

## Al-Fe-Ni (Aluminum-Iron-Nickel)

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This technologically-important system has been investigated experimentally a number of times. Many reviews and updates have also appeared in the last two decades [1988Ray, 1992Bud, 1994Rag, 1995Vil, 2005Cac, 2005Rag, 2006Ele, 2006Rag, 2008Rag]. For a recent detailed summary, the reader is referred to the updated version of [1992Bud] by [2005Cac], the review of [2006Ele] and the update by [2008Rag] on the work of [2007Chu1]. The present update will be limited to a review of three recent publications, featuring isothermal sections at 1200, 1100, 1000 and 900 °C for the Al-lean alloys by [2008Chu], two isothermal sections at 850 and 627 °C for Al-rich alloys by [2008Zha], and a vertical section along Fe<sub>3</sub>Al-Ni<sub>3</sub>Al join by [2007Chu2, 2008Chu].

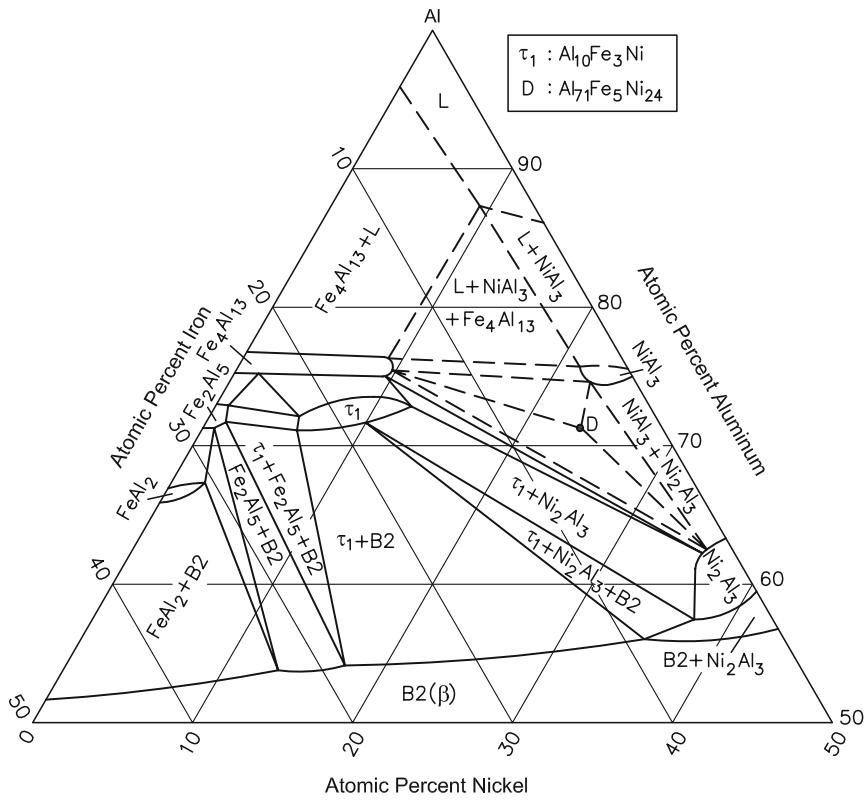
$B2$  and  $D0_3$  forms. Apart from the high temperature phase  $\epsilon$ , there are three other intermediate phases in this system:  $\text{FeAl}_2$  (triclinic),  $\text{Fe}_2\text{Al}_5$  (70-73 at.% Al, orthorhombic), and  $\text{FeAl}_3$  (or  $\text{Fe}_4\text{Al}_{13}$ ) (74.5-76.6 at.% Al, monoclinic). The Al-Ni phase diagram [1993Oka] shows five intermediate phases:  $\text{NiAl}_3$  ( $D0_{11}$ ,  $\text{Fe}_3\text{C}$ -type orthorhombic),  $\text{Ni}_2\text{Al}_3$  ( $D5_{13}$ -type hexagonal),  $\text{NiAl}$  ( $B2$ ,  $\text{CsCl}$ -type cubic, denoted  $\beta$ ),  $\text{Ni}_5\text{Al}_3$  ( $\text{Ga}_3\text{Pt}_5$ -type orthorhombic), and  $\text{Ni}_3\text{Al}$  ( $L1_2$ ,  $\text{AuCu}_3$ -type cubic, denoted  $\gamma'$ ). The Fe-Ni phase diagram [1993Swa] is characterized by a very narrow solidification range with a peritectic reaction at 1514 °C between bcc  $\delta$  and liquid that yields the Fe-based fcc solid solution. A continuous fcc solid solution denoted  $\gamma$  is stable over a wide range of temperature. At 517 °C, an ordered phase  $\text{FeNi}_3$  forms congruently from  $\gamma$ .

# Binary Systems

The Al-Fe phase diagram [1993Kat] shows that the face-centered cubic (fcc) solid solution based on Fe is restricted by a  $\gamma$  loop. The body-centered cubic (bcc) solid solution exists in the disordered  $A2$  form ( $\alpha$ ), as well as the ordered

## Ternary Phase Equilibria

With starting metals of 99.999% Al, 99.99% Fe and 99.95% Ni, [2008Zha] arc-melted 19 ternary alloys under Ar atm. The final anneal was at 850 °C for 18 d or at



**Fig. 1** Al-Fe-Ni isothermal section at 850 °C for Al-rich alloys [2008Zha]

## Section II: Phase Diagram Evaluations

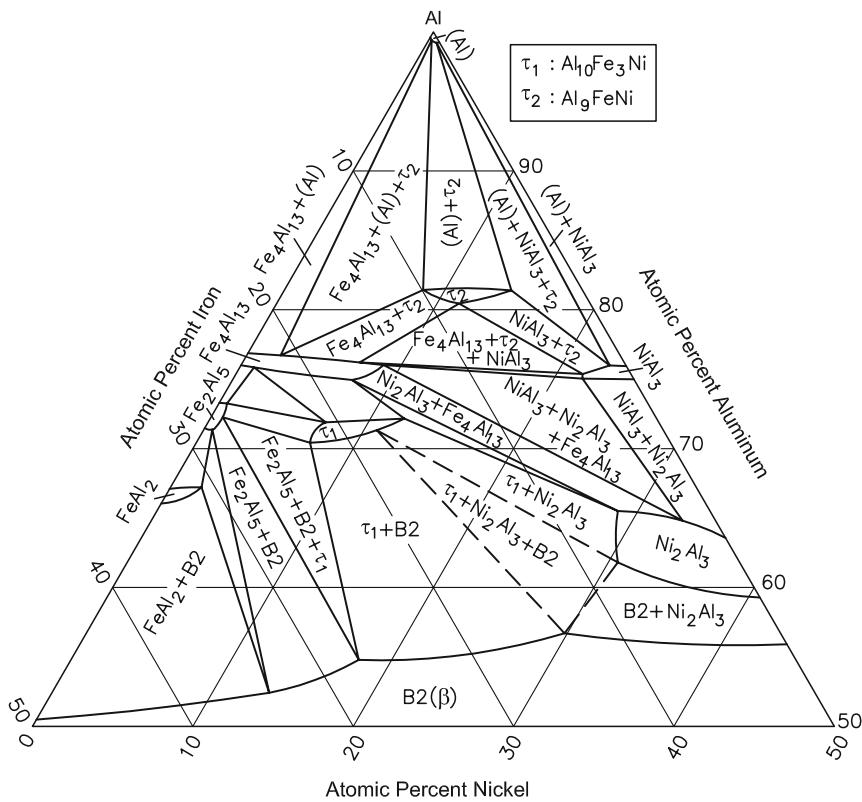


Fig. 2 Al-Fe-Ni isothermal section at 627 °C for Al-rich alloys [2008Zha]

627 °C for 50 d, followed by water quenching. The phase equilibria were studied by optical microscopy and x-ray powder diffraction. The compositions of coexisting phases were measured with energy dispersive x-ray analysis on a scanning electron microscope or by electron probe microanalysis (EPMA) and listed. The isothermal sections constructed by [2008Zha] at 850 and 627 °C are shown in Fig. 1-2.

At 850 °C (Fig. 1), the ternary phase  $\tau_1$  (denoted as  $\tau_2$  by [2008Zha]) and the quasicrystalline decagonal phase D (denoted  $\tau_3$  by [2008Zha]) are present. The homogeneity range of  $\tau_1$  is 5.7-13.1 at.% Ni and 14.4-22.2 at.% Fe. The binary phases  $\text{FeAl}_2$ ,  $\text{Fe}_2\text{Al}_5$ , and  $\text{Fe}_4\text{Al}_{13}$  dissolve 2.1, 1.3 and 9.6 at.% Ni respectively.  $\text{Al}_2\text{Ni}_3$  dissolves up to 4.9 at.% Fe. At 850 °C, the D phase is shown to be in four-phase equilibrium with  $\text{Fe}_4\text{Al}_{13}$ ,  $\text{NiAl}_3$ , and  $\text{Ni}_2\text{Al}_3$ . It decomposes at this temperature through a ternary eutectoid reaction. This section at 850 °C by [2008Zha] is remarkably similar to that of [2007Chu1] reviewed by [2008Rag], the two versions appearing as independent investigations.

At 627 °C (Fig. 2), the ternary phases  $\tau_1$  (denoted  $\tau_2$  by [2008Zha]) and  $\tau_2$  (denoted  $\tau_1$  by [2008Zha]) are present. The homogeneity range of  $\tau_1$  is 7.2-12.4 at.% Ni and 15.6-22.5 at.% Fe. The range of  $\tau_2$  is 8.7-14.3 at.% Ni and 4.5-10.0 at.% Fe. The solubility of Ni in  $\text{FeAl}_2$ ,  $\text{Fe}_2\text{Al}_5$  and  $\text{Fe}_4\text{Al}_{13}$  is 2.0, 1.5, and 8.9 at.% respectively.  $\text{Ni}_2\text{Al}_3$  dissolves up to 3.6 at.% Fe.

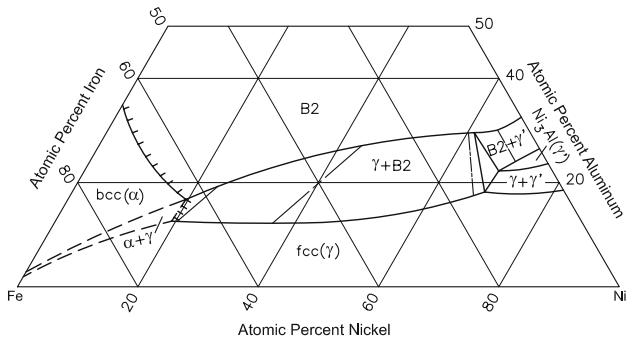
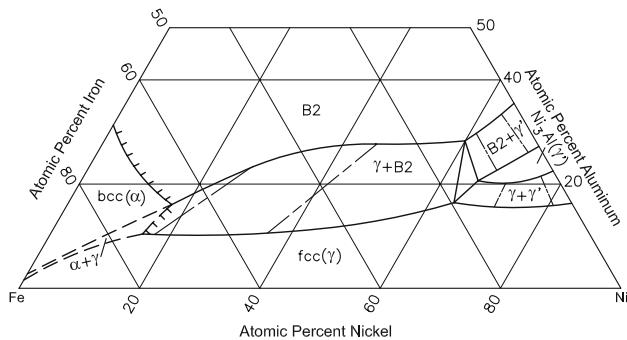
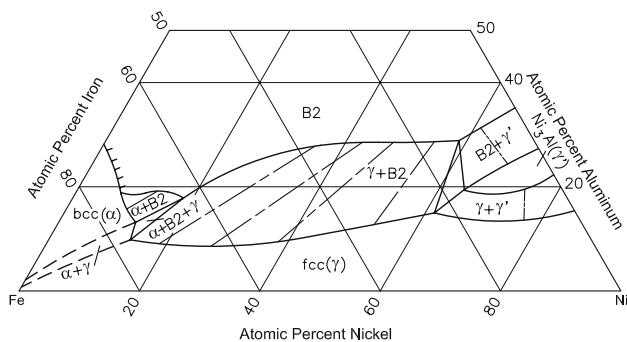


Fig. 3 Al-Fe-Ni isothermal section at 1200 °C for Al-lean alloys [2008Chu]

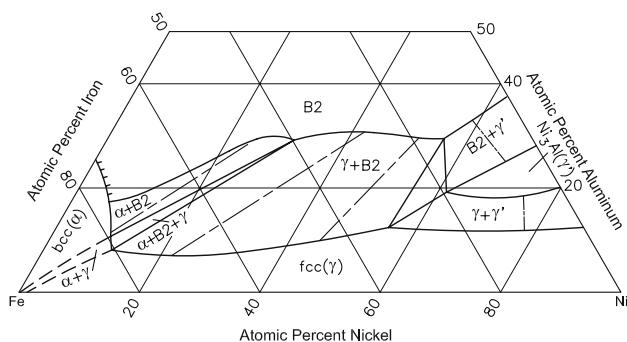
The phase relationships between the ordered and disordered forms of bcc Fe in the ternary region were clarified in the recent study of [2008Chu]. With starting metals of 99.999% Al, 99.99% Fe and 99.99% Ni, [2008Chu] arc-melted under Ar atm 28 ternary alloys containing up to 28 at.% Al. The samples were annealed at 1200, 1100, 1000 and 900 °C for ½, 4, 14 and 28 d respectively and quenched in water. The phase equilibria were studied with x-ray powder diffraction and electron probe microanalysis. The partial isothermal sections constructed by [2008Chu] at 1200, 1100, 1000, and 900 °C are shown in Fig. 3-6. At 1200 °C (Fig. 3) and 1100 °C



**Fig. 4** Al-Fe-Ni isothermal section at 1100 °C for Al-lean alloys [2008Chu]



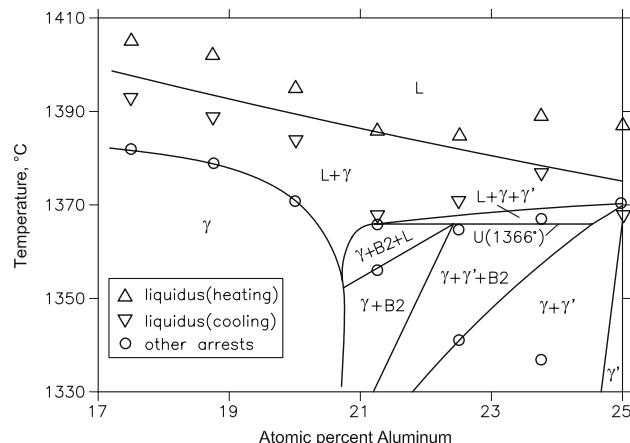
**Fig. 5** Al-Fe-Ni isothermal section at 1000 °C for Al-lean alloys [2008Chu]



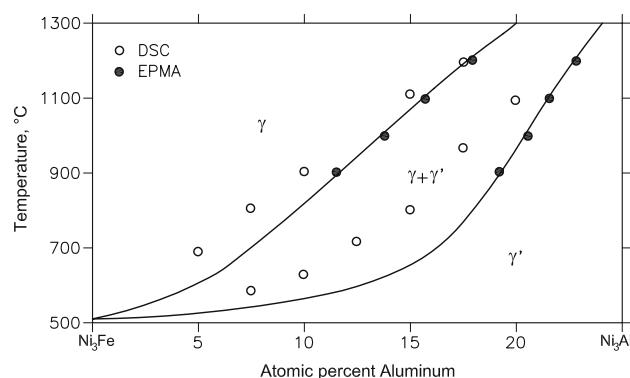
**Fig. 6** Al-Fe-Ni isothermal section at 900 °C for Al-lean alloys [2008Chu]

(Fig. 4), the  $bcc \leftrightarrow B2$  transition is fully second order. At 1000 and 900 °C (Fig. 5 and 6), the transition is partially second-order and partially first-order. The ( $bcc + B2$ ) two-phase region gives rise to a three-phase ( $fcc + bcc + B2$ ) equilibrium.

[2006Rag] reviewed the vertical section along the Ni<sub>3</sub>Fe-Ni<sub>3</sub>Al join from the studies of [2005Him], omitting the region near the Ni<sub>3</sub>Al end, where the binary Ni<sub>3</sub>Al forms through a peritectic reaction. The details of this region were studied recently by [2007Chu2]. Their results are shown in Fig. 7. An invariant U-type transition reaction:  $L + \gamma' \leftrightarrow \gamma + B2$  occurs at 1366 °C, with the composition (in at.%) of



**Fig. 7** Al-Fe-Ni invariant horizontal near the Ni<sub>3</sub>Al-end of the vertical section along the Ni<sub>3</sub>Fe-Ni<sub>3</sub>Al join [2007Chu2]



**Fig. 8** Al-Fe-Ni pseudobinary region of the vertical section along the Ni<sub>3</sub>Fe-Ni<sub>3</sub>Al join [2008Chu]

the participating phases being L (5Fe73Ni22Al),  $\gamma'$  (0.5Fe75Ni24.5Al),  $\gamma$  (4Fe76Ni20Al), and B2 (0.5Fe71.5-Ni28Al). Towards the Ni<sub>3</sub>Fe-side, the vertical section is pseudobinary in nature. [2008Chu] found that the location of the phase boundaries in this region is influenced by the experimental technique used, as shown in Fig. 8. The experimental points from differential scanning calorimetry (DSC) used by [2005Him] are consistently higher than those found with EPMA used by [2008Chu]. The EPMA analysis of equilibrated samples is to be considered more accurate, as compared to those obtained from the dynamic method of DSC.

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