

# DSC Investigations of the Phase Transitions of $[\text{M}(\text{NH}_3)_6](\text{ClO}_4)_2$ and $[\text{M}(\text{NH}_3)_6](\text{BF}_4)_2$ , where $\text{M} = \text{Co}$ and $\text{Cd}$

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Z. Naturforsch. **54 a**, 590–594 (1999); received August 13, 1999

Solid polymorphism of four compounds of the type  $[\text{M}(\text{NH}_3)_6](\text{XY}_4)_2$ , where  $\text{M} = \text{Co}^{2+}$  or  $\text{Cd}^{2+}$ , and  $\text{XY}_4 = \text{ClO}_4^-$  or  $\text{BF}_4^-$  has been studied at 100 - 300 K by DSC. One or two phase transitions of the investigated compounds have been found. For the compounds with  $\text{M} = \text{Co}$  the phase transitions have not yet been described in the literature. For the compounds with  $\text{M} = \text{Cd}$  the phase transition-temperature is in good agreement with the results obtained by NMR and EPR. Generally, for  $[\text{M}(\text{NH}_3)_6](\text{BF}_4)_2$  compounds ( $\text{M} = \text{Mg}, \text{Fe}, \text{Co},$  or  $\text{Ni}$ ) the phase transition temperature  $T_{\text{C1}}$  is lower than that for the corresponding  $[\text{M}(\text{NH}_3)_6](\text{ClO}_4)_2$ , but for compounds with  $\text{M} = \text{Cd}$  it is higher. However, the enthalpy and entropy changes at the  $T_{\text{C1}}$  phase transitions of  $[\text{M}(\text{NH}_3)_6](\text{BF}_4)_2$  are always lower than those for  $[\text{M}(\text{NH}_3)_6](\text{ClO}_4)_2$ . Moreover, for the compounds of this type a correlation between the transition temperature  $T_{\text{C1}}$  and the crystal lattice parameter  $a$  has been found.

*Key words:* Phase Transitions; DSC Method; Chlorate(VII) and Tetrafluoroborate of Hexaaminacobalt(II) and Hexaaminacadmium(II).