#### PROCEEDINGS OF THE 4th E.C.F. CONFERENCE

AN ANALYTICAL MODEL OF INTERGRANULAR CREEP CRACKING.
APPLICATION TO PRECIPITATION-HARDENED ALL ALLOYS EMBRITTLED BY Pb.

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An analytical model has been developed to account for the kinetic evolution of crack populations in creeping alloys. Its application to measured crack distributions in precipitation hardened Al-base alloys embrittled by Pb leads to simple equations for crack growth v(x) and nucleation rate g(t) as functions of time t and crack length x:

 $v(x) = v_1 x$   $g(t) = g_0[1-exp \alpha(t-t_0)]$ 

These equations allow the influence of microstructural and compositional variables and of creep conditions on intergranular damage to be analysed through their effects on the parameters  $\mathbf{v}_1$  and  $\mathbf{g}_{\circ}.$ 

#### INTRODUCTION

Creep damage has generally been thought of as occurring by either of two distinct processes, wedge-type cracks and cavities. Moreover, the majority of microscopic models for creep cracking only take into account a single, semi-infinite, plane crack, ignoring deflections at triple junctions, and those for cavitation consider the evolution of either a single cavity or that of a uniformly distributed population of identical cavities.

Recent work has shown, however that these two types of damage often correspond to two different scales or stages of a single damage process, since cavitation was observed to develop at the tip of advancing cracks (1-6), and models were proposed to explain this behaviour (7-10). However these microscopic models do not take into account the experimental fact that creep damage takes the form of a non-uniform distribution of defects, cavities and/or cracks, which may therefore develop unevenly according to their size and the local values of other parameters.

In view of the complexity of these relationships, phenomenological models can be of some help in describing the kinetic evolution of a defect population, concurrently with physical models which shed light on the elementary mechanisms but overlook the statistical aspect of creep damage.

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#### EXPERIMENTAL DETAILS

In a study (6,11,12) of creep embrittlement of precipitation-hardened 6060-type Al-Mg-Si alloys by an unsoluble impurity, Pb, a quantitative experimental analysis of intergranular cracking was carried out by optical microscopy on four heats containing up to 164 ppm Pb.

The partial densities of cracks were empirically defined as the numbers  $\rho_{\Lambda}^{\Lambda}$  of cracks per unit area of longitudinal sections having a length in the interval  $[n\Delta - \frac{\Lambda}{2}, n\Lambda + \frac{\Lambda}{2}]$ , where n was an integer and  $\Delta$ , the width of all crack classes, corresponded to one division of the micrometer,  $\Delta = 5.75~\mu m$ . It was considered that the detection limit of the technique was about  $\Delta/2$ , which is the size of large inclusions and large intragranular cavities (6) with which the incipient cracks might be mistaken. Defects shorter than  $\Delta/2$  were thus ignored in the analysis.

Although a drastic decrease in creep life with increasing applied stress and Pb content was straightforwardly observed (6) no simple relationship could be establishedbetween creep cracking characteristics and macroscopic rupture properties ( $\hat{\epsilon}_{m}$ ,  $\sigma_{r}$ ,  $t_{r}$ ,  $\epsilon_{r}$ , etc) or Pb content (11,12). An analytical model of cracking, based on statistically significant data ( $_{\sim}120~000$  cracks counted and classified over 76 testpieces) was therefore developed in order to isolate the incidences of the various parameters.

#### GENERAL MODEL

# The observed micro-mechanism of crack growth

A phenomenological model of creep-crack distribution evolution has already been developed by Lindborg (13), based on the observation that the majority of intergranular cracks were arrested at triple junctions at both ends, which implied that growth was virtually instantaneous along individual grain boundary facets, and was controlled essentially by re-nucleation beyond triple junctions. Lindborg's probabilistic rationale, well suited to that discontinuous type of growth, in not compatible with the mode of crack growth in the present Al-base alloys, where the cracks advandiately ahead (i.e. within a few  $\mu m$ ) of their tip (6). This microscopic mechanism explains that crack growth is continuous at the scale of grain facets and is not basically controlled by arrest at triple junctions, resulting in the fact that more than 2/3 of the observed cracks have at least one of their ends between two maxima when the crack length, expressed in terms of number of grain boundary facets, is an integer, and the absolute maximum is for cracks smaller than one facet. Conversely, the crack density continuously decreases with increasing crack length when the latter is measured in absolute (metric) units, fig. 1-a, starting from a minimum value  $\Delta$  related to the detection limit  $\Delta/2$ . Thus, only the latter type of distribution has been considered, and a continuum model had to be developed to account for the optical microscopy observations.

## Hypotheses of the model

a - Coalescence of, and interaction between, individual cracks are neglected, which is reasonable in view of the small absolute density of cracks even at rupture (11).

b - The growth rate v of an individual crack in a given alloy for given heat treatment and testing conditions is essentially a function of its length x. In most of the microscopic models of crack growth, v is found to increase as  $K^m$ , where K is the applied stress intensity factor, proportional to  $x^{1/2}$ , and m is a constant. Therefore v(x) is assumed to have the form:

 $v(x) = \frac{dx}{dt} = v_q \ x^q \qquad \qquad (\underline{growth} \ equation)$  where q = m/2. When vacancy diffusion is responsible for crack growth and crack tip stress concentrations are not neglected as in ref. 14,15,  $2 \le m \le 4$  (8,9,10) and thus  $1 \le q \le 2$ . When crack extension is controlled solely by plastic deformation, m becomes equal to the stress component of the creep law  $\varepsilon = A\sigma^m$  and can therefore be much larger (7). Experimental values ranging between 3 and 30 have been obtained (16).

c - The partial crack density at time t,  $\rho(x,t)$ , is an analytical function of x and t whose values at  $x = n\Delta$  are related to the experimentally measured  $\rho_n^{\Delta}$  by the equation :

$$\rho(n\Delta, t) = \rho \frac{\Delta}{n} / \Delta \tag{2}$$

d - Only an empirical acceptation of nucleation is implied here: the "nucleation rate"  $g^{\Delta}\left(t\right)$  is considered as the number of cracks which become detectable, i.e. longer than  $\Delta/2$ , per unit time. The same convention, reminded by the superscript  $\Delta$ , applies to the total density  $\rho^{\Delta}(t)$  and integrated length  $\mathscr{L}^{\Delta}\left(t\right)$  of cracks per unit area at time t.

#### Basic equations of the model

The possibility of crack sintering being outruled,  $v\left(x\right)\,\geqslant\,0$ and the ranking of existing cracks according to their length and the ranking of existing cracks according to their length remains unchanged during growth. Therefore the number of cracks having their lengths in the interval  $[x-\frac{\Delta}{2}, x+\frac{\Delta}{2}]$  at time t, will be found unchanged in the interval  $[(x-\frac{\Delta}{2})+v(x-\frac{\Delta}{2}).dt, (x+\frac{\Delta}{2})+v(x+\frac{\Delta}{2}).dt]$  at time t+dt. This yields the distribution equation, which is formally identical to a conservation equation:

Similarly, the "nucleation" rate is equal to the number of cracks whose lengths grow over  $\Delta/2$  per unit time:

$$g^{\Delta}(t) = v(\frac{\Delta}{2}) \cdot \rho(\frac{\Delta}{2}, t)$$
 (4)

The incubation time to is that at which the <u>first</u> crack becomes detectable (i.e.  $x = \Delta/2$ ). Under the above assumptions, this crack will remain the longest one throughout creep life. Its length xmax at time t is related to t by the equation :  $t-t_0 = \int_{\Delta/2}^{X_{max}} [v(x)]^{-1} dx$ 

$$t - t_o = \int_{\Lambda/2}^{x_{\text{max}}} \left[ v(x) \right]^{-1} dx \tag{5}$$

The total density and length of cracks can be calculated by:
$$\rho^{\Delta}(t) = \int_{\Delta/2}^{X_{\text{max}}} \rho(x,t) dx \tag{6}$$

$$\mathcal{Z}^{\Delta}(t) = \int_{\Delta/2}^{X_{\text{max}}} x \rho(x,t) dx \tag{7}$$
Eleven boundary conditions can be written:

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Eleven boundary conditions can be written:

$$\forall t, \ t_{o} \leq t < t_{r} :$$

$$g^{\Delta}(t) > 0 \ ; \ \rho^{\Delta}(t) > 0 \ ; \ \mathcal{L}^{\Delta}(t) > 0$$
(8)

$$\forall t, t_{o} \leq t < t_{r} ; \forall x, \frac{\Delta}{2} \leq x \leq x_{max} :$$

$$\rho \geq 0 ; \frac{\partial \rho}{\partial t} \geq 0 ; \frac{\partial \rho}{\partial x} \leq 0 ; \frac{\partial^{2} \rho}{\partial x^{2}} \geq 0$$
(9)

$$\rho\left(\frac{\Delta}{2}, t_{o}\right) = \rho^{\Delta}\left(t_{o}\right) = \mathcal{L}^{\Delta}\left(t_{o}\right) = 0 \tag{10}$$

$$\rho\left(\mathbf{x}_{\text{max}},\mathsf{t}\right) = 0 \tag{11}$$

Among these, conditions (9) are imposed to comply with the shape and time dependence of the  $\rho$  vs. x curves, fig. 1-b. In particular the condition on curvature is strictly valid for large x only.

### Resolution of the equations

Integration of the growth equation 1 yields :

$$x = (\frac{\Delta}{2}) \cdot \exp[v_q(t - t_i)/q]$$
(12)

where  $t_i$  is the time at which a given crack is first detected i.e.  $x_i = \Delta/2$ . For the first crack  $t_i = t_o$  and  $x = x_{max}$ , eqn. 5.

A general solution of eqn. 3 with separated variables has been obtained (11), using the above expression of v(x), eqn. 1:

obtained (II), using the above expression of 
$$V(x,y)$$
 equal 
$$\rho(x,t) = (g_o/v_q) x^{-q} \left\{1 - \exp[\alpha(t-t_o) - C\beta(x)]\right\}$$
(13)

where  $g_0$ ,  $\alpha$  and C are constants,

$$C = (\alpha/v_q) (\Delta/2)^{1-q}$$
(14)

and 
$$q \neq 1$$
  $\beta(x) = (1-q)^{-1} [(2x/\Delta)^{1-q} - 1]$  (15-a)  
 $q = 1$   $\beta(x) = \ln(2x/\Delta)$  (15-b)

The "nucleation" rate and the total density of cracks as defined by eqns. 4 and 6 have the expressions:

$$g^{\Delta}(t) = g_{0} \{1 - \exp[\alpha(t - t_{0})]\}$$
 (16)

$$\rho^{\Delta}(t) = g_{0}(t - t_{0}) + (g_{0}/\alpha)\{1 - \exp[\alpha(t - t_{0})]\}$$
 (17)

Condition (11) leads to the time dependence of  $x_{\text{max}}$ :

$$\alpha(t-t_o) = C\beta(x_{max})$$
 (18)

The whole set of conditions (8-11) essentially reduces to (11) :

$$\begin{array}{c|c} a.g. < 0 \\ \hline g > 1 & |C| \leqslant 3q \end{array}$$
 (19-a)

The basic condition eqn.19-adecomposes into two cases, fig.2: a - If  $\alpha>0$  and  $g_o<0$ , the "nucleation" rate and crack density increase exponentially with time, eqns. 16,17, and the majority of cracks are formed at the end of creep life. This behaviour is expected in case of essentially strain-controlled creep damage, which dominates at high stresses and creep rates. b - If  $\alpha<0$  and  $g_o>0$ ,  $g^{\Delta}(t)$  tends asymptotically towards a constant value  $g_o$  which then has the simple physical meaning of a steady state "nucleation rate". Accordingly, the total density of cracks  $\rho^{\Delta}(t)$  tends to become linear with time :

$$\rho^{\Delta}(t) \rightarrow g_{\circ}(t-t_{\circ}) \tag{20}$$

The time necessary to establish this steady state regime is of the order of  $1/\alpha\,.$ 

#### APPLICATION OF THE MODEL TO Al-Mg-Si ALLOYS

The observed variations of total crack density with time, fig. 3-a, being virtually linear (for 65 and 164 ppm Pb), at least after a certain time (for 29 ppm Pb), imply that  $\underline{\alpha}<0$  and  $\underline{g}_o>0$  in the alloys and creep conditions considered  $(T=373~K\simeq0.45~T_m\,;~\sigma=140~to~210~MPa)$ . These linear plots also imply that the exponential in eqn. 17 has become negligible, and, since  $\beta$  (x) increases with x and C is negative (as  $\alpha$ ), eqn. 13 is equivalent to :

$$\rho(x, t) \simeq (g_{\circ}/v_{q})x^{-q} = B_{q} x^{-q}$$
 (21)

for short cracks. Similarly, according to eqn. 2:

 $\rho_n^\Delta(n,\,t) \simeq B_q \Delta^{1-q} n^{-q}$  (22) Considering experimentally determined histograms of  $\rho_n^\Delta$ , fig. 4, it appears that q=1 is the best value consistent with the results for the lower values of n, i.e. in the range where the approximation of eqn. 21 applies. Only a few values of  $\rho_n^\Delta$  appear somewhat erratic, due to the difficulty in identifying the smallest cracks. These conclusions hold for all the results obtained with 76 specimens of 4 alloys in 3 heat treatment conditions, and creep times ranging from 6 to 1 800 h (11,12).

The value q = 1 is consistent with the predictions of models of crack advance involving vacancy diffusion-controlled growth of cavities ahead of crack tip (8,9), i.e.  $v_{\alpha}\kappa^2$  for the larger stress intensity factors, and with our experimental observations which reveal the importance of diffusion processes in the creep-rupture of these alloys (6,11,12) :

- cavities are themselves divided into smaller vacuoles bounded by simple cristallographic planes, as occurs in the surface diffusion-controlled formation of equilibrium cavities; this causes the surfaces of freshly formed cracks to appear pitted at an extremely fine scale ( $\leqslant$  0.1  $\mu m$ ), see fig. 5 of ref. 6 ;

faces of freshly formed cracks to appear pitted at an extremely fine scale ( $\leqslant$  0.1  $\mu m)$ , see fig. 5 of ref. 6; 
– later on, these fine features smooth out by surface diffusion; 
– the embrittling effect of Pb is enhanced as the creep rate is decreased, i.e. as time dependent mechanisms are allowed to prevail; 
– Auger electron spectroscopy has demonstrated that Pb rapidly covers fracture surfaces by surface diffusion even at low temperatures; this is probably the basic mechanism of embrittlement.

The influence of microstructural and compositional variables upon "nucleation" and growth can be assessed by considering their effects on the parameters  $v_1$  and  $g_0$  of the model.  $g_0$  is the slope of the straight lines in fig. 3-a, and  $B_1$  can be graphically determined in fig. 4 with the help of eqn. 22. Then  $v_1 = g_0/B_1$ . The loglog plots of  $v_1$  and  $g_0$  vs. the minimum creep rate  $\hat{\epsilon}_m$ , fig. 5, reveal that remarkable correlations of the type

$$v_1 = v_1^{\circ} (\dot{\epsilon}_{\mathfrak{m}})^{\mathsf{r}} \qquad g_{\circ} = g_{\circ}^{\circ} . (\dot{\epsilon}_{\mathfrak{m}})^{\mathsf{S}}$$
 (23)

hold over 4 orders of magnitude of  $\dot{\epsilon}_m$ , which emphasizes the importance of plastic creep deformation process in crack nucleation and growth. The parameters  $v_1^{\circ}$ ,  $g_{\circ}^{\circ}$ , r and s are characteristic of the material condition. Both  $v_1^{\circ}$  and  $g_{\circ}^{\circ}$  increase with increasing Pb content (11,12), which implies that both "nucleation" and growth are enhanced by the embrittling impurity. A decrease in grain size strongly increases  $B_1$  by favoring "nucleation" ( $g_{\circ}$  increases) and hindering growth ( $v_1$  decreases). In the precipitation-hardened conditions,  $v_1$ ,  $g_{\circ}$  and  $g_{\circ}$  are virtually independent of aging time and Mg content, which implies that, for a given type of micro-

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structure, crack nucleation and growth are not affected by hardness intrinsically but only via its effect on creep rate.

In the case q=1, simple equations relating  $v_1$  and  $g_0$  $x_{max}$  can be established by combining eqns. 14, 15-a and 18:

(24)(cf. eqn. 12)  $v_1(t-t_0) = ln(2x_{max}/\Delta)$ (25) $g_o(t-t_o) = B_1 \ln(2x_{max}/\Delta)$ 

Numerical values of  $v_1$  and  $g_0$  can be consistently deduced from  $\ln\left(x_{max}\right)$  vs. t plots, but this often leads to negative values of to, as expected from fig. 4. In the present model, negative incubation times allow to account formally for the existence of defect nuclei prior to the onset of creep, as can be induced by cold work (17), or for a regime of intense nucleation during loading or primary creep (18), as depicted schematically in fig. 6.

### CONCLUDING REMARKS

Due to its general character, the model is applicable to a variety of materials, provided a number of crack distributions are known of materials, provided a number of crack distributions are known experimentally. These are necessary to identify the parameters, beginning with the sign of  $\alpha$ , which allows to choose between type a ("tertiary" cracking) or type b ("continuous" cracking) behaviours as depicted in fig. 2, and the value of q which may depend on the material type and on the ranges of temperature and applied stress. Even though they appear quite intricate in the raw experimental results, the influences of the various microstructural-compositional parameters and of creep conditions upon the cracking and thus on the rupture-life behaviours, lend themselves to a more straightforward analysis when the model equations and their parameters  $(g_{\circ},\,v_{q})$  and probably  $\alpha$  in other cases) are used, despite meters (g.,  $v_q$  and probably  $\alpha$  in other cases) are used, despite the simplicity of the initial assumptions of the model.

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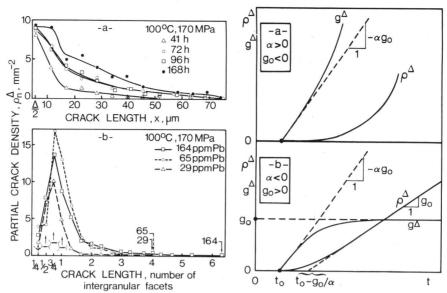


Figure 1 Examples of crack distribution 
Figure 2 Schematic of predicted  $\rho^{\Lambda}(t)$  in terms of : a - absolute length ; b - and  $g^{\Lambda}(t)$  curves with the condition in terms of : a - absolute length; b - number of cracked intergranular facets.

 $\alpha.g_{\circ} < 0.$ 

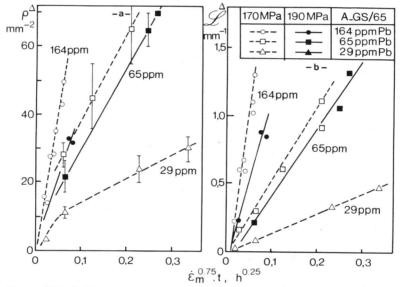


Figure 3 Evolution of total crack density  $\rho^\Delta$  and length  $\mathcal{L}^\Delta$  in peakaged alloys with various Pb contents. Time has been corrected for variations in creep rate  $\dot{\epsilon}_{m}$ .

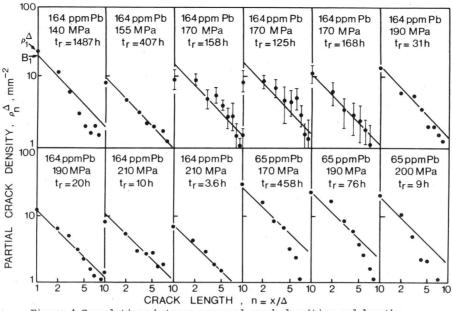


Figure 4 Correlations between measured crack densities and lengths (the slope of all solid lines is q=1).

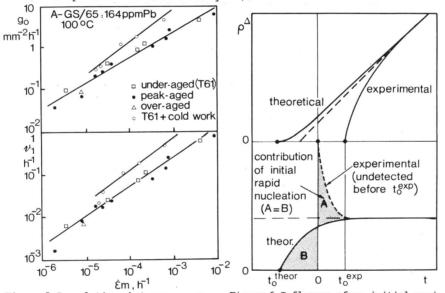


Figure 6 Influence of an initial regime of rapid nucleation superimposed on the solution of the model equations.