

Al-Er-Fe (Aluminum-Erbium-Iron)

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A partial isothermal section for this ternary system at 500 °C was determined by [2001Yan] and was reviewed by [2004Rag]. Recently, [2010Jem] determined an isothermal section for this ternary system at 800 °C covering the entire composition range, which depicts six ternary compounds.

Binary Systems

The Al-Er phase diagram [Massalski2] depicts the following intermediate compounds: ErAl₃ (*L*₁₂, AuCu₃-type cubic), ErAl₂ (*C*₁₅, MgCu₂-type cubic), ErAl (ErAl-type orthorhombic), Er₃Al₂ (Zr₃Al₂-type tetragonal), and Er₂Al (*C*₂₃, Co₂Si-type orthorhombic). In the Fe-Al phase diagram, the γ solid solution based on face-centered cubic (fcc) Fe is restricted by a loop. The solid solution based on the body-centered cubic (bcc) Fe exists in both the disordered *A*₂ and ordered *B*₂ and *D*₀₃ forms. Apart from the high temperature phase ϵ (*D*₈₂, Cu₅Zn₈-type cubic), there are three intermediate phases in the system with restricted ranges of homogeneity: FeAl₂ (triclinic), Fe₂Al₅ (orthorhombic), and FeAl₃ or Fe₄Al₁₃ (monoclinic). The Er-Fe phase diagram [Massalski2] shows four line compounds: Fe₁₇Er₂ (Ni₁₇Th₂-type hexagonal), Fe₂₃Er₆ (*D*₈_a, Mn₂₃Th₆-type

cubic), Fe₃Er (Be₃Nb-type rhombohedral), and Fe₂Er (*C*₁₅, MgCu₂-type cubic).

Ternary Compounds

Five ternary compounds, all with a range of homogeneity, were identified in this system at 800 °C by [2010Jem]. The structural details as determined by [2010Jem] are listed in Table 1. ErFe_{12-*x*}Al_{*x*} ($5 \leq x \leq 8$) (denoted here as τ_1) has the ThMn₁₂-type tetragonal structure. Er₂Fe_{17-*x*}Al_{*x*} ($4.75 \leq x \leq 5.7$) (τ_2) has the Cu₇Tb-type hexagonal structure. Er₂Fe_{17-*x*}Al_{*x*} ($5.7 \leq x \leq 9.5$) (τ_3) has the Th₂Zn₁₇-type rhombohedral structure. ErFe_{3-*x*}Al_{*x*} ($0.5 \leq x \leq 1$) (τ_4) has the DyFe₂Al-type hexagonal structure. ErFe_{1+*x*}Al_{1-*x*} ($-0.2 \leq x \leq 0.25$) (τ_5) is of MgZn₂-type hexagonal. In addition, a sixth compound of fixed composition ErFe₂Al₁₀ (τ_6) with orthorhombic symmetry was earlier reported by [1998Thi] (see Table 1).

Isothermal Section

With starting metals of 99.99% Al, 99.8% Er, and 99.9% Fe, [2010Jem] arc-melted under Ar atm 90 alloy samples.

Table 1 Al-Er-Fe crystal structure and lattice parameter data [2010Jem]

Phase	Composition, atomic %	Pearson symbol	Space group	Prototype	Lattice parameter, nm
ErFe _{12-<i>x</i>} Al _{<i>x</i>} (τ_1) ($5 \leq x \leq 8$)	38.5-61.5 Al 53.8-30.8 Fe 7.7 Er	<i>tI</i> 26	<i>I</i> 4/ <i>mmm</i>	ThMn ₁₂	<i>a</i> = 0.8594-0.8704 <i>c</i> = 0.4981-0.5037
Er ₂ Fe _{17-<i>x</i>} Al _{<i>x</i>} (τ_2) ($4.75 \leq x \leq 5.7$)	25.0-30.0 Al 64.5-59.5 Fe 10.5 Er	<i>hP</i> 8	<i>P</i> 6/ <i>mmm</i>	Cu ₇ Tb	<i>a</i> = 0.4956 (a) <i>c</i> = 0.4229
Er ₂ Fe _{17-<i>x</i>} Al _{<i>x</i>} (τ_3) ($5.7 \leq x \leq 9.5$)	30.0-50.0 Al 59.5-39.5 Fe 10.5 Er	<i>hR</i> 19	<i>R</i> $\bar{3}$ <i>m</i>	ThZn ₁₇	<i>a</i> = 0.8695 (b) <i>c</i> = 1.2692
ErFe _{3-<i>x</i>} Al _{<i>x</i>} (τ_4) ($0.5 \leq x \leq 1$)	12.5-25 Al 62.5-50 Fe 25 Er	<i>hP</i> 24	<i>P</i> 6 ₃ / <i>mmc</i>	DyFe ₂ Al	<i>a</i> = 0.5156 (c) <i>c</i> = 1.6522
ErFe _{1+<i>x</i>} Al _{1-<i>x</i>} (τ_5) ($-0.2 \leq x \leq 0.25$)	40-25 Al 26.7-41.7 Fe 33.3 Er	<i>hP</i> 12	<i>P</i> 6 ₃ / <i>mmc</i>	MgZn ₂	<i>a</i> = 0.5332 (d) <i>c</i> = 0.8683
ErFe ₂ Al ₁₀ (τ_6)	76.9 Al 15.4 Fe 7.7 Er	<i>oC</i> 52	<i>Cmcm</i>	YbFe ₂ Al ₁₀	<i>a</i> = 0.8948 <i>b</i> = 1.0136 <i>c</i> = 0.8988

(a) at $x = 5.2$, (b) at $x = 9$, (c) at $x = 1$, and (d) at $x = 0$

