

FINITE ELEMENT SIMULATION OF STRAIN LOCALIZATION IN INELASTIC SOLIDS BY ENERGY RELAXATION

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Abstract. A new approach for the treatment of strain localization in inelastic material is proposed. It is based on energy minimization principles associated with micro-structure developments. Shear bands are treated as micro-shearing of rank-one laminates. It is assumed that the thickness of the shear band represented by its volume fraction tends to zero, and the energy inside the shear band is a function of the norm of the strain field. The existence of shear bands in the structure leads to an ill-posed problem which can be solved by means of energy relaxation. The performance of the proposed concept is demonstrated through numerical simulation of tension test under plane strain conditions. Numerical results show that mesh sensitivity can be completely removed.

Keywords: Energy relaxation, shear band, strain localization.

1. INTRODUCTION

Regions of high strain localization by intense shearing are referred to as shear bands. The formation of shear bands is accompanied by a softening response, characterized by a decrease in strength of the material with accumulated inelastic strain, often leading to complete failure [1, 2]. Predictions of the onset and evolution of shear bands play an important role in determining the safety of structures, improving mechanical properties of material and designing material microstructure.

In recent years a new methodology based on *energy relaxation* has been developed to simulate the development of material microstructures [3]-[9]. The advantage of this theory, when applied to the problem at hand, is the natural formation of shear bands based on the energy minimization principles associated with micro-structure developments.

In this paper a theoretical framework for the treatment of shear localizations in inelastic materials at small strains is developed. The theory is based on energy minimization principles associated with micro-structure developments under the assumption of a shear band of zero thickness and of a special form of the energy within the shear band. Shear bands are treated as laminates of first order. The problem of the non-convex energy arising due to the formation of shear bands is solved by energy relaxation in order to ensure

that the corresponding problem is well-posed. Numerical simulations of a tension test are shown in order to evaluate the performance of the proposed concept.

2. SHEAR BANDS AS SPECIAL LAMINATES

In this paper localization phenomena are regarded as micro-structure developments associated with nonconvex potentials. We assume the micro-structures to consist of two domains: a low-strain domain and a high-strain one. Let us consider a so-called RVE (representative volume element) obtained by zooming in on the region around the point A as shown in Fig. 1. The RVE is split into two volume fractions: the volume fraction ξ of the low-strain domain and the volume fraction $(1 - \xi)$ of the high-strain domain.

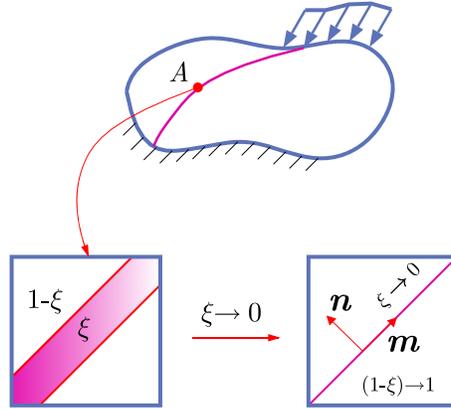


Fig. 1. Shear band is treated as the micro-shearing of a rank-one laminate

When strain localization occurs, the following assumptions related to the volume fraction ξ are made:

(1) After the onset of localization, the width of the shear band represented by the volume fraction ξ tends to zero

$$\xi \rightarrow 0. \quad (1)$$

(2) Most of the deformation is concentrated parallel to the band

$$\mathbf{m} \cdot \mathbf{n} = 0 \quad (2)$$

where \mathbf{m} and \mathbf{n} are two unit vectors giving the direction of shear band evolution.

(3) The potential (energy) inside the shear band W_2 is assumed to satisfy

$$W_2(\xi \boldsymbol{\varepsilon}) = |\xi| W_2(\boldsymbol{\varepsilon}). \quad (3)$$

For example $W_2(\boldsymbol{\varepsilon})$ may be taken in the following form

$$W_2(\boldsymbol{\varepsilon}) = A \|\boldsymbol{\varepsilon}\| \quad (4)$$

where A is a material parameter.

The first assumption is based on the observations that the width of shear bands is normally very small relative to the overall dimension of the system under consideration

[10, 11]. This assumption corresponds to the concept of strong discontinuities proposed in [12, 13] among others.

The second assumption based on the results of the experiment by *Finno et al, 1996* [14]. In their experiment, they showed that the normal movements are much smaller than those in the tangent direction inside the shear band. This assumption implies the two vectors \mathbf{m} and \mathbf{n} being orthogonal.

According to the third assumption, the potential inside a shear band is positive homogeneous of first-order in the strain field (3). We will see later on, that only for this very form of the potential as given in eq. (4), it has the desired property leading to strong discontinuities.

In the following considerations we develop a simple approach to the treatment of shear localization based on these assumptions.

3. TWO-DIMENSIONAL PROBLEM

3.1. Micro-strain

In the two-dimensional problem the micro-strains $\boldsymbol{\varepsilon}_1$ and $\boldsymbol{\varepsilon}_2$ can be written as

$$\boldsymbol{\varepsilon}_1 = \boldsymbol{\varepsilon} - \xi(\mathbf{a} \otimes \mathbf{n})^s \quad (5)$$

$$\boldsymbol{\varepsilon}_2 = \boldsymbol{\varepsilon} + (1 - \xi)(\mathbf{a} \otimes \mathbf{n})^s = \boldsymbol{\varepsilon} - \xi(\mathbf{a} \otimes \mathbf{n})^s + (\mathbf{a} \otimes \mathbf{n})^s \quad (6)$$

where $(\mathbf{a} \otimes \mathbf{n})^s = \frac{1}{2}(\mathbf{a} \otimes \mathbf{n} + \mathbf{n} \otimes \mathbf{a})$.

Let us define s by

$$\xi \mathbf{a} = s \mathbf{m}, \quad (7)$$

where $\|\mathbf{m}\| = 1$. Herein \mathbf{m} and \mathbf{n} are two unit vectors giving the direction of shear band evolution, s is an appropriately rescaled variable.

On inserting eq. (7) into eqs. (5) and (6), we have

$$\boldsymbol{\varepsilon}_1 = \boldsymbol{\varepsilon} - s(\mathbf{m} \otimes \mathbf{n})^s, \quad (8)$$

$$\boldsymbol{\varepsilon}_2 = \boldsymbol{\varepsilon} - s(\mathbf{m} \otimes \mathbf{n})^s + \frac{s}{\xi}(\mathbf{m} \otimes \mathbf{n})^s. \quad (9)$$

As ξ tends to zero $\frac{s}{\xi}$ will grow out of bounds. Thus, eq. (9) can be simplified as

$$\boldsymbol{\varepsilon}_2 \approx \frac{s}{\xi}(\mathbf{m} \otimes \mathbf{n})^s. \quad (10)$$

3.2. Relaxed potential

The mixed potential of the two domains can be written in the following form

$$W^{mix}(\boldsymbol{\varepsilon}) = W_1(\boldsymbol{\varepsilon} - s(\mathbf{m} \otimes \mathbf{n})^s) + |s| W_2((\mathbf{m} \otimes \mathbf{n})^s). \quad (11)$$

The relaxed potential is computed in order to ensure the well-posedness of the boundary-value problem governing the emergence of a shear band. It is obtained by the minimization procedure

$$W_R(\boldsymbol{\varepsilon}) = \inf \{ W^{mix}(\boldsymbol{\varepsilon}) \mid s, \mathbf{m}, \mathbf{n} : \|\mathbf{m}\| = \|\mathbf{n}\| = 1 \}. \quad (12)$$

Table 1. Minimization problem: $\inf_s W(s)$

Scalar minimization problem:	$\inf_s W(s)$
Potential:	$W(s) = as^2 + bs + c s $ with $c > 0, a > 0$
Solution:	$\inf_s W(s) = -\frac{1}{4a}(b - c)_+^2$
Minimizer:	$s = -\frac{1}{2a}(b - c)_+ \text{sign}(b)$
Abbreviations:	$(b - c)_+ = \begin{cases} 0 & \text{for } b \leq c \\ b - c & \text{for } b > c \end{cases}$ $\text{sign}(b) = \frac{ b }{b}$ for $b \neq 0$

Let us consider two specific potentials representing the behaviour at small and large strains, respectively

$$W_1(\boldsymbol{\varepsilon}) = \frac{1}{2} \boldsymbol{\varepsilon} : \boldsymbol{\mathcal{C}} : \boldsymbol{\varepsilon}, \quad (13)$$

$$W_2(\boldsymbol{\gamma}) = A \|\boldsymbol{\gamma}\|, \quad (14)$$

where $\boldsymbol{\mathcal{C}}$ is symmetric fourth-order, positive definite tensor. Substituting (13) and (14) into (11), one gets

$$W^{mix}(\boldsymbol{\varepsilon}) = \frac{1}{2} (\boldsymbol{\varepsilon} - s\boldsymbol{\gamma}) : \boldsymbol{\mathcal{C}} : (\boldsymbol{\varepsilon} - s\boldsymbol{\gamma}) + A |s| \|\boldsymbol{\gamma}\| \quad (15)$$

where $\boldsymbol{\gamma} = (\boldsymbol{m} \otimes \boldsymbol{n})^s$.

On inserting (2) into (15) we have the simplified form of the mixed energy

$$W^{mix}(\boldsymbol{\varepsilon}) = \frac{1}{2} (\boldsymbol{\varepsilon} - s\boldsymbol{\gamma}) : \boldsymbol{\mathcal{C}} : (\boldsymbol{\varepsilon} - s\boldsymbol{\gamma}) + \frac{A}{\sqrt{2}} |s|. \quad (16)$$

Using the results in Table 1, minimization of (16) with respect to s yields

$$s = \frac{\text{sign}(\boldsymbol{\varepsilon} : \boldsymbol{\mathcal{C}} : \boldsymbol{\gamma})}{(\boldsymbol{\gamma} : \boldsymbol{\mathcal{C}} : \boldsymbol{\gamma})} \left[|\boldsymbol{\varepsilon} : \boldsymbol{\mathcal{C}} : \boldsymbol{\gamma}| - \frac{A}{\sqrt{2}} \right]_+, \quad (17)$$

and the potential (16) with s given by (17) takes the form

$$\inf_s W^{mix}(\boldsymbol{\varepsilon}) = \frac{1}{2} \boldsymbol{\varepsilon} : \boldsymbol{\mathcal{C}} : \boldsymbol{\varepsilon} - \frac{1}{2} \frac{\left[|\boldsymbol{\varepsilon} : \boldsymbol{\mathcal{C}} : \boldsymbol{\gamma}| - \frac{A}{\sqrt{2}} \right]_+^2}{\boldsymbol{\gamma} : \boldsymbol{\mathcal{C}} : \boldsymbol{\gamma}}. \quad (18)$$

Herein $a = \frac{1}{2} \boldsymbol{\gamma} : \boldsymbol{\mathcal{C}} : \boldsymbol{\gamma}$, $b = -\boldsymbol{\varepsilon} : \boldsymbol{\mathcal{C}} : \boldsymbol{\gamma}$, $c = \frac{A}{\sqrt{2}}$. Easily one can recognise that a is positive due to the positive definiteness of the fourth-order tensor $\boldsymbol{\mathcal{C}}$.

3.3. Computation of stress and the tangent operator

The stress and the tangent operator are derived from the direct derivative of the relaxed potential (12). The derivative of (12) reads

$$\frac{\partial W_R}{\partial \boldsymbol{\varepsilon}} = \frac{\partial W^{mix}}{\partial \boldsymbol{\varepsilon}} + \frac{\partial W^{mix}}{\partial s} \frac{\partial s}{\partial \boldsymbol{\varepsilon}} + \frac{\partial W^{mix}}{\partial \mathbf{m}} \frac{\partial \mathbf{m}}{\partial \boldsymbol{\varepsilon}} + \frac{\partial W^{mix}}{\partial \mathbf{n}} \frac{\partial \mathbf{n}}{\partial \boldsymbol{\varepsilon}}. \quad (19)$$

It is observed that the three last terms in eq. (19) vanish due to the stationarity condition of the minimization problem (12). Thus, the relaxed stress which is an appropriate average of the two micro-stresses has the form

$$\boldsymbol{\sigma} = \frac{\partial W^{mix}}{\partial \boldsymbol{\varepsilon}}. \quad (20)$$

Considering the form of the potential (18) we obtain

$$\boldsymbol{\sigma} = \mathbf{C} : \boldsymbol{\varepsilon} - s \mathbf{C} : \boldsymbol{\gamma}. \quad (21)$$

The tangent operator is given by

$$\mathbf{A} = \frac{\partial^2 W_R}{\partial \boldsymbol{\varepsilon}^2} = \frac{\partial \boldsymbol{\sigma}}{\partial \boldsymbol{\varepsilon}} = \mathbf{C} - (\mathbf{C} : \boldsymbol{\gamma}) \otimes \frac{\partial s}{\partial \boldsymbol{\varepsilon}} - s \frac{\partial (\mathbf{C} : \boldsymbol{\gamma})}{\partial \boldsymbol{\varepsilon}}. \quad (22)$$

3.4. Localization criterion

In the derivations above a central role is played by the quantity

$$L = \left[|\boldsymbol{\varepsilon} : \mathbf{C} : \boldsymbol{\gamma}| - \frac{A}{\sqrt{2}} \right]. \quad (23)$$

As the process of deformation progresses, L may be negative, zero or positive. A positive value in turn signals the onset of localization. A criterion that can be shown to be equivalent to the well-known notion of loss of ellipticity.

i. $L \leq 0$: we have $s = 0$. The relaxed potential $W_R(\boldsymbol{\varepsilon})$ reduces to the elastic strain energy $W_1(\boldsymbol{\varepsilon})$.

ii. $L > 0$: we have $s \neq 0$. A shear band starts to develop. The homogeneous deformation $\boldsymbol{\varepsilon}$ decomposes into the two micro-strains $\boldsymbol{\varepsilon}_1$ and $\boldsymbol{\varepsilon}_2$. The nonconvex potential energy W^{mix} is replaced with the approximated rank-one convexification $W_R(\boldsymbol{\varepsilon})$ to ensure well-posedness of the problem.

4. APPLICATION OF RELAXATION THEORY TO LINEAR ISOTROPIC MATERIAL

The fourth-order isotropic elastic tensor \mathbf{C} takes the form

$$C_{ijkl} = \lambda \delta_{ij} \delta_{kl} + \mu (\delta_{ik} \delta_{jl} + \delta_{il} \delta_{jk}), \quad (24)$$

or in the tensor notation

$$\mathbf{C} = \lambda \mathbf{I} \otimes \mathbf{I} + \mu (\mathbf{I} + \bar{\mathbf{I}}), \quad (25)$$

where λ and μ are Lamé constants.

The simplified form of the mixed energy is obtained by substituting (24) into eq. (16)

$$W^{mix}(\boldsymbol{\epsilon}) = \frac{\lambda}{2}(\text{tr}\boldsymbol{\epsilon})^2 + \mu \left(\|\boldsymbol{\epsilon}\|^2 - 2s \mathbf{m} \cdot \boldsymbol{\epsilon} \mathbf{n} + \frac{s^2}{2} \right) + \frac{A}{\sqrt{2}} |s|. \quad (26)$$

Based on the minimization procedure the relaxed potential (12) reads

$$W_R = \inf_{s, \mathbf{m}, \mathbf{n}} W^{mix} = \frac{\lambda}{2} \text{tr}(\boldsymbol{\epsilon})^2 + \mu \|\boldsymbol{\epsilon}\|^2 - 2\mu \left[|\epsilon_{mn}| - \frac{\alpha}{2} \right]_+^2, \quad (27)$$

where $\alpha = \frac{A}{\sqrt{2}\mu}$, $|\epsilon_{mn}|$ is the maximum shear strain.

Solution s yields

$$s = (2|\epsilon_{mn}| - \alpha)_+ \text{sign}(\epsilon_{mn}) = \begin{cases} 0 & \text{if } |\epsilon_{mn}| < \alpha/2 \\ (2|\epsilon_{mn}| - \alpha) \text{sign}(\epsilon_{mn}) & \text{if } |\epsilon_{mn}| \geq \alpha/2 \end{cases} \quad (28)$$

The relaxed stress is obtained from eqs. (21) and (27)

$$\boldsymbol{\sigma} = \lambda \text{tr}\boldsymbol{\epsilon} \mathbf{I} + 2\mu\boldsymbol{\epsilon} - 2\mu s (\mathbf{m} \otimes \mathbf{n})^s. \quad (29)$$

On inserting (27) into (22) we get the tangent operator

$$\mathcal{A} = \lambda \mathbf{I} \otimes \mathbf{I} + \mu(\mathcal{I} + \bar{\mathcal{I}}) - 2\mu \left[(\mathbf{m} \otimes \mathbf{n})^s \otimes \frac{\partial s}{\partial \boldsymbol{\epsilon}} + s \frac{\partial (\mathbf{m} \otimes \mathbf{n})^s}{\partial \boldsymbol{\epsilon}} \right]. \quad (30)$$

The capability of the proposed model is demonstrated through numerical simulation of a tension test in the next Section.

5. TENSION TEST

In this section the model presented in Section 4 are implemented into the finite element code FEAP [15]. In what follows we investigate a tension test under plane strain conditions. Different kinds of elements are used in this example as summarized in Fig. 2.

Element type	Method	Number of Gauss points	Notation
	MES ¹	2x2	MES
	EAS ²	2x2	EAS
	QM6 ³	2x2	QM6
	Disp ⁴	2x2	Q4

Fig. 2. Notation of element type: MES¹ - the mixed enhanced strain method [16], EAS² - the enhanced assumed strain method [17], QM6³ - the method of incompatible mode [18], Disp⁴ - the displacement method [19]

A sample is subjected to a prescribed vertical displacement under plane strain conditions. The geometry of the specimen and the boundary conditions imposed on the displacements as well as the material parameters are given in Fig. 3. In order to trigger the shear band formation a geometrical imperfection along the height of specimen is introduced.

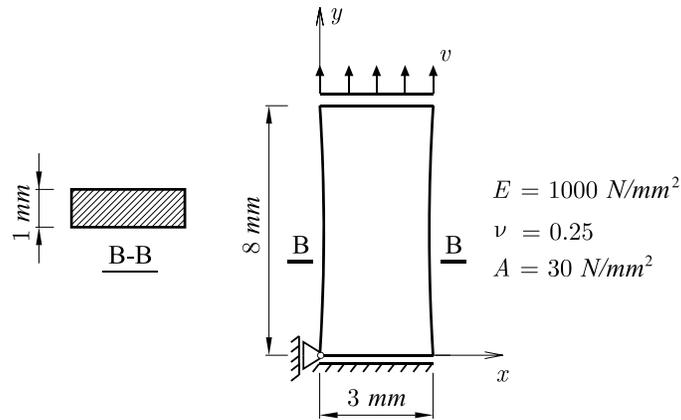


Fig. 3. Localization in tension. Geometry and boundary conditions

Four discretizations of the domain are considered: 3x8, 7x18, 14x36, 21x54 elements. The load-displacement curves behave identically for $v < 0.3126\text{mm}$ as shown in Figs. 4a and 5a. The localization is delayed until the maximum shear strain is equal to $\frac{\alpha}{2}$ corresponding to $v \approx 0.3126\text{mm}$, then the performance of diverse finite element methods starts to differ. The localization band is straight, at 45° with the direction of principal strain.

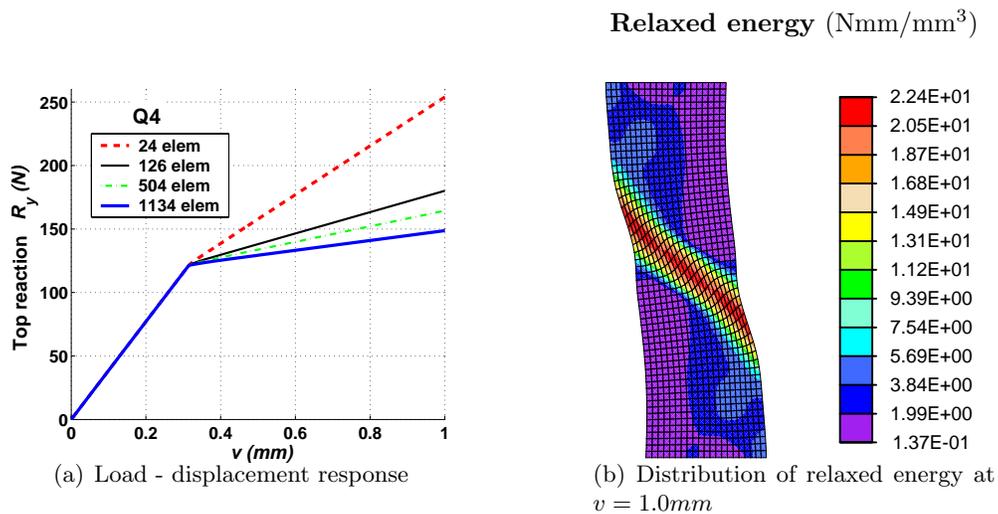


Fig. 4. Localization in tension by use of Q4 element (the displacement method)

The displacement method is unable to capture the localization and shows hardening behaviour as depicted in Fig. 4a and 4b.

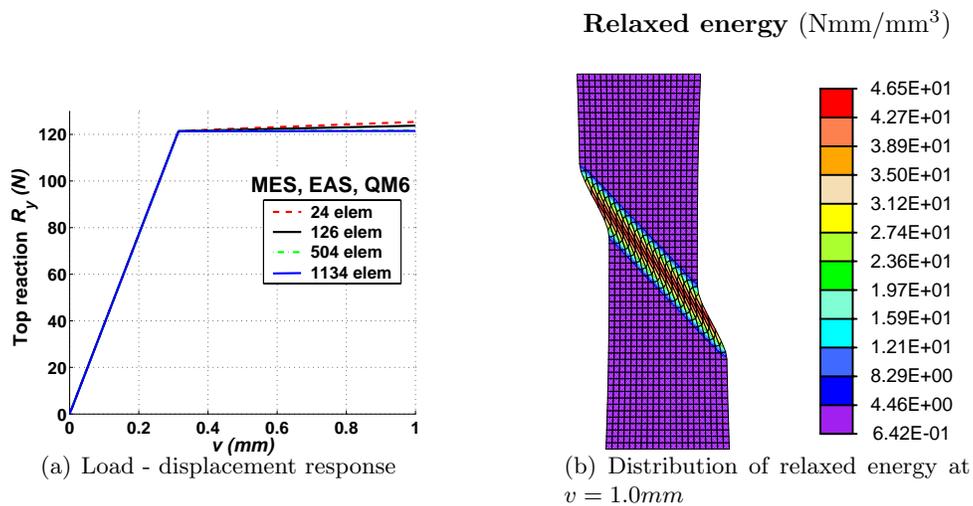


Fig. 5. Localization in tension by use of MES element (the mixed enhanced strain method), EAS element (the enhanced assumed strain method) and QM6 element (the incompatible mode method)

As we expected in Fig. 5a, the solutions obtained by using MES, EAS and QM6 elements are practically coincident in term of vertical reactions, thus showing mesh independence of the proposed relaxed theory. The distribution of the relaxed energy shown in Fig. 5b points out the localized elements.

The behaviour of the relaxed potential as well as the relaxed vertical stress σ_y of the element 356 at the first gauss point inside the shear band is depicted in Fig. 6. After the

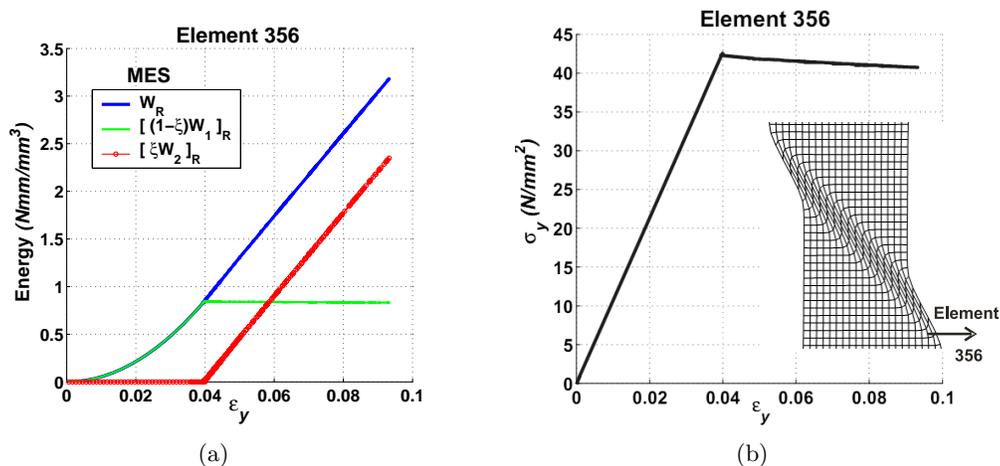


Fig. 6. Relaxed energy and relaxed normal stress σ_y by use of MES element (the mixed enhanced strain method) at the first Gauss point of the element 356

bifurcation point, the relaxed potential of the small-strain domain approaches asymptotically a constant value, whereas the relaxed potential of the large-strain domain is active inside the shear band only and increases continuously with the prescribed displacement v . As a result, the relaxed stress approaches nearly a constant value.

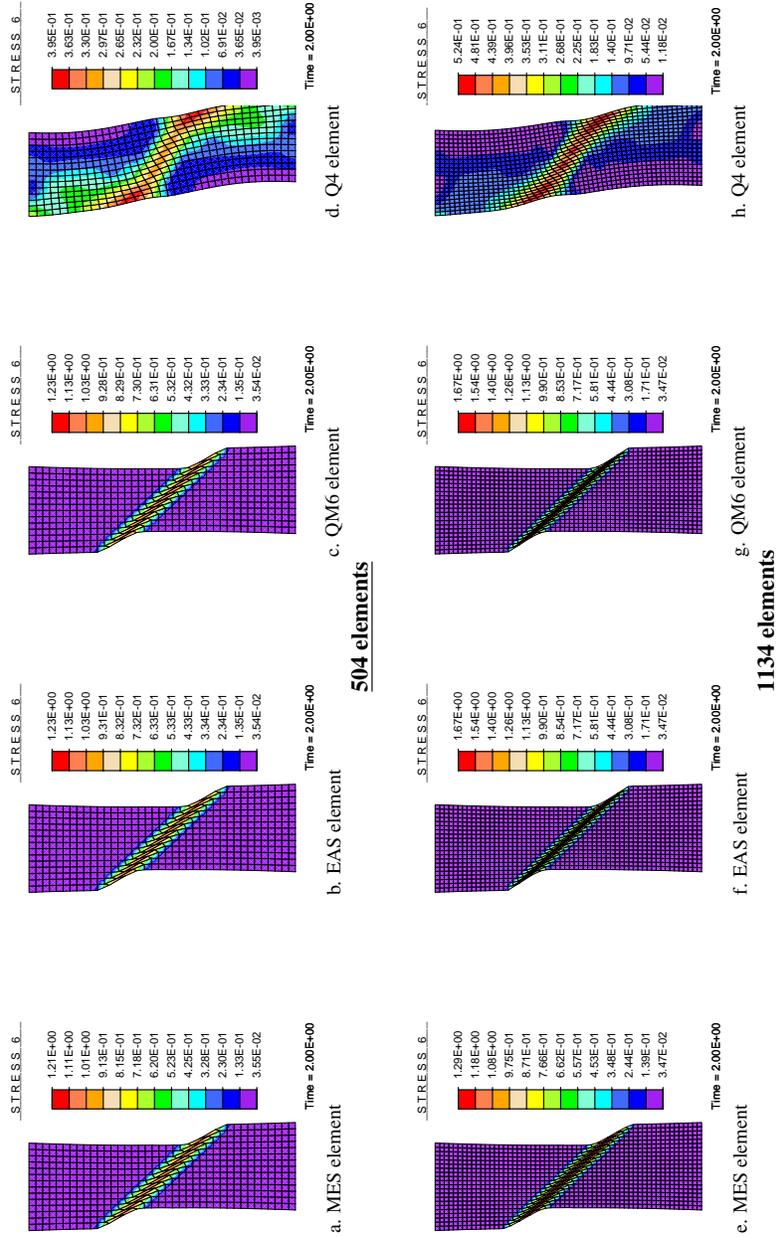


Fig. 7. Localization in tension. Comparison of the different finite element methods: Distribution of shear strain at $v = 1.0 \text{ mm}$

It is very interesting to see the performance of the different finite element formulations by considering the deformation shape of sample (Fig. 7). It is observed that the solution obtained by use of the Q4 element fails to capture effects of strain localization.

6. CONCLUSIONS

The paper focuses on a new approach for the treatment of strain localization in solid materials. The theory is based on minimization principles associated with micro-structure developments under the assumptions of a shear band of a zero thickness and the presence of a mixed potential inside the shear band.

Localization phenomena are regarded as micro-structure developments associated with nonconvex potentials. The nonconvexity of the mixed potential occurring due to the formation of strain localization is resolved by relaxation in order to ensure the well-posedness of the associated boundary value problem. The relaxed potential is obtained via local minimization problem of the mixed potential. The onset of localization is detected through the proposed optimization process. Numerical results show clearly mesh-independent behavior in the sense that shear bands are as narrowly as the mesh-resolution allows, while all other features of the solution are independent of the chosen discretization.

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