

# Modelling of active transformation of microstructure of two-phase Ti alloys during hot working

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Being very strong and of limited ductility, Ti alloys require special techniques to manufacture the parts with complex shapes. Many of these technologies are based on superplastic and near-to-superplastic deforming. In these processes the transformation of the microstructure of the material can be very significant and can lead to changes in the mechanical properties of the material during deformation. Because of this an appropriate description of the correlation between mechanical loading, changes in microstructure and mechanical behavior of material is required. A phenomenological scalar model with an internal variable based on a statistical description of microstructure is proposed and used for simulation of the high temperature deformation of Ti-6Al-4V alloy. The problems of obtaining the parameters of the model and sensitivity of the model to the accuracy of this process are discussed.

Keywords: constitutive modelling, simulation, superplasticity, microstructure transformation.

## 1. Introduction

It is established [1,2] that almost all polycrystalline materials and even ceramics are capable of showing under certain conditions a special kind of mechanical behavior called «superplasticity» (SP). Thus even materials that are of limited ductility in normal conditions can be deformed to hundreds and sometimes a few thousand percent frequently without occurrence of a localization of deformation which accelerates the process of failure. The conditions required for such behavior of materials, are usually determined experimentally and thus by nature empirical. For example, for the majority of alloys, for the maintenance of the superplastic condition it is necessary that the temperature ( $T$ ) is in the order  $0.4T_m$  or greater (where  $T_m$ - melting temperature on absolute scale), and the strain-rate lies in the range of  $10^{-5} \text{ s}^{-1}$  to  $10^{-1} \text{ s}^{-1}$  and the microstructure of the material is fine-grained (the average size of grains  $\leq 10\mu\text{m}$ ) and equi-axed. However, due to the approximate nature of these constraints, it is practically impossible to say exactly when the deformation of the material becomes superplastic, and at what point during the deformation it ceases to be so. For this reason, the design of technological processes using the phenomenon of superplasticity is generally done in such a way that the entire process of deformation is carried out in the «optimal superplasticity» regime. This regime is experimentally the set of conditions, in which parameter of strain-rate sensitivity [3–5] reaches the maximum.

$$m = \left( \frac{\partial \log \sigma}{\partial \log \dot{\epsilon}} \right)_{T, \epsilon = \text{const}} \quad (1)$$

Here  $\sigma$  and  $\dot{\epsilon}$  are the dimensionless stress and strain rate respectively (hence it only refers to uniaxial loading). There is even an approach of interactive change of loading conditions to maintain the maximum rate sensitivity during the

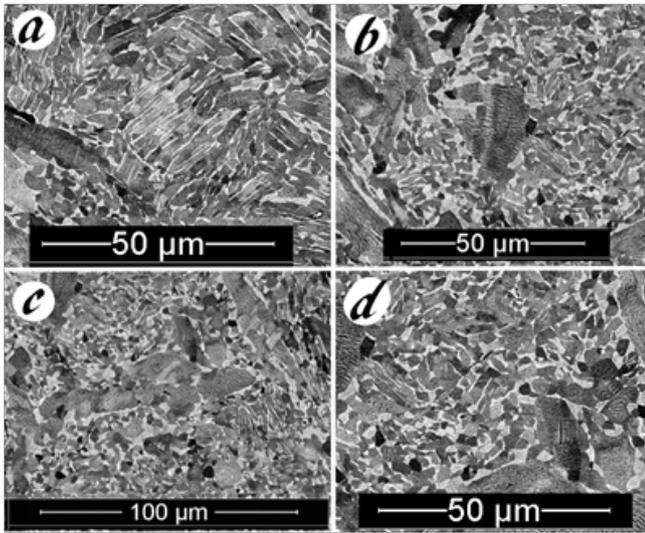
deformation process [6]. During optimum superplasticity, the material behaves like a viscous liquid and the flow stress can be considered to be a function of strain rate, temperature and initial grain size:

$$\sigma = f(\dot{\epsilon}, T, d_0). \quad (2)$$

The change in microstructure of the material in this case is not significant, and its transformation to a first approximation can be neglected, taking into account only the state of the material before deformation (e.g. its initial average grain size).

In practice, maintaining the optimal conditions of superplastic deformation (SPD) in industrial processes is sometimes difficult or even impossible and, moreover, it is economically inefficient. Indeed, it is quite difficult and costly to implement the process of isothermal deformation for a long time, especially for large dimensions of the workpiece and to prepare the required original fine-grained microstructure. Low strain-rates typical for SPD make the forming process time consuming and inefficient. Due to these reasons the cases of terminating the optimal regime are coming to the stage more and more frequently. Going beyond the limits of the optimum SP happens occasionally, but currently it is often being done deliberately e.g. with increasing strain rate to speed up the process at some stage (the so-called «high strain-rate SPD»), with lowering the temperature to reduce power consumption («low-temperature SP»), or refining of an unprepared microstructure during manufacturing of the product. Such processes, where optimum superplastic conditions are violated, but the material deforms almost homogeneously without visible flow localizations or fractures till relatively high strains (compared to a conventional high- temperature deformation), are classified as «near-to-superplastic deformation».

The main difference between these processes from the optimal SP is an active transformation of microstructure



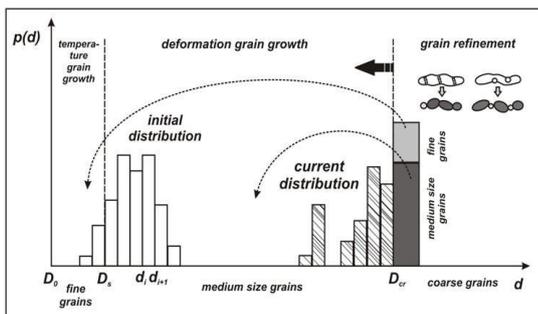
**Fig. 1.** The SEM snaps of Ti-6Al-4V alloy with initial Widmanstätten microstructure after deformation of  $\varepsilon=0.5$ ,  $\dot{\varepsilon} = 10^{-4} s^{-1}$   $T=850^{\circ}C$  illustrating main mechanisms of microstructure transformations.

occurring during deformation. It means that mechanical properties of material will significantly depend on the history of deformation (e.g. see Fig. 1). From the mathematical point of view it leads to a functional nature of constitutive models describing these processes. This fact was proved by many experiments showing the dependence of mechanical behavior of the material on the history of deformation [2, 4, 9, 10]. For the ability of numerical simulation (especially FEM) of such deformation processes a special model is proposed. The purpose of this model is to consider the active transformation of the microstructure during deformation and its effect on the mechanical properties of the material. This model might approximately describe the behavior of other Titanium alloys, but the mechanisms behind most Titanium alloys are too different to be predicted by it.

## 2. Constitutive modelling

### Description of the model

The proposed model, unlike a number of physical models [2], has a phenomenological character and will remain within continuum mechanics. To take into account active transformation of a microstructure during deformation, an additional characteristic variable will be assigned to a representative volume. This variable ( $\eta$ ) is used to represent



**Fig. 2.** The scheme of microstructure transformation used in the model.

the current state of the microstructure. The choice of this parameter is difficult, as microstructure transformation happens at different scale levels and mapping this is not trivial [11]. So to begin with, we use the most obvious and easily experimentally measurable parameter — the average grain size. If we are going to involve in our consideration not only a uniform fine-grained equiaxial microstructure, but also non-perfect ones like bimodal, lamellar or Widmanstätten, the direct kinetic equation for the average grain size will have only an indirect physical sense. If the microstructure simultaneously contains a mixture of grain sizes possibly with different morphology, then the micro-level mechanisms of deformation working in these grains may be also different. For each grade of grains a specific micro-model should be used. This approach will provide for the change of the average grain size some functional nature. The instantaneous change of this parameter will depend not only on current strain rate, stress, strain and temperature, but also on the current state of the microstructure.

Mathematically the approach described above will be represented using statistical methods by introducing the function of the distribution of the grain sizes. It is assumed that the representative volume contains grains of various sizes; each grain is identified only by one size — the characteristic size  $d$ . The distribution function  $p(d)$ , specifies the relative volume ( $V_i$ ) occupied by the grains which belong to a size range  $d_i$  ( $d_i - \frac{\Delta}{2} \leq d < d_i + \frac{\Delta}{2}$ ). Here all parameters of the model are dimensionless.

The parameter characterising a microstructure, has the physical sense of the average grain size and is defined as a probabilistic mean of distribution:

$$\eta = \sum_i p(d_i) \cdot d_i \cdot \quad (3)$$

As it was mentioned above it is possible that within the volume element several mechanisms of deformation may be operating simultaneously this being dependent upon the grain size distribution within the material. Using the distribution function provides us the ability to take all of them into account. Here each physical law is applied only to the grain size for which it is relevant [1]. For example, if grains are small and globular, the basic mechanism of their growth will be slow diffusion growth (governed by temperature). For medium size grains deformation growth will be a dominating mechanism (depending on the stress and individual size of grain). When the accumulated energy of deformation increases, the grains which are bigger than some critical value (depending on the value of the strain energy) start to refine (by dislocation creep and recrystallization). This scheme is shown in the fig.2. To take this into account in the model, the modification of the size of the grains of each type is calculated with a different mathematical law. The slow temperature-dependent growth of small grains is described by the equation:

$$\frac{\partial d_i}{\partial t} = k \cdot \exp\left(-\frac{Q}{RT}\right), \quad d_i \leq D_s, \quad (4)$$

where  $t$  is the time. The limitation  $d_i \leq D_s$  selects fine grains smaller than parameter  $D_s$ ; a dimensionless combination

$(Q/RT)$  involves temperature  $T$  (absolute scale), universal gas constant  $R$ , energy of activation  $Q$ , and material constant  $k$ .

Mechanism of deformation growth dominating for the medium size grains ( $D_s < d_i < D_{cr}$ ) was studied by researchers and a number of models are proposed. Here the model of Perevesentsev is used [12], for which:

$$\frac{\partial d_i}{\partial t} = \left( \frac{C_1 \sigma^4 + C_2 \sigma^5}{1 + C_3 \sigma^4 d_i} \right) \cdot \exp\left(-\frac{Q}{RT}\right), \quad (5)$$

$$D_s < d_i < D_{cr}$$

where  $\delta$  is dimensionless stress;  $C_1, C_2, C_3$  — parameters, which in initial model [12] represent combinations of physical constants of a material, but can be also considered as phenomenological and found from experiments.

As soon as the increments of the grain sizes are calculated, distribution function is modified accordingly:

$$p(d_i)|_{t+\Delta t} = \sum_{d_j + \Delta d_j \in (d_i)} p(d_j)|_t, \quad (6)$$

$$d_i < D_{cr}$$

Grains grow only until they reach some critical size  $D_{cr}$ , as soon they exceed it a refining process will commence. However, experimental results show, that even long lamellar grains remain unrefined if the deformation is small (e.g. in the centre of the specimen under torsion), and relatively small grains can get recrystallized at large deformations. It was also observed that at high stresses refinement is more active than at low ones. That means that the critical size  $D_{cr}$  is not a constant value. Here it is assumed to depend on accumulated strain energy:

The flow chat of the algorithm used for simulation of microstructure transformation.

$$D_{cr} = D^* \left[ \exp(-q \cdot \int \sigma \cdot d\varepsilon) \right] \quad (7)$$

where  $D^*$  and  $q > 0$  — constants.

Refinement of large grains can happen due to two basic physical mechanisms: **rearrangement of dislocations into dislocation walls and recrystallization**. In the first case, sometimes referred to as ‘strain assisted continuous recrystallization’ dislocations initially form low angle boundaries, which gradually convert to high angle boundaries probably facilitated by grain boundary sliding/rotation. In the second case a nucleus of a new, undeformed, grain arises in some part of the border of a larger grain, then it grows and divides the pre-existing grain into several parts. The elementary mathematical model of refinement used here is based on the assumption that all grains that exceed a critical size are refined into small and medium grains in proportion to  $\chi/(1-\chi)$ .

$$V_{ref} = \sum p(d_i), \quad d_i + \Delta d_i \geq D_{cr} \quad (8)$$

$$p(D_0) := p(D_0) + \chi \cdot V_{ref} \quad (9)$$

$$p(D_m) := p(D_m) + (1-\chi) \cdot V_{ref}$$

where  $V_{ref}$  is the total relative volume of «large» grains with the sizes greater than  $D_{cr}$ . Parameters  $D_0$  and  $D_m$  are mean sizes of small and medium grains produced by refinement correspondingly.  $D_0$  is a fixed size for new recrystallized grains; it is assumed to be almost always the same. The size of the medium grains changes in the process of deformation because  $D_{cr}$  decreases and the original sizes of the grains which undergo refinement also become smaller. So  $D_m$  can be assumed to be equal to either the total average grain size from the previous time step, or as some value depending on  $D_{cr}$ .

To complete the phenomenological model the relation between macroscopic parameters of loading such as stress and strain rate are required. It can be taken in the form of two serial basic elements: viscoelastic and «superplastic». As a viscoelastic element the generalized body proposed by Maxwell [2] modified for the temperature can be taken, whereas the superplastic element can be taken in standard form, but involving the internal parameter  $\eta$ :

$$\dot{\varepsilon} = \dot{\varepsilon}_e + \dot{\varepsilon}_v + \dot{\varepsilon}_{sp}, \quad (10)$$

$$\dot{\varepsilon}_e + \dot{\varepsilon}_v = \left( \frac{\dot{\sigma}}{E} + \lambda \cdot \sigma \right) \cdot \exp\left(-\frac{Q}{RT}\right), \quad (11)$$

$$\dot{\varepsilon}_{sp} = A \sigma^{1/m} \eta^{-k} \cdot \exp\left(-\frac{Q}{RT}\right), \quad (12)$$

where  $E, \lambda, A, m$  and  $k$  are the constants of model.

## Simulation

Numerical simulation of the uniaxial (tensile) loading with constant strain rate and under constant temperature was done in MATLAB. The flow chart of the algorithm used is shown in the fig.3 and the results of simulation in the fig.4.

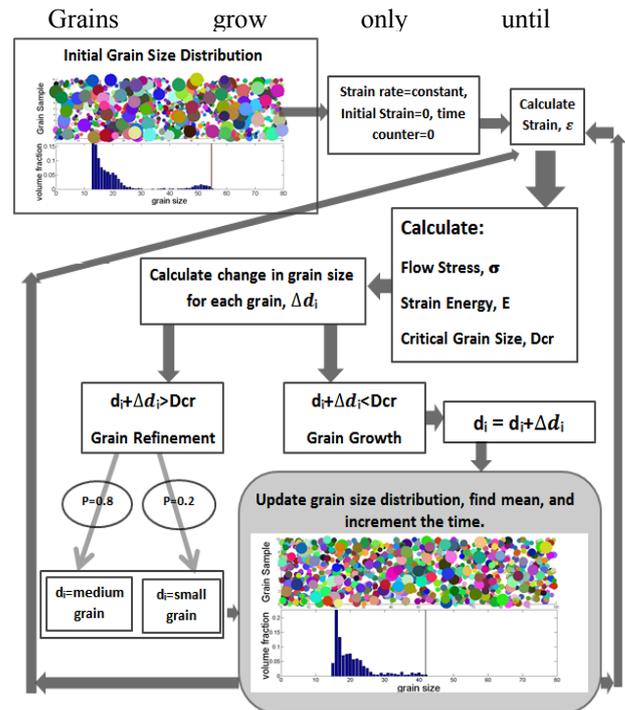


Fig. 3. The flow chat of the algorithm used for simulation of microstructure transformation.

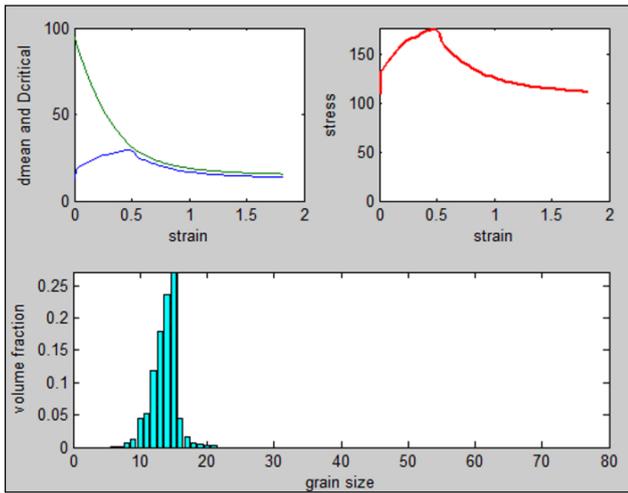


Fig. 4. The example of the results of simulation.

It can be seen from the results of simulation that the stress-strain curve obtained has hardening in the first part followed by softening with a gradually decreasing slope. Such mechanical behavior is typical for Ti alloys deformed in near-to-superplastic regimes [15]. However for practical use of this model in real technological simulations the question of experimentally obtaining the parameters of the model must be solved.

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### 3. Finding the parameters of the model

Two parameters of the model:  $m$  — strain rate sensitivity and  $Q$  — activation energy, are the most important for the accuracy of calculations because the results of simulation are very sensitive to them. There are two main problems with obtaining these parameters from experimental data. It is assumed that they are constants for a given material, but in reality both of them strongly depend on the strain rate and temperature. The second problem is related with the

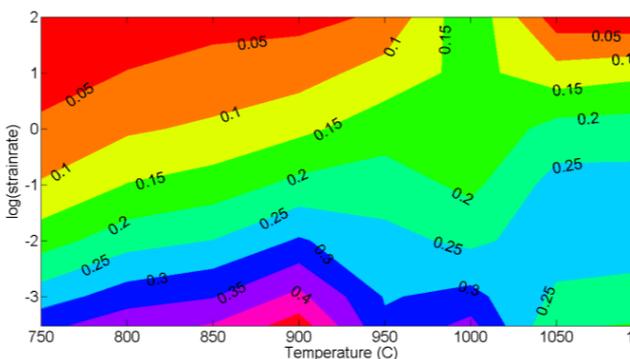


Fig. 6. Contour plot for the strain rate sensitivity  $m$  (3<sup>rd</sup> order).

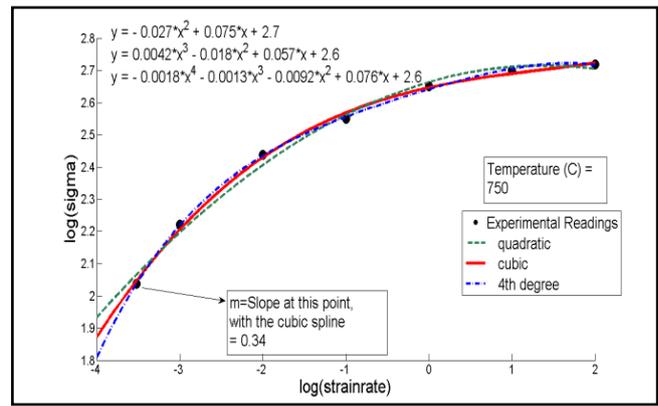


Fig. 5. Approximation of the experimental data with splines of different orders for differentiation required for obtaining strain rate sensitivity.

need to extrapolate experimental data from several different experiments.

$$m = \left( \frac{\partial \log \sigma}{\partial \log \dot{\epsilon}} \right)_{T, \epsilon = const} \quad (13)$$

$$Q = \left( \frac{\partial \log \sigma}{\partial (1/T)} \right)_{\dot{\epsilon}, \epsilon = const} \cdot \frac{R}{m} \quad (14)$$

If experimental data is taken from different tests, and in our case several experiments with different strain rates and several experiments with different temperatures are required, the amount of data is very limited and contains certain inaccuracies.

To illustrate the problem of finding  $m$  and  $Q$  and the robustness of this process, the experimental data published in [16] were used. These experiments were conducted for the Ti alloy Ti6Al4V with a fine globular microstructure for 7 constant strain rates (0.0003, 0.001, 0.1, 1, 10 and 100s<sup>-1</sup>) and 8 temperatures (750, 800, 850, 900, 950, 1000, 1050 and 1100°C). This range of conditions covers both superplastic as well as neighboring hot visco-plastic regimes of deforming.

To find required parameters, the approximation of available experimental data with polynomial splines of different order was made. The example of it is shown in the fig. 5. The values of strain rate sensitivity  $m$ , obtained with interpolation of different orders are given in the tab.1. It can be seen that from 4<sup>th</sup> order the values do not change much,

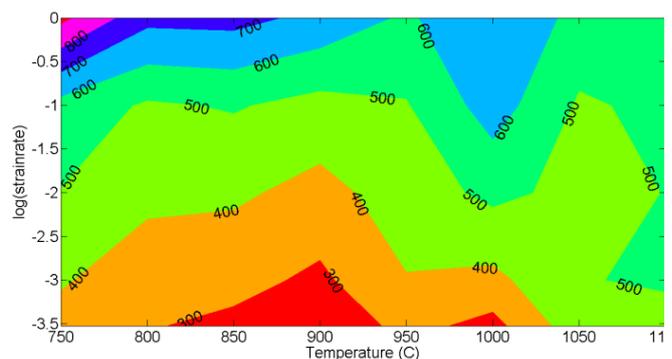


Fig. 7. Contour plot for the parameter  $Q$ , kJ mol<sup>-1</sup> (4<sup>th</sup> order).

Table 1

The dependance of strain rate sensitivity parameter to the order of polynomial interpolation and accuracy of input data

Degree of Interpolation	Original Value (v1)	Value with random-error in log(sigma) (v2)	v1-v2 (Sensitivity)	Value with random error in log(sigma) (v3)	v1-v3 (Sensitivity)
<i>Quadratic</i>	0.25	0.26	0.01	0.211	0.039
<i>Cubic</i>	0.30	0.29	0.01	0.273	0.027
<i>4th Degree</i>	0.35	0.28	0.07	0.385	0.035
<i>5th Degree</i>	0.36	0.16	0.2	0.410	0.050
<i>6th Degree</i>	0.35	0.18	0.17	0.415	0.065

Table 2

Values of strain rate sensitivity  $m$  (3<sup>rd</sup> order)

$\dot{\epsilon} / T$	0.0003	0.001	0.01	0.1	1	10	100
<b>750</b>	0.34	0.27	0.18	0.10	0.06	0.03	0.04
<b>800</b>	0.38	0.32	0.23	0.15	0.09	0.05	0.03
<b>850</b>	0.40	0.35	0.25	0.17	0.11	0.06	0.04
<b>900</b>	0.48	0.41	0.31	0.21	0.14	0.08	0.03
<b>950</b>	0.31	0.30	0.26	0.23	0.18	0.12	0.06
<b>1000</b>	0.37	0.32	0.24	0.19	0.16	0.16	0.19
<b>1050</b>	0.20	0.24	0.28	0.27	0.22	0.12	0.02
<b>1100</b>	0.20	0.23	0.27	0.27	0.22	0.14	0.01

Table 3

Values of effective activation energy  $Q$  (in KJ mol<sup>-1</sup>) (3<sup>rd</sup> order)

$\dot{\epsilon} / T$	0.0003	0.001	0.01	0.1	1	10
<b>750</b>	190	225	240	300	786	950
<b>800</b>	275	320	350	370	520	780
<b>850</b>	310	360	490	620	920	1300
<b>900</b>	265	320	455	610	910	1400
<b>950</b>	375	435	520	585	710	900
<b>1000</b>	260	365	510	625	670	560
<b>1050</b>	320	400	340	335	360	505
<b>1100</b>	160	300	210	200	165	150

which may imply that highest-possible order (6<sup>th</sup>) would lead to the most accurate results. But higher order polynomial approximations are more sensitive to the inaccuracy of input data and results of calculations are thus less robust.

To check this, randomized error of 10% was introduced to initial experimental data (to the value of stress), this depicts the actual accuracy limit of majority of high temperature mechanical tests. Obtained results and their deviation from the original ones are also listed in tab.1. These results show that increasing the order of polynomial order from 4 to 6 increases the sensitivity to inaccuracies almost twice. So for all further calculations 4th order polynomial approximations were used.

Using these experimental data [16] and approximations which were found to be optimal, the values of strain rate sensitivity  $m$  and activation energy  $Q$  were calculated for different temperatures and strain rates. Obtained results are given in the tabs. 2 and 3, and the fields of these values are show in the fig.6.

This clearly shows that both parameters significantly depend on temperature and strain rate. In FEM simulation of real technological processes temperature and strain rate can vary from one point of a workpiece to another. For equation (8) and the total model to give correct results, the parameters  $m$  and  $Q$  must be continuously verified and adjusted according to local conditions. This can be done with the help of the fields shown in figs.6 and 7.

Other parameters of the model are also important but less critical because the model is much less sensitive to them than to the above described ones. Nevertheless, the analysis of the appropriate choice of physical models and the possibilities of obtaining the parameters describing micro-mechanisms has to be investigated systematically. In this paper due to limited space these questions are not discussed.

#### 4. Conclusion

The importance of the consideration of microstructure in simulation of processes of hot forming including ones utilizing superplasticity was analyzed and a constitutive model able to take into account active transformation of microstructure was proposed.

Being based on a statistical approach this model can deal with several micro-mechanisms at a time. This feature is very important for two-phase alloys, especially for titanium alloys, which can have variable morphologies of microstructure.

Proposed model involves grain growth as well as refinement of the microstructure that gives the ability to describe both hardening and softening of the material. The transformation of the microstructure at every step of the simulation gives the ability to trace functional (history dependent) character of the process as well as the effect of initial microstructure on material properties.

Proposed model can be suitable for the family of  $\alpha+\beta$  Ti alloys with similar type of microstructure and similar mechanisms of its transformation. However for other titanium alloys, e.g. the intermetallic titanium aluminides, having a fundamentally different microstructure, the choice of the microstructure parameters as well as kinetic models of their behavior must be different.

The procedure of obtaining parameters of the model from experimental data and robustness of results is discussed and an optimal method suggested.

The values of strain rate sensitivity and activation energy were calculated for different temperatures and strain rates. It was shown that these parameters significantly depend on these conditions and this must be taken into account in simulation.

The reason for higher inaccuracies pertaining to polynomials of greater degree is purely mathematical. The higher the degree of a function, the greater number of stationary points it will contain.

Simulation with the model using calculated parameters approximately matched the experimental results.

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