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# Microstructural simulation of friction stir welding using a cellular automaton method: a microstructure prediction of AZ91 magnesium alloy

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## Abstract

**Background:** Recently, some researchers have simulated FSW using FEM and studied the influence of process parameters and tool geometry on material flow, welding force, and temperature and strain distributions during friction stir processing. Additionally, in terms of microstructure modeling, various approaches such as the Cellular Automaton (CA) model have been developed to simulate microstructural evolution during plastic deformation processes.

**Method:** In this work, a finite element model (FEM) is established to study the microstructure evolution during friction stir welding (FSW) of AZ91 magnesium alloy. To this aim, first, the hot compression tests at different temperatures and strain rates were carried out to achieve the flow stress curves. Then, the hardening parameter, the recovery parameter and the strain rate sensitivity were calculated according to flow stress results and using the Kocks–Mecking model. Next, a continuum based thermo-mechanically coupled rigid-viscoplastic FEM model was proposed in Deform-3D software to simulate the FSW of AZ91 magnesium alloy. To evaluate microstructure of the weld zone a model is proposed based on the combination of Cellular Automaton and Laasraoui-Jonas models.

**Results:** Temperature history, strain distribution and welding force are achieved through thermomechanical model and microstructure and grain size distribution are achieved by microstructure evolution model. The effects of rotational and traverse speeds on the grain size and microstructure of weld zone are considered.

**Conclusion:** There is a good agreement between results of numerical models and experiments in the aspects of welding forces, temperature history and grain size. Additionally, the proposed microstructure evolution model can simulate accurately the dynamic recrystallization (DRX) process during FSW and its resulted microstructure.

**Keywords:** FSW simulation; Microstructural evolution; DRX; Cellular automaton; Laasraoui-Jonas model

## Background

AZ91, a magnesium alloy, is one of the most commercially and commonly used magnesium alloys. This alloy, containing 9 wt% Al, 1 wt% Zn, and 0.2–0.3 wt% Mn as major alloying elements, contains a good combination of castability, mechanical strength, and ductility (Suresh et al. 2009). This has made AZ91 a popular light metal alloy especially among automotive industries whose aim is manufacturing lightweight vehicles (Srinivasan et al.

2010). However, the use of AZ91 in different industries is not yet extended comparing to its competitors such as aluminum alloys and plastics, partially due to the difficulty in controlling its microstructure (Asadi et al. 2010a).

Friction stir welding (FSW) as a relatively new welding technique has gained wide applications in different industries such as aerospace, automotive, and maritime. It has been utilized to weld and process different aluminum (Heidarzadeh et al. 2015), Mg (Asadi et al. 2012; 2010b; Motaleb-nejad et al. 2014; Faraji and Asadi 2011), and Cu (Farrokhi et al. 2013) alloys, some of

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which are classified as practically unweldable alloys in use of conventional welding methods.

Recently, some researchers have simulated FSW using a finite element model (FEM). Buffa et al. (2013; 2012; 2006) simulated friction stir welding using a 3D finite element method. Their model effectively determines the relationships between the tool forces and process parameters. Shojaeefard et al. (2013) studied the influence of pin profile and shoulder diameter on material flow, welding force, temperature, and strain distributions. Marzbanrad et al. (2014) investigated the effect of tool pin profile, and Asadi et al. (2011a; Tutunchilar et al. 2012a) studied the effect of the process parameters on material flow, temperature, and strain distributions during friction stir processing.

It is clear that the grain size in the weld zone has a great influence on the mechanical properties of weld such as hardness, tensile strength, plasticity, and toughness properties, and therefore, fine-grain structure could enhance these properties (Asadi et al. 2010b; Farrokhi et al. 2013; Heidarzadeh et al. 2014). Since it is difficult and time consuming to investigate experimentally the microstructure of weld, numerical simulations could be very applicable in different manufacturing processes (Liu et al. 2013; Wang et al. 2010). In terms of microstructure modeling, various approaches such as the cellular automaton (CA), the Monte Carlo model, and the phase field model have been developed to simulate microstructural evolution during processes (Liu et al. 2013). Although all these models successfully simulate microstructural evolution, most of the CA model is employed because of its length scale calibrations and straightforward time. Discrete spatial and temporal evolution of complex systems via applying local or global deterministic or probabilistic transformation rules to the location of a lattice is the main algorithm of the CA method. Many researchers have shown that CA offers a computationally efficient framework for simulation of microstructural evolution (Liu et al. 2013).

Timoshenkov et al. simulated the microstructure evolution in steel using CA for thermo-mechanical treatment. Tsai et al. (2010) predicted the morphologies in the solidification process for Cu-0.6Cr (mass fraction, %) alloy and Wang et al. (2010) simulated the dynamic recrystallization (DRX) characteristic in hot compression of steel using the CA method. They stated that the CA model can simulate the nucleation and growth kinetics of dynamically recrystallized grains in hot working process. Besides these advantages, this method could not consider solely the effects of the process parameters on DRX and the relationship between the nucleation sites and the distribution of dislocation density (Liu et al. 2013).

In fact, dislocation density plays a crucial role in nucleation and therefore, in microstructural evolution during DRX (Qian and Gou 2004). In order to evaluate the dislocation density, many models have been developed,

such as the Laasraoui-Jonas (LJ) model (Laasraoui and Jonas 1991), the Kocks-Mecking (KM) model (Mecking and Kocks 1981), and the Estrin-Mecking (EM) model (Yazdipour et al. 2007). These models are internal variable dislocation density models and deal with calculation of flow stress and evaluation of dislocation density during hot deformation processes. Additionally, the modified LJ model (Gourdet and Montheullet 2003) considers the effects of grain boundary migration on dislocation density. Therefore, the modified LJ model grants a more realistic evolution for dislocation density. Li et al. (2012) used the LJ model to simulate the microstructural evolution during hot extrusion of Mg-Al-Ca-based alloy. Similarly, Liu et al. (2013) simulated the DRX for hot compression of AZ31 magnesium alloy.

In the present study, a model is developed to simulate the microstructural evolution of AZ91 magnesium alloy during FSW. To this aim, first, the flow stress curves are obtained via the hot compression tests carried out at different temperatures and strain rates. Then, the hardening parameter, the recovery parameter, and the strain rate sensitivity are calculated according to flow stress results. Next, a continuum based thermo-mechanically coupled rigid-viscoplastic FEM model for the FSW process is proposed in Deform-3D software coupled with the combination of CA and LJ models for microstructural evolution and dislocation density computation. The relationship between the nucleation rate and the dislocation density as well as between critical strain, critical dislocation density, and DRX are investigated. Next, the optical-microscope images from weld zone and base metal are used to validate the microstructure predicting model. Furthermore, nucleation and grain growth are shown by the micro-images extracted from simulation. Finally, the effect of the process parameters on grain size is studied.

## Methods

The methods of this research include some experimental investigations and tests, numerical model for the process simulation, and the microstructure evolution formulation for modeling of dynamic recrystallization. All these methods are described in below.

### Experimental setup

AZ91 magnesium alloy plates with the chemical composition shown in Table 1 and in 5-mm thickness were friction stir welded in different conditions. The rotational and traverse speeds were varied between 710–1400 rpm and

**Table 1** Chemical composition of the AZ91 magnesium alloy used in this study

Element	Al	Zn	Mn	Si	Cu	Ni	Fe	Mg
Weight percentage	9.1	0.68	0.21	0.085	0.0097	0.001	0.0029	Bal.



**Fig. 1** FSW tool used in this study

25–100 mm/min, respectively. The FSW tool was made of 2344 hot working steel with a circular pin in a 5-mm diameter and 4.8-mm height (Fig. 1). The tool shoulder diameter was 18 mm and the tool tilt angle was 3°. To track the temperature variations during the process, an infrared thermometer with  $\pm 3$  °C accuracy was employed. Meanwhile, a one-component Kistler dynamometer was utilized to measure the axial force.

The welded specimens were cut in transverse sections and prepared by standard metallographic techniques. The etching solution was composed of 5 mL acetic acid, 5 mL HCl, 6 g picric acid, 7 mL nitric acid, 100 mL ethanol, and 10 mL water for 1–2 s. Microstructural

observations of the samples were carried out by optical microscopy (OM), and a linear intercept method was employed to measure the average grain size ( $d = 1.74 L$ ;  $L$  is the linear intercept size).

In order to acquire the flow stress of the AZ91 Mg alloy, hot compression test was conducted at different temperatures among 200–450 °C and strain rates of 0.001, 0.01, 0.1 and 1 s<sup>-1</sup>. The hot compression test specimens were 15 mm in height and 10 mm in diameter.

#### Numerical model of friction stir welding

Due to the ability of accurate modeling of severe plastic deformation (SPD) processes, Deform-3D™ software is employed to simulate the FSW (Tutunchilar et al. 2012a). To simplify the problem, the tool was selected as a rigid material and workpiece was supposed to be a rigid-viscoplastic material.

The Arrhenius equation was chosen to determine the relationship between the strain rate, the flow stress, and the temperature (Asadi et al. 2011a).

The tool and workpiece were meshed non-uniformly in about 25,000 and 58,000 tetrahedral elements, and an automatic remeshing system was utilized. Finer elements in 0.5 mm mean size were placed under the tool pin and shoulder. A shear friction factor of 0.4 was selected due to the best agreement between the experimental and simulated results for temperature history. Figure 2 illustrates the simulated model for welding of AZ91 magnesium alloy, and Fig. 3, comparing the temperature history and axial force results, shows a good agreement between the experimental and predicted forces.

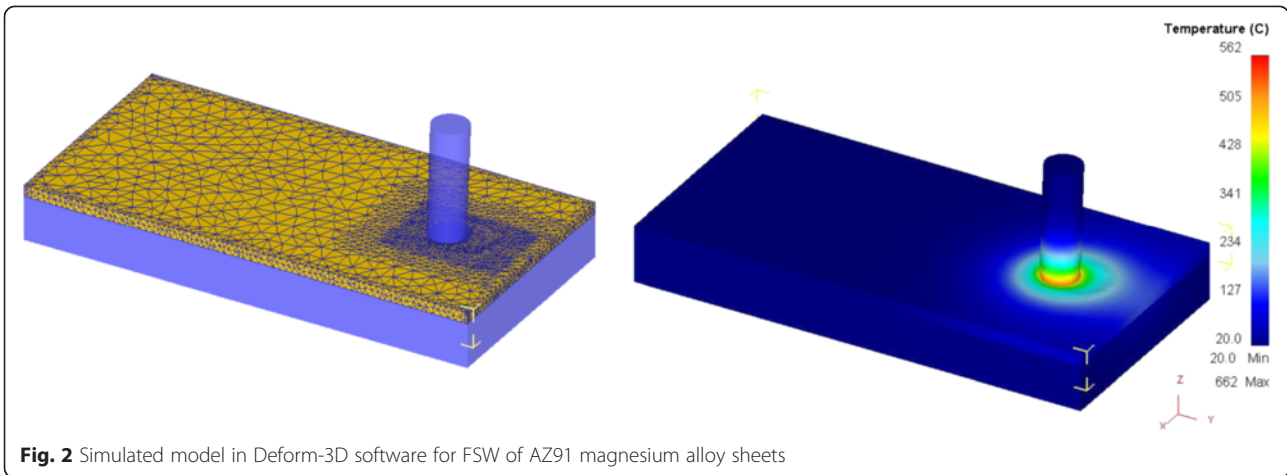
#### Microstructure evolution model

The microstructural evolution during FSW of AZ91 magnesium alloy is simulated using the combination of CA and LJ methods in Deform-3D software which are coupled to the thermo-mechanically, rigid-viscoplastic FEM model of FSW.

Evidently, the dislocation density is the most effective factor on nucleation and microstructure evolution during hot deformation processes (Liu et al. 2013; Qian and Gou 2004). Therefore, besides the temperature history, calculation of dislocation density is crucial in microstructure evolution of the FSW process. The temperature history will be achieved via the FSW FEM model. However, the dislocation density calculation requires coupled models (such as CA and LJ).

#### Dislocation density calculation

The main factors affecting the dislocation density during plastic deformation are evidently these three concurrent phenomena: (1) work hardening (WH), (2) dynamic recovery (DRV), and (3) DRX (Liu et al. 2013). In the



present research, the modified LJ model was employed to illustrate the evolution of dislocation density during DRX in the FSW process:

$$d\rho_i = (h-r\rho_i)d\varepsilon - \rho_i dV \tag{1}$$

In this equation,  $\rho_i$  is the dislocation density in the  $i$ th

grain;  $h$ , the strain hardening parameter;  $r$ , the recovery coefficient parameter;  $\varepsilon$ , the strain; and  $dV$ , the volume swept by grain boundary movement. To precisely calculate the amount of parameters in LJ model (Eq. 1) the KM model is also used:

$$\dot{\rho}_i = (K_1\sqrt{\rho_i} - K_2\rho_i)\dot{\varepsilon}_{eff}^p \tag{2}$$

where  $K_1$  is the constant denoting the work hardening and  $K_2$  is the softening parameter for dislocation annihilation.

#### Nucleation and growth during DRX

Several nucleation models are developed for DRX, but Qian and Gou (2004), by proposing Eq. 2, showed that the nucleation rate is proportional linearly to strain rate and exponentially to temperature.

$$\dot{n} = C\dot{\varepsilon}e^{-\frac{\Delta H}{RT_{abs}}} \tag{3}$$

where  $\dot{n}$  represents the nucleation rate;  $\Delta H$ , the activation energy, acquirable by the flow curves;  $R$ , the gas constant;  $T_{abs}$ , absolute temperature; and  $C$ , a constant. The relationship reveals that by increasing the process temperature and strain rate, the nucleation number raises.

In the metals with medium to low stacking fault energy such as Mg, the DRX onsets when the dislocation density reaches a critical value (Liu et al. 2013; Roberts and Ahlblom 1978). Generally, nuclei predominantly nucleate along the grain boundaries in which sufficient stored energy is present by a critical dislocation density:

$$\rho_{cd} = \frac{20\gamma_b\dot{\varepsilon}}{3blM\tau^2} \tag{4}$$

where  $b$  represents the Burger's vector;  $l$ , the dislocation free path;  $M$ , the grain boundary mobility;  $\tau$ , the dislocation line energy; and  $\gamma_b$ , the grain boundary energy, which can be computed by Eq. 5 (Xiao et al. 2008):

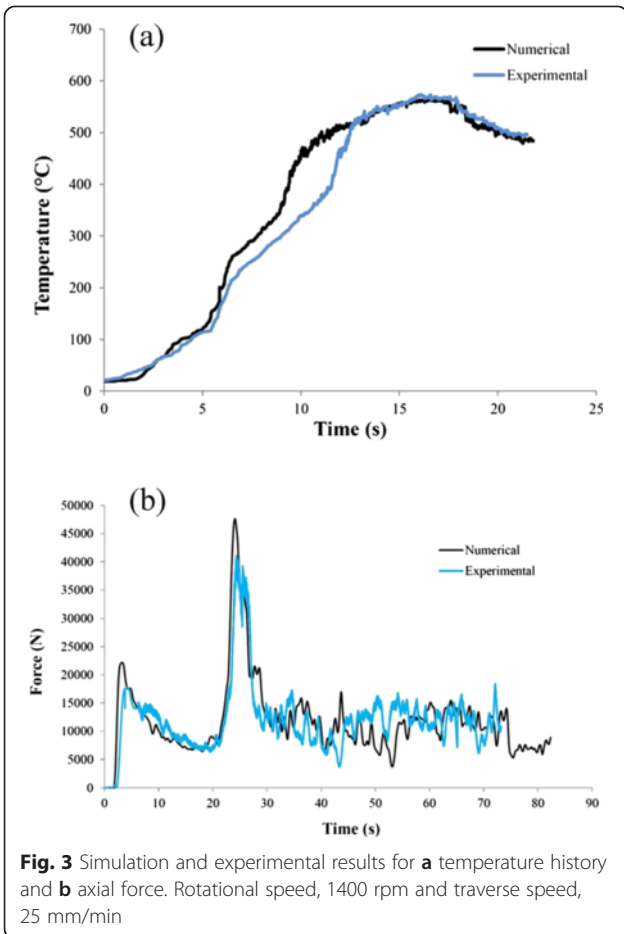


Fig. 3 Simulation and experimental results for a temperature history and b axial force. Rotational speed, 1400 rpm and traverse speed, 25 mm/min



$$\gamma_i = \begin{cases} \gamma_m \frac{\theta_i}{\theta_m} \left( 1 - \ln \left( \frac{\theta_i}{\theta_m} \right) \right), & \theta_i \geq 15^\circ \\ \gamma_m \frac{\theta_i}{\theta_m} \left( 1 - \ln \left( \frac{\theta_i}{\theta_m} \right) \right), & \theta_i < 15^\circ \end{cases} \quad (5)$$

Where  $\gamma_m$  is the high angle grain boundary energy which can be described by Equation 6;  $\theta_i$ , is the misorientation among the  $i$ th recrystallized grain and its neighbor; and  $\theta_m$ , the misorientation between high angle boundaries.

$$\gamma_m = \frac{\mu b \theta_m}{4\pi(1-\nu)} \quad (6)$$

where,  $\nu$  is the Poisson ratio. The grain boundary migration velocity is influenced by the net pressure on the boundary and can be described as:

$$v_i = \frac{MF_i}{4\pi r_i^2} \quad (7)$$

where  $v_i$  represents the grain boundary migration velocity;  $M$ , the grain boundary mobility;  $r_i$ , the  $i$ th recrystallized grain radius; and  $F_i$ , the driving force is attainable through Eq. 7.

$$F_i = 4\pi r_i^2 \tau \rho_i - 8\pi r_i \gamma_i \quad (8)$$

### Cellular automaton method

A representative volume in  $1000 \times 1000 \mu\text{m}$  is used in the CA model and discretized using 250,000 tetragonal cells. Each cell involves four variables: one variable for orientation showing the orientation of grain and calculates the energy of grain boundary by Eqs. 5 and 6; one variable for status indicating which grain is recrystallized; one for dislocation density determining the site energy by Eqs. 1 and 2; and one for color exhibiting different grains. The orientation is randomly selected to be between 0 and  $180^\circ$  for base metal and recrystallized grains, and dislocation density is set to be zero for recrystallized grains. When the dislocation density goes beyond the critical value, obtained by Eq. 4, the DRX starts. Next, the nucleation rate can be calculated by Eq. 3 for each time step. In terms of recrystallized grain growth, the recrystallized grain can continuously grow until the driving force ( $F_i$ ) is positive. The growth velocity ( $v_i$ ) can be obtained through Eqs. 7 and 8 for each time step. The radius of the  $i$ th recrystallized grain represents the volume swept by the grain boundary. Then, the dislocation density can be calculated by Eq. 1 for each time step. These computations are terminated, when the pre-set strain is reached. The cells in which recovery happens are random (Peckzak and Luton 1993), and to select certain number of lattices  $N_r$  for recovery occurrence in each time step, the following equation is selected:

$$N_r = \frac{2N_1N_2}{K} h(d\varepsilon)^{1-2m_0} \quad (9)$$

where  $N_1$  and  $N_2$  represent the number of rows and columns in cells and  $K$  is the material constant (Peckzak and Luton 1993).

### Work hardening and recovery parameters

The true stress-strain curves, obtained from the hot compression test, at different temperatures and strain rates are illustrated in Fig. 4. It is evident that the flow curves involve four stages of (1) work hardening, (2) transition, (3) softening, and (4) steady, up to breaking down. The single peak in the flow curves implies the occurrence of DRX (Asadi et al. 2011a; Xin et al. 2010). It can be observed from Fig. 4a–d that increase in temperature results in a reduction in the incubation period. This phenomenon is mostly affected by the critical strain and dislocation accumulation rates (Song et al. 2014). The work hardening rate can be calculated from the slope of true stress-strain curves in the work hardening stage (Liu et al. 2013). The average strain hardening parameter ( $h$ ) and the recovery coefficient parameter ( $r$ ) can be calculated through Eqs. 10 and 11 (Gourdet and Montheullet 2003):

$$h = h_0 \dot{\varepsilon}^m \exp\left(\frac{m\Delta H_b}{RT_{\text{abs}}}\right) \quad (10)$$

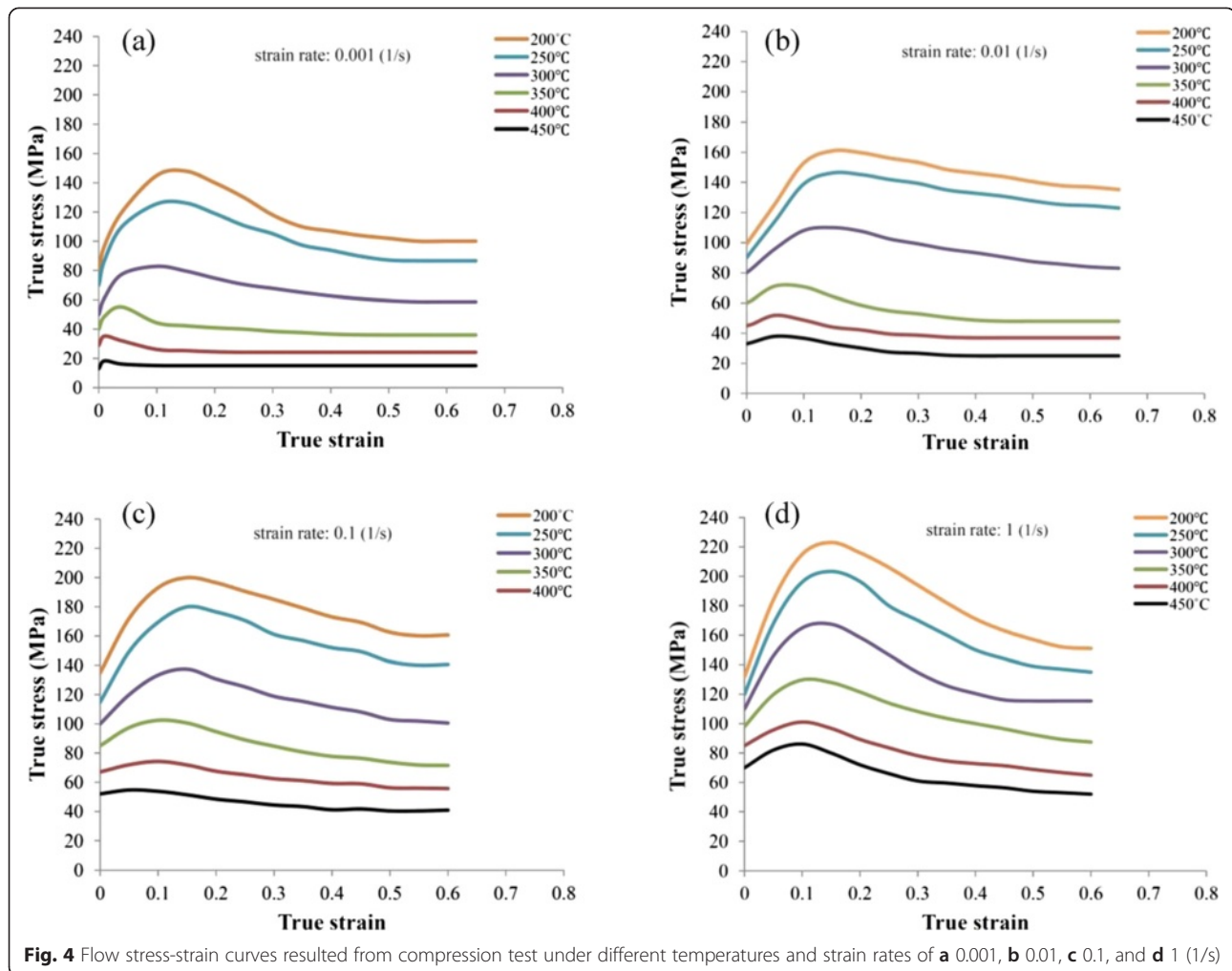
$$r = r_0 \dot{\varepsilon}^{-m} \exp\left(\frac{-m\Delta H_b}{RT_{\text{abs}}}\right) \quad (11)$$

where  $\dot{\varepsilon}$  represents the strain rate;  $m$ , the strain rate sensitivity;  $\Delta H_b$ , the activation energy for self-diffusion;  $h_0$ , the hardening parameter (in Eq. 10); and  $r_0$ , the recovery parameter (in Eq. 11).

By applying regression analysis for flow stress results, the values for  $h_0$ ,  $r_0$ , and  $m$  are achieved as: 1.34e13, 15.6, and 0.2, respectively (for more details see ref. (Liu et al. 2013; Guangyin et al. 2001; Unigovski 2009)). The material parameters used in the simulation process are summarized in Table 2.

### Results and discussion

As cited before, the temperature and strain distributions in the weld zone are crucial factors in determining the grain size of the weld nugget. Therefore, the temperature history and strain distribution during FSW will be reported in this paper. Then, the microstructural properties of the weld zone such as microstructural image, grain size, and nucleation sites will be considered as well as the effect of the process parameters on these properties.



#### Temperature history and strain distribution during FSW

It is well-accepted that frictional heat and plasticizing action formed by the FSW tool over the process causes the joining of two parts. Generation of a certain amount

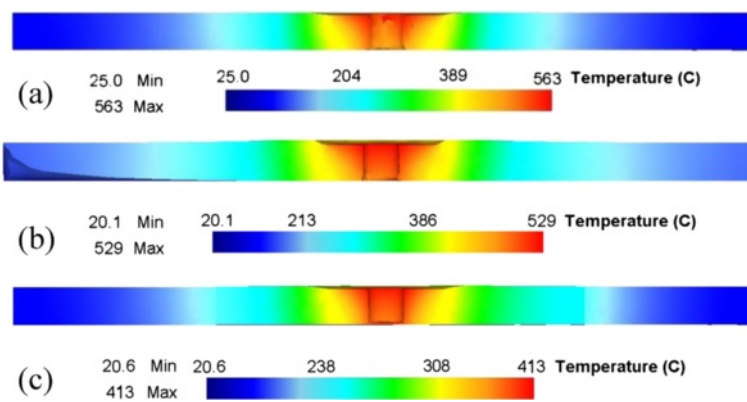
**Table 2** Materials parameters for AZ91 magnesium alloy used in simulation model

Material parameter	Value
Primary dislocation density, $\rho_0 (\mu\text{m}^{-2})$	0.01 (Liu et al. 2013)
Shear modulus, $G(\text{MPa})$	17,000 (Guangyin et al. 2001; Knezevic et al. 2010)
Activation energy for self-diffusion, $\Delta H_b \frac{\text{kJ}}{\text{mol}}$	140 (Guangyin et al. 2001)
Activation energy, $\Delta H \frac{\text{kJ}}{\text{mol}}$	147
Material constant in Eq. 9, $K$	6030 (Liu et al. 2013)
Shear friction factor, $m$	0.4
Hardening parameter, $h_0$	1.34e13
Recovery parameter, $r_0$	15.6
Strain rate sensitivity, $m$	0.2

of frictional heat along with the presence of a high hydrostatic pressure along the joint line are the crucial requirements during the FSW to keep the processing material in a well-plasticized region with a suitable temperature, and finally to form a sound weld.

Figure 5 shows temperature distribution in the cross section of welds produced by different rotational and traverse speeds. It is obvious that, for all welding conditions, temperature profile is nearly symmetric about the weld line, since the tool rotational speed is a dominating factor in heat generation rather than the tool traverse speed (Buffa et al. 2006). Furthermore, the figure reveals that the peak temperature grows as an increase in rotational speed or a decrease in traverse speed takes place. In summary, it can be mentioned that by increasing the rotational speed to traverse speed ratio ( $w/v$ ) the heat generation will be increased.

The strain distribution in the cross section of weld is shown in Fig. 6 for welds produced by different rotational and traverse speeds. It is obvious that the strain profiles are asymmetric about the weld line and the maximum



**Fig. 5** Temperature distribution in the weld zone **a** 1400 rpm, 25 mm/min; **b** 1400 rpm, 50 mm/min; and **c** 710 rpm, 50 mm/min

strain is biased towards the advancing side because of the positive combination of the traverse and rotational speeds which leads to a higher plastic deformation and strain rate (Asadi et al. 2011a; Tutunchilar et al. 2012a).

It can be seen that an increase in the rotational speed or a decrease in the traverse speed results in an increase in the amount of strain. Indeed by increasing the rotational speed to traverse speed ratio ( $w/v$ ), the tool pin rotation rate in a certain distance increases, leading to a rise in strain of material under processing.

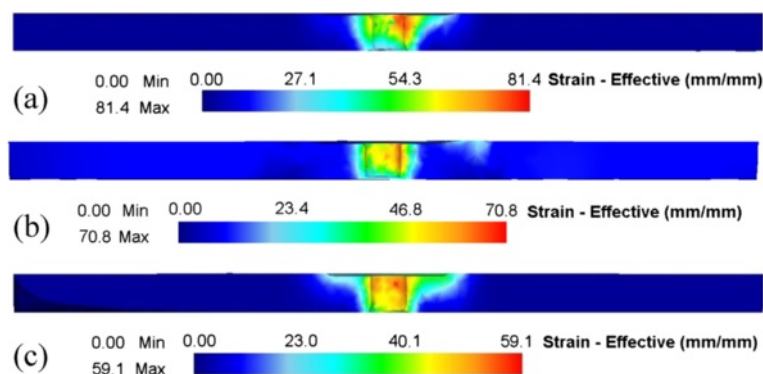
#### Microstructure evolution during FSW

It is well-known that DRX during FSW will cause the generation of fine equiaxed grains in the stir zone. Process parameters, tool geometry, welding material composition, temperature history, vertical pressure, and active cooling are the effective factors on the microstructure and grain size of the weld (Besharati Givi and Asadi 2014). In this paper, the effects of FSW on the microstructure of base metal as well as the effects of process parameters on the

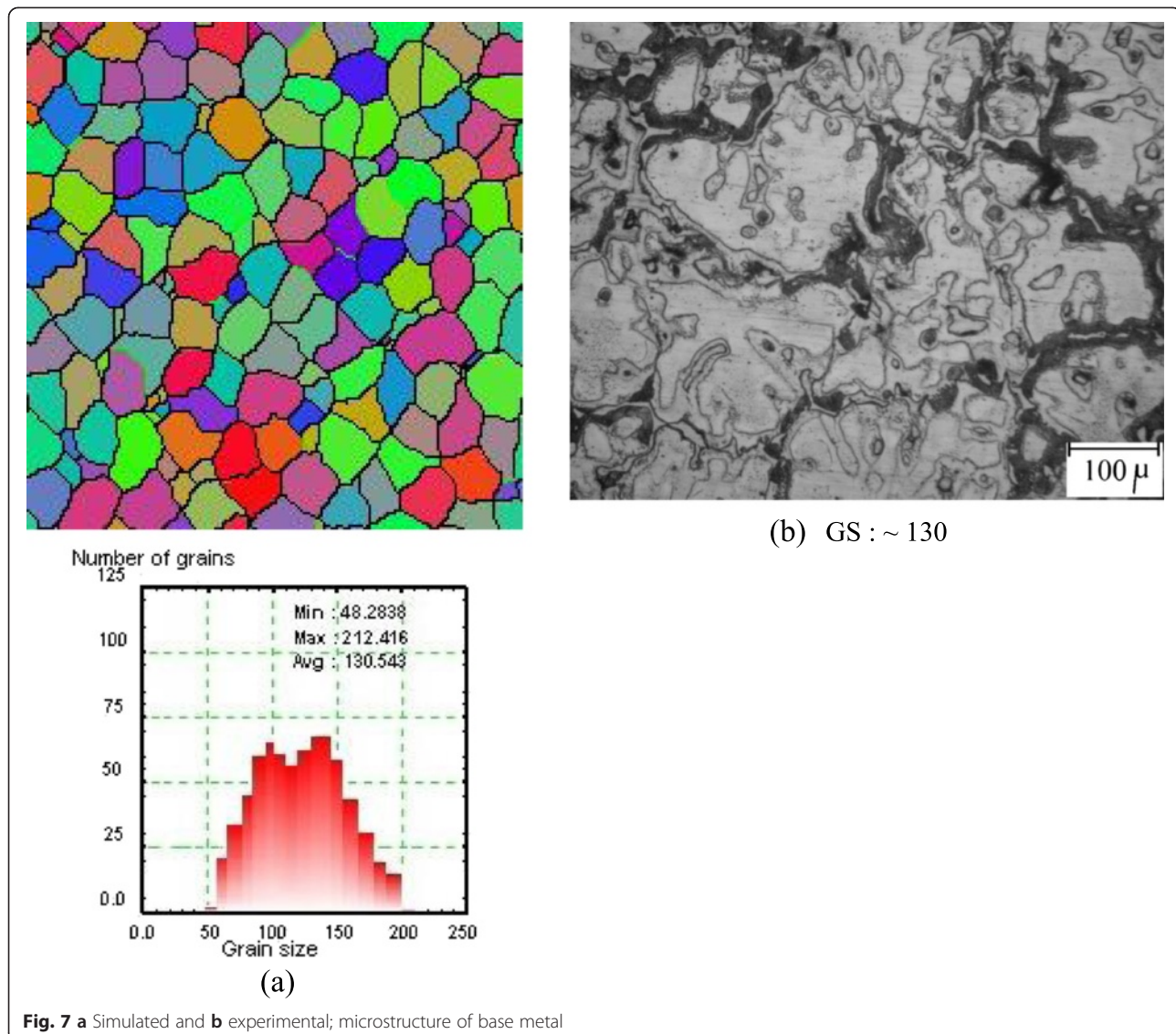
microstructure of stir zone are discussed by simulation results.

Song et al. (2014) simulated the microstructure evolution during FSW of titanium alloy using CA model and reported that their established model is more suitable for hot compression process. However, they did not compare the grain size with an experimental microstructure and did not report any simulated microstructure of the weld zone.

As mentioned before, the combination of LJ and KM models is employed to analyze the microstructural evolution of AZ91 magnesium alloy during the FSW. Fig. 7 shows the experimental and simulated microstructures for the base metal. In fact, according to the average grain size of the experimental microstructure, the number of grains in a certain area of CA cell ( $1000 \times 1000 \mu\text{m}$ ) can be obtained. Next, in this area a determined number of nuclei are distributed randomly and are then allowed to grow. The microstructure obtained by simulation for base metal can be seen in Fig. 7a, in which the different grains are represented by different colors while the black



**Fig. 6** Effective strain contours in rotational and traverse speeds of **a** 1400 rpm, 25 mm/min; **b** 1400 rpm, 50 mm/min; and **c** 710 rpm, 50 mm/min



lines show the grain boundaries. The grain size distribution in this figure indicates the average grain size of  $\sim 130 \mu\text{m}$ . Grain size in the experimental base metal (Fig. 7b), calculated by linear intercept, is also about  $130 \mu\text{m}$ .

To verify the ability of presented model in predicting the microstructure of weld zone, the simulated microstructures are compared with the experimental ones in Fig. 8 for the specimens produced by rotational and traverse speeds of  $710 \text{ rpm} \times 50 \text{ mm/min}$  and  $1400 \text{ rpm} \times 25 \text{ mm/min}$ . Comparing the average grain size for simulated and experimental microstructures reveals a good agreement between them. It demonstrates that the model presented for simulation of recrystallization during FSW of AZ91 magnesium alloy works precisely and can be employed to predict the microstructure and grain size in the stir zone.

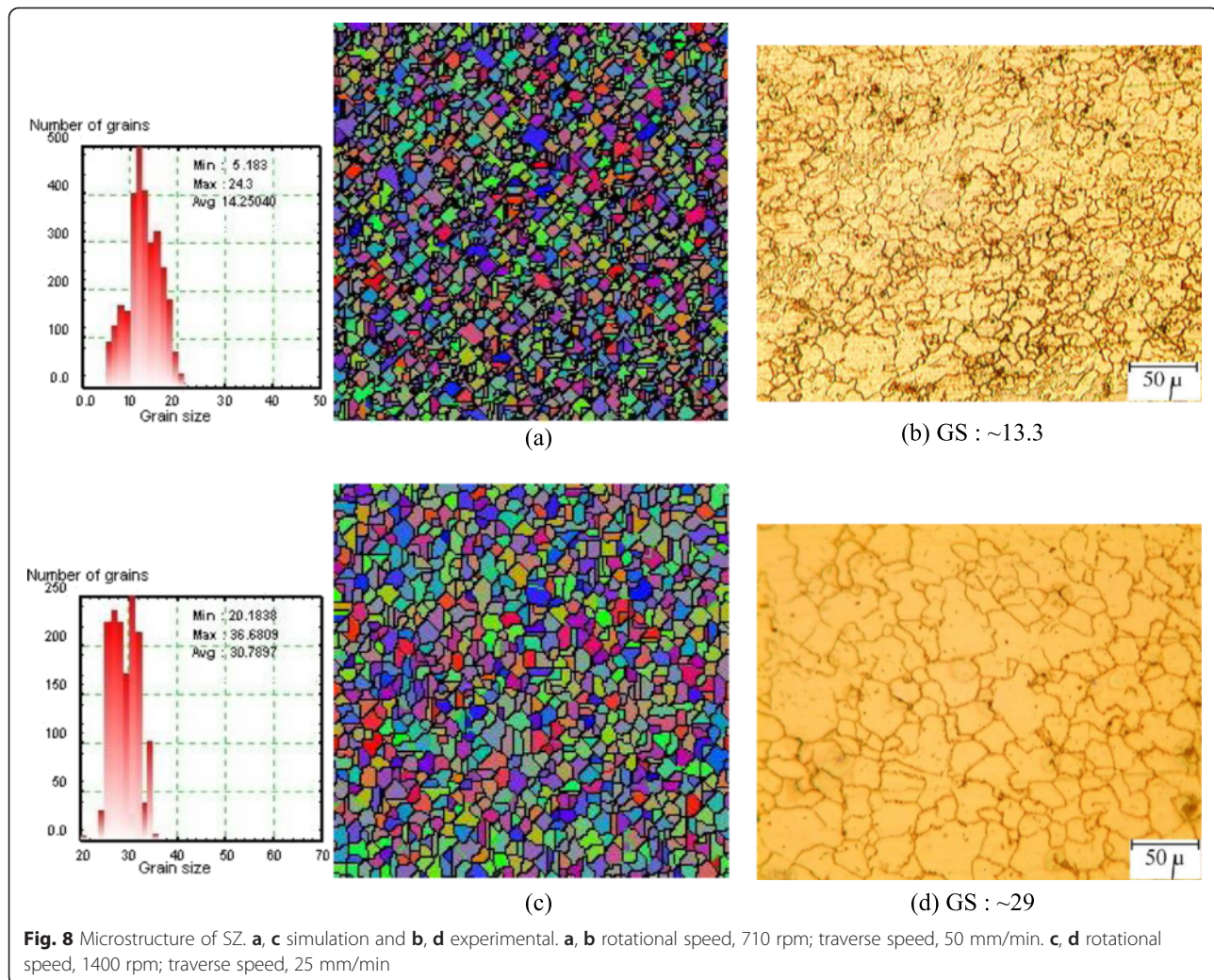
As cited before, by increasing the  $w/v$ , the heat generation is increased and therefore the grain growth and

grain size will be increased. The  $w/v$  ratio for Fig. 8a, b is  $\frac{710}{50} = 14.2$ , and it is  $\frac{1400}{25} = 56$  for Fig. 8c, d. In this case, the grain size is increased from  $13.3 \mu\text{m}$  in Fig. 8b to  $29 \mu\text{m}$  in Fig. 8d.

Comparing the microstructures for base metal and weld zone (Figs. 7 and 8), and as reported widely in literature, the grain size is reduced drastically by FSW (from  $130 \mu\text{m}$  in base metal to  $10\text{--}30 \mu\text{m}$  in the weld zone). The combination of high temperature, as a result of frictional heat, and high strain rate, due to the severe plastic deformation, leads to generation of fine recrystallized grains in the stir zone (Fig. 8) (Asadi et al. 2012; 2010b). Therefore, the grain refinement occurs by the DRX including nucleation and growth. The presented model can precisely reveals the changes in microstructure during the DRX process.

Figure 9 shows the nucleation and grain growth steps simulated by the presented model. This figure clearly





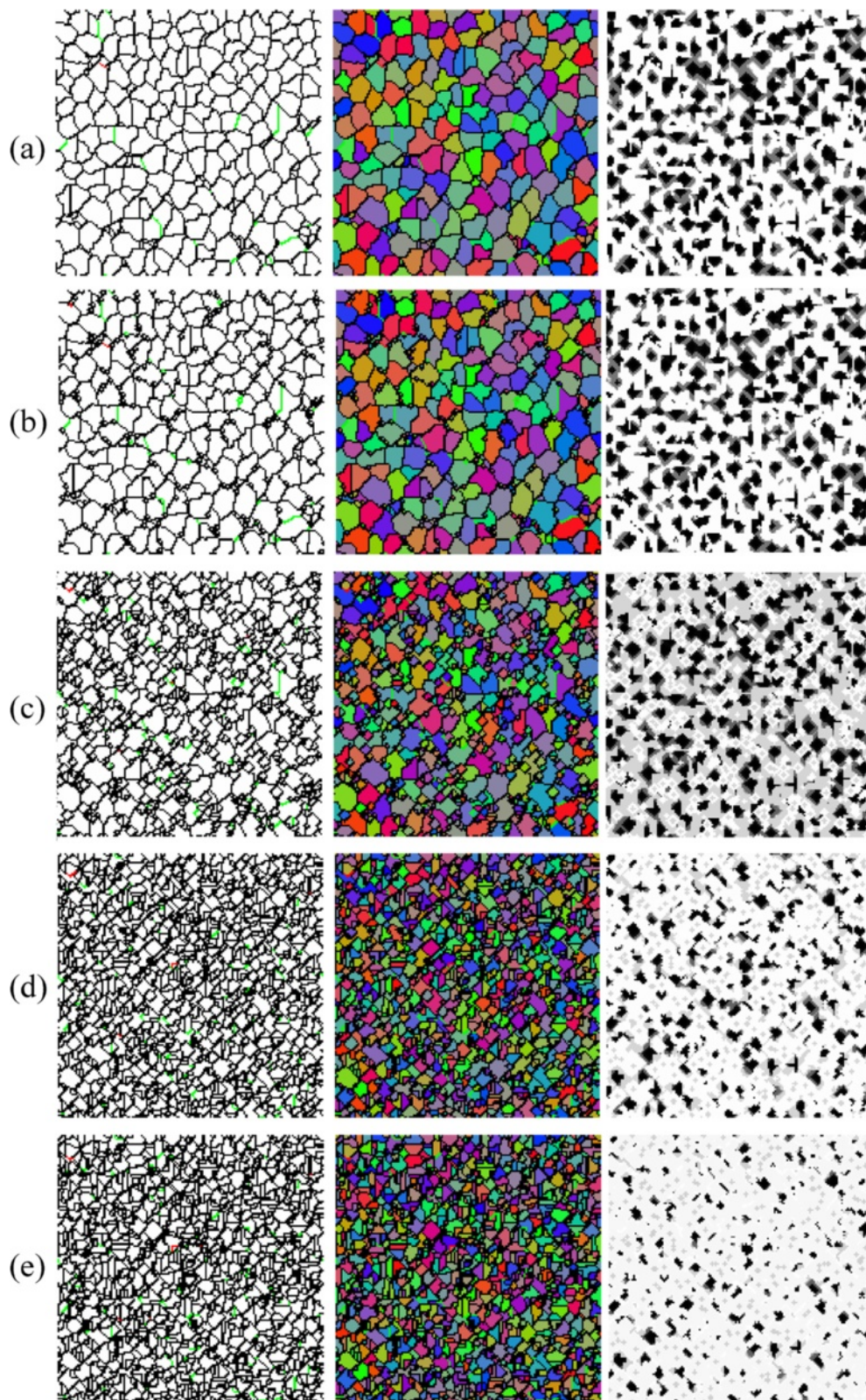
illustrates that the new grains nucleate from the old grains' boundary (Fig. 9b) and then start to grow (Fig. 9c). Indeed, the nucleation of new grains occurs in the preferential sites, and then, due to the presence of enough heat and the dislocation density difference, the grain growth begins and continues as far as the driving force for boundary migration is positive. Since in the center of stir zone, there are enough heat and strain rate, the microstructure is filled all by the recrystallized fine grains (Fig. 9e).

In each step (Fig. 9a–e), the figures on the left side and middle shows the microstructure of stir zone and the figure on the right side illustrates the dislocation accumulation map of that step. The darker points in the dislocation accumulation map represent the areas with higher dislocation density. Noting to the dislocation accumulation maps at different steps, it can be concluded that, by nucleation and growth, white points form inside the dark areas in the map and onset to consume them. Therefore, little by little, the

dark areas reduce demonstrating the reduction of dislocation density.

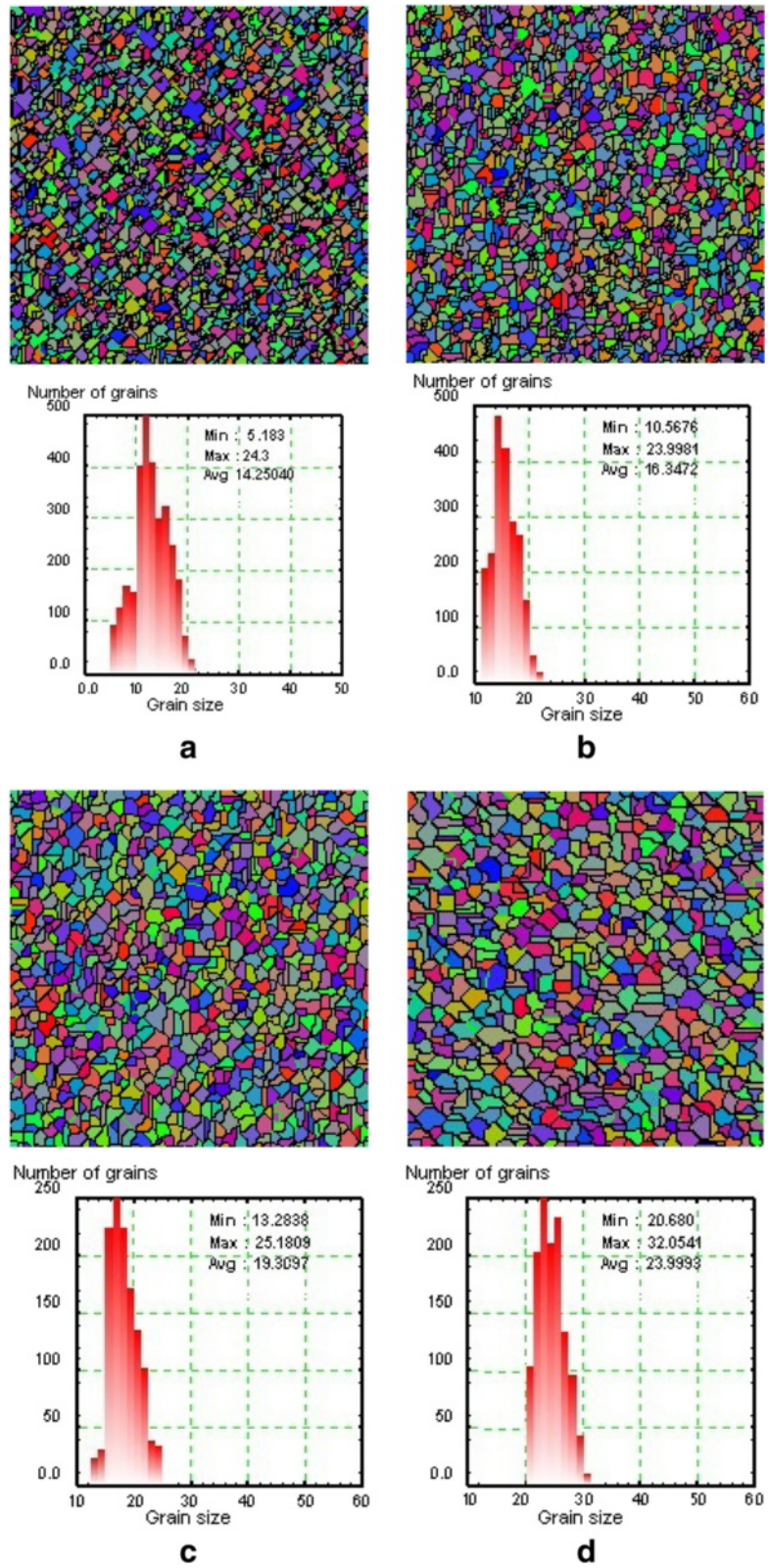
It should be stated that by deformation progress, the dislocation density goes beyond a critical value and the DRX initiates. Then, the new grains with zero dislocation density start to grow. On the other side, the simultaneous work hardening generates dislocation density inside the DRX grains and thus, the driving force for grain boundary migration drops down as far as the growth stops. On the other words, the driving force for the grain growth reduces gradually as its dislocation density inside the grain increases with deformation progress and finally the grain growth ceases. Also in another case, if the DRX grain sticks into another, both grains cease to grow at the clash point, while free sections of their boundaries can continue the growth. Additionally, in very high deformation rates, the DRX grain's boundaries again can become nucleation sites, if its dislocation density reaches the critical value for nucleation (Liu et al. 2013; Song et al. 2014).



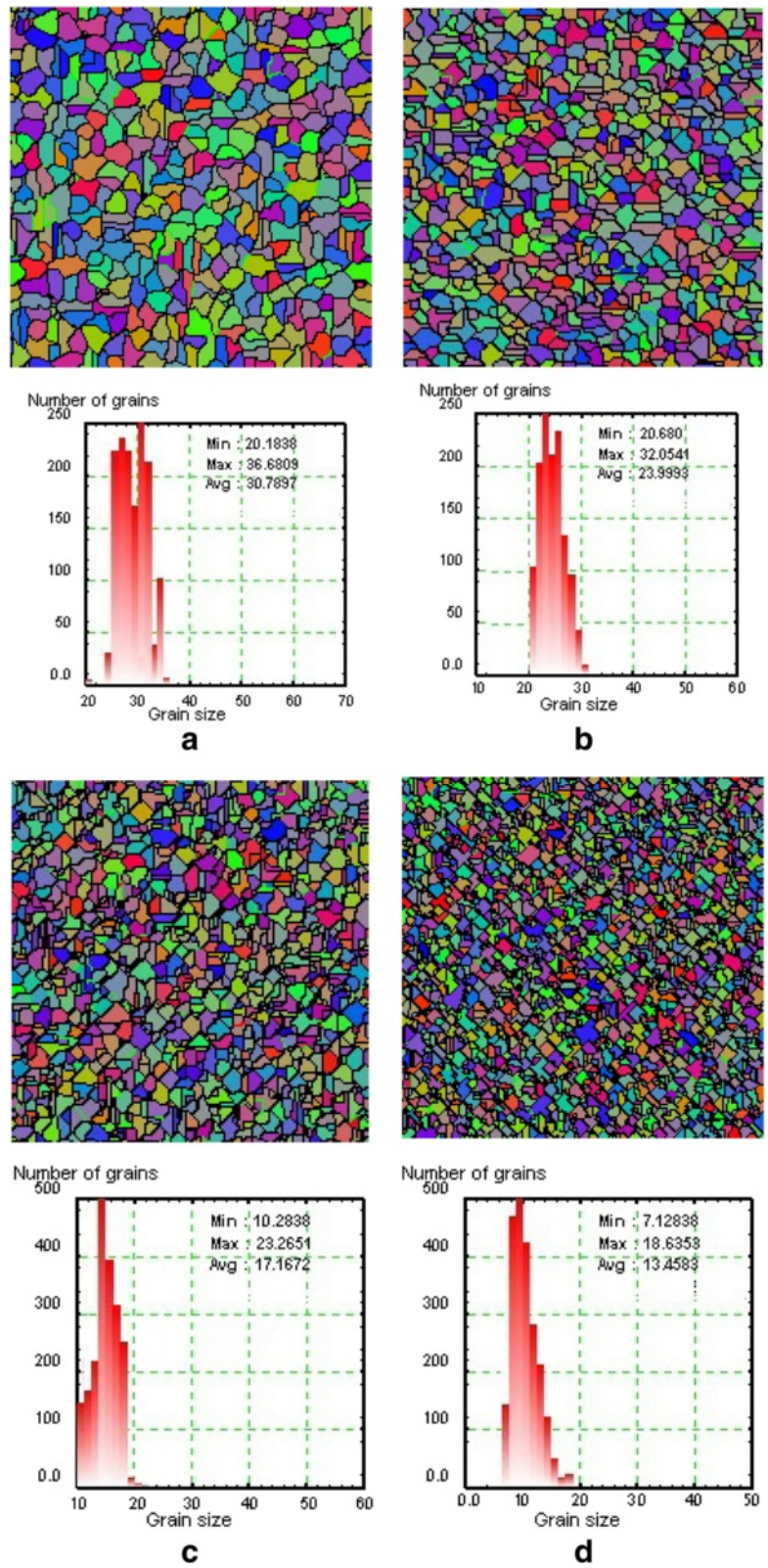


**Fig. 9 a–e** Nucleation and grain growth steps simulated by the presented model to predict the microstructure in stir zone. *Left and middle:* snapshot of microstructure of the weld zone during DRX, and *right:* dislocation accumulation





**Fig. 10** Microstructural of stir zone in the weld produced in a traverse speed of 50 mm/min and rotational speed of **a** 710, **b** 900, **c** 1120, and **d** 1400 rpm



**Fig. 11** Microstructure of stir zone in the weld produced in rotational speed of 1400 rpm and traverse speed of **a** 25, **b** 50, **c** 80, and **d** 100 mm/min



Figure 10, illustrating the microstructure and grain size distribution for welds simulated in different rotational speeds, reveals that the average grain size in the stir zone rises as the tool rotational speed increases. This result is widely reported by experimental works (Besharati Givi and Asadi 2014; Asadi et al. 2011b). Similarly, Fig. 11 shows the microstructure and grain size distribution for the welds simulated in different traverse speeds where the rotational speed is 1400 rpm. It is clear that by increasing the traverse speed, the average grain size reduces. Indeed, increase in  $w/v$  results in a rise in the amount of generated heat and peak temperature, leading to acceleration of grain growth step, and thus, the final microstructure will contain larger grains (Salekrostam et al. 2010; Tutunchilar et al. 2012b).

## Conclusions

In this work, a finite element model is established to study the microstructure evolution during FSW of AZ91 magnesium alloy. The hardening parameter, the recovery parameter, and the strain rate sensitivity, required for the model, are calculated according to flow stress results. A continuum based thermo-mechanically coupled rigid-viscoplastic FEM model is proposed in Deform-3D software to simulate the FSW of AZ91. To evaluate microstructure of the weld zone, a model is proposed based on the combination of cellular automaton and Laasraoui-Jonas models. Results show that the simulated microstructure of the weld zone has a good agreement with that of the experiments. The proposed model can simulate the dynamic recrystallization process during friction stir welding and predict the grain size and microstructure of the weld zone precisely. The simulated grain size under different process parameters reveals that by increasing the  $w/v$  parameter, the grain size increases.

## Competing interests

The authors declare that they have no competing interests.

## Authors' contributions

PA carried out all the experimental tests. PA and MA performed the simulation of process and microstructural evolution method and formulation. MK and BG edited the manuscript as the supervisor of the project. All authors read and approved the final manuscript.

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