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K. Takegahara, A. Yanase, T. Kasuya

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Erratum

Unusual anisotropy mechanism in Ce-pnictides due to virtual valence fluctuation.

K. TAKEGAHARA, A. YANASE and T. KASUYA


Figures 1 and 2 are incorrect because there is an error in the numerical calculations of equations (4), (5) and (6). The correct results are shown in figures 1 and 2. Therefore, the easy direction is always in [001] and the hard direction [111] in good agreement with the experiments. The anisotropy energy in case (ii) is smaller than that of case (i) and about a half of the value listed in the paper.

Fig. 1. — Calculated energy shift \( U \) for case (i). Equations (4) and (5) give the same result.

Fig. 2. — Calculated energy shift \( U \) for case (ii).