

# Al-Cu-Zn (Aluminum-Copper-Zinc)

V. Raghavan

The previous update on this ternary system by [2007Rag] reviewed mainly the thermodynamic assessment of [1998Lia]. Very recently, [2009Ren] determined a partial isothermal section at 360 °C for Cu-lean alloys and found that Cu stabilizes the miscibility gap of the Al-Zn fcc phase.

## Binary Systems

The Al-Cu phase diagram [2004Ria] depicts the following intermediate phases: CuAl<sub>2</sub> (C16-type tetragonal, denoted  $\theta$ ), CuAl(HT) ( $\eta_1$ , orthorhombic) CuAl(LT) ( $\eta_2$ , monoclinic), Cu<sub>5</sub>Al<sub>4</sub>(HT) ( $\zeta_1$ , orthorhombic, space group *Fmm*2), Cu<sub>5</sub>Al<sub>4</sub>(LT) ( $\zeta_2$ , orthorhombic, space group *Imm*2),  $\varepsilon_1$ (HT) (cubic?),  $\varepsilon_2$ (LT) (*B8*<sub>1</sub>, NiAs-type hexagonal), Cu<sub>3</sub>Al<sub>2</sub> (rhombohedral), Cu<sub>9</sub>Al<sub>4</sub>(HT) ( $\gamma_0$ , *D8*<sub>2</sub>, Cu<sub>5</sub>Zn<sub>8</sub>-type cubic), Cu<sub>9</sub>Al<sub>4</sub>(LT) ( $\gamma_1$ , *D8*<sub>3</sub>, Cu<sub>9</sub>Al<sub>4</sub>-type cubic), and Cu<sub>3</sub>Al ( $\beta$ , bcc). The Al-Zn phase diagram [1993Che] contains no intermediate phases. A miscibility gap occurs in the Al-based face centered cubic (fcc) solid solution below 351 °C, where the fcc phase splits into fcc<sub>1</sub> and fcc<sub>2</sub>. The monotectoid reaction fcc<sub>2</sub> ↔ fcc<sub>1</sub> + (Zn) follows at 277 °C. The Cu-Zn phase diagram [1993Kow, Massalski2] is characterized by a series of peritectic reactions, which yield CuZn ( $\beta$ , bcc), Cu<sub>5</sub>Zn<sub>8</sub> ( $\gamma$ , *D8*<sub>2</sub>-type cubic), CuZn<sub>3</sub> ( $\delta$ , *B2*, CsCl-type cubic), and CuZn<sub>4</sub> ( $\varepsilon$ , cph). Zn (cph) has a *c/a* axial ratio much larger than  $\varepsilon$  and the two coexisting cph phases are modeled separately with different interaction parameters [1993Kow]. The  $\beta$  phase orders to a CsCl-type *B2* phase ( $\beta'$ ) through a second-order transition below ~460 °C.

## Ternary Phases

A ternary phase with rhombohedral symmetry and with the nominal composition Al<sub>4</sub>Cu<sub>3</sub>Zn (denoted  $\tau$ ) is known in this system [Pearson3]. The homogeneity range of  $\tau$  and its temperature dependence are not clearly defined. The structurally-related, low-temperature form  $\tau'$  was found to be stable by [2005Hao] between 400 °C and room temperature.

## Ternary Phase Equilibria

With starting metals of 99.999% purity, [2009Ren] prepared diffusion couples of Cu-Zn master alloys with pure Al. The diffusion couples were annealed at 360 °C for 48 h and quenched in water. The coexisting phase compositions were measured by electron probe microanalysis and listed. The isothermal section at 360 °C constructed by

[2009Ren] is shown in Fig. 1. In the Al-Zn system, the miscibility gap in the fcc phase closes at 351 °C. The addition of Cu stabilizes this gap, which is stable at 360 °C, even with as low as 0.1 at.% addition of Cu. The gap increases in width with increasing Cu content. In an earlier study, [2004Ren] determined tie-lines between fcc<sub>1</sub> and fcc<sub>2</sub> at 340 and 320 °C for small additions of Cu and these are shown in Fig. 2. The Cu addition shifts the fcc<sub>1</sub>/(fcc<sub>1</sub> + fcc<sub>2</sub>) and (fcc<sub>1</sub> + fcc<sub>2</sub>)/fcc<sub>2</sub> phase boundaries towards the Al-rich corner. Also, Cu segregates in the fcc<sub>2</sub> phase in preference to fcc<sub>1</sub>.

The computed results of [2009Dai] indicate that the stabilizing effect of Cu on the Al-Zn miscibility gap increases with increasing Cu content, culminating in a metastable miscibility gap in the Al-Cu binary system. These calculations, however, ignore the intervening effect of the binary and ternary compounds.

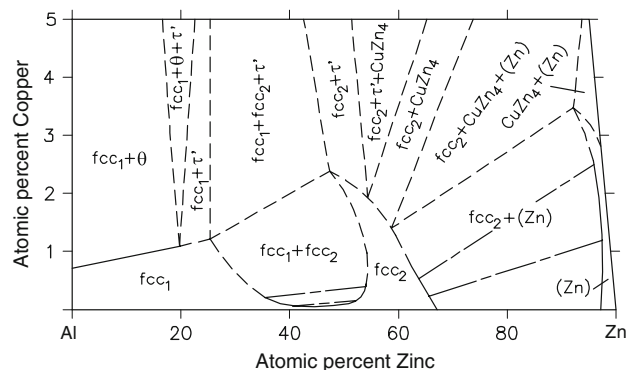


Fig. 1 Al-Cu-Zn partial isothermal section at 360 °C [2009Ren]

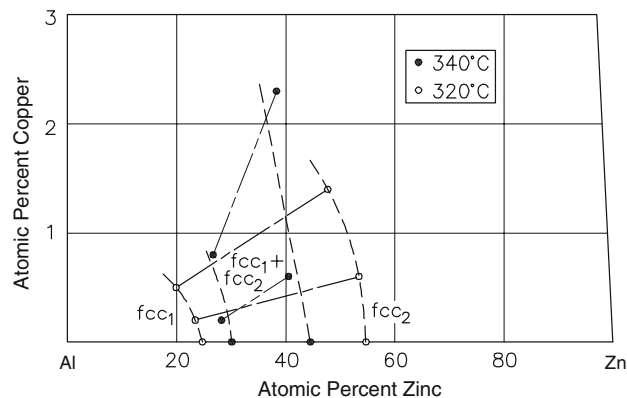


Fig. 2 Al-Cu-Zn tie-lines between fcc<sub>1</sub> and fcc<sub>2</sub> at 340 and 320 °C [2004Ren]

## Section II: Phase Diagram Evaluations

### References

- 1993Che:** S.L. Chen and Y.A. Chang, A Thermodynamic Analysis of the Al-Zn System and Phase Diagram Calculation, *CALPHAD*, 1993, **17**(2), p 113-124
- 1993Kow:** M. Kowalski and P.J. Spencer, Thermodynamic Reevaluation of the Cu-Zn System, *J. Phase Equilibria*, 1993, **14**(4), p 432-438
- 1998Lia:** H. Liang and Y.A. Chang, A Thermodynamic Description for the Al-Cu-Zn System, *J. Phase Equilibria*, 1998, **19**(1), p 25-37
- 2004Ren:** Y.P. Ren, H. Chen, K. Wang, S.M. Hao, and D.J. Dong, Effect of Cu Addition on the ( $\alpha_1 + \alpha_2$ ) Miscibility Gap of the Al-Zn System, *Scripta Mater.*, 2004, **51**, p 267-270
- 2004Ria:** P. Riani, L. Arrighi, R. Marazza, D. Mazzone, G. Zanicchi, and R. Ferro, Ternary Rare-Earth Aluminum Systems with Copper: A Review and a Contribution to Their Assessment, *J. Phase Equilib. Diffus.*, 2004, **25**(1), p 22-52
- 2005Hao:** S.M. Hao, H.X. Li, H. Chen, and Y.P. Ren, Phase Transformation of Al-Zn(Cu) Alloys and the Phase Diagram of Al-Zn-Cu System, *Cailiao Yu Yejin Xuebao (J. Mater. Metall.)*, 2005, **4**(2), p 92-101
- 2007Rag:** V. Raghavan, Al-Cu-Zn (Aluminum-Copper-Zinc), *J. Phase Equilib. Diffus.*, 2007, **28**(2), p 183-188
- 2009Dai:** L.L. Dai, H.X. Li, and Y.P. Ren, Thermodynamic Calculation on the Miscibility Gap of fcc Al-Based Solid Solution in the Al-Zn-Cu System, *J. Alloys Compd.*, 2009, **478**, p 144-146
- 2009Ren:** Y.P. Ren, G.W. Qin, W.L. Pei, and S.M. Hao, The ( $\alpha_1 + \alpha_2$ ) Miscibility Gap of the Al-Zn-Cu System at 360 °C, *Scripta Mater.*, 2009, **61**, p 36-39