

Co-Fe-Pr-Sm (Cobalt-Iron-Praseodymium-Samarium)

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An isothermal section at 800 °C at a constant Sm/Pr ratio of 1 was determined recently by [2001Wan] for this quaternary system for compositions up to 33.3 at.% of rare earth metal content.

Binary Systems

The Co-Fe phase diagram [1984Nis] is characterized by an extremely narrow solidification range. The face-centered-cubic (fcc) Fe forms a continuous solid solution γ with α Co over a wide range of temperature. The $\gamma \rightarrow (\alpha\text{Fe})$ body-centered-cubic (bcc) transformation temperature is initially raised by the addition of Co, reaching a maximum of 985 °C at 45 at.% Co. At 730 °C, the bcc phase of equiatomic composition orders to a CsCl type $B2$ structure. The updated Co-Pr phase diagram [2001Oka] depicts nine intermediate phases: $\text{Pr}_2\text{Co}_{17}$, PrCo_5 , $\text{Pr}_5\text{Co}_{19}$, Pr_2Co_7 , PrCo_3 , PrCo_2 , Pr_4Co_3 , Pr_5Co_2 , and Pr_3Co . See [2001Oka] for a summary of the crystal structure data. The Co-Sm phase diagram [2000Cam] depicts eight intermediate phases: $\text{Sm}_2\text{Co}_{17}$, SmCo_5 , $\text{Sm}_5\text{Co}_{19}$, Sm_2Co_7 , SmCo_3 , SmCo_2 , Sm_9Co_4 , and Sm_3Co . Among these, only $\text{Sm}_2\text{Co}_{17}$ and SmCo_5 show small homogeneity ranges at high temperatures. The crystal structure data on the Co-Sm phases are summarized by [1992Rag2]. The Fe-Pr phase diagram [1999Zha] depicts the $\text{Th}_2\text{Zn}_{17}$ type rhombohedral phase

$\text{Pr}_2\text{Fe}_{17}$. Both crystalline forms of PrFe_2 ($C14$ and $C15$) are metastable. The Fe-Sm phase diagram [1982Kub] depicts three line compounds, $\text{Sm}_2\text{Fe}_{17}$, SmFe_3 , and SmFe_2 . They all form peritectically, with the final eutectic solidification of Sm-rich alloys at 720 °C. The Pr-Sm phase diagram [Massalski2] shows no intermediate phases. βPr and γSm (both bcc) form a continuous solid solution.

Ternary Systems

No phase diagram information is available for the Co-Fe-Pr system. The lattice parameter variation of the solid solutions $\text{Pr}(\text{Co,Fe})_2$ and $\text{Pr}_2(\text{Co,Fe})_{17}$ was summarized by [1992Rag1]. The review of the Co-Fe-Sm system by [1992Rag2] gave a schematic liquidus surface, a reaction scheme and two isothermal sections at 1200 and 800 °C. No ternary compounds were found. There appear to be no reports on the phase equilibria of the Co-Pr-Sm and Fe-Pr-Sm systems.

The Quaternary Phase Equilibria

With starting metals of purity of 99.9% Co, 99.8% Fe, 99.9% Pr, and 99.9% Sm, [2001Wan] melted 45 alloy compositions at $\text{Sm/Pr} = 1$ and $(\text{Sm} + \text{Pr}) \leq 33.3$ at.% in an arc furnace under Ar atm. The alloy samples were

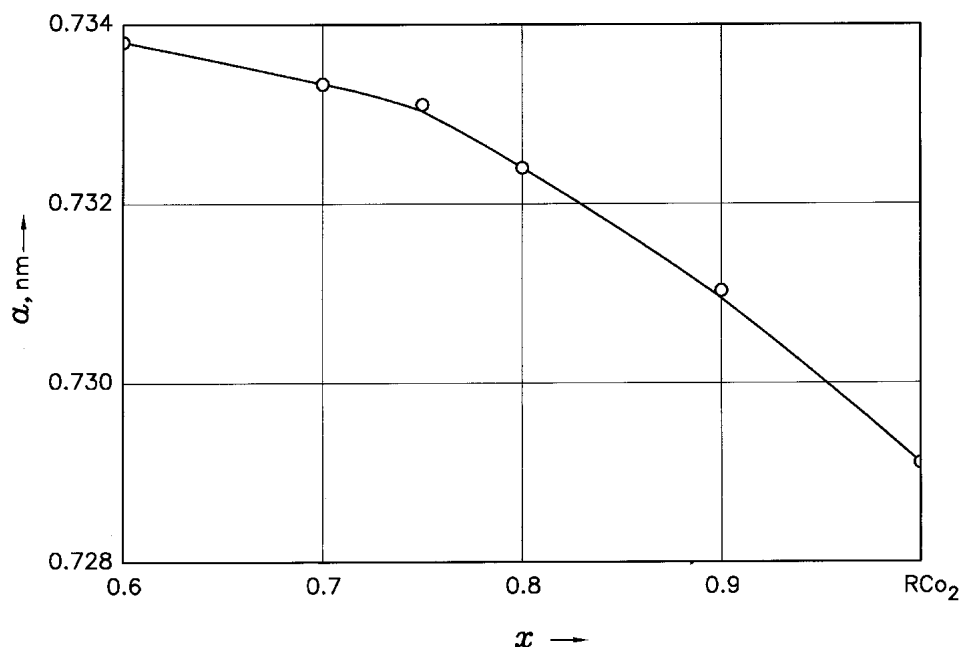


Fig. 1 Co-Fe-Pr-Sm lattice parameter variation of $\text{R}(\text{Fe}_{1-x}\text{Co}_x)_2$ alloys. $\text{R} = \text{Sm}_{0.5}\text{Pr}_{0.5}$ [2001Wan].

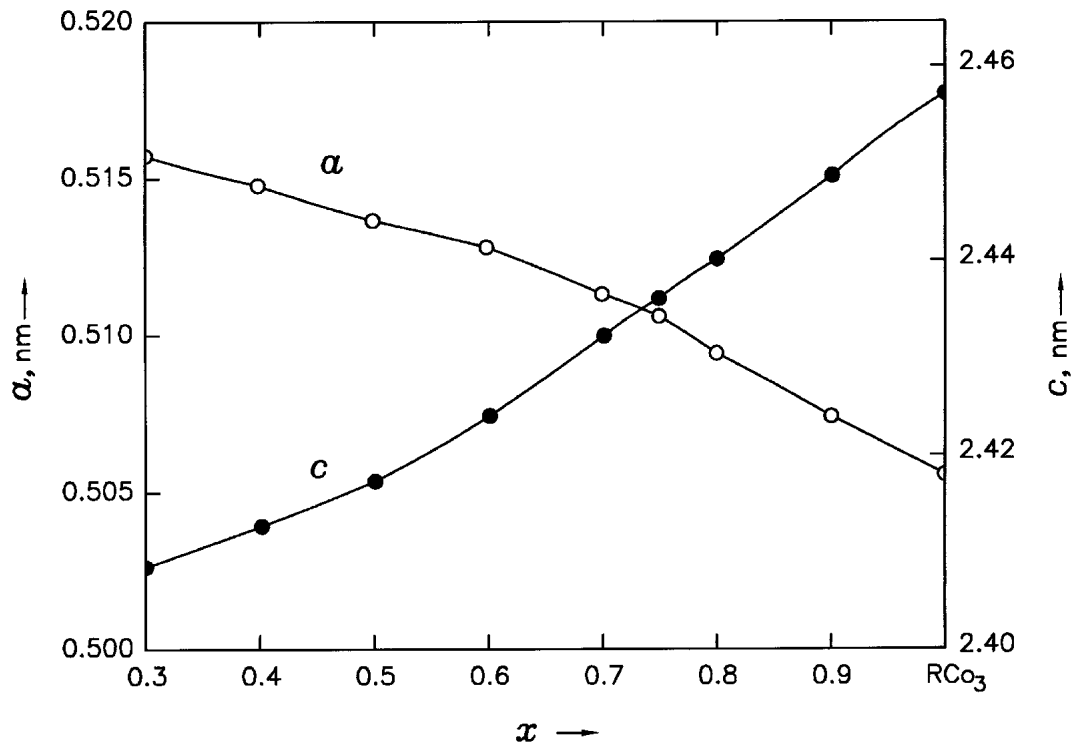


Fig. 2 Co-Fe-Pr-Sm lattice parameter variation of $R(\text{Fe}_{1-x}\text{Co}_x)_3$ alloys. $R = \text{Sm}_{0.5}\text{Pr}_{0.5}$ [2001Wan]

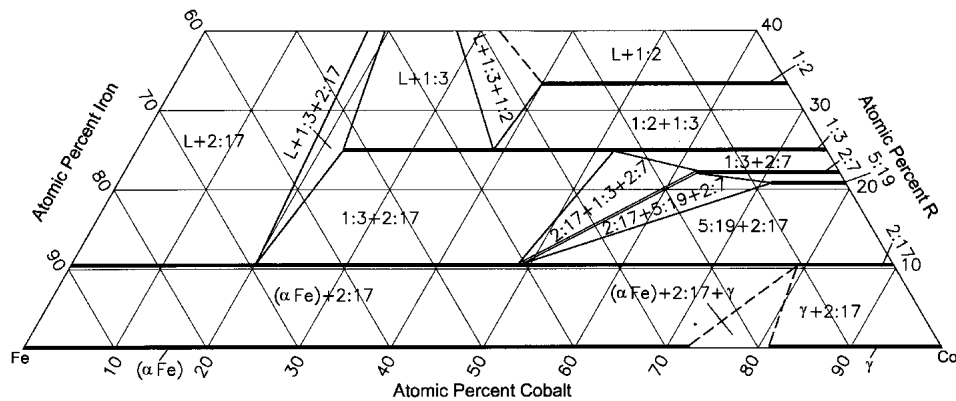


Fig. 3 Co-Fe-Pr-Sm isothermal section at 800 °C for $\text{Sm}/\text{Pr} = 1$ ($R = \text{Sm}_{0.5}\text{Pr}_{0.5}$) [2001Wan]

given a final anneal at 800 °C for 15–20 days and quenched in water. The phase equilibria were studied by optical microscopy, x-ray powder diffraction, and electron probe microanalysis.

In the composition range studied, five quaternary solid solutions based on binary compounds were identified by [2001Wan]: $(\text{Sm}, \text{Pr})(\text{Fe}, \text{Co})_2$, $(\text{Sm}, \text{Pr})(\text{Fe}, \text{Co})_3$, $(\text{Sm}, \text{Pr})_2(\text{Fe}, \text{Co})_7$, $(\text{Sm}, \text{Pr})_5(\text{Fe}, \text{Co})_{19}$, and $(\text{Sm}, \text{Pr})_2(\text{Fe}, \text{Co})_{17}$. Defining $(\text{Sm}_{0.5}\text{Pr}_{0.5}) = R$, the C15 type cubic phase $R(\text{Fe}_{1-x}\text{Co}_x)_2$ (denoted 1:2) is stable for $0.6 \leq x \leq 1.0$. The lattice parameter variation with x is shown in Fig. 1 [2001Wan]. The Be_3Nb type rhombohedral phase $R(\text{Fe}_{1-x}\text{Co}_x)_3$ (denoted 1:3) is stable for $0.3 \leq x \leq 1.0$. The

lattice parameter variation in this range is shown in Fig. 2 [2001Wan]. The Ce_2Ni_7 type hexagonal structure $R_2(\text{Fe}_{1-x}\text{Co}_x)_7$ (denoted 2:7) is stable for $0.8 \leq x \leq 1.0$. The $\text{Ce}_5\text{Co}_{19}$ type rhombohedral structure $R_5(\text{Fe}_{1-x}\text{Co}_x)_{19}$ (denoted 5:19) dissolves 8 at.% Fe at constant R content. The $\text{Th}_2\text{Zn}_{17}$ type rhombohedral structure $R_2(\text{Fe}_{1-x}\text{Co}_x)_{17}$ (denoted 2:17) forms a continuous solid solution, i.e., for all values of x from 0 to 1. There is no stable phase corresponding to the formula $R\text{Co}_5$ at 800 °C. No true ternary or quaternary compound was found by [2001Wan]. The isothermal section constructed by [2001Wan] at 800 °C for Sm/Pr ratio of 1 is redrawn in Fig. 3 to agree with the accepted binary data.

Section II: Phase Diagram Evaluations

References

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