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Persistent slip bands: The bowing and passing model revisitedLadislav Kubin^{a*} and Maxime Sauzay^b^aLaboratoire d'Étude des Microstructures, CNRS-ONERA (UMR 104 CNRS),
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91191 Gif-sur-Yvette Cedex, France**Abstract**

Brown's bowing and passing model for persistent slip bands (PSBs) was extended in Cu and Ni to the whole range of temperatures in which a saturation plateau is observed. Advantage was taken of the similitude relation to rewrite in dimensionless form the equations of the original model and of a more accurate revisited version. Input quantities for computing the solutions were taken from a previous study of experimental results; all unknown quantities could then be directly calculated without any assumption or approximation. The comparison between experimental results and the predictions of the revisited model confirms the basic assumptions of the bowing and passing model, according to which the thermally activated annihilation of screw dipoles is governing the channel widths, the Orowan stresses and the critical stresses in the channels. All other assumptions and numerical predictions are perfectly confirmed, save for the occurrence of small resistive stresses in the channels. In addition, a better understanding of the complex behavior of PSB walls under stress is necessary in order to accurately determine the plastic strain amplitude of PSBs. Mesoscale and atomistic simulations are needed for further modeling of the wall properties and the screw dipole annihilations.

Keywords: Persistent slip bands, Saturation stress, Similitude, Modeling, Screw annihilations.

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1 Introduction

Persistent slip bands (PSBs) are formed in ductile materials cycled in single slip at imposed plastic strain amplitudes per cycle typically between 10^{-4} and 10^{-2} . Since their discovery in the 1950s these localized bands of plastic strain have drawn a lot of attention for practical reasons related to fatigue damage as well as for fundamental reasons. Indeed, they constitute a unique example of mechanical response exhibiting a constant saturation stress. Hence, there is a wealth of experimental and theoretical literature on the relation between their almost periodic wall-and-channel dislocation structure and their properties [1-4].

The PSB channels contain a constant density of screw dislocations, which shuttle forward and backward during cycling. These dislocations mutually annihilate by cross-slip after having travelled a certain mean free path and deposited edge segments along the PSB walls. A small fraction of these edge segments is able to cross the walls; it emerges and expands in the neighboring channels where it produces fresh screw dislocations of both signs. Hence, the saturation of the screw dislocation density results from an exact balance between creation and annihilation mechanisms. The PSB walls contain a large density of short edge segments, mostly in the primary plane but also in the cross-slip plane. The microstructure was found to be rather complex [5] and, as yet, former models for saturation in the walls were not revisited. The walls are partly permeable to edge dislocations, but screw dislocation motion in the channels is producing the largest part of the imposed plastic strain amplitude.

In its simplest form, the similitude relation expresses that the characteristic dimension of a dislocation microstructure, λ , is proportional to the average length of the dislocation segments, $\bar{\ell}$, whereas the flow stress τ_{\square} is inversely proportional to it. Thus, the flow stress is inversely proportional to the characteristic dimension and is written $\tau_{\square} = K\mu b/\lambda$, where K is a

similitude constant, μ is the shear modulus and b is the magnitude of the Burgers vector [7].

The saturation stress of PSBs is strongly temperature dependent, so that investigations on similitude require measurements performed on microstructures formed at different temperatures.

In the recent study [8], the similitude properties of PSB channels and their consequences were investigated in full detail from experimental results in copper and nickel crystals cycled at saturation. The domains of temperature investigated encompassed the whole domain of occurrence of PSBs. It was found that both PSB channels and PSBs follow similitude relations. This may seem paradoxical in the case of PSBs since the wall-widths do not depend on temperature; it is so because of a compensating term arising from PSB channels, in which the critical stress for screw dipole annihilations is smaller than the saturation stress (see the discussion in Section 3 of [8]). The obtained results were confronted to the predictions of Brown's bowing and passing model [9], according to which the critical flow stress of PSB channels is governed by the annihilation of screw dipoles by cross-slip. This critical stress is defined as the stress corresponding to the maximum height of stable dipoles, that is, to their passing stress. As a whole, the predictions of the bowing and passing model were found to be in reasonable agreement with experimental results. However, as the model does not account for any other interaction than the ones between screw dislocations, the critical stresses in the channels were found smaller than the saturation stresses by a few MPa.

An alternative model, the composite model, was developed over the years by Mughrabi and co-workers, essentially at room temperature (see [4] for a full review). This model emphasizes the occurrence of significant internal stresses in the walls and channels and assumes that the annihilations of screw dipoles occurs spontaneously without the help of thermal activation.

Till now, debates about the modeling of the saturation stress essentially focused on the properties of screw dislocations in PSB channels at 300 K. The motivation of the present work is to provide an extension to a wide range of temperatures. The most basic equation of the bowing and passing model [9] is rather intricate and full predictions cannot be expressed in analytical form. Thus, the model was cast into a more transparent form with the help of a few simplifying assumptions, which were validated at room temperature. In the present work, the basic equation is solved numerically. This makes it possible to predict without any approximation the values of all the relevant quantities as a function of flow stress and temperature. For this purpose, use is made of the results drawn in [8] from experimental studies on copper and nickel.

In Section 2, the basic equations of the model are set in dimensionless forms, which incorporate the similitude constants of PSB channels determined in [8]. The results presented in Section 3 are concerned with the temperature and critical stress dependencies of the critical screw dipole heights and of the coefficient determining the respective contributions of the bowing and passing stresses to the critical stress. Except for the small differences between critical and saturation stresses mentioned above, a striking agreement is found between the model predictions and experimental results on copper and nickel. A few other predictions, in particular on the intrinsic plastic shear amplitudes of PSBs, are discussed in Section 4 and concluding remarks highlighting the major results are presented in Section 5.

2. The bowing and passing model revisited

This part recalls first the original bowing and passing model and next the way it is expanded to a wide range of temperatures. Further, a revisited version is established in order

to verify whether or not an approximation made at room temperature is valid at all temperatures.

2.1. The original bowing and passing model

In its original form, Brown's bowing and passing model [9] has brought an improved analytical answer to comments by Mughrabi and Pschenitzka [10] on a previous attempt [11] to estimate the saturation stress in PSB channels.

The discussion of the model is based on Eq. (10*)¹. This equation yields the critical stress of the channels τ_c as a linear combination of the Orowan stress τ_{Or} , which sets screw dislocations in motion, and the passing stress, τ_{pass} , which corresponds to the maximum critical height, h_c , for the annihilation of screw dipoles

$$\tau_c = \alpha_B \tau_{Or} + \tau_{pass} = \alpha_B \frac{2E_{edge}}{bd_{ch}} + \frac{\mu b}{4\pi h_c} \quad (1)$$

In the assumed absence of any other stress contribution, the flow stress τ_c is assumed to be the stress at saturation. The constant coefficient α_B accounts for the contribution of the Orowan stress to the flow stress. In this Orowan stress, E_{edge} is the line energy of edge segments (see Supplementary Section S1 for line tensions and line energies) trailed by the bowing screws in channels of width d_{ch} ². The contribution of the passing stress is expressed in terms of the critical dipole height, h_c . In the present context, it is convenient to introduce similitude in Eq. (1) by rewriting it in the form

¹ In what follows, starred equations are those of the original bowing and passing model [9].

² In Brown's model [9], the channel widths (or wall spacings) are denoted by d . In what follows, d denotes the periodicity of a PSB, $d = d_{ch} + d_w$, where d_w is the wall thickness.

$$K_{ch} = \frac{\tau_c d_{ch}}{\mu b} = \frac{2\alpha_B E_{edge}}{\mu b^2} + \frac{1}{4\pi} \left(\frac{d_{ch}}{h_c} \right), \quad (2)$$

where the first equality expresses the similitude relation in the channels and K_{ch} is the related similitude slope. In Eqs. (1) or (2) the channel widths and wall thicknesses are known from experiment. The value usually quoted in the literature for the critical annihilation distance of copper at room temperature is $h_c \approx 50$ nm for copper at room temperature [1,12]. The value of the coefficient α_B is critical too because it contributes to the magnitude of resistive stresses in the channels, $\delta\tau_{ch} = \tau_{PSB} - \tau_c$, where τ_{PSB} is the saturation stress. Such resistive stresses were found in the previous study [8] but, as they are not included in the bowing and passing model, they do not appear in its equations.

In the model, the coefficient α_B is drawn from a rather complex master equation describing the critical configuration of a dipole of flexible screw dislocations at the passing stress. For this purpose, a small transition zone is defined close to the PSB walls. It connects the straight edge lines deposited on the walls by the motion of the screws to the two interacting screw lines with large critical radii in the central part of the channels. In the transition zone, the characters of the interacting lines go from nearly edge to nearly screw and the curvature radii are quite small. The equilibrium condition for this critical configuration is estimated using an effective value of the line tension, T_{eff} , which is taken for line orientations of 45° (see Sections S1 and S2.2). As a result, one obtains the master equation given by Eq. (9*):

$$\tau_c = \frac{2E_{edge}}{bd_{ch}} + \frac{\mu b}{4\pi h_c} \left(1 - \frac{b^2 d_{ch}^4 \left(\tau_c - \frac{\mu b}{4\pi h_c} \right)^2}{640 h_c^2 T_{screw}^2} \right) k, \quad (3)$$

where $k = (1 - 2 T_{\text{eff}} / \tau_c d_{ch} b)$ and T_{screw} is the line tension of screw dislocations (Eqs. S1-2 and S2-3). At room temperature k does not differ much from unity, so that it is assumed that $k = 1$ at all temperatures. Eq. (3) is then rewritten in terms of the coefficient α_B as given by Eqs. (1) or (2). This leads to Eq. (11*)

$$1 - \alpha_B - \alpha_B^2 \frac{\mu b^2}{2E_{\text{edge}}} \frac{1}{640\pi} \left(\frac{d_{ch}}{h_c} \right)^3 \left(\frac{E_{\text{edge}}}{T_{\text{screw}}} \right)^2 = 0 \quad (4)$$

The derivation of Eq. (4) is reproduced at end of Section S2.2. The value of α_B is obtained by introducing into Eq. (4) a plausible value $d_{ch}/h_c \approx 30$ in copper at 300 K. This yields $\alpha_B \approx 0.5$, which is adopted for further developments of the original model. According to Eq. (1), the internal stresses calculated by Mughrabi and Pschenitzka [10] should then correspond to a smaller value of α_B .

2.2. The revisited bowing and passing model

In the original model, Eq. (3) is simplified by taking $k = 1$. The equations of the revisited model are obtained by fully treating the term k and solving numerically Eq. (3) for the critical dipole height h_c . This allows Eq. (4) to be further solved for the coefficient α_B . For this purpose, and as was done for Eq. (1), Eqs. (3) and (4) are rewritten in such a way as to incorporate the similitude coefficient of the channels, K_{ch} . The expanded forms of Eqs. (3) and (4) are derived in Section S2.2. The revisited form of Eq. (3) is written

$$F\left(\frac{d_{ch}}{h_c}\right) = 0 = -K_{ch} + \frac{k}{4\pi} \left(\frac{d_{ch}}{h_c}\right) + \frac{2E_{\text{edge}}}{\mu b^2} - k \left(\frac{d_{ch}}{h_c}\right)^3 \frac{\left(K_{ch} - \frac{1}{4\pi} \left(\frac{d_{ch}}{h_c}\right)\right)^2}{4\pi 640 (T_{\text{screw}} / \mu b^2)^2} \quad (5)$$

This polynomial equation of fifth order in d_{ch}/h_c has a single-valued solution in the domain of temperatures where PSBs are observed. Hence, the temperature dependence of h_c can be determined unequivocally. The revisited Eq. (4) is written

$$(1 - \alpha_B) - \alpha_B^2 \frac{\mu b^2}{2E_{\text{edge}}} \frac{k}{640\pi} \left(\frac{d_{ch}}{h_c} \right)^3 \left(\frac{E_{\text{edge}}}{T_{\text{screw}}} \right)^2 - \frac{(1-k)\mu b^2}{8\pi E_{\text{edge}}} \left(\frac{d_{ch}}{h_c} \right) = 0 \quad (6)$$

Equation (6) has a single positive root in the interval $0 \leq \alpha_B \leq 1$. At this step, the critical stress in the channels can be obtained directly from Eq. (2) and other properties can be determined by further developing the bowing and passing model. As shown in Section S2.2, setting $k = 1$ into Eq. (5) yields Eq. (3) of the original model. One can also easily verify that Eq. (6) reduces to Eq. (4) of the original model.

3. The expanded bowing and passing model: numerical results

This part reports on the predictions drawn from the original and revisited models expanded as a function of temperature. The experimental data on Cu and Ni are recalled first; they provide input values for solving the two models. The quantities successively determined are the critical dipole heights h_c , the coefficient α_B , the critical stresses τ_c and the dependence of critical dipole heights on critical stresses. All computations were performed using Mathematica v. 6 (Wolfram Research).

3.1. Experimental data

The experimental data used in the present study are taken from reference [8], which can be consulted for more detail. For copper, two sets of experimental measurements were lumped together, one by Basinski and Basinski [3] between 4.2 and 295 K and the other by Holzwarth and Essmann [13,14] between 77 and 430 K. The highest investigated temperatures correspond to the vanishing of the saturation plateau. In what follows, these two data sets are referred to as Cu_{BB} and Cu_{HE} . Although the channel widths exhibit large fluctuations [14], they could be measured with a reasonably good accuracy. The wall thicknesses are constant with $d_w = 0.11 \pm 0.01 \mu\text{m}$ and, as they were not measured in [3], it was assumed that the value of d_w is unchanged below 77 K.

In nickel, the measurements by Tippelt et al. [5] and Hähner et al. [15] include four data points between 77 to 600 K. A data point at 750 K was removed in [8] because of the coexistence of dislocation cells and PSBs in the microstructure. The wall thicknesses are also constant in nickel with $d_w = 0.155 \mu\text{m} \pm 10\%$ [6], but no scatter is available for the channel widths .

The additional data given in Table 1 are drawn from [8] and will be used in what follows. One may notice, as was done in [8], that the similitude properties and the saturation stresses at 0 K of Cu and Ni differ from each other.

Table 1 about here

The similitude properties of PSBs as a whole were drawn by plotting τ_{PSB}/μ as a function of b/d and taking the slope of the regression line $K_{PSB} = \tau_{PSB}d/\mu$. The corresponding determination coefficients were always found larger than $R^2 = 0.99$. For Cu and Ni, the regression lines pass through the origin within quite small stress intercepts, respectively, -5×10^{-2} MPa and 1.02 MPa. As far as the channels are concerned, the situation is

different as the stress intercepts are much larger (Table 1). The positive values indicate that the flow stresses of the channels are smaller than the saturation stresses. Similar stress intercepts were measured by Bretschneider et al. [16] in Ni and drawn from several experimental data by Brown (see Section 4 of [9]).

As previously done in [8], all the numerical calculations presented in what follows employ the approximation proposed by Scattergood and Bacon for the shear moduli and Poisson's ratios (ν) of Cu and Ni (Section S1). The temperature dependence of the elastic stiffnesses was taken into account (Section S3). In what follows, the main source of error arises from the measurements of channel widths. The measurements on Cu_{HE} are more accurate ($\pm 4\%$) than the ones on Cu_{BB} ($\pm 7\%$). The error on wall thicknesses is comparatively smaller³ but it is included.

3.2. The critical dipole heights in the channels

The ratio d_{ch}/h_c was calculated from the revisited Eq. (5), in which the input quantities are the similitude slope K_{ch} of the channels, the channel widths d_{ch} and the wall thicknesses. The values at $T = 0$ K were included. The critical dipole heights were directly obtained from the product $h_c = d_{ch}(d_{ch}/h_c)^{-1}$. According to the original model, the ratio d_{ch}/h_c should be about 30 or perhaps a bit less. In Table 2, the values of the coefficient k , the ratio d_{ch}/h_c and the critical dipole heights obtained from the revisited and original versions of the model are compared at three temperatures.

Table 2 about here

³ The wall thickness only appears in the logarithmic term of the line energy of edge dislocations (Eq. S1-1).

In the revisited model, the values of the coefficient k decrease in a non-negligible manner with increasing temperature. Hence, it seems that the approximation $k = 1$ is rather rough. Whereas the original model yields ratios d_{ch}/h_c that decrease with increasing temperature, these ratios are almost constant in the revisited model. This arises from a compensating effect of the term k . In the revisited model, the critical screw dipole heights at room temperature are close to the most frequently quoted value of 50 nm for copper [1,12], specifically 44.8 ± 3 nm for a saturation stress of 29 MPa in Cu_{BB} and 49.3 ± 2 nm for a saturation stress of 28 MPa in Cu_{HE}. For nickel h_c (293 K) \approx 48 nm.

Figure 1a shows revisited similitude plots of the scaled saturation stresses as a function of the scaled critical dipole heights in Cu and Ni. The similitude slopes are almost the same and the positive stress intercepts reveal the occurrence of resistive stresses, which will be further discussed in Section 3.3. Fig. 1b shows the temperature dependence of the revisited critical dipole heights.

Figure 1ab about here

One can see from overlapping data in Fig. 1b that the dipole heights are almost identical in Cu and Ni. The only significant differences reside in the critical stresses at 0 K (Table 1) and in the maximum temperatures at which a saturation plateau was obtained. The same features were experimentally observed on the temperature dependence of the scaled stresses τ_{II} in monotonic deformation (see [17] p. 160 and Fig. 28 of [6]). Hence, in spite of their similarities in an intermediate temperature domain, the stress dependences of the activation energies for screw dipole annihilations in Cu and Ni are expected to be different. This is not surprising since the scaled stacking fault energies $\gamma/\mu b$ are also different, 4.0×10^3 for Cu and 6.0×10^3 for Ni (see Table B1 of [18]).

In copper, the critical dipole heights at 0 K are $h_{co}(\text{Cu}) \approx 16$ nm and $h_{co}(\text{Ni}) \approx 20$ nm. The simulations by Pauš et al. [19] shown in Fig. 1b yield similar values for Cu at all temperatures, with $h_c(\text{Cu}) = 23$ nm at 4.2 K. Similar values were also found for nickel at 77 and 300 K (see Table 1 of [19]). Mitchell [17] also compiled extrapolated resolved stresses for the onset of the dynamic recovery stage at 0 K in monotonic deformation⁴. From his data and the similitude relations of the critical dipole heights (Fig. 1a), one draws $h_{co}(\text{Cu}) = 9 - 17$ nm and $h_{co}(\text{Ni}) = 12 - 35$ nm. These comparisons show that the present results are compatible with experiment and confirm the thermally activated character of screw dipole annihilations.

Although atomistic simulations also confirmed that screw dislocation cross-slip is thermally activated [20], only few of them were devoted to the annihilation of screw dipoles (e.g., [20-22]). Unfortunately, they did not incorporate the contribution of the externally applied stress, so that the critical dipole heights at 0 K were too small and the critical stresses too high. In such stress-free conditions, it is not possible to obtain a realistic insight into the bowing and passing mechanism.

3.3. The coefficient α_B and the critical stresses in the channels

As the values of k and d_{ch}/h_c are known, the coefficient α_B can be calculated from Eqs. (4) and (6) for the two versions of the expanded model. The results are given in Table 3.

Table 3 about here

The original model yields α_B values that increase with increasing temperature. The stabilizing effect of the term k manifests itself again in the revisited model. In Cu, one has α_B

⁴ In metals with low stacking fault energy, twinning may interfere with dynamic recovery.

(Cu) = 0.503 ± 0.007 over the whole temperature range. Hence, the value estimated by Brown [9] at 300 K, $\alpha_B \approx 0.5$, is actually temperature-independent. For Ni, $\alpha_B(\text{Ni}) = 0.532 \pm 0.04$.

As the values of d_{ch}/h_c and α_B are known, the critical stresses τ_c can be directly estimated from Eq. (2). In the two versions of the model, the similitude coefficient of the channels, K_{ch} , is an input property. For both Cu and Ni, the critical stresses are found smaller than the saturation stresses. Hence, the stress difference $\delta\tau_{ch} = \tau_{PSB} - \tau_c$ represents additional resistive stresses. The stresses $\delta\tau_{ch}$ and τ_c yielded by the original model are obtained as follows. The similitude relation in the channels is written $K_{ch} = \tau_c d_{ch}/\mu b$ and the resistive stresses are obtained from

$$\delta\tau_{ch} = \tau_{PSB} - \tau_c = \tau_{PSB} - \frac{K_{ch}\mu b}{d_{ch}}, \quad (7)$$

where all numerical values are known from experiment in the last term at the right-hand side. The original model yields almost exactly the same results as the comparison with experiment performed in [8]. For instance, the average values, 5.33 MPa for Cu and 5.97 MPa for Ni, are very close to the ones previously determined by another method (Table 1). The same holds for the individual resistive stresses. In the revisited model, the term k slightly modifies the similitude relation because it contains a drifting logarithmic term (Eq. S2-3). As a result, the average resistive stresses $\delta\tau_{ch}$ decrease to 5.0 MPa in Cu and 4.84 MPa in Ni.

All the stresses discussed above, plus the bowing stress, are plotted as a function of temperature in Fig. 2a. The resistive stresses are obtained through a difference between two large quantities. Thus, they exhibit large fluctuations, at least $\pm 8\%$ for Cu_{HE} and $\pm 14\%$ for Cu_{BB} , owing to the accumulation of known and unknown experimental uncertainties. In addition, they do not seem to exhibit a well-marked temperature dependence.

Fig. 2ab about here

Figure 2b shows a similitude plot of the critical stresses in the revisited model as a function of the critical dipole heights in Cu and Ni. The similitude coefficients (K_{hc}) only slightly differ from each other and from the ones obtained for the whole PSBs in the revisited model (Fig. 1a). The stress intercepts are negligibly small and well below experimental uncertainties. The similitude properties of the channels in the revisited model are only slightly modified with respect to the ones previously obtained in Ref. [8]. For Cu and Ni, the coefficients $K_{hc} = \tau_{c, rev.} d_{ch}/\mu b$ respectively differ by 2% and 0.5%.

In summary the revisited model is more accurate than the original one. It yields constant values for the coefficient α_B and the ratio d_{ch}/h_c as well as critical stresses in the channels that are smaller than the saturation stresses by a few MPa. The critical screw dipole heights are thermally activated, they follow a similitude relation and their values at 300 K is $h_c \approx 50$ nm.

4. Additional predictions of the revisited bowing and passing model

Brown's model [9] contains a few predictions for persistent slip, which do not derive directly from the complex Eq. (5) and are drawn from various other considerations.

4.1. Equilibrium of PSB walls and similitude

The basic Eq. (1) is rewritten in Eq. (12*) as a sum of the Orowan stress and a complementary passing stress term τ_p

$$\tau_c = \tau_{Or} + \tau_p = \tau_{Or} + \left(\frac{\mu b}{4\pi h_c} - (1 - \alpha_B) \tau_{Or} \right), \quad (8)$$

The original model [9] is further developed on the basis of Eq. (8) by making an hypothesis about the equilibrium of PSB walls. Considering that the motion of screws under the Orowan stress τ_{Or} tends to push too closely spaced walls back in place, whereas the formation of screw dipoles tends to pull the walls together, the equilibrium of wall spacings requires that $\tau_{Or}/\tau_p = 1$ or, equivalently, $\tau_c = 2\tau_{Or}$. By importing these two relations in Eq. (8) one obtains Eq. (14*) of the original model, $\tau_c = \mu b/2\pi(2-\alpha_B)h_c$. This last equation is rewritten in the form of a similitude relation

$$K_{hc} = \frac{\tau_c h_c}{\mu b} = \frac{2\tau_{Or} h_c}{\mu b} = \frac{1}{2\pi(2-\alpha_B)} \approx \frac{1}{3\pi} \quad (9)$$

The relations drawn from the equilibrium hypothesis were tested on the revisited version of the model. For Cu and Ni, the ratios τ_{Or}/τ_p are almost independent of temperature with average values larger than unity, 1.200 ± 0.003 in Cu and 1.36 ± 0.03 in Ni. The ratios τ_c/τ_{Or} behave in the same manner with average values smaller than two, 1.83 ± 0.03 in Cu and 1.73 ± 0.01 in Ni. By reproducing the simple calculation performed above without attributing values to the two ratios and to α_B , one obtains the following corrected form of Eq. (9)

$$K_{hc} = \frac{\tau_c h_c}{\mu b} = \frac{(\tau_c / \tau_{Or}) \tau_{Or} h_c}{\mu b} = \frac{(\tau_c / \tau_{Or})}{4\pi(1 + (\tau_p / \tau_{Or}) - \alpha_B)} = 0.109 \text{ (Cu); } 0.114 \text{ (Ni)} \quad (10)$$

This relation reproduces almost perfectly the two values of the similitude coefficients given in Fig. 2b. Thus, the two stress ratios τ_{Or}/τ_p and τ_c/τ_{Or} do have constant values.

The measured channel widths, which are input quantities, are all fixed; in copper they are all given with error bars representing the amplitude of their fluctuations. Eq. (10) shows that in such conditions, the hypothesis based on the equilibrium of the walls is by construction embedded into the bowing and passing model. Thus, no assumption is needed because the constancy of the two stress ratios simply derives from the fact that the critical dipole heights follow a similitude relation.

4.2. *The intrinsic plastic strain amplitude of PSBs*

The intrinsic plastic strain amplitude of PSBs is defined as the shear strain γ_{PSB} per quarter of cycle under the saturation stress τ_{PSB} . Reliable experimental values of this quantity are rather scarce. Here, use is made of the data by Holzwarth and Essmann [14] in copper, which agree well with other published values. The plastic strain amplitude was found to depend linearly on the saturation stress. Four values of γ_{PSB} were determined at different temperatures from direct measurements of the active volume fraction of PSBs.

The plastic strain amplitude γ_{PSB} was estimated by Brown in 2004 [11]. The probability for complete annihilations of screw dislocations was calculated as a function of the probability p that an edge dislocation crosses a wall (Eq. (20) of [11]). In the improved model [9], the introduction of the coefficient α_{B} does not modify the previous result. As all experimental results were plotted as a function of τ_{PSB} the critical stress τ_{c} (Eq. (15*) of [9]) is replaced here by τ_{PSB} . A linear relation is then obtained between γ_{PSB} and τ_{PSB}

$$\gamma_{\text{PSB}} = \frac{\pi(1+p)^2}{(1-p)^2} \frac{3d_{\text{ch}}}{3d_{\text{ch}} + d_{\text{w}}} \frac{\tau_{\text{PSB}}}{\mu} = 4\pi \frac{3d_{\text{ch}}}{3d_{\text{ch}} + d_{\text{w}}} \frac{\tau_{\text{PSB}}}{\mu} \approx 4\pi \frac{\tau_{\text{PSB}}}{\mu} \quad (11)$$

It was shown in [11] that $p \approx 1/3$, so that the term containing the probability p is equal to 4π . The next term is a tentative correction close to unity, which accounts for screw dipole annihilations in the walls. The prediction drawn from Eq. (11) holds for the two versions of the model since the involved quantities are the same in both cases. It is compared to the experimental results on Cu_{HE} in Fig. 3a. The experimental and predicted linear dependences on saturation stress are somehow different. An attempt was made to fit the experimental data by adjusting the value of the coefficient p . The best result was obtained by taking $p = 0.39$.

Fig. 3ab about here

Brown [11] proposed a simplified alternative model for γ_{PSB} based on the annihilations of rectangular edge-screw loops expanding through walls and channels (see Section S4 for more detail on the model). The intrinsic plastic shear strain amplitude of PSBs is given by

$$\gamma_{\text{PSB}} = b/2Fh_c, \quad (12)$$

where F is a free statistical parameter. With $F = 1.43$, Eq. (12) reproduces rather well the experimental results (Fig. 3b).

Figure 3 shows that the predictive ability of Eqs. (11) and (12) may be improved by adjusting in each case the value of a single coefficient. In the two equations it is assumed that the channels and walls deform under an uniform saturation stress. This cannot be since the flow stress of the channels is smaller than the saturation stress, so that the flow stress of the walls is necessarily larger than τ_{PSB} . Furthermore, the real microstructure of the walls that was described by Tippelt et al. [5] is much more complex than the traditional ones that were used to derive Eqs. (11) and (12). In particular the walls contain a large density of small prismatic dislocation loops lying in the cross-slip plane, which should strongly interact with incoming

dislocations. Thus, the loops expanding through the walls must exhibit quite complicated shapes.

5. Concluding remarks

The bowing and passing model for PSBs [9] was extended to the whole domain of temperatures in which a saturation plateau is observed. Advantage was taken of the similitude relation to rewrite in dimensionless form the equations of the original model and of a revisited model, which is more accurate. Input quantities for the computation of solutions were taken from a previous study of experimental results on copper and nickel. In such conditions, all unknown coefficients and variables were directly calculated without any assumption or approximation.

The comparison between experimental results and the predictions of the original and revisited bowing and passing models were discussed in the previous sections. This part emphasizes the major results obtained from the revisited model and their consequences.

- Two important quantities were found to be constant irrespective of temperature. i) – The ratio d_{ch}/h_c of the channel width to the critical dipole height is about 29 in Cu, close to the value assumed by Brown [9] at 300 K, and about 25 in Ni. ii) – The coefficient α_B governs the contribution of the Orowan stress to the critical stress and the magnitude of the resistive stress in the channels. Its values are α_B (Cu) = 0.5, as was assumed by Brown [9] at 300 K, and α_B (Ni) = 0.53.
- The critical dipole heights, h_c , were computed as a function of temperature in the whole domain of existence of PSBs. At room temperature, the revisited model yields values close to 50 nm in copper, in agreement with the experimental estimate given in [1,12], and in nickel. At 0 K, the critical dipole heights are h_{co} (Cu) = 16 nm and h_{co} (Ni) = 20 nm. Although

atomistic simulations confirmed that cross-slip is a thermally activated phenomenon, they did not provide realistic predictions for the critical dipole heights. The present results may serve as a benchmark test for further studies, in which an external loading is applied to the interacting screw dislocations.

- The critical dipole heights are related to the critical stresses in the channels, τ_c , through a perfect similitude relation passing through the origin (Fig. 2b). The channels widths are also related to the critical stresses through another perfect similitude relation (Fig. 6 of [8]). This arises because the ratio d_{ch}/h_c is a constant. The ratio τ_c/τ_{Or} of the critical stress to the Orowan stress in the channels was assumed in [9] to be a constant from considerations on the equilibrium of the walls. The value of this ratio is actually rooted in the similitude relation between the critical dipole heights and the critical stresses (Section 3.1). Finally, the critical stresses at 0 K are similar to the critical stresses for the onset of the dynamic recovery stage III in monotonic deformation (Section 3.2). These results are of prime importance because they confirm again the basic assumption of the bowing and passing model, according to which the thermally activated annihilation of screw dipoles is governing the channel widths, the Orowan stresses and the critical stresses in the channels. In contrast, the central role played by the critical dipole height challenges a basic assumption of the composite model according to which the annihilation of screw dipoles is only stress-assisted.
- In this study no new information could be drawn about the origin of the small resistive stresses in the channels. It is only confirmed that the critical stresses in the channels are slightly smaller than the saturation stresses by a few MPa and that they do not exhibit a well-marked temperature dependence [8]. In this previous study the critical stresses were directly obtained from experiment through the stress intercepts of the similitude relation in the channels. It was shown in Section 3.1 of [8] that the use of the line tension proposed by Scattergood and Bacon [S3] would slightly improve the results. Nevertheless, the magnitudes

of the resistive stresses in the channels are now reasonably well known. Their existence was not considered in the bowing and passing model; taking them into account allowed all other predictions of the model to be confirmed. According to Mughrabi and Pschneitzka [10], the critical stress τ_c should be about 20% larger than the largest of the contributions from the Orowan and dipole bowing stresses. In copper at room temperature, we obtained from the present data a resistive stress of 11 MPa, whereas the revisited model yields a resistive stress around 4 MPa, and a coefficient $\alpha_B \approx 0.2$ instead of 0.5. Thus the model [10] overestimates the resistive stresses.

- The first dislocation dynamics simulations on channels of persistent slip bands [23–25] were intended to compare the predictions of Brown's original model [9] and the model by Mughrabi and Pschneitzka [10]. No real consensus was obtained because these simulations, as well as the model, strongly depended on simplifying assumptions made about the complex microstructure of the walls. More recent DD simulations (e.g. [19]) and mass simulations that are now in progress will certainly contribute to clarify the situation.
- The models proposed by Brown [9,11] for the intrinsic plastic strain amplitude of PSBs take into account the penetration and crossing of PSB walls by edge dislocations. However, to obtain accurate predictions, it is necessary to account in a more realistic manner for the microstructure of the walls. This modeling task should be performed in parallel with dislocation dynamics simulations and atomistic simulations for short-range interactions.

In summary, the numerical solutions of Brown's expanded and revisited model for bowing and passing [9] are in excellent agreement with its basic assumptions and predictions. The only exception is the occurrence of small resistive stresses in the channels. In addition, much remains to be understood about the behavior of the walls. Analytical approaches may be extremely complex and only mesoscale mass simulations and atomistic simulations can help

going further. Finally, it may be interesting to examine as a function of temperature the compatibility of the composite model [4] with the bowing and passing model.

Appendix A. Supplementary data

Supplementary data associated with this article can be found at

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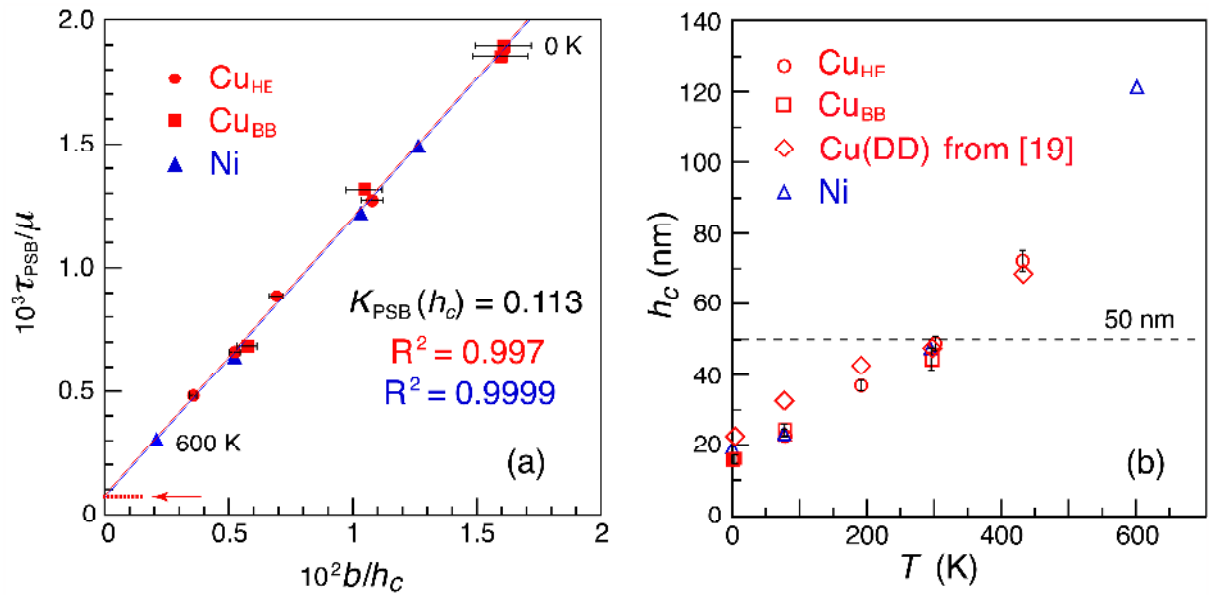


Fig. 1. Critical dipole heights, h_c , in Cu and Ni as yielded by the revisited model. The values at 0K and available error bars are included ($\pm 8\%$ for Cu_{HE} and $\pm 14\%$ for Cu_{BB}). (a)

Similitude plot of the scaled saturation stress τ_{PSB}/μ vs. h_c/b . The arrow near the origin points at small intercepts with the stress axis. (b) Temperature dependence of the critical dipole heights. The data Cu_{DD} is taken from dislocation dynamics simulations of the bowing and passing model performed by Pauš et al. [19].

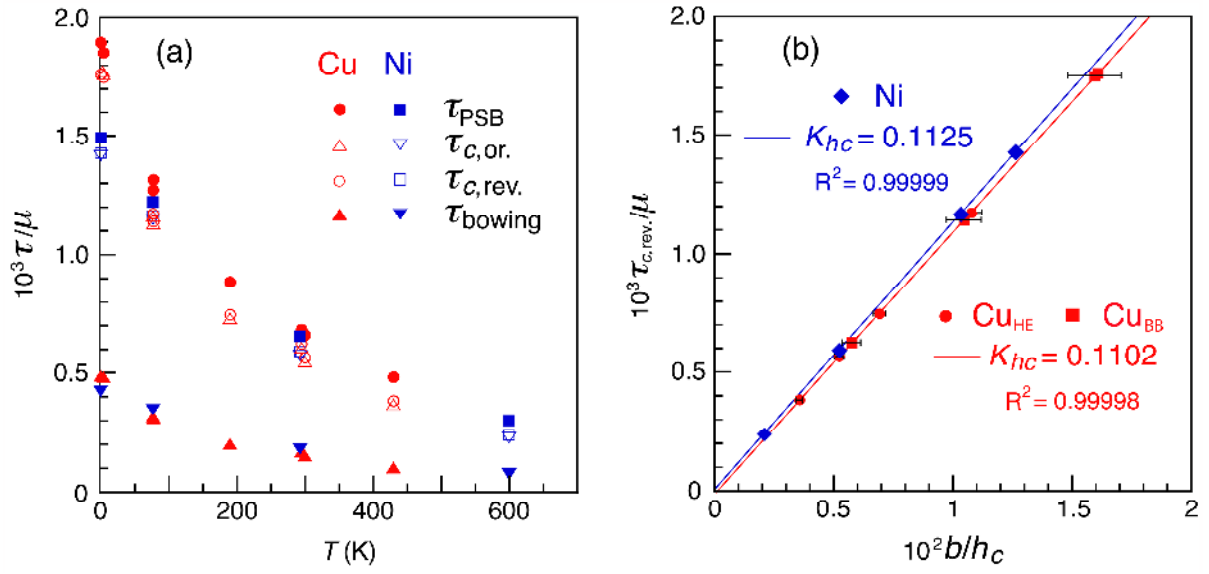


Fig. 2. (a) Temperature dependence of several stresses in Cu and Ni. τ_{PSB} is the saturation stress, $\tau_{\text{c, or.}}$ and $\tau_{\text{c, rev.}}$ are the critical stresses in the channels in the original and revisited versions of the model and τ_{bowing} is the bowing stress in the revisited model. (b) Similitude plot of the scaled critical stresses $\tau_{\text{c, rev.}} / \mu$ as a function of b/h_c . The stress values at 0 K are included in both figures. The error bars are $\pm 8\%$ for Cu_{HE} and $\pm 14\%$ for Cu_{BB}.

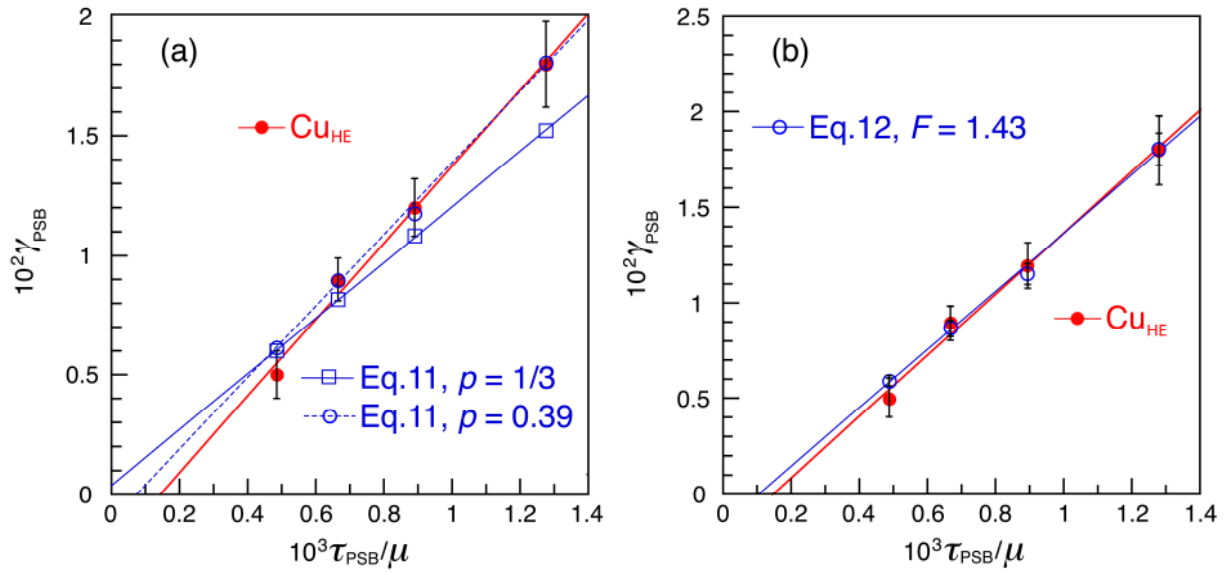


Fig. 3.

The plastic strain amplitude of PSBs, γ_{PSB} , as a function of the scaled critical stress τ_{PSB} in Cu_{HE} (after [14]) compared to the predictions of the bowing and passing model. (a)

Predictions drawn from Eq. (11), with $p = 1/3$ and from a fit with $p = 0.39$. (b) Prediction

obtained from Eq. (12) for the revisited model with a value $F = 1.43$ for the free-parameter.

The error bars of the measured and predicted values are included.

Table 1

Data drawn from experimental results on Cu (Cu_{BB} and Cu_{HE} merged together) and Ni. From left to right: – the extrapolated saturation stresses at 0 K scaled by the shear modulus at 0 K, – the similitude slopes for PSBs (K_{PSB}) and PSB channels (K_{ch}) with, in parentheses, the stress intercepts (in MPa) and the correlation coefficients, – the average values of the moderately temperature-dependent resistive stresses in the channels (in MPa). The stresses at 0 K are included in the last three sets of results.

	τ_0/μ_0	K_{PSB}	K_{ch}	$\langle \delta\tau_{\text{ch}} \rangle$
Cu	$1.90 \cdot 10^3$	4.06, (-0.05 – 0.992)	3.10, (5.40 – 0.992)	5.31
Ni	$1.55 \cdot 10^3$	3.58, (1.02 – 0.999)	2.78, (5.53 – 0.998)	5.96

Table 2

Values of k , d_{ch}/h_c and the critical dipole heights h_c in Cu and Ni at 0 K, room temperature (RT) and the maximum temperature of the saturation plateau ($T_{\max} \approx 430$ K for Cu and ≈ 600 K for Ni). The calculations were performed on the expanded original (or.) and revisited (rev.) versions of Brown's model [9]. The room temperature values are 295 K for Cu_{BB} (left), 300 K for Cu_{HE} (right) and 293 K for Ni.

$T(K)$	0	300		T_{\max}	
k_B	1	1		1	or.
k_B (Cu)	1.0	0.94	0.93	0.90 ₅	rev.
k_B (Ni)	0.98	0.92		0.86	rev.
d_{ch}/h_c (Cu)	28.6	26.6	26.4	25.4	or.
	28	29.0	29.1	29.5	rev.
d_{ch}/h_c (Ni)	24	23.6		18.8	or.
	24.6	25.0		24.7	rev.
h_c (Cu) nm	15.9 ₅	48.9 ₅		84.1	rev.
h_c (Ni) nm	19.8	47.9		121.5	rev.

Table 3

Rounded values of α_B in Cu and Ni at 0 K, room temperature (RT) and the maximum temperature ($T_{\max} = 430$ K for Cu and 600 K for Ni) in the original (or.) and revisited (rev.) versions of Brown's model [9]. The room temperature values are 295 K for Cu_{BB} (left), 300 K for Cu_{HE} (right) and 293 K for Ni.

$T(\text{K})$	0	300	T_{\max}		
α_B (Cu)	0.49	0.57	0.58	0.62	or.
	0.50	0.51	0.50	0.51	rev.
α_B (Ni)	0.53	0.63	0.74	0.74	or.
	0.52	0.53	0.56	0.56	rev.