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C. Engwer, I.S. Pop, T. Wick

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Dynamic and weighted stabilizations of the L -scheme applied to a phase-field model for fracture propagation

Christian Engwer, Iuliu Sorin Pop, Thomas Wick

Abstract We consider a phase-field fracture propagation model, which consists of two (nonlinear) coupled partial differential equations. The first equation describes the displacement evolution, and the second is a smoothed indicator variable, describing the crack position. We propose an iterative scheme, the so-called L -scheme, with a dynamic update of the stabilization parameters during the iterations. Our algorithmic improvements are substantiated with two numerical tests. The dynamic adjustments of the stabilization parameters lead to a significant reduction of iteration numbers in comparison to constant stabilization values.

1 Introduction

This work is an extension of [3] in which an L -type iterative scheme (see [5, 8]) with stabilizing parameters for solving phase-field fracture problems was proposed. In [3], the stabilization parameters were chosen as constants throughout an entire computation. With these choices, the convergence of the scheme has been proven rigorously. The resulting approach performs well in the sense that an unlimited number of iterations compared to a truncated scheme yields the same numerical solution. The results were validated by investigating the load-displacements curves.

Christian Engwer
Institut für Numerische und Angewandte Mathematik Fachbereich Mathematik und Informatik der
Universität Münster Einsteinstrasse 62 48149 Münster, Germany, e-mail: christian.engwer@uni-
muenster.de

Iuliu Sorin Pop
Universiteit Hasselt, Faculty of Sciences, Agoralaan Gebouw D - B-3590 Diepenbeek, Belgium,
e-mail: sorin.pop@uhasselt.be

Thomas Wick
Leibniz Universität Hannover, Institut für Angewandte Mathematik, Welfengarten 1, 30167 Han-
nover, Germany, e-mail: thomas.wick@ifam.uni-hannover.de

Moreover, the robustness of the scheme w.r.t. spatial mesh refinement was shown. Nonetheless, the iteration numbers (for an unlimited number of iterations) remained high.

In this work, we propose and compare two extensions of the aforementioned scheme. First, we update the L scheme parameters dynamically. Second, we use an adaptive weight depending on the fracture location inside the domain. For the latter idea, we use the phase-field variable to weight L locally.

The outline of this work is as follows: In Section 2 the model is stated whereas Section 3 presents the dynamic choice of the stabilization parameters. In Section 4, we present two numerical tests to study the performance of the proposed scheme.

2 The phase-field fracture model

We consider an elliptic problem stemming from the crack propagation model proposed in [3]. $\Omega \subset \mathbb{R}^d$ is a d -dimensional, polygonal and bounded domain. We use the spaces $W^{1,\infty}(\Omega)$, containing functions having essentially bounded weak derivatives in any direction, and $H_0^1(\Omega)$ containing functions vanishing at the boundary of Ω (in the sense of traces) and having square integrable weak derivatives. (\cdot, \cdot) stands for the $L^2(\Omega)$ inner product. For the ease of writing we use the notations $V := (H_0^1(\Omega))^d$ and $W := W^{1,\infty}(\Omega)$. The vector-valued displacements are denoted by u . For modeling fracture propagation in Ω , a phase field variable φ is used. This approximates the characteristic function of the intact region of Ω . Written in weak form, we solve the following problems iteratively

- **Problem 1ⁱ**: Given $(u^{i-1}, \varphi^{i-1}) \in V \times W$, find $u^i \in V$ s.t. for all $v \in V$

$$a_u(u^i, v) := L_u(u^i - u^{i-1}, v) + \left(g(\varphi^{i-1}) \sigma^+(u^i), \mathbf{e}(v) \right) + \left(\sigma^-(u^i), \mathbf{e}(v) \right) = 0. \quad (1)$$
- **Problem 2ⁱ**: Given $(\varphi^{i-1}, u^i, \bar{\varphi}) \in W \times V \times W$, find $\varphi^i \in W$ s.t. for all $\psi \in W$

$$a_\varphi(\varphi^i, \psi) := L_\varphi(\varphi^i - \varphi^{i-1}, \psi) + G_c \varepsilon(\nabla \varphi^i, \nabla \psi) - \frac{G_c}{\varepsilon} (1 - \varphi^i, \psi) + (1 - \kappa)(\varphi^i \sigma^+(u^i) : \mathbf{e}(u^i), \psi) + (\Xi + \gamma[\varphi^i - \bar{\varphi}]^+, \psi) = 0. \quad (2)$$

In case of convergence, the first terms in the above are vanishing, and the limit pair $(u, \varphi) \in V \times W$ solves a time discrete counterpart of the model in [3], if $\bar{\varphi}$ is interpreted as the phase field at the previous time step. In this context, with $\Xi \in L^2(\Omega)$ and $\gamma > 0$, the last term in (2) is the augmented Lagrangian penalization proposed in [9] for the irreversibility constraint of the fracture propagation.

Furthermore, in the above, ε is a (small) phase-field regularization parameter, $G_c > 0$ is the critical elastic energy restitution rate, and $0 < \kappa \ll 1$ is a regularization parameter used to avoid the degeneracy of the elastic energy. The latter is similar to replacing the fracture with a softer material. Next, $g(\varphi) := (1 - \kappa)\varphi^2 + \kappa$ is the degradation function, and $\mathbf{e} := \frac{1}{2}(\nabla u + \nabla u^T)$ is the strain tensor.

The stress tensor in the above is split into a tensile and compressive part,

$$\boldsymbol{\sigma}^+ := 2\mu_s \mathbf{e}^+ + \lambda_s [\text{tr}(\mathbf{e})]^+ I, \quad \boldsymbol{\sigma}^- := 2\mu_s (\mathbf{e} - \mathbf{e}^+) + \lambda_s (\text{tr}(\mathbf{e}) - [\text{tr}(\mathbf{e})]^+) I,$$

where $[\cdot]^+$ stands for the positive cut of the argument. Further, $\mathbf{e}^+ = \mathbf{P}\boldsymbol{\Lambda}^+\mathbf{P}^T$, with \mathbf{P} being the matrix containing the unit eigenvectors corresponding to the eigenvalues of the strain tensor \mathbf{e} . In particular, for $d = 2$ one has $\mathbf{P} = [v_1, v_2]$ and

$$\boldsymbol{\Lambda}^+ := \boldsymbol{\Lambda}^+(u) := \begin{pmatrix} [\lambda_1(u)]^+ & 0 \\ 0 & [\lambda_2(u)]^+ \end{pmatrix}.$$

3 The L -scheme with dynamic updates of the stabilization parameters

The iteration (1)-(2) is essentially the scheme proposed in [3], in which the stabilization parameters L_u and L_φ are taken constant. To improve the convergence behaviour of the scheme, we propose a dynamic update of these parameters.

Dynamic update at each iteration / constant in space: The iteration discussed in [3] uses constant parameters L_u and L_φ . With this choice, the convergence has been proved rigorously. However, the number of iterations can remain high. High iteration numbers for phase-field fracture problems were also reported in [4, 10]. To improve the efficiency, we suggest in this work to update L_u and L_φ at each iteration i :

$$L_i = a(i)L_{i-1}, \quad \text{where } L_i := L_{u,i} = L_{\varphi,i}.$$

Inspired by numerical continuation methods in e.g. [1], one would naturally choose a large L_0 and $a(i) := a < 1$ to obtain a decreasing sequence $L_0 > L_1 > L_2 > \dots$, updated until a lower bound L_- is reached. However, this seems not to be a good choice in phase-field fracture since the system does not have a unique solution. Consequently, with increasing i the iterations would oscillate in approaching one or another solution, and the algorithm convergence deteriorates. For this reason, we propose the other way around: the closer the iteration is to some solution, the larger the stabilization parameters is chosen, so that the iterations remain close to this solution. We choose $a(i) := a > 1$, yielding $L_0 < L_1 < L_2 < \dots$ up to a maximal L_* .

On the specific choice of the parameters: A possible choice for a is $a(i) := 5^i$ ($i = 0, 1, 2, \dots$), while $L_0 := 10^{-10}$. This heuristic choice and may be improved by using the solution within the iteration procedure, or a-posteriori error estimates for the iteration error. Moreover, $a(i) := 5^i$ is motivated as follows. Higher values greater than 5 would emphasize too much the stabilization. On the other hand, too low values, do not lead to any significant enhancement of the convergence behaviour. We substantiate these claims by also using $a(i) = 10^i$ and $a(i) = 20^i$ in our computations.

Dynamic update using the iteration: An extension of the strategy is to adapt the L -scheme parameters in space by using the phase-field variable $\varphi^{n,i-1}$. We still take $L_i = aL_{i-1}$, but now $a := a(i, \varphi^{n,i-1})$. Away from the fracture, we have $\varphi \approx 1$ and

essentially only the elasticity component (2) is being solved. On the other hand, the stabilization is important in the fracture region, for which we take

$$L_i = a(i, \varphi^{i-1})L_{i-1}, \quad \text{with } a(i, \varphi^{i-1}) := (1 - \varphi^{i-1})a.$$

Recalling that the fracture is characterised by $\varphi \approx 0$, it becomes clear that the stabilization parameters are acting mainly in the fracture region. Finally, to improve further the convergence behaviour of the scheme we adapt Ξ at each iteration. In this case we take $\Xi_i = \Xi_{i-1} + \gamma[\varphi^{i-1} - \bar{\varphi}]^+$.

Algorithm 1 Dynamic variant of the L-scheme for a phase-field fracture

Choose $\gamma > 0$, $a > 1$, as well as Ξ^0 and L_0 . Set $i = 0$.

repeat

 Let $i = i + 1$;

 Solve the two problems, namely

 Solve the nonlinear elasticity problem in (1)

 Solve the nonlinear phase-field problem in (2)

 Update $L^i = aL^{i-1}$

 Update $\Xi^i = \Xi^{i-1} + \gamma[\varphi^i - \bar{\varphi}]^+$

until

$\max\{\|a_u(u^i, v)\|, \|a_\varphi(\varphi^i, \psi)\|/v \in V, \psi \in W\} \leq \text{TOL}$,

The final algorithm: The algorithm is based on the iterative procedure for phase-field fracture originally proposed in [9]. Therein, the inequality constraint is realized by an augmented Lagrangian iteration. Within this loop we update the L scheme parameters too. The resulting is sketched in Algorithm 3, in which $\text{TOL} = 10^{-6}$ is taken, and $L = L_u = L_\varphi$.

Remark 1 For the solution of both nonlinear subproblems (1) and (2), we use a monotonicity-based Newton method (details see e.g., in [10]) with the tolerance 10^{-8} . Inside Newton's method, we solve the linear systems with a direct solver.

4 Numerical tests

We consider two test examples. Details for the first test can be found in [7]. The setup of the second test can be found for instance in [6]. Both examples were already computed in [3] and the results therein are compared to the ones obtained here. The scheme is implemented in a code based on the deal.II library [2].

Single edge notched shear test: The configuration is shown in Figure 1. Specifically, we use $\mu_s = 80.77 \text{ kN/mm}^2$, $\lambda_s = 121.15 \text{ kN/mm}^2$, and $G_c = 2.7 \text{ N/mm}$. The crack growth is driven by a non-homogeneous Dirichlet condition for the displacement field on Γ_{top} , the top boundary of B . We increase the displacement on Γ_{top} over time, namely we apply non-homogeneous Dirichlet conditions:

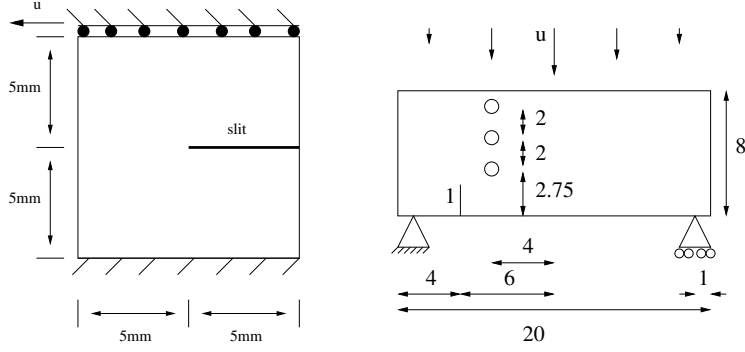


Fig. 1 Examples 1 and 2. The following conditions are prescribed: on the left and right boundaries, $u_y = 0$ mm and traction-free in x -direction. On the bottom part, $u_x = u_y = 0$ mm. On Γ_{top} , $u_y = 0$ mm and u_x is as stated in (3). Finally, the lower part of the slit is fixed in y -direction, i.e., $u_y = 0$ mm. Right: Asymmetric notched three point bending test. The three holes have each a diameter of 0.5. All units are in mm .

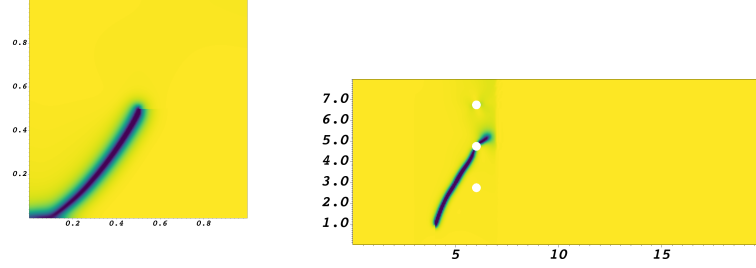


Fig. 2 Examples 1 and 2. Numerical solutions on the finest meshes and at the end time. The cracks are displayed in dark blue color.

$$u_x = t\bar{u}, \quad \bar{u} = 1 \text{ mm/s}, \quad (3)$$

where t denotes the current loading time. Furthermore, we set $\kappa = 10^{-10}$ [mm] and $\varepsilon = 2h$ [mm]. We evaluate the surface load vector on the Γ_{top} as

$$\tau = (F_x, F_y) := \int_{\Gamma_{\text{top}}} \sigma(u) \nu ds, \quad (4)$$

with normal vector ν , and we are particularly interested in the shear force F_x . Three different meshes with 1024 (Ref. 4), 4096 (Ref. 5) and 16384 (Ref. 6) elements are observed in order to show the robustness of the proposed schemes. The results are shown in Figure 6.

Our findings are summarized in Figure 3. The numerical solutions for all four different strategies for choosing L are practically identical, only the number of iterations being different. Here, $L = 0$ and $L = 1e - 2$ denote tests in which $L = L_u = L_\varphi$ are taken constant throughout the entire computation. The newly

proposed dynamic versions are denoted by *L dynamic* and *L dyn. weighted*. We observe a significant reduction in the computational cost when using the dynamic *L*-schemes. The maximum number of iterations is 21 for both the weighted version and the spatially-constant *L*-scheme. This number is reduced to 12 iterations using $a = 20$ while the accuracy only slightly changes.

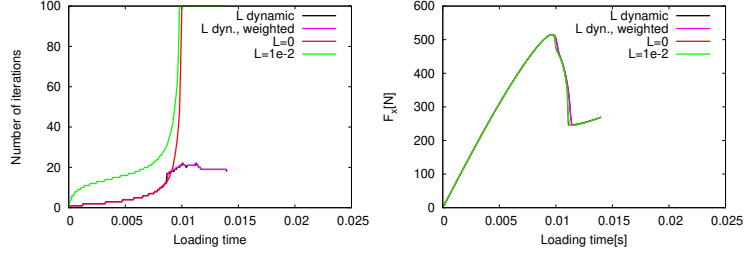


Fig. 3 Example 1. Comparison of dynamic *L* updates, the weighted version, and constant *L*. Left: number of iterations. Right: load-displacement curves.

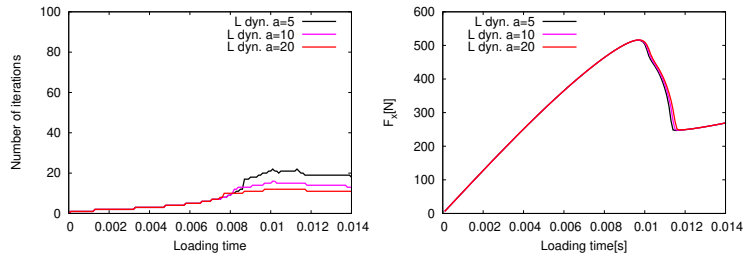


Fig. 4 Example 1. Comparison of different α for the dynamic *L* scheme.

Asymmetrically notched three point bending test: The configuration is shown in Figure 1 (right). The initial mesh is 3, 4 and 5 times uniformly refined, yielding 3904, 15616 and 62464 mesh elements with the minimal mesh size parameter $h_3 = 0.135$, $h_4 = 0.066$ and $h_5 = 0.033$. As material parameters, we use $\mu_s = 8 \text{ kN/mm}^2$, $\lambda_s = 12 \text{ kN/mm}^2$, and $G_c = 1 \times 10^{-3} \text{ kN/mm}$. Furthermore, we set $k = 10^{-10} h [\text{mm}]$ and $\varepsilon = 2h$.

Figure 5 presents the number of iterations and the load-displacement curves. The number of iterations is decreasing from 500 (in the figures cut to 100) for the classical L-scheme, to a maximum of 25 when using the dynamic updates. The choice of weighting does not seem to have a significant influence on the number of iterations though. The crack starts growing a bit later when using the dynamic updates, which can be inferred from the right plot in Figure 5. Thus, the stabilization parameters have a slight influence on the physical solution. This can be explained in the following way. In regions where $\varphi = 0$ the solution component u is not uniquely defined. This leads to a sub-optimal convergence behaviour of the L-scheme. With the dynamic L-scheme we regain uniqueness, but at the cost of a slightly modified physical problem.

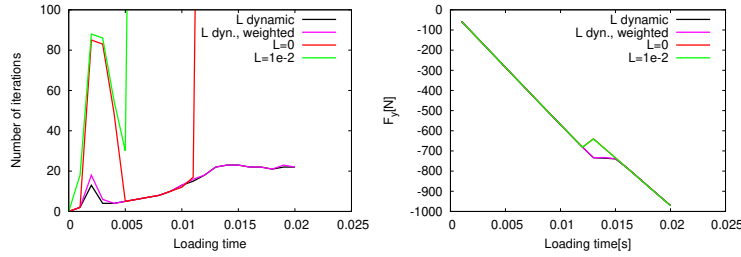


Fig. 5 Example 2: Left: The number of iterations for the different schemes; the results for $L = 0$ and $L = 1e - 2$ are taken from [3]. Right: The load-displacement curves; a slight difference can be observed in the results, indicating that the dynamic updates lead to a slight delay in the prediction of the starting time for the fracture growth.

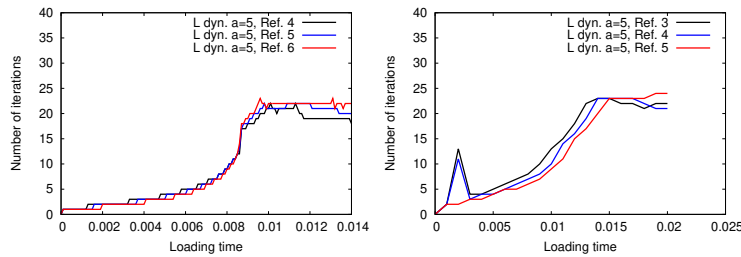


Fig. 6 Examples 1 and 2 for the dynamic L scheme using $\alpha = 5$; three different mesh levels are used in order to verify the robustness of the proposed scheme. The results indicate that the mesh size does not influence the number of the iterations.

Remark 2 Noteworthy, the number of iterations for the dynamic L-scheme is robust with respect to the mesh refinement, as shown in Figure 6. This is in line with the analysis in [3, 5, 8], where it is proved that the convergence rate does not depend on the spatial discretization.

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