

Comment on Al-Mn (Aluminum-Manganese)

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Fig. 1 shows the Al-Mn phase diagram calculated by [92Jan]. The diagram is somewhat different from [Massalski2] (Fig. 2), which was redrawn from [87Mca] (with phase names changed for easier comparison).

The primary concern of the editor with regard to Fig. 1 is the for of Al_8Mn_5 phase field extended to include the δ phase field. Many ternary phase diagrams, for example Al-Co-Mn and Al-Cr-Mn, clearly indicate that δ and (δMn) belong the same bcc phase field. Therefore, bcc δ and orthorhombic Al_8Mn_5 should be differentiated, as in Fig. 2. In this regard, however, the liquidus and solidus of (δMn) and δ in Fig. 2 should be smoothly continuous in the metastable range, which is not the case in [87Mca]. The absence of λ in [92Jan] may be justifiable because λ with a composition very similar to Al_4Mn may be metastable [93Oka].

The Al-Mn crystal structure data in [87Mca] are updated in Table 1.

Cited References

- 30Wes:** A. Westgren, *Z. Metallkd.*, 22, 372 (1930).
38Hof: W. Hofmann, *Aluminium*, 20, 865-872 (1938).
58Bla: J.A. Bland, *Acta Crystallogr.*, 11, 236-244 (1958).
58Kon: H. Kono, *J. Phys. Soc. Jpn.*, 13(12), 1444-1451 (1958).
60Sch: K. Schubert, S. Bahn, W. Burkhardt, R. Gohle, H.G. Meissner, M. Potzschke, and E. Stolz, *Naturwissenschaften*, 47(13), 303 (1960).
61Tay: M.A. Taylor, *Acta Crystallogr.*, 14, 84 (1961).
75Bar: J.G. Barlock and L.F. Mondolfo, *Z. Metallkd.*, 66, 605-611 (1975).
75Oni: T. Onishi and Y. Nakatani, *Keikinzoku*, 25(7), 253-258 (1975).
87Mca: A.J. McAlister and J.L. Murray, *Bull. Alloy Phase Diagrams*, 8(5), 438-447 (1978).
89Sho: C.B. Shoemaker, D.A. Kessler, and D.P. Shoemaker, *Acta Crystallogr. B*, 45, 13-20 (1989).
92Jan: A. Jansson, *Metall. Trans. A*, 23(11), 2953-2962 (1992).
93Oka: H. Okamoto and T.B. Massalski, *J. Phase Equilibria.*, 14(3), 316-335 (1993).

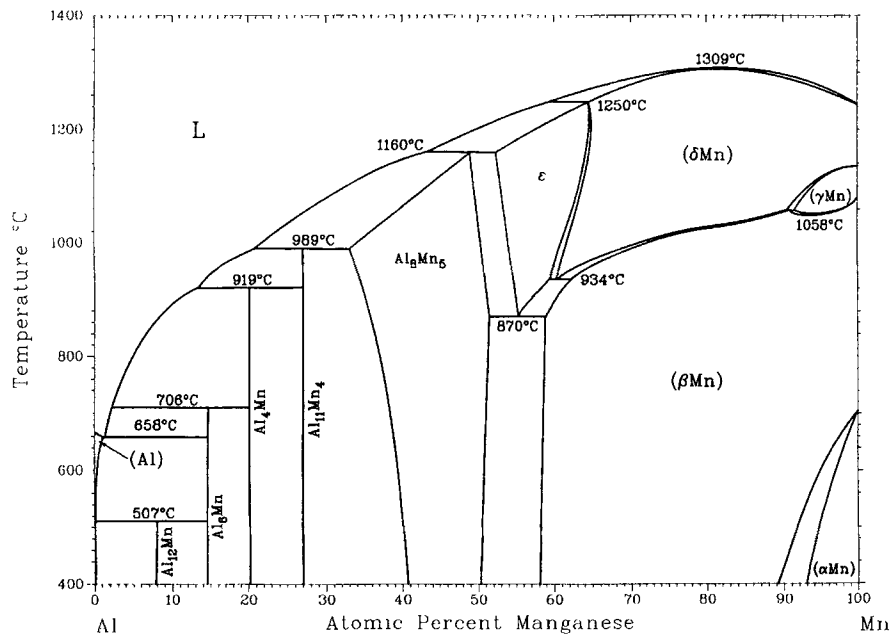


Fig. 1 Al-Mn phase diagram calculated by [92Jan].

Section III: Survey of Current Literature

Table 1 Al-Mn Crystal Structure Data

Phase	Composition, at.% Mn(1)	Pearson symbol	Space group	Strukturbericht designation	Prototype	Reference
(Al)	0 to 0.2	<i>cF4</i>	<i>Fm$\bar{3}m$</i>	A1	Cu	...
Al ₁₂ Mn	7.7	<i>cI27</i>	<i>Im$\bar{3}$</i>	[75Bar]
Al ₆ Mn	14.3	<i>oC28</i>	<i>Cmcm</i>	<i>D2_h</i>	Al ₆ Mn	[38Hof]
δ	?	<i>oP60</i>	<i>Pn\bar{m}</i>	[76Oni]
Al ₄ Mn	20	<i>hP574</i>	<i>P6₃/mmc</i>	[89Sho]
β Al ₁₁ Mn ₄	?	<i>oP160</i>	<i>Pn$\bar{m}a$</i>	[61Tay]
α Al ₁₁ Mn ₄	26.7	<i>aP30</i>	<i>P$\bar{1}$</i>	[58Bla]
γ	?
Al ₈ Mn ₅	33 to 50	<i>hR26</i>	<i>R$\bar{3}m$</i>	<i>D8₁₀</i>	Al ₈ Cr ₅	[60Sch]
δ	35 to 52	<i>cI2</i>	<i>Im$\bar{3}m$</i>	A2	W	[30Wes]
ϵ	52 to 64	<i>hP2</i>	<i>P6₃/mmc</i>	A3	Mg	[58Kon]
(δ Mn)	60 to 100	<i>cI2</i>	<i>Im$\bar{3}m$</i>	A2	W	...
(γ Mn)	91 to 100	<i>cF4</i>	<i>Fm$\bar{3}m$</i>	A1	Cu	...
(β Mn)	59 to 100	<i>cP20</i>	<i>P4₃2</i>	A13	β Mn	...
(α Mn)	? to 100	<i>cI58</i>	<i>I$\bar{4}3m$</i>	A12	α Mn	...

(1) Primarily according to [92Jan].

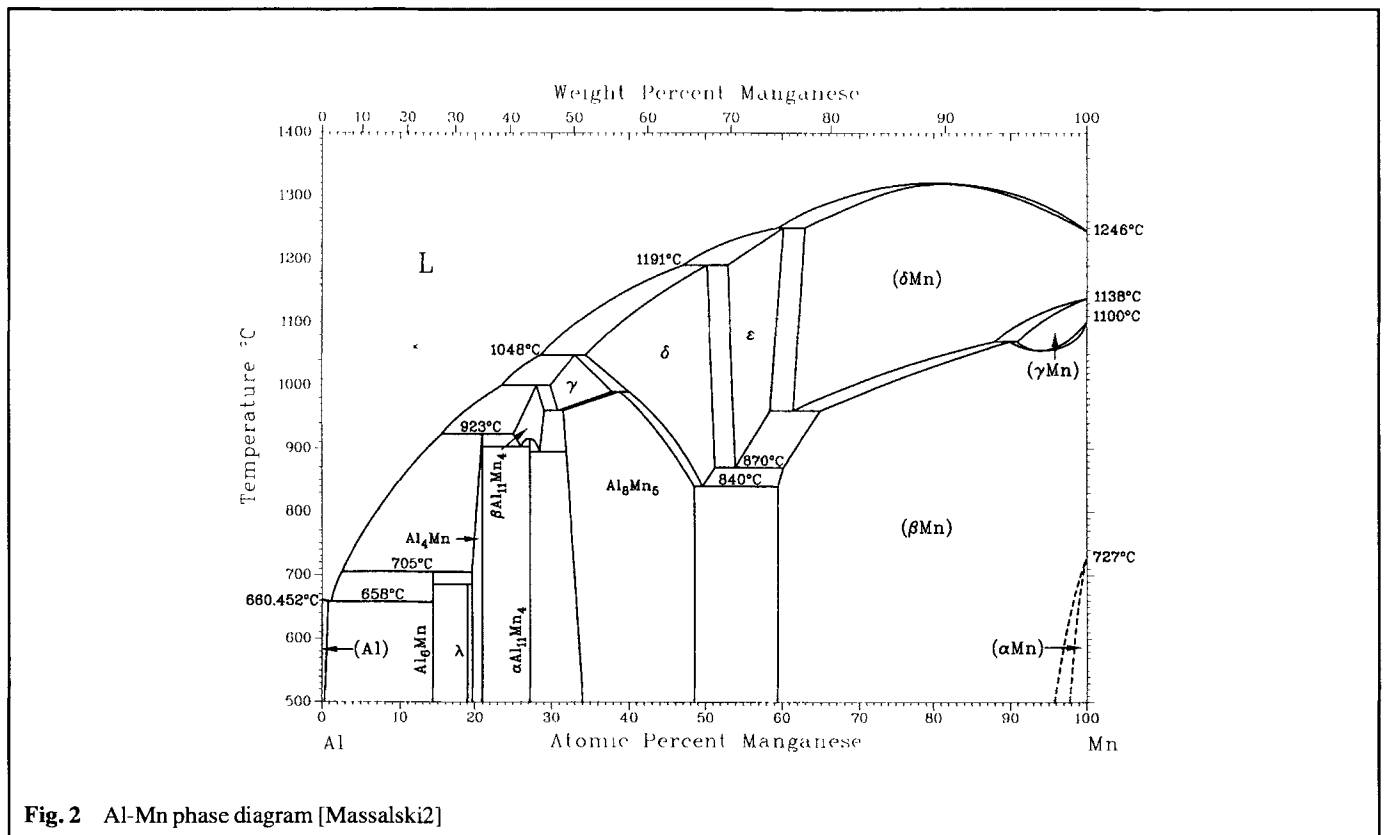


Fig. 2 Al-Mn phase diagram [Massalski2]