## Fe-P-Y (Iron-Phosphorus-Yttrium)

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A partial isothermal section for this system at 800 °C was determined recently by [2002Ori]. The section depicts two ternary compounds.

## **Binary Systems**

A partial phase diagram is known for the Fe-P system [1982Kub]. The intermediate compound Fe<sub>3</sub>P forms through a peritectic reaction at 1166 °C between liquid and Fe<sub>2</sub>P. Fe<sub>2</sub>P forms congruently at 1370 °C. Fe<sub>3</sub>P is body centered tetragonal with the Ni<sub>3</sub>P-type structure. Fe<sub>2</sub>P has the hexagonal C22 structure. The other intermediate phases at higher P contents are FeP (orthorhombic MnP type), FeP<sub>2</sub> [orthorhombic FeS<sub>2</sub> (marcasite) type], and FeP<sub>4</sub> (monoclinic). The Fe-Y phase diagram [1992Zha] depicts four intermediate phases with limited or no homogeneity ranges:  $Y_2Fe_{17}$ ,  $Y_6Fe_{23}$ ,  $YFe_3$ , and  $YFe_2$ . See [1992Zha] for crystal structure data. The Y-P phase diagram is not known. The compound YP has the cubic NaCl type structure.

## **Ternary Isothermal Section**

With starting materials of 99.95 Fe, 99.98 P, and 99.5 Y (all at.%), Orishchin et al. [2002Ori] melted 17 ternary alloy compositions in an arc furnace under Ar atm. The alloys were finally annealed at 800 °C for 500 h and quenched in water. The phase equilibria were studied mainly by x-ray

powder diffraction, supplemented by metallography and electron probe microanalysis. The isothermal section at 800 °C determined by [2002Ori] up to 50 at.% P is redrawn in Fig. 1 to agree with the accepted binary data. Two ternary compounds are stable at this temperature.  $Y_2Fe_{12}P_7$  (denoted  $\tau_1$  here) has the hexagonal  $Zr_2Fe_{12}P_7$  type of structure. The crystal structure of  $YFe_5P_2$  (the experimentallydetermined composition is  $YFe_{4.8}P_{1.8}$ ), denoted  $\tau_2$  in Fig. 1, was not resolved. The ternary compound  $YFe_5P_3$  [1984Jei] was not found at this temperature. The compound YP forms tie lines with the two ternary compounds as well as with all the binary compounds (except  $Fe_3P$ ). The solubility of the third component in the binary phases is negligible.  $Y_2Fe_{17}$ , which has two crystal modifications, was found to have the hexagonal Th<sub>2</sub>Ni<sub>17</sub> type structure at 800 °C [2002Ori].

## References

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Fig. 1 Fe-P-Y isothermal section at 800 °C [2002Ori]. Narrow two-phase regions around tie-triangles are omitted