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## Fatigue Crack Growth Rate Model for Metallic Alloys

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### Abstract

A model has been created to allow the quantitative estimation of the fatigue crack growth rate in steels as a function of mechanical properties, test–specimen characteristics, stress–intensity range and test–frequency. With this design, the remarkable result is that the method which is based on steels, can be used without modification, and without any prior fatigue test, to estimate the crack growth rates in nickel, titanium and aluminium alloys. It appears therefore that a large proportion of the differences in the fatigue crack growth rate of metallic alloys can be explained in terms of the macroscopic tensile properties of the material rather than the details of the microstructure and chemical composition.

**Keywords:** Fatigue Crack Growth, Neural Network, Steel, Superalloys, Titanium Alloys, Aluminium Alloys

### 1 Introduction

It is understood that fatigue crack growth is a consequence of the accumulation of damage by deformation in the plastic zone at the crack tip. At low loads the deformation is governed by the cyclic variation in the stress-intensity range  $\Delta K$ . The crack extension per cycle (da/dN) becomes measurable at a threshold  $\Delta K_{th}$ , followed by the slower extension rate in the Paris Law regime [1–4] described by the proportionality

$$\log \Delta K \propto \log \left\{ \frac{da}{dN} \right\}^m \tag{1}$$

where da/dN is the average crack advance per cycle, and m is known as the Paris exponent. The equation can be interpreted in terms of a variety of physical mechanisms [5, 6], in which case the proportionality constant (C) becomes a function of the Young's Modulus E, the Poisson's ratio  $\nu$ , and the yield and ultimate tensile strengths  $\sigma_Y$  and  $\sigma_U$  respectively. Based on the possible

mechanisms consistent with the Paris Law, attempts have been made to generally interpret fatigue crack growth data on the basis of just the mechanical properties of the material [5, 6].

Elber modified the relation with an effective stress intensity range  $\Delta K_{eff}$  to allow for variable amplitude loading, arguing that cracks grow only when their tips are open [7]:

$$\frac{da}{dN} = C_0 (\Delta K_{eff})^m \quad \text{with} \quad C_0 = C/(0.7^m) \tag{2}$$

These equations do not explicitly contain material properties; Duggan [8] expressed the crack growth rate in terms of the elastic modulus, toughness, and ductility:

$$\frac{da}{dN} = \left(\frac{\pi}{32}\right)^{\frac{1}{2\eta}} \frac{1}{\eta} \left[\frac{2}{\epsilon_f E(K_{Ic} - K_{max})} \left(1 - \frac{K}{K_{Ic}}\right)\right]^{\frac{1}{\eta}} K^{\frac{2}{\eta}}$$
(3)

where  $\eta$  and  $\epsilon$  are the fatigue ductility exponent and coefficient respectively, E is the elastic modulus,  $K_{Ic}$  is the critical stress intensity for fracture. Ramsamooj and Shugar have also accounted for toughness, yield strength and modulus, but not for frequency and the analysis is presumably limited to mode I loading [9]. Their model is interesting in that it generalises against iron, aluminium and titanium alloys; it does however require a prior knowledge of the threshold stress intensity range for fatigue crack growth.

The aim here was to exploit published fatigue crack growth data to create a model based on physical variables which are readily measured in a tensile test, rather than rely on inputs which depend on fatigue testing, and further, to include variables which account for test-specimen parameters. The model uses neural network analysis; although there are physically based models available in the literature, for example, [10], they require fitting parameters; a neural network is the most general way of achieving fitting without making prior assumptions about the relationship to which the data are fitted [11]. There have been other attempts to use neural networks for this purpose [12] but they do not adequately treat the uncertainties of modelling so it is not possible to properly assess the predictions made. The original intention here was to study steels, but as will be seen later, the model was, without modification, found to generalise to other alloy systems.

#### 2 Method

A versatile method for treating empirical data is the neural network in a Bayesian framework. The theory behind practical Bayesian networks has been described in [13, 14] and the background information theory is available in a seminal textbook on the subject [15]. In addition, this method been reviewed thoroughly [11], as have been its applications [16]. Indeed, there have been diverse applications which lead to useful and verifiable predictions in the context of low-cycle fatigue [17], the estimation of bainite plate thickness [18], the calculation of ferrite number in stainless steels [19], the estimation of tensile strength [20, 27], impact strength [21, 26], the effect of processing parameters on marageing steels [22], the modelling of strain induced martensitic transformation [23], and the reduction in steel varieties [28], to name but a few. There has even been an assessment of procedures needed to design networks which are well-assessed in their performance [24]. Given this plethora of literature, only specific points of relevance are introduced here.

With neural networks, the input data  $x_j$  are multiplied by weights, but the sum of all these products forms the argument of a flexible mathematical function (known as the *transfer function*), here a hyperbolic tangent. The output y is therefore a non-linear function of  $x_j$ . The exact shape of the hyperbolic tangent can be varied by altering the weights. Further degrees of non-linearity can be introduced by combining several of these hyperbolic tangents, so that the network is able to model highly non-linear relationships. The nature of the transfer functions and the weights define a reproducible mathematical function which represents the empirical data.

The network just described is essentially a non-linear regression method which, because of its flexibility, is able to capture complicated data, whilst at the same time avoiding overfitting. There are a number of interesting outputs other than the coefficients which help recognise the significance of each input. First, there is the *noise* in the output, associated with the fact the input set is unlikely to be comprehensive – i.e., a different result is obtained from identical experiments. Secondly, there is the *uncertainty of modelling* because many mathematical functions may be able to adequately represent known data but which behave differently when extrapolated. A knowledge of this uncertainty helps make the method less risky in extrapolation. This uncertainty can be expected to be large in regions of the input domain where data are sparse or exceptionally noisy.

### 3 The Variables

Published [25] fatigue crack growth data for tests done in ordinary air, at room temperature, were digitised, covering steels with chemical compositions in the range presented in Table 1. Traces element concentrations (Ti, Al, V, S, P) together with the details of heat treatment can be found in the original compilation [25]. The properties of a steel depend on the composition and heat treatment, but fatigue crack propagation should depend to a large extent on macroscopic mechanical properties. It was deliberately decided to focus on easily measured properties obtained from a tensile test, rather than use inputs such as the threshold stress intensity which would defeat the purpose of modelling since a fatigue test would be required before a prediction could be made. The dimensions of the test specimens and the test conditions are also important in this respect and were included in the analysis. The advantage of this approach also is that a large quantity of data are available with each of the input variables listed in Table 1. Data for both axial mode I and an in-plane bending mode II were incorporated; mode III data were not available.

The plots in Fig. 1 illustrate the distribution of data, but clearly cannot represent multidimensional dependencies. However, the neural network method used here is based on a Bayesian framework [13, 15] so that the predictions are associated with a *modelling uncertainty* whose magnitude depends on the position in the input domain where a calculation is done. As pointed out previously, the details of the neural network and Bayesian framework used have been fully described elsewhere so only the essential points are included in this paper.

#### 3.1 Training the Model

The data were randomly and equally divided into the training and testing sets, and normalised [11]. One hundred networks were trained, with hidden units ranging from one to twenty and five seeds in each case. This is in order to select a committee of models which gives the best generalisation on unseen data [11, 13, 14, 29]. The performance of the optimum committee accompanies by  $\pm 1\sigma$  modelling uncertainties is illustrated in Fig. 2. Of the total of 12807 data, only 158 can be classified as mild outliers which are more than  $3\sigma$  from the measured values. The noise in the output of the committee model was found to be  $\sigma = \pm 4\%$ , which is a constant additional error to the modelling uncertainties plotted in subsequent graphs. The network perceived significances, which indicate the ability of an input to explain the variation in the output (akin a partial correlation coefficient) are shown in Fig. 3. The elongation, ultimate tensile strength and proof stress are significant in influencing da/dN but it is natural that the stress intensity range  $\Delta K$  should have the greatest effect. Although it is expected in a valid test that specimen size should not influence da/dN [30], it is likely that true plane strain conditions do not exist in all the cases studied, and hence a specimen size effect is perceived in Fig. 3. Such behaviour has been reported previously, with the crack growth rate increasing as plane–strain conditions are approached [31].

One way of assessing a model is by making predictions, in this case on a bearing steel of relevance in our other research. The steel of interest is variously known as SUJ2, AISI 52100 and En31 in different countries and has the approximate composition 1C, 0.3–1.1Mn, 1.2–1.4Cr, 0.2–0.4Si wt%. The inputs required were obtained from [32]: 5% elongation, 2030 MPa 0.2% proof stress, 2240 MPa tensile strength, loading mode 2, specimen length 80 mm, specimen thickness 2 mm, pre-crack size 3 mm, frequency 2 Hz and stress ratio 0.

Fig. 4 shows the outcome, with the model not only capturing the trend in the variation of da/dN versus  $\Delta K$  over several orders of magnitude, and both for the threshold and Paris regions of the curve, but giving also a reasonable absolute prediction accuracy.

Although all of the data used to create the model were from experiments on steels [25], the inputs include only mechanical and test parameters. It was imagined that the model should therefore apply without modification to other alloys.

#### 3.2 Ni, Ti and Al Alloys

Calculations for three nickel-base superalloys Udimet 700, Inconel 718 and Waspaloy; their detailed compositions can be found in [33–37]. Fig. 5 compares the model and the experimental data using the inputs listed in Table 2. The calculations are represented with the uncertainty range and the reported measurements [33–37] as points. The results are fascinating since the model correctly estimates the Paris slopes, although it marginally overestimates the fatigue behaviour (we have checked that this overestimation is not explained by modulus variations between the different materials). A similar level of agreement was found for titanium Ti-6Al-4V, 7075 aluminium alloy were made and compared against published measurements [38, 39], Fig. 6. The model nicely captured the slope for both the titanium and aluminium alloys, and it again slightly overestimates the fatigue crack growth rates.

In order to further test the model a colleague from industry supplied input data (last two columns, Table 2) without revealing the alloy type for the purpose of making blind predictions, for which the crack growth rates would be revealed after the calculations are made. Fig. 7 shows calculations, and the subsequent experimental data on the same Ti6Al4V alloy with two different heat treatments. The agreement obtained is good.

### 4 Conclusions

- 1. It has been possible to design a neural network model for fatigue crack growth in steels, which relies on inputs that consist only of properties that can be obtained from a simple tensile test, and using information about specimen geometry and testing parameters.
- 2. Given the nature of the inputs, it has been demonstrated that although the model is based entirely on data from steels, it can be applied without modification to nickel, titanium and aluminium alloys.
- 3. Given recent work where a similar approach has been used in modelling the hot-tensile strength [40] and stretch flangeability [41], it becomes evident that the neural network method has enormous potential for creating models for complex mechanical properties on the basis of simple experiments, such as the data obtained during tensile testing.

The computer program associated with this work can be downloaded freely from:

http://www.msm.cam.ac.uk/map/mapmain.html

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Element	Range	Element	Range
Carbon	0.1 – 0.8	Copper	0 – 0.2
Chromium	$0\!-\!5$	Manganese	0 - 2
Molybdenum	$0\!-\!2$	Nickel	$0\!-\!2$
Silicon	0 - 2		
Variable	Range	Variable	Range
Elongation / $\%$	0.2 - 61.8	0.2% Proof stress / MPa	122 - 1735
Tensile strength / MPa	270 - 2206	Specimen length / mm $$	13 - 260
Specimen thickness / mm	1.2 - 134	Pre-crack length / mm $$	1 - 52
Stress ratio	-1-1	Frequency / Hz	1 - 150
$\Delta K$ / ${\rm MPa}{\rm m}^{1/2}$	2.5 - 142	$\frac{da}{dN}$ / mm cycle <sup>-1</sup>	$9.82{\times}10^{-10}{-}4.86{\times}10^{-1}$

Table 1: Chemical composition range (wt%) of the steels studied. The lower half of the table shows the variables actually included in the analysis.

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Table 2:	The inputs	for the	prediction	ns in	Figs	5,	6, 7	covering	nickel,	titaniun	ı and	alum	inium
alloys.													
Variable	9		Figure number										
				- 1	-		~ 1	-	- 0	0 0		_	-1

Figure number									
5a	5b	5c	5d	5e	5f	6a	6b	7a	$7\mathrm{b}$
5	15	20	20	27	33	14	8	20	14
1020	1172	1113	1113	1076	921	930	524	1172	940
1520	1404	1373	1373	1441	1351	970	464	1440	998
72.5	63.5	50.8	31.8	62.5	5	155	155	7	7
12.5	25.4	12.7	8.89	25	3	40	40	7	7
12.5	18.3	6.4	5.3	17.5	0.4	9	9	0.5	0.5
0.1	0.1	0.05	0.05	0.5	0.5	-1	0.5	0.1	0.5
40	20	0.667	0.667	20	100	20	20	0.25	100
	5a $5$ $1020$ $1520$ $72.5$ $12.5$ $12.5$ $0.1$ $40$	5a5b515102011721520140472.563.512.525.412.518.30.10.14020	5a5b5c5152010201172111315201404137372.563.550.812.525.412.712.518.36.40.10.10.0540200.667	5a         5b         5c         5d           5         15         20         20           1020         1172         1113         1113           1520         1404         1373         1373           72.5         63.5         50.8         31.8           12.5         25.4         12.7         8.89           12.5         18.3         6.4         5.3           0.1         0.1         0.05         0.05           40         20         0.667         0.667	Figure nu5a5b5c5d5e515202027102011721113111310761520140413731373144172.563.550.831.862.512.525.412.78.892512.518.36.45.317.50.10.10.050.050.540200.6670.66720	5a5b5c5d5e5f515202027331020117211131113107692115201404137313731441135172.563.550.831.862.5512.525.412.78.8925312.518.36.45.317.50.40.10.10.050.050.50.540200.6670.66720100	Figure number5a5b5c5d5e5f6a51520202733141020117211131113107692193015201404137313731441135197072.563.550.831.862.5515512.525.412.78.892534012.518.36.45.317.50.490.10.10.050.050.50.5-140200.6670.6672010020	Figure number5a5b5c5d5e5f6a6b515202027331481020117211131113107692193052415201404137313731441135197046472.563.550.831.862.5515515512.525.412.78.89253404012.518.36.45.317.50.4990.10.10.050.050.50.5-10.540200.6670.667201002020	Figure number5a5b5c5d5e5f6a6b7a5152020273314820102011721113111310769219305241172152014041373137314411351970464144072.563.550.831.862.55155155712.525.412.78.892534040712.518.36.45.317.50.4990.50.10.10.050.050.51.51550.140200.6670.6672010020200.25



Figure 1: Distribution of data used to create the model.



Figure 2: Performance of the committee of models on the entire dataset of 12807 experiments.



Figure 3: Perceived significance of the inputs in the committee model. Both the mean significance and the upper and lower limits from the members of the committee are shown.



Figure 4: Calculations for a bearing steel. The points represent experimental data from [32], whereas the uncertainty range illustrated is calculated.



Figure 5: Predictions represented by the uncertainty range, and experimental data presented as points, for nickel based superalloys. (a) Udimet 700, data from [33]. (b) Inconel 718 with data dues to [34]. (c,d) Inconel 718, data from [35]. (e) Waspaloy, data from [36]. (f) Waspaloy, data from [37].



Figure 6: Calculations (uncertainty ranges) for titanium and aluminium alloys compared with measurements (points) due to [38, 39].



Figure 7: Blind predictions for Ti 6/4 forging material.