

# Al-Fe-La (Aluminum-Iron-Lanthanum)

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The previous review of this system by [1992Rag] presented an isothermal section at 500 °C, which depicts six ternary compounds. Recently, [1995Tan] determined an isothermal section at ~25 °C with five ternary compounds. The known crystal structures of three of these compounds are in agreement in the two studies. However, the homogeneity ranges of the compounds differ significantly.

## Binary Systems

The Al-Fe phase diagram reviewed by [1993Kat] shows that the fcc solid solution based on Fe is restricted by a  $\gamma$  loop. The bcc solid solution  $\alpha$  exists in both ordered and disordered forms. Apart from the high-temperature phase  $\varepsilon$ , there are three intermediate phases in the system with restricted ranges of homogeneity:  $\text{FeAl}_2$ ,  $\text{Fe}_2\text{Al}_5$ , and  $\text{FeAl}_3$ . The Al-La phase diagram reviewed by [1988Gsc] shows six intermediate phases:  $\text{Al}_{11}\text{La}_3$ ,  $\text{Al}_3\text{La}$ ,  $\text{Al}_x\text{La}$ ,  $\text{Al}_2\text{La}$ ,  $\text{AlLa}$ , and  $\text{AlLa}_3$ . Recently, [1996Sac] studied the La-rich region and reported that  $\text{AlLa}_3$  decomposes at 400 °C to  $\text{AlLa}$  and ( $\beta\text{La}$ ). The Fe-La phase diagram reviewed by [1982Kub] shows no intermediate phases. For crystal structure data on the above binary compounds, see [Pearson3].

## Ternary Compounds

A phase of unknown structure  $\tau_1$  is present at 500 °C on the 33.3 at.% La line between 20 and 33.3 at.% Al [1992Rag].

[1995Tan] found a phase of unknown structure at 36 at.% La and 20 at.% Al at ~25 °C. This phase is denoted  $\tau_1$  here.  $\text{La}_2\text{Fe}_7\text{Al}_{10}$  ( $\tau_2$ ) has the  $\text{Th}_2\text{Zn}_{17}$ -type rhombohedral structure [1995Tan]. The same phase at 500 °C [1992Rag] extends from 52.6 to 57.9 at.% Al at constant La content.  $\text{LaFe}_{1.2}\text{Al}_{7.8}$  ( $\tau_3$ ) is of unknown structure [1992Rag, 1995Tan]. The  $\text{ThMn}_{12}$ -type tetragonal structure  $\text{LaFe}_4\text{Al}_8$  ( $\tau_4$ ) (the 1:12 phase) occurs at the same composition both at 500 and 25 °C. The  $\text{NaZn}_{13}$ -type cubic structure  $\text{La}(\text{Fe},\text{Al})_{13}$  ( $\tau_5$ ) has a much wider range of Al content at 25 °C than at 500 °C.  $\text{LaFe}_2\text{Al}_{10}$  found at 500 °C [1992Rag] is not present at 25 °C.  $\text{La}(\text{Al},\text{Fe})_2$  is an  $\text{MgCu}_2$ -type Laves phase extending from  $\text{Al}_2\text{La}$  into the ternary region at 25 °C (not at 500 °C). The structural details of these compounds determined by [1995Tan] are listed in Table 1.

## Ternary Isothermal Section

With starting metals of purity 99.9%, [1995Tan] melted 105 alloy compositions in an arc furnace under Ar atm. The alloys were homogenized between 900 and 500 °C and furnace cooled to room temperature. It is assumed here that the phase equilibria correspond to room temperature (~25 °C). The phase identification was carried out by x-ray powder diffraction analysis. [1995Tan]'s isothermal section is redrawn in Fig. 1 to agree with the accepted binary data. All

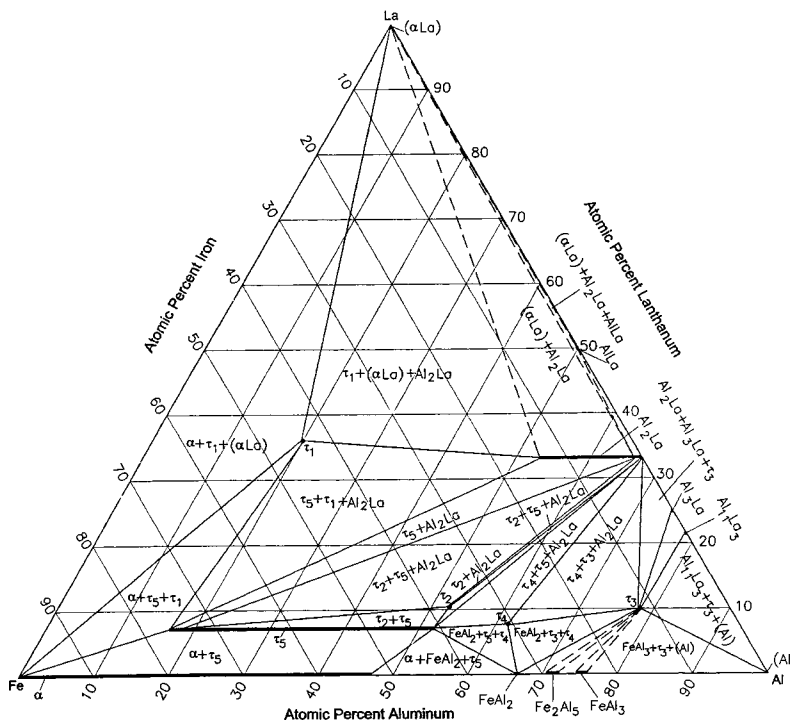


Fig. 1 Al-Fe-La isothermal section at ~25 °C [1995Tan]. The thin two-phase fields around tie-triangles are omitted

**Table 1 Al-Fe-La crystal structure and lattice parameter data**

Phase	Composition, at.%		Pearson symbol	Space group	Strukturbericht designation	Prototype	Lattice parameter, nm
	La	Al					
La <sub>9</sub> Fe <sub>11</sub> Al <sub>5</sub> ( $\tau_1$ )	36.0	20.0					
La <sub>2</sub> Fe <sub>7</sub> Al <sub>10</sub> ( $\tau_2$ )	10.5	52.6	<i>hR19</i>	$R\bar{3}m$		Th <sub>2</sub> Zn <sub>17</sub>	$a = 0.8962$ $c = 1.298$
LaFe <sub>1.2</sub> Al <sub>7.8</sub> ( $\tau_3$ )	10.0	78.0					
LaFe <sub>4</sub> Al <sub>8</sub> ( $\tau_4$ ) (1:12)	7.7	61.5	<i>tI26</i>	<i>I4/mmm</i>	<i>D2<sub>b</sub></i>	ThMn <sub>12</sub>	$a = 0.8900$ $c = 0.5075$
La(Fe,Al) <sub>13</sub> ( $\tau_5$ )	7.1	16.7–52	<i>cF112</i>	$Fm\bar{3}c$	<i>D2<sub>3</sub></i>	NaZn <sub>13</sub>	$a = 1.1668\text{--}1.1983$
La(Al,Fe) <sub>2</sub>	33.3	66.7–53.3	<i>cF24</i>	$Fd\bar{3}m$	<i>C15</i>	MgCu <sub>2</sub>	$a = 0.8147\text{--}0.8111$

the ternary phases listed in Table 1 are present. Due to lack of data in the ternary region, the ordered and disordered forms of bcc  $\alpha$  are not shown separately in Fig. 1. With the exception of Al<sub>2</sub>La, none of the Al-La or Fe-Al compounds show any solubility for the third component. AlLa<sub>3</sub> (shown by [1995Tan]) is omitted in Fig. 1, in line with the findings of [1996Sac].

#### References

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