

Al-Sb-Y (Aluminum-Antimony-Yttrium)

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In the previous work on this ternary system, two isothermal sections have been reported: at 500 °C by [1971Mur] and at 527 °C by [2003Zen]. The latter study was reviewed briefly by [2008Rag]. Recently, [2009Zha] reported a thermodynamic assessment of this system.

Binary Systems

The Al-Sb phase diagram depicts the stoichiometric compound AlSb (*B3*, Sphalerite-type cubic). The Al-Y phase diagram [Massalski2] has the following intermediate phases: α Al₃Y (*D0₁₉*, Ni₃Sn-type hexagonal), β Al₃Y (*BaP₃*-type rhombohedral), Al₂Y (*C15*, MgCu₂-type cubic), AlY (*B_f*, CrB-type orthorhombic), Al₂Y₃ (*Al₂Zr₃*-type tetragonal), and AlY₂ (*C23*, Co₂Si-type orthorhombic). The Sb-Y phase diagram [Massalski2] depicts the following intermediate phases: SbY₃ (*PTi₃*-type tetragonal), Sb₃Y₅ (*D8₈*, Mn₅Si₃-type hexagonal), Sb₃Y₄ (*D7₃*, Th₃P₄-type cubic, stable between 2120 and 1660 °C), and SbY (*B1*, NaCl-type cubic).

Computed Ternary Phase Equilibria

[2009Zha] described the liquid, fcc, bcc, cph and rhombohedral phases as substitutional solutions. The binary compounds were treated as stoichiometric phases. The solubility of Al up to 27 at.% in Sb₃Y₅ was taken into account. The binary descriptions were used from the literature data. The ternary interaction parameters were set to zero. The optimized parameters were listed. Two isothermal sections at 527 and 500 °C and a liquidus projection were computed by [2009Zha]. The computed isothermal section at 527 °C agrees with that determined experimentally by [2003Zen]. The computed isothermal section at 500 °C is shown in Fig. 1. The triangulations of (Al) + α Al₃Y + SbY, (Al) + AlSb + SbY and AlSb + SbY + (Sb) agree with those found by [1971Mur], who studied the system up to 33.3 at.% Y. No experimental data are

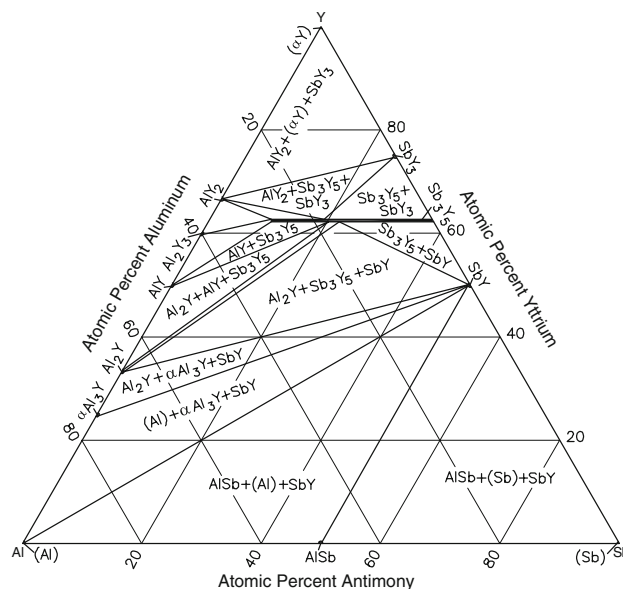


Fig. 1 Al-Sb-Y computed isothermal section at 500 °C [2009Zha]. Narrow two-phase regions are omitted

available for comparison with the computed liquidus projection of [2009Zha] (not shown here).

References

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