

# Al-Cr-Ni-W (Aluminum-Chromium-Nickel-Tungsten)

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Recently, [2003Bur] determined isothermal sections of this quaternary system at 71 and 76 at.% Ni and at 1000 and 900 °C. The previous work on this system by [1961Ali] presented three isothermal sections at 1100 °C and at constant weight ratios of W:Al = 3:1, 1:1, and 1:3 respectively. Also, [1989Cha] gave partial isothermal sections at 1250 and 1000 °C at a constant Ni content of 75 at.%.

## Binary Systems

For the Al-Cr, Al-W, Cr-Ni, Cr-W, and Ni-W, binary systems, [Massalski2] can be used as a starting source. A partial list of updates on these systems is [2000Mah] (Al-Cr) and [1997Oka] (Cr-Ni). For a brief update of the Al-Ni system, see the Al-Nb-Ni update in this issue.

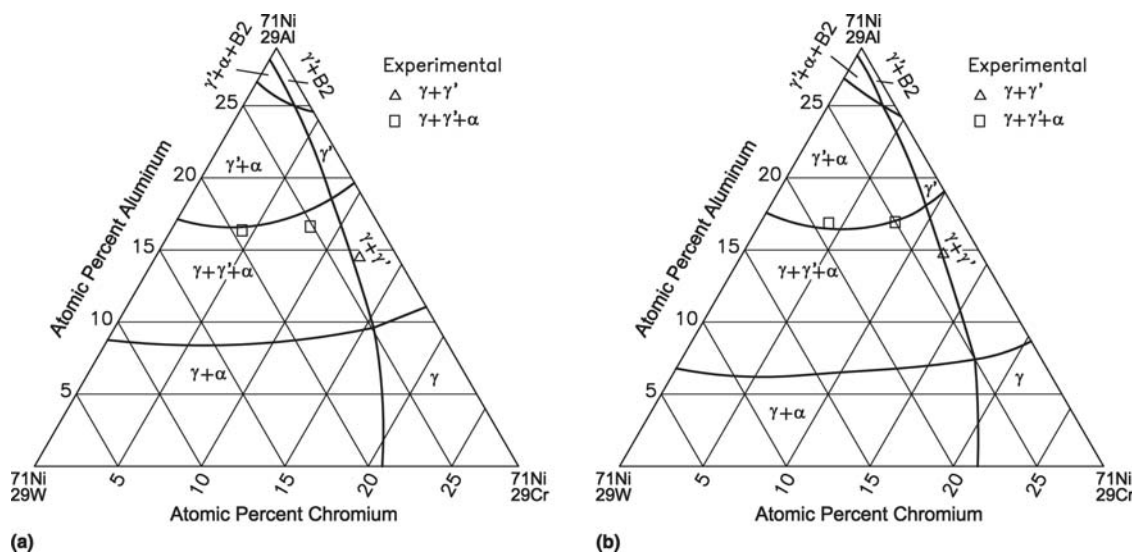


Fig. 1 Al-Cr-Ni-W computed isothermal sections at 71 at.% Ni (a) 1000 °C and (b) 900 °C [2003Bur]

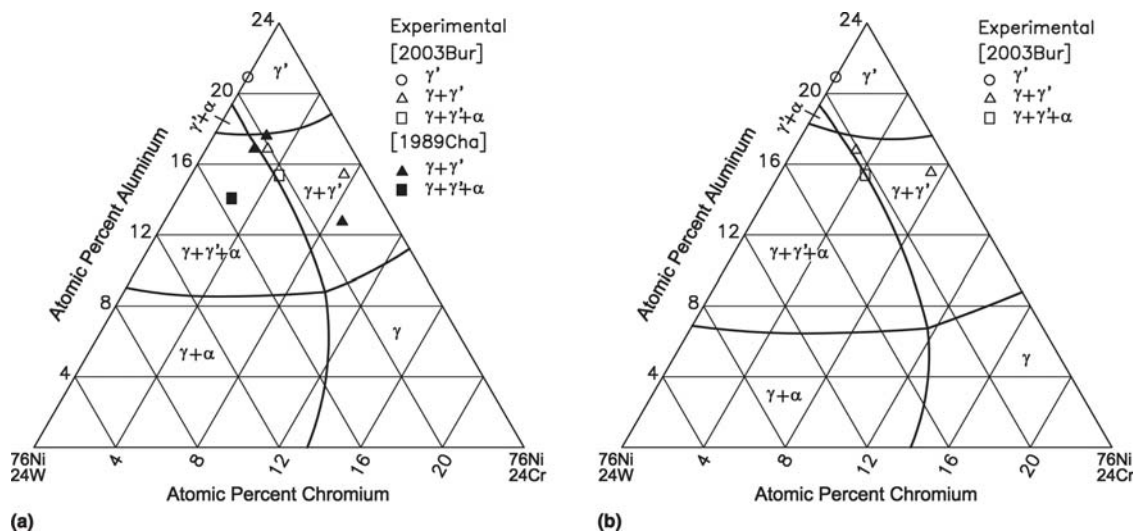


Fig. 2 Al-Cr-Ni-W computed isothermal sections at 76 at.% Ni (a) 1000 °C and (b) 900 °C [2003Bur]

## Ternary Systems

Compilations of data on the Al-Cr-Ni, Al-Ni-W, and Cr-Ni-W systems can be found in [1995Vil1], [1995Vil2], and [1995Vil3], respectively. No phase diagram data appear to be available on the Al-Cr-W system. In the Ni-rich alloys reviewed here, the Al-Cr-W system is not relevant.

## Quaternary Phase Equilibria

With starting metals of purity of 99.9 to 99.99%, [2003Bur] induction-melted seven Ni-rich alloys with compositions in the range of 14.5 to 21.0 wt.% Al, 0 to 12.5 wt.% Cr, and 0.8 to 7.6 wt.% W. The samples were annealed at 1000 °C for 176 h or at 900 °C for 672 h, followed by water quenching. The phase equilibria were studied by metallography and selected area electron diffraction in a transmission electron microscope. Compositions of the co-existing phases were measured by energy-dispersive x-ray spectroscopy and listed. The phase equilibria were computed using the CALPHAD approach. To obtain a better fit with the experimental data, the interaction parameters for the  $\gamma$  (fcc) and  $\gamma'$  ( $L1_2$ ) phases of the Al-Ni-W system were reoptimized and used in conjunction with the literature values of the other parameters in the calculations. The computed isothermal sections at 1000 and 900 °C at 71 and 76 at.% Ni are compared in Fig. 1 and 2 with the limited

experimental data. In addition to  $\gamma$  and  $\gamma'$  phases, the W-based bcc solid solution denoted  $\alpha$  is present.

## References

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