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Errata to Landolt-Börnstein subvolume III/15a

- p. 125, Fig. 98. Rb the ordinate lable $(\varrho - \varrho_0)$ should read $10 \cdot (\varrho - \varrho_0)/T$ and in the figure caption $(\varrho - \varrho_0)$ vs. T should read $(\varrho - \varrho_0)/T$ vs. T
- p. 186, Table 1 the host metal **Mo** is omitted to the left of the solute Co (third line). The solutes Co, Fe, Nb, Re, Ta, W and Zr should refer to the host metal **Mo**: **MoCo**, **MoFe**, **MoNb**, **MoRe**, **MoTa**, **MoW** and **MoZr**. The entries for **MoW**, 0.72 and 0.9 $\mu\Omega\text{cm}/\text{at\%}$ should read 0.072 and 0.09 $\mu\Omega\text{cm}/\text{at\%}$, respectively
- p. 227, Fig. 191. **InSn** the concentration of Sn, c , is given in at%
- p. 233, Fig. 224. **MoW** the unit of ϱ , $\mu\Omega\text{cm}$ should read $10^{-1} \cdot \mu\Omega\text{cm}$ the slope $0.72 \mu\Omega\text{cm}/\text{at\%}$ should read $0.072 \mu\Omega\text{cm}/\text{at\%}$
- p. 286 80 Me 1 should read 80 My 1
- p. 293, T_K^* the Kondo temperature T_K^* is defined by Eqs. (11), (17), and (21)
- p. 296, Table 1 **YTmSe** should read **YSeTm**
- p. 297, Table 1 **PdCrSi** should read **PdSiCr**
PdMnSi should read **PdSiMn**
PdFeSi should read **PdSiFe**
PdCoSi should read **PdSiCo**
- p. 298, Table 1 **CeAl₃** should read **Al₃Ce**
- p. 302, Table 3, **RhFe** 71 L 2 should read 71 L 3
74 R 4 should read 74 R 3 (note: Rusky should read Rusby)
For additional information on **RhFe** see Coles, B.R.: Phys. Lett. **8** (1964) 243 and Oliveira, jun., N.S., Foner, S.: Phys. Lett. **34A** (1971) 15
- p. 309, Table 4 $Y_x Tm_{(1-x)}$ should read $Y_x Tm_{(1-x)} Se$
- p. 315, Table 5 in the column "Remarks": Tm maximum temperature at 50 K should read ...at 5 K
- (La_(1-x)Ce_x)Al₂ **TiV** should read **(Ti, V)₂O₃**
- p. 334, Table 7 74 R 4 should read 74 R 3 (note: Rusky should read Rusby)
- p. 363, Fig. 105 Rusky should read Rusby
- p. 381, 74 R 1 74 R 3