

Al-Cu-Fe (Aluminum-Copper-Iron)

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The review of this ternary system by [1992Rag] presented liquidus projections for Al- and Cu-rich alloys from [1953Phi] and [1941Yut], respectively, and an isothermal section at 600 °C from [1971Pre]. In an update, [2005Rag] reviewed the more-recent studies of [1998Wan] and [2003Zha1, 2003Zha2] and presented a liquidus projection and several isothermal sections between 1200 and 560 °C. The system was also reviewed by [2005Wan]. Most recently, [2009Che] investigated the region above 40 at.% Al at 600 °C and carried out a detailed thermodynamic analysis. This work will be reviewed here briefly.

In the Fe-Al phase diagram [Massalski2], the solid solution γ based on face-centered cubic (fcc) Fe is restricted by a loop. The solid solution based on the body-centered cubic (bcc) Fe (α) exists in both the disordered $B2$ and ordered $D0_3$ forms. The high-temperature phase ϵ is stable between 1232 and 1102 °C and has the $D8_2$, Cu_5Zn_8 -type cubic structure, with $a = 0.89757$ nm [2010Ste]. There are three other intermediate phases in the system with restricted ranges of homogeneity: FeAl_2 (triclinic), Fe_2Al_5 (orthorhombic), and FeAl_3 or $\text{Fe}_4\text{Al}_{13}$ (monoclinic). There are no intermediate phases in the Cu-Fe system. A metastable liquid miscibility gap is known.

Binary Systems

The Al-Cu phase diagram [2004Ria] depicts the following intermediate phases: CuAl_2 ($C16$ -type tetragonal, denoted θ), $\text{CuAl}(\text{HT})$ (η_1 , orthorhombic), $\text{CuAl}(\text{LT})$ (η_2 , monoclinic), $\text{Cu}_5\text{Al}_4(\text{HT})$ (ζ_1 , orthorhombic, space group $Fmm2$), $\text{Cu}_5\text{Al}_4(\text{LT})$ (ζ_2 , orthorhombic, space group $Imm2$), $\epsilon_1(\text{HT})$ (bcc), $\epsilon_2(\text{LT})$ ($B8_1$, NiAs-type hexagonal), Cu_3Al_2 (rhombohedral), $\text{Cu}_9\text{Al}_4(\text{HT})$ (γ_0 , $D8_2$, Cu_5Zn_8 -type cubic), $\text{Cu}_9\text{Al}_4(\text{LT})$ (γ_1 , $D8_3$, Cu_9Al_4 -type cubic), and $\text{Cu}_3\text{Al}(\beta$, bcc).

Ternary Phases

Three established ternary phases are known in this system [1992Rag, 2009Che]. FeCu_2Al_7 (denoted ω by [2009Che] and τ_2 by [1992Rag]) is tetragonal, space group $P4/mnc$, with lattice parameters $a = 0.63417 - 0.63499$ nm and $c = 1.48498 - 1.48377$ nm [2009Che]. FeCu_2Al_6 (denoted τ_1 by [2009Che] and as I here) is the stable icosahedral quasicrystalline phase [1987Tsa]. [2004Tur] studied

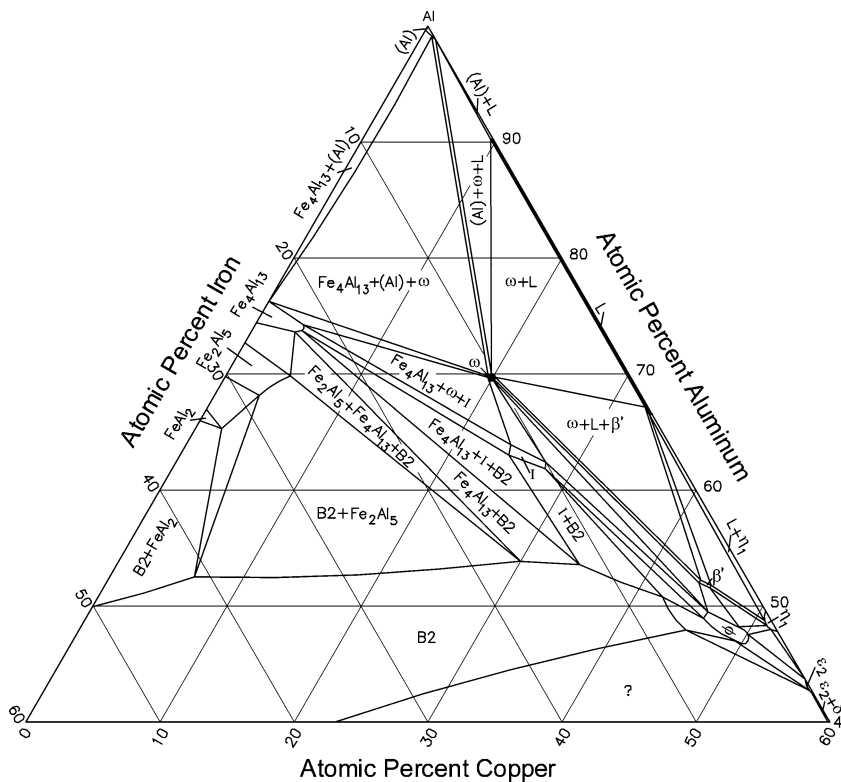


Fig. 1 Al-Cu-Fe experimental isothermal section at 600 °C [2009Che]

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the effect of pressure up to ~ 1.6 GPa on the formation of I. The formation temperature was between 750 and 800 °C and changed little with pressure [2004Tur]. $\text{FeCu}_{10}\text{Al}_{10}$ (denoted φ by [2009Che] and as τ_3 by [1992Rag]) forms in the solid state and is present at 600 °C. The presence of a fourth phase $\text{Fe}_4\text{CuAl}_{23}$ (labeled α in earlier reports and as τ_1 by [1992Rag]) was discounted by [2009Che] and was not included in their thermodynamic analysis.

Ternary Phase Equilibria

With starting metals of 99.99% Al, 99.99% Cu, and 99.99% Fe, [2009Che] arc-melted under Ar atm 14 Al-rich ternary alloys. The alloys were annealed at 600 °C for 35 days and quenched in water. The phase equilibria were studied with x-ray powder diffraction, scanning electron microscopy, and energy dispersive x-ray spectroscopy. Differential thermal analysis was carried out at a heating rate of 5 °C per min. The identified phases and their compositions were listed. The isothermal section at 600 °C for Al-rich alloys constructed by [2009Che] is shown in Fig. 1. The three ternary phases ω , I, and φ are present. In addition, [2009Che] found a phase β' , which appears to arise from a miscibility gap in the B2 solid solution.

In their thermodynamic modeling, [2009Che] treated liquid, fcc and bcc (ε_1 , Cu_3Al , and $\alpha\delta\text{Fe}$) as substitutional solid solutions. The magnetic contribution to the Gibbs energy of fcc and bcc phases was taken into account. To describe the ordered bcc (B2) phase, an ordering energy term was added to the description of the disordered bcc phase. A two sublattice model was used to describe the

Fe-Al binary compounds to allow for the solubility of Cu, which mainly substitutes for Al. The Al-Cu phase γ_1 was described similarly to account for the Fe solubility. In cases where no ternary solubility has been reported, the binary phases θ , η , ζ , δ , and ε_2 were described as pure compounds. The ternary compounds FeCu_2Al_6 (I) and $\text{FeCu}_{10}\text{Al}_{10}$ (φ)

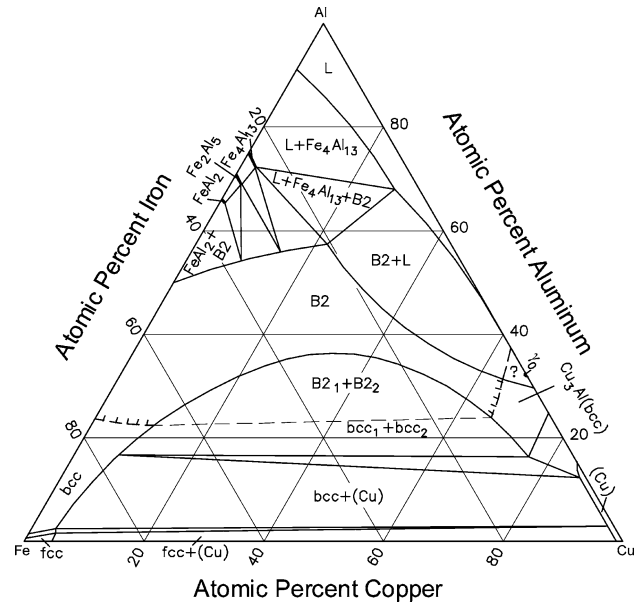


Fig. 3 Al-Cu-Fe computed isothermal section at 1000 °C [2009Che]

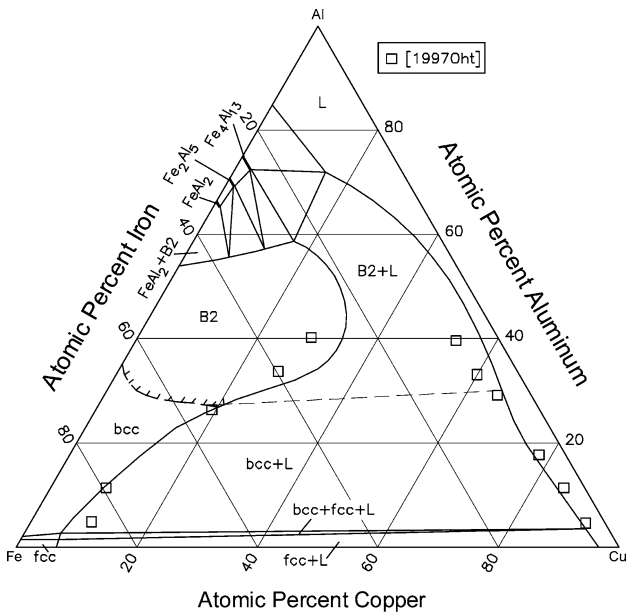


Fig. 2 Al-Cu-Fe computed isothermal section at 1100 °C [2009Che]

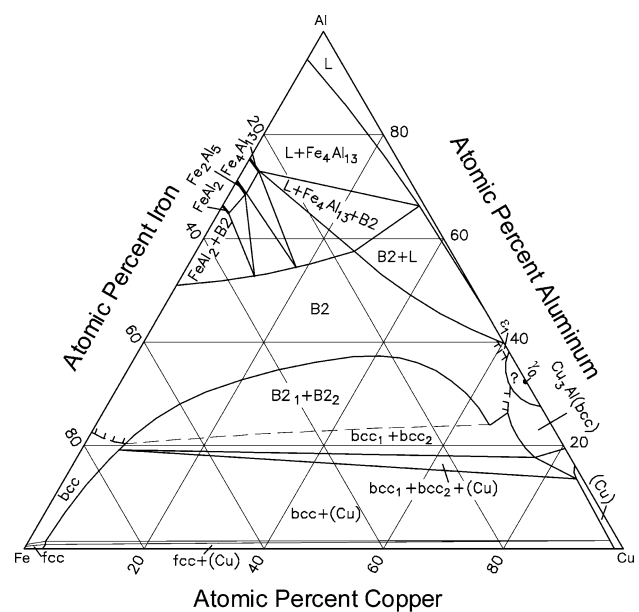


Fig. 4 Al-Cu-Fe computed isothermal section at 900 °C [2009Che]

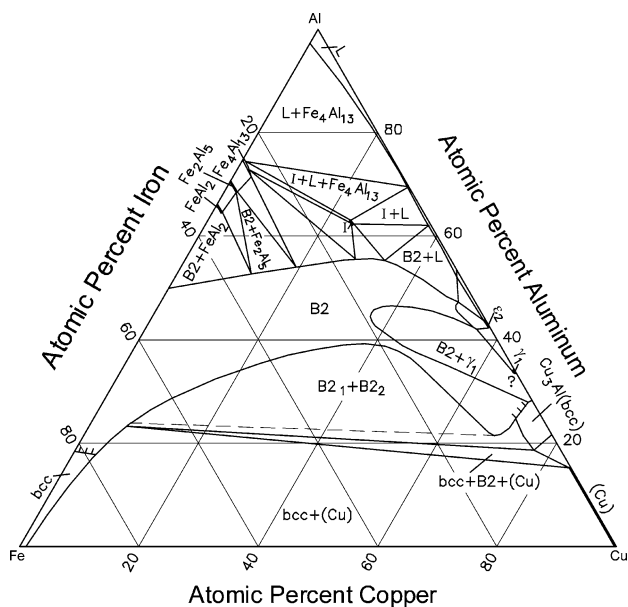


Fig. 5 Al-Cu-Fe computed isothermal section at 800 °C [2009Che]

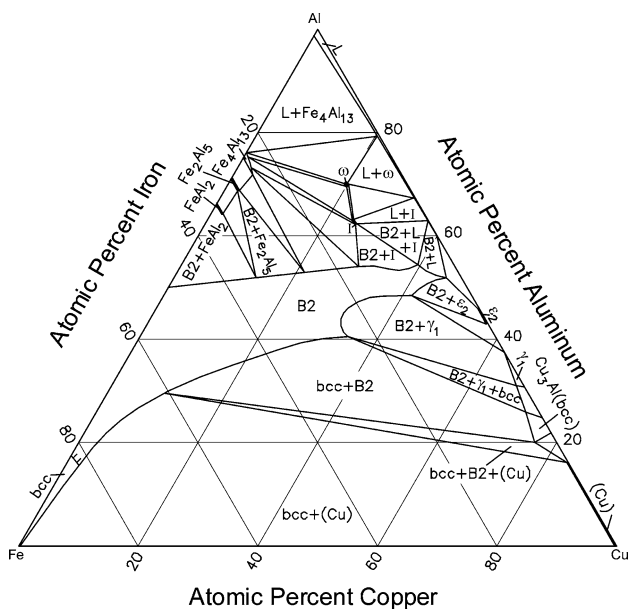


Fig. 6 Al-Cu-Fe computed isothermal section at 700 °C [2009Che]

were assumed to have a homogeneity range at constant Fe content, whereas FeCu_2Al_7 (ω) was treated as a stoichiometric compound. The binary descriptions were taken from the literature. The experimental data of [1941Yut], [1997Oht], [1998Wan], [2003Zha1], [2003Zha2], and [2009Che] were used as inputs. The optimized interaction parameters were listed.

[2009Che] computed eight isothermal sections at 100 °C intervals between 1300 and 600 °C. The computed sections were found to be in general agreement with the experimental data. Here, the computed isothermal section at 1100 °C is compared with the experimental data of [1997Oht] in Fig. 2. The computed isothermal sections at 1000, 900, 800, and 700 °C are shown in Fig. 3-6. Due to the low solubility of Fe in some of the Al-Cu intermediate phases and the proximity of the ordering $\text{bcc} \leftrightarrow \text{B2}$ transitions, the phase relationships near the Al-Cu side are not clearly defined in Fig. 3-6. The quasicrystalline phase I, which forms peritectically at 882 °C [2003Zha2], is present at 800 °C (Fig. 5). The ω phase, which also forms peritectically at 740 °C [2003Zha2] (762 °C [2009Che]), is present additionally at 700 °C (Fig. 6). The computed isothermal section at 600 °C (not shown here) agrees with the experimental results shown in Fig. 1. The invariant equilibria computed by [2009Che] were found to have some differences with the liquidus projection and reaction scheme presented by [2003Zha1] and reviewed by [2005Rag].

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