# AI-Cr-Ni (Aluminum-Chromium-Nickel)

V. Raghavan

The earlier reviews of this ternary system were those of [1984Mer], [1991Rog], and [1995Vil]. [2004Vel] updated the data reviewed by [1991Rog]. An update by [2006Rag] presented a computed liquidus projection, compared several isothermal sections of [1999Hua] and [2001Dup] and gave vertical sections along the NiAl-Cr and Ni<sub>3</sub>Al-'Ni<sub>3</sub>Cr' joins. An isothermal section at 1150 °C for Al-lean alloys determined by [2006Kit] was reviewed briefly by [2008Rag]. The ternary phases in the Al-rich region were recently investigated by [2008Gru1], who presented isothermal sections at 1000 and 900 °C. At 1000 °C, two ternary phases labeled  $\zeta$  and  $\varepsilon$  were found; at 900 °C, only  $\zeta$ was found. Very recently, for the Al-rich region, [2008Wei] determined a liquidus projection and an isothermal section at 700 °C. In addition to ζ, two other ternary phases of unresolved structure were identified at 700 °C.

## **Binary Systems**

The Al-rich end of the Al-Cr phase diagram investigated recently by [2008Gru2] is shown in Fig. 1. The intermediate phases in the system are:  $CrAl_7$  ( $V_7Al_{45}$ -type monoclinic, denoted  $\theta$ ),  $Cr_2Al_{11}$  ( $CrAl_5$ -type monoclinic, denoted  $\eta$ ),



Fig. 1 Al-Cr Phase diagram in the Al-rich region [2008Gru2]

 $CrAl_4$  (hexagonal,  $P6_3/mmc$ , denoted  $\mu$ ),  $CrAl_3$  ( $\nu$ ),  $Cr_2Al$ (MoSi<sub>2</sub>-type tetragonal) and an unconfirmed low-temperature phase X at 75 at.% Cr. Between 30 and 41 at.% Cr, five phases were reported earlier:  $\alpha Cr_4 A_9$ ,  $\beta Cr_4 Al_9$ ,  $\gamma Cr_4 Al_9$ ,  $\alpha Cr_5Al_8$  and  $\beta Cr_5Al_8$ , with no well-established phase boundaries between them. The work of [2008Gru2] shows only two phases in this region. The high-temperature phase denoted  $\gamma_1$  is cubic (Cu<sub>5</sub>Zn<sub>8</sub>-type) and transforms on cooling via a second-order transition to  $\gamma_2$ , the transition temperature decreasing from 1140 to 1060 °C with increasing Al content. The  $\gamma_2$  phase is rhombohedral (Cr<sub>5</sub>Al<sub>8</sub>-type). The Al-Ni phase diagram [1993Oka] shows five intermediate phases: NiAl<sub>3</sub> (Fe<sub>3</sub>C-type orthorhombic), Ni<sub>2</sub>Al<sub>3</sub>  $(D5_{13}$ -type hexagonal, denoted  $\delta$ ), NiAl (B2, CsCl-type cubic, denoted  $\beta$ ), Ni<sub>5</sub>Al<sub>3</sub> (Ga<sub>3</sub>Pt<sub>5</sub>-type orthorhombic), and Ni<sub>3</sub>Al ( $L1_2$ , AuCu<sub>3</sub>-type cubic, denoted  $\gamma'$ ). The Cr-Ni phase diagram is of the simple eutectic type, with Ni dissolving up to  $\sim 50$  at.% Cr and Cr dissolving up to ~32 at.% Ni.

## **Ternary Phase Equilibria**

About 20 alloys were induction-melted under Ar atm by [2008Gru1]. Alloys containing >60 at.% Al were annealed at 1000 °C for 90-120 h or at 900 °C for 110-142 h. Alloys with <60 at.% Al were annealed at 1150 °C for 49 h, at 1025 °C for 400 h or at 1000 °C for 382-686 h. All alloys were quenched in water after annealing. The phase equilibria were studied by x-ray powder diffraction, scanning and transmission electron microscopy and differential thermal analysis (DTA) at heating and cooling rates of 10-50 °C per min. The local phase composition was determined by the energy dispersive x-ray analysis.

The isothermal sections constructed by [2008Gru1] for Al-rich alloys at 1150 and 1025 °C (Fig. 2 and 3) are in broad agreement with earlier results reviewed by [2006Rag]. The isothermal sections of [2008Gru1] at 1000 and 900 °C are redrawn in Fig. 4 and 5. At 1000 °C (Fig. 4), two ternary phases were found. A hexagonal phase labeled  $\zeta$  by [2008Gru1] (called  $\kappa$  or  $\rho_2$  in earlier literature) was found to have a composition range between Al<sub>79.0</sub> Ni<sub>3.0</sub>Cr<sub>18.0</sub> and Al<sub>71.5</sub>Ni<sub>9.0</sub>Cr<sub>19.5</sub> [2008Gru1], with lattice parameters of a = 1.7674 nm, and c = 1.2516 nm. The DTA experiments indicated its formation at  $\sim 1030$  °C. The second ternary phase at 1000 °C labeled  $\varepsilon$  (see Fig. 4) has orthorhombic symmetry with  $a \approx 1.26$  nm,  $b \approx 3.48$  nm, and  $c \approx 2.02$  nm and occurs around the composition Al<sub>76 5</sub>Ni<sub>2 0</sub>Cr<sub>21 5</sub>. Interpretation in the earlier literature of  $\varepsilon$ as the binary phase  $CrAl_4$  ( $\mu$ ) was discounted by [2008Gru1]. The third component solubility in the binary phases  $\mu$ ,  $\gamma_2$ , and  $\delta$  phases was found to be about 1 at.% Ni, 3 at.% Ni, and 2.5 at.% Cr respectively. At 900 °C (Fig. 5),

### Section II: Phase Diagram Evaluations



Fig. 2 Al-Cr-Ni isothermal section at 1150 °C [2008Gru1]



Fig. 3 Al-Cr-Ni isothermal section at 1025 °C [2008Gru1]

the ternary phase  $\epsilon$  was not found. The homogeneity range of  $\zeta$  at 900 °C becomes wider and extends between  $Al_{81}Ni_3Cr_{16},~Al_{76.5}Ni_3Cr_{20.5},~Al_{76.5}Ni_9Cr_{14.5},~and~Al_{71.5}~Ni_9Cr_{19.5}$ 

With starting metals of 99.99% purity, [2008Wei] arcmelted more than 30 Al-rich ternary alloys under Ar atm. The alloys were annealed between 950 and 600 °C for 2 weeks and quenched in water. The phase equilibria were studied with x-ray powder diffraction, metallography, energy dispersive x-ray analysis on a scanning electron microscope, and DTA at a heating and cooling rate of 5 °C/min. The homogeneity range of  $\zeta$  (designated  $\tau_1$  by [2008Wei]) at 700 °C was found to match with that found by [2008Gru1] at 900 °C. The second ternary phase labeled  $\tau_2$  had an x-ray pattern characteristic of decagonal phases. The third ternary phase  $\tau_3$  had a complex pattern. The



Fig. 4 Al-Cr-Ni isothermal section at 1000 °C [2008Gru1]



Fig. 5 Al-Cr-Ni isothermal section at 900 °C for Al-rich alloys [2008Gru1]

structures of  $\tau_2$  and  $\tau_3$  were not resolved. The isothermal section at 700 °C constructed by [2008Wei] is shown in Fig. 6. The binary phase field designated  $\gamma_2 Cr_5 Al_8 - \gamma_3 Cr_4 Al_9$  by [2008Wei] is shown as  $\gamma_2$ . The binary phase Ni<sub>2</sub>Al<sub>3</sub> ( $\delta$ ) was found to dissolve up to 12 at.% Cr, which is much higher than 2.5 at.% found by [2008Gru1]. CrAl<sub>3</sub> (v) shown tentatively stable up to 750 °C by [2008Gru2] was not found by [2008Wei] at 700 °C.

The liquidus projection constructed by [2008Wei] is shown in Fig. 7, with some tentative modifications. [2008Wei] adopted three phases in the  $\gamma$ -region of the



Fig. 6 Al-Cr-Ni isothermal section at 700 °C in the Al-rich region [2008Wei]

Al-Cr system designated as  $\gamma_1(Cr,Al)$  or  $\gamma_1Cr_5Al_8$ ,  $\gamma_2Cr_5Al_8$ and  $\gamma_3Cr_4Al_9$ . In the Al-Cr phase diagram accepted here, only  $\gamma_1$  (cubic) and  $\gamma_2$  (rhombohedral) are recognized. Accordingly, the liquid-solid reactions involving  $\gamma_3Cr_4Al_9$ are omitted in Fig. 7. Only one binary peritectic reaction  $L + (Cr) \leftrightarrow \gamma_1$  is shown. [2008Wei] admitted that the issues regarding  $\gamma_2Cr_5Al_8-\gamma_3Cr_4Al_9$  are not entirely resolved. Clearly, there is need for a general consensus on the phases present in the  $\gamma$  field of the Al-Cr system. A reaction scheme developed by [2008Wei] (not shown here) will need modification, if the tentative view accepted here regarding the Al-Cr system prevails.

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Fig. 7 Al-Cr-Ni liquidus projection in the Al-rich region [2008Wei]

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