

## Magnetic anisotropy and magnetic transitions in RFe<sub>10</sub>Mo<sub>2</sub>

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Compounds of the composition RFe10 Mo2 with R = Y, Nd, Sm, Gd, Tb, Dy, Ho, Er, Tm and Lu were studied by means of the ac-initial susceptibility (x) and the singular point detection (SPD) technique in the temperature range from 4.2 to 300 K. Spin reorientation transitions were detected for the compounds RFe10Mo2 with R = Nd and Dy. For the compounds RFe10Mo2 with R = Y, Ho, Er, Tm and Lu the ac-initial susceptibility data indicate a magnetic transition of unidentified nature. For SmFe 10 Mo2 a first-order magnetization process (FOMP) is observed below 170 K.

The ThMn<sub>12</sub> structure with RFe<sub>12</sub> was found to be unstable. However, compounds of the composition  $RFe_{12-x}M_x$  (M = Ti, V, Cr, Si, Mo; x = 1 or 2) with ThMn<sub>12</sub> structure can be formed for various rare earths. Such compounds are interesting for permanent magnet applications because of their high Fe concentration (causing possibly a high saturation magnetization) and their uniaxial crystal structure (causing possibly high magnetocrystalline anisotropy which is basic for a high coercivity). Since 1987, compounds of the type RFe<sub>11</sub>Ti, RFe<sub>10</sub>V<sub>2</sub> have been studied intensively [1-4]. In this contribution experimental investigations on RFe10 Mo2 are reported.

Samples of RFe<sub>10</sub>Mo<sub>2</sub> with R = Y, Nd, Sm, Gd, Tb, Dy, Ho, Er, Tm and Lu were prepared by arc melting and were subsequently homogenized at 850°C. The homogenized samples have been checked by X-ray diffraction to be single phase with tetragonal structure. In order to detect a possible temperature-induced magnetic transition, all samples were subjected to measurements of the temperature dependence of the ac-initial susceptibility in an ac magnetometer in the temperature range from 4.2 to 300 K using an ac-field of 40 A/m and a frequency of 125 Hz. The spin reorientation transitions, described as a change of the easy direction of magnetization (EDM) from one crystallographic direction to another with varying temperature, were detected for RFe10 Mo2 with R = Nd (at 180 K) and R = Dy (at 143 and 63 K). No spin reorientation transition was traced for TbFe10Mo2. Figure 1 shows the temperature dependence of the ac-initial susceptibility ( $\chi$ ) and its first derivative ( $d\chi/dT$ ) for DyFe10 Mo2. It follows from this figure that two peaks are evident in the curve of the  $d\chi/dT$  versus T. Similar to the case of DyFe11Ti [5], the following explanation can be given for DyFe10 Mo2. The peak at the higher temperature (at about 143 K) indicates a change of the EDM from the c-axis to a cone, whereas the

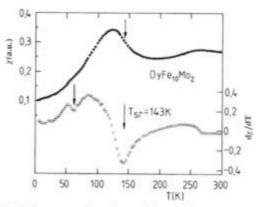


Fig. 1. Temperature dependence of the ac-initial susceptibility  $\chi$  (solid circles) and its first derivative  $d\chi/dT$  (open circles) of DyFe10Mo2.

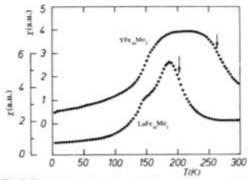


Fig. 2. Temperature dependence of the ac-initial susceptibility x of YFe10Mo2 and LuFe10Mo2.

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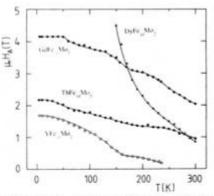


Fig. 3. Temperature dependence of the anisotropy field H<sub>A</sub> determined by the SPD technique for YFe<sub>10</sub>Mo<sub>2</sub>. GdFe<sub>10</sub>· Mo<sub>2</sub>. DyFe<sub>10</sub>Mo<sub>2</sub> and TbFe<sub>10</sub>Mo<sub>2</sub>.

peak at the lower temperature (at about 63 K) indicates a change of the EDM from the cone to the basal plane. Figure 2 shows the temperature dependence of the ac-initial susceptibility of YFe10Mo2 and LuFe10 Mo2. A very pronounced peak can be observed for YFe10Mo2 at 266 K and for LuFe10Mo2 at 204 K. A similar phenomenon was also found for compounds  $RFe_{10}Mo_2$  with R = Ho (231 K), Er (219 K) and Tm (204 K). All these experimental facts together with those of the singular point detection (SPD) measurements (see YFe10Mo2 in fig. 3) force us to the conclusion that a temperature-induced magnetic phase transition occurs in the Fe sublattice of RFe10Mo2 compounds. According to this assumption the magnetic transition observed in HoFe10 Mo2, ErFe10 Mo2 and TmFe<sub>10</sub>Mo<sub>2</sub> can be understood as being due to a competition between the Fe sublattice anisotropy and the very low anisotropy of the Ho-, Er- and Tm sublattice. The nature of this unidentified transition of the Fe sublattice has been explained as the temperatureinduced competition of the anisotropy among the different Fe sites.

The SPD theory predicts a singularity in the  $d^2M/dH^2$  versus H curve at  $H=H_A$  for an uniaxial compound for which the external field is applied perpendicular to the EDM. Figure 3 shows the temperature dependence of the anisotropy field  $H_A$  for YFe<sub>10</sub>Mo<sub>2</sub>, GdFe<sub>10</sub>Mo<sub>2</sub>, DyFe<sub>10</sub>Mo<sub>2</sub> and TbFe<sub>10</sub>Mo<sub>2</sub>. From the results of YFe<sub>10</sub>Mo<sub>2</sub> and GdFe<sub>10</sub>Mo<sub>2</sub>, it is concluded that the Fe sublattice anisotropy of RFe<sub>10</sub>Mo<sub>2</sub> is uniaxial. The anisotropy of TbFe<sub>10</sub>Mo<sub>2</sub> is also uniaxial, even down to 4.2 K, whereas the anisotropy of DyFe<sub>10</sub>Mo<sub>2</sub> is uniaxial only above 143 K, which is in agreement with the susceptibility measurement (see fig. 1). Figure 4 shows the temperature dependence of the ac-initial susceptibility  $\chi$  and its

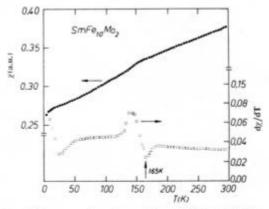


Fig. 4. Temperature dependence of the ac-initial susceptibility  $\chi$  (solid circles) and its first derivative  $d\chi/dT$  (open squares) of SmFe<sub>10</sub>Mo<sub>2</sub>.

first derivative  $d_X/dT$  for SmFc<sub>10</sub>Mo<sub>2</sub>. Due to the large uniaxial anisotropy of the Sm sublattice, no spin reorientation transition is predicted for SmFc<sub>10</sub>Mo<sub>2</sub>. However, for SmFc<sub>10</sub>Mo<sub>2</sub> a change of the shape of the  $\chi(T)$  curve (resulting in a peak in the curve of the  $d_X/dT$  versus T) is evident (see fig. 4). According to our previously experience on Nd<sub>2</sub>Fc<sub>14</sub>B [6], this change might give an evidence for the onset temperature of a first-order magnetization process. In order to verify this prediction, a direct measurement of the magnetization process at various temperatures from 4.2 to 300 K was performed for SmFc<sub>10</sub>Mo<sub>2</sub> in a pulsed field system by the SPD technique. Figure 5 shows the measured results. From this figure it is evident that, as

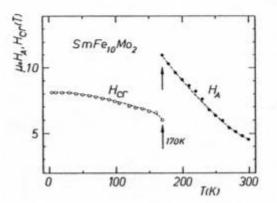


Fig. 5. Temperature dependence of the anisotropy field H<sub>A</sub> and critical FOMP field H<sub>cr</sub> of SmFe<sub>10</sub>Mo<sub>2</sub>.

predicted, a type-II FOMP occurs indeed in  $\rm SmFe_{10}$  Mo  $_2$  below 170 K.

## References

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