

# Cu-Fe-Zn (Copper-Iron-Zinc)

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In their early work, [1934Bau] investigated Cu-rich alloys containing a few mass% of Fe and presented a number of isothermal and vertical sections. Later, [1992Rag] reviewed and presented full isothermal sections at 1000 and  $\sim 700$  °C from the studies of [1980Bud]. In a recent update, [2008Rag] briefly reviewed an isothermal section at 460 °C near the Zn corner from the work of [2008Ave]. A thermodynamic description applicable to Cu-rich alloys was developed by [2008Mie].

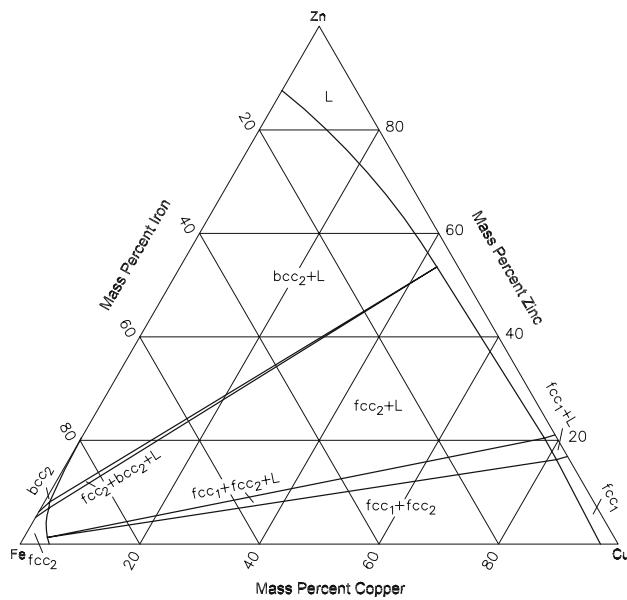
## Binary Systems

There are no intermediate phases in the Cu-Fe system. A metastable liquid miscibility gap has been experimentally measured in this system. The Cu-Zn phase diagram [Massalski2] is characterized by a series of peritectic reactions, which yield CuZn ( $\beta$ , bcc), CuZn ( $\beta'$ , CsCl-type cubic), Cu<sub>5</sub>Zn<sub>8</sub> ( $\gamma$ ,  $D_{8_2}$ -type cubic), CuZn<sub>3</sub> ( $\delta$ ,  $B_2$ , CsCl-type cubic), and CuZn<sub>4</sub> ( $\epsilon$ , cph). The intermediate phases in the Fe-Zn system [Massalski2] are:  $\Gamma$  (Fe<sub>3</sub>Zn<sub>10</sub>; Cu<sub>5</sub>Zn<sub>8</sub>-type cubic),  $\Gamma_1$  (Fe<sub>11</sub>Zn<sub>40</sub>; cubic, space group  $F\bar{4}3m$ , 408 atoms/cell),  $\delta$  (FeZn<sub>10</sub>; FeZn<sub>10</sub>-type hexagonal), and  $\zeta$  (CoZn<sub>13</sub>-type monoclinic).

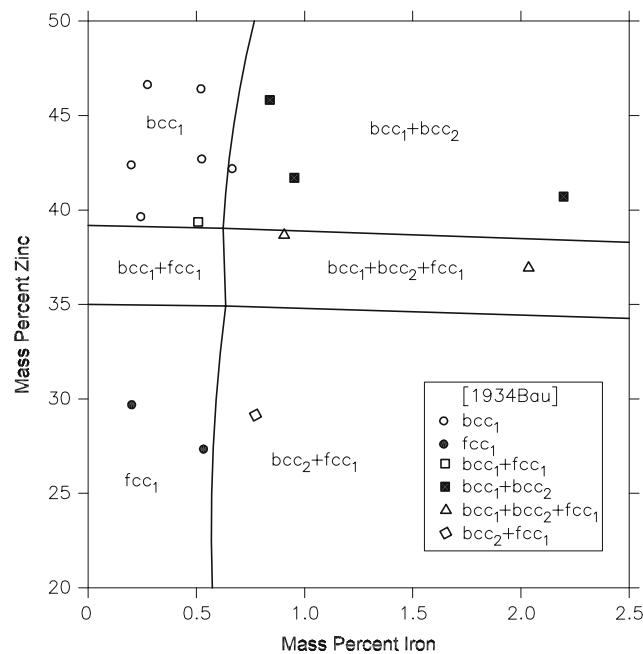
## Computed Ternary Phase Equilibria

In the thermodynamic description, [2008Mie] modeled the liquid, face centered cubic (fcc), body centered cubic (bcc), close packed hexagonal (cph), and the cubic  $\gamma$  phase of the Cu-Zn binary system as substitutional solutions. Ternary interaction parameters were introduced for the liquid, fcc, and bcc phases. It is doubtful whether a single lattice description is applicable to the  $\gamma$  phase, which has the Cu<sub>5</sub>Zn<sub>8</sub>-type of structure with a homogeneity range. The Fe-Zn binary compounds were treated as stoichiometric compounds, with nil solubility of the third component. Experimental data on phase equilibria from [1934Bau] and [1952Haw] and data on the Zn activity in liquid solutions from [1987Ryc] were used in the optimization.

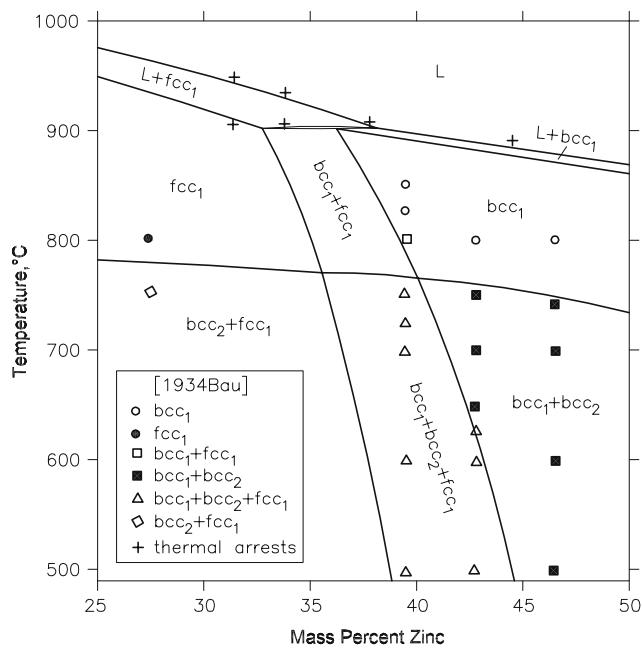
A liquidus projection, two full isothermal sections at 1200 and 1000 °C, six isothermal sections for Cu-rich alloys at 900, 800, 700, 672, 600, and 500 °C were computed. Also, four vertical sections at 0.5 and 0.9 mass% Fe and at 64 and 75 mass% Cu, respectively, were computed. The computed sections were compared mainly with the experimental data of [1934Bau]. As examples, the full isothermal section at 1000 °C, the Cu-rich isothermal section at 800 °C and the vertical section at 0.5 mass% Fe are shown in Fig. 1-3. The agreement with the data of [1934Bau] (Fig. 2 and 3) is satisfactory.



**Fig. 1** Cu-Fe-Zn computed isothermal section at 1000 °C [2008Mie]



**Fig. 2** Cu-Fe-Zn computed isothermal section at 800 °C for Cu-rich alloys [2008Mie]



**Fig. 3** Cu-Fe-Zn computed vertical section at 0.5 mass% Fe [2008Mie]

## References

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