

Fe-Ge-La (Iron-Germanium-Lanthanum)

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Recently, [2008Zhu] determined an isothermal section of this system at 500 °C, which depicts three ternary compounds.

Binary Systems

In the Fe-Ge phase diagram, the intermediate phases found stable at 500 °C by [2008Zhu] are: Fe₃Ge (*L1*₂, AuCu₃-type cubic), Fe₅Ge₃ (*B8*₁, NiAs-type hexagonal), Fe₆Ge₅ (Fe₆Ge₅-type monoclinic, space group *C2/m*), FeGe, and FeGe₂ (*C16*, CuAl₂-type tetragonal). The Fe-La phase diagram [1997Zha] shows no intermediate phases. In the Ge-La system [Masslski2], the following seven compounds were found by [2008Zhu] at 500 °C: La₃Ge (tetragonal), La₅Ge₃ (*D8*₈, Mn₅Si₃-type hexagonal), La₄Ge₃ (*D7*₃, Th₃P₄-type cubic), La₅Ge₄ (Ge₄Sm₅-type orthorhombic), LaGe (*B27*, FeB-type orthorhombic), La₂Ge₃ (orthorhombic, space group *Imam*), and βLaGe_{2-x} (orthorhombic). The phase La₂Ge₃ is not seen in the phase diagram of [Masslski2], but is listed in [Pearson3] and found by [2008Zhu].

Ternary Phases

[2008Zhu] prepared a series of samples covering the composition range LaFe_xGe₂ ($x = 0.6-1$) and found that the compositions at the previously reported compounds LaFe_{0.69}Ge₂ and LaFeGe₂ consisted of two-phase mixtures of LaFe_{0.6}Ge₂ and LaFe₂Ge₂. The compound LaFe_{0.6}Ge₂

(denoted as τ₃ in Table 1 and Fig. 1 and as C by [2008Zhu]) has the CeNiSi₂-type orthorhombic structure [1990Fra]. LaFe₂Ge₂ (denoted τ₁ here and A by [2008Zhu]) has the Al₄Ba-type tetragonal structure. LaFeGe₃ (denoted τ₂ here and B by [2008Zhu]) has the BaNiSn₃-type tetragonal structure [1995Yam], see Table 1 for details.

Ternary Isothermal Section

With starting metals of 99.9% Fe, 99.999% Ge, and 99.9% La, [2008Zhu] arc-melted 126 alloy compositions under Ar atm. The alloys were given a final anneal at 500 °C for 10 days and quenched in liquid nitrogen. The phase equilibria were studied by optical microscopy, x-ray powder diffraction and differential thermal analysis at a heating rate of 10 °C/min. The isothermal section at 500 °C constructed by [2008Zhu] is shown in Fig. 1. The three ternary compounds τ₁, τ₂, and τ₃ are present. Whereas τ₁ and τ₂ are stoichiometric, LaFe_{0.6}Ge₂ (τ₃) shows a small homogeneity range of 16.7-18.0 at.% Fe ($x = 0.60-0.66$) [2008Zhu]. The equilibrium involving the ordered forms of bcc Fe (*B2* and *D0*₃) was not investigated by [2008Zhu] and these phases are not shown separately from bcc in Fig. 1. In a study of a series of alloy samples of La(Fe_{1-x}Ge_x)₁₃ ($x = 0-0.2$) at 500 °C, [2008Zhu] did not find the compound with the NaZn₁₃-type of structure reported by several authors in the related La(Fe_{1-x}Si_x)₁₃ alloys. Also, [2008Zhu] found that two other reported compounds La₃Fe₂Ge₄ and La₁₅FeGe₉ to be two-phase mixtures at 500 °C. The structural details of La₃Fe₂Ge₄ [Pearson3] and La₁₅FeGe₉ [1996Gul] are listed tentatively at the end of Table 1.

Table 1 Fe-Ge-La crystal structure and lattice parameter data [2008Zhu, Pearson3]

Phase	Composition, at.%	Pearson symbol	Space group	Prototype	Lattice parameter, nm
LaFe ₂ Ge ₂ (τ ₁ or A)	40 Fe 40 Ge 20 La	<i>tI10</i>	<i>I4/mmm</i>	Al ₄ Ba	$a = 0.4110$ $c = 1.0581$
LaFeGe ₃ (τ ₂ or B)	20 Fe 60 Ge 20 La	<i>tI10</i>	<i>I4mm</i>	BaNiSn ₃	$a = 0.4368$ $c = 0.9985$
LaFe _{0.6-0.66} Ge ₂ (τ ₃ or C)	16.7-18 Fe 55.6-54.6 Ge 27.8-27.3 La	<i>oC16</i>	<i>Cmcm</i>	CeNiSi ₂	$a = 0.4347$ $b = 1.683$ $c = 0.4210$
La ₃ Fe ₂ Ge ₄	22.2 Fe 44.4 Ge 33.3 La	<i>hP3</i>	<i>P6/mmm</i>	AlB ₂	$a = 0.4216$ $c = 0.4236$
La ₁₅ FeGe ₉	4 Fe 36 Ge 60 La	<i>hP?</i>	<i>P6₃/mcm</i>	Mn ₅ Si ₃	$a = 1.54810$ $c = 0.68768$

Section II: Phase Diagram Evaluations

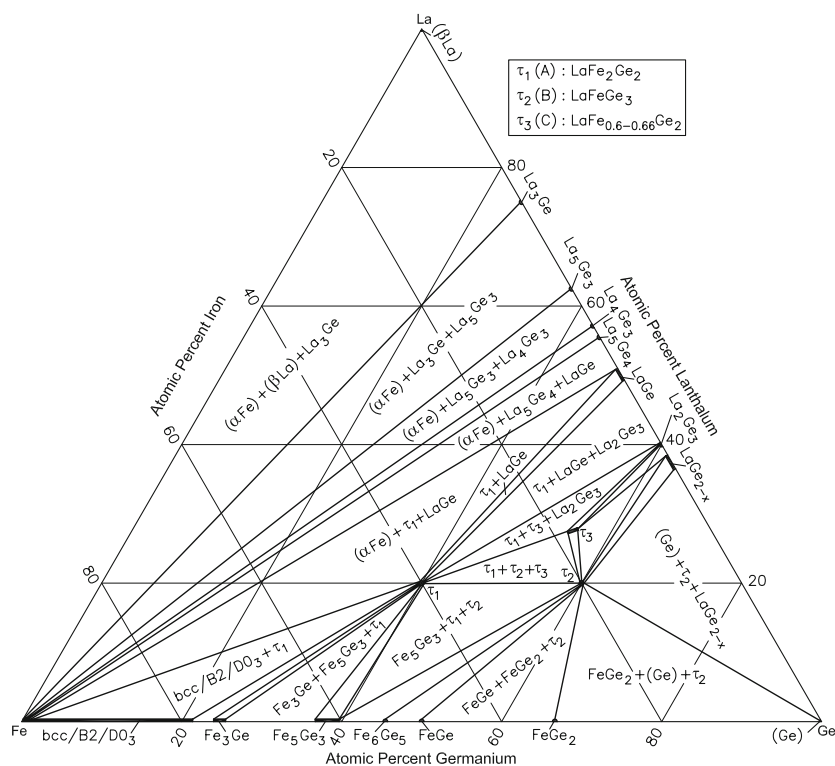


Fig. 1 Fe-Ge-La isothermal section at 500 °C [2008Zhu]. Narrow two-phase regions are omitted

References

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