Weiss behaviour ( $\mu_{\text{eff}} = 4.70 \mu_{\text{B}}$ ,  $\Theta = -1.3$  K).

*Key words:* Chromium, Tetrachloroaluminate, Crystal Structure, Jahn-Teller Distortion,  
Group-Subgroup Relations