Staggered fracture-elasticity solver

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1 Introduction

This document describes the weak form and associated linearized equations for the staggered fracture-elasticity solver implemented in *IFEM*, based on the papers by Borden *et al.* [1], and Gerasimov and De Lorenzis [2].

2 Energy functionals

The energy functional for the quasi-static brittle fracture problem, using a phase-field to represent the crack geometry is in [1] given as follows (see Equation (17) therein):

$$E_{c}(\boldsymbol{u},c) = \int_{\Omega} \left[g(c) \Psi_{0}^{+}(\boldsymbol{\epsilon}(\boldsymbol{u})) + \Psi_{0}^{-}(\boldsymbol{\epsilon}(\boldsymbol{u})) \right] dV$$

$$+ G_{c} \int_{\Omega} \left(\frac{1}{4\ell_{0}} (1-c)^{2} + \ell_{0} \boldsymbol{\nabla} c \cdot \boldsymbol{\nabla} c \right) dV$$

$$(1)$$

where c is the phase field describing the crack, i.e., it has the value 1.0 where the material is undamaged and equal to 0.0 for fully cracked material. $g(c) = (1 - k)c^2 + k$ is the stress degradation function used to scale down the tensile part of the strain energy density in the elasticity equation. $k \ge 0.0$ is a (small) stability parameter that can be used to improve the conditioning of the resulting linear equation system.

In [2], an alternative formulation is used, where d=1-c is used as the unknown phase field instead of c, i.e., it has the value 0.0 in the undamaged material and 1.0 in the fully cracked material. Moreover, a penalty term is introduced to ensure crack irreversibility, and the length scale parameter ℓ_0 is replaced by $\ell=2\ell_0$. The energy functional therefore reads as follows (see also Equation (3) in [2]):

$$E_{d}(\boldsymbol{u}, d) = \int_{\Omega} \left[g(d) \Psi_{0}^{+}(\boldsymbol{\epsilon}(\boldsymbol{u})) + \Psi_{0}^{-}(\boldsymbol{\epsilon}(\boldsymbol{u})) \right] dV$$

$$+ G_{c} \int_{\Omega} \left(\frac{1}{2\ell} d^{2} + \frac{\ell}{2} \boldsymbol{\nabla} d \cdot \boldsymbol{\nabla} d \right) dV + \frac{1}{2\gamma} \int_{CR_{l-1}} (1 - d)^{2} dV$$
(2)

where the stress degradation function now reads $g(d) = (1-k)(1-d)^2 + k$.

The directional derivatives of E_c and E_d read, respectively:

$$E'_{c}(\boldsymbol{u}, c; \boldsymbol{v}, w) := \frac{\partial E_{c}}{\partial \boldsymbol{u}} \cdot \boldsymbol{v} + \frac{\partial E_{c}}{\partial c} w$$

$$= \int_{\Omega} \left[g(c) \frac{\partial \Psi_{0}^{+}}{\partial \boldsymbol{\epsilon}} (\boldsymbol{\epsilon}(\boldsymbol{u})) + \frac{\partial \Psi_{0}^{-}}{\partial \boldsymbol{\epsilon}} (\boldsymbol{\epsilon}(\boldsymbol{u})) \right] : \boldsymbol{\epsilon}(\boldsymbol{v}) \, dV$$

$$+ \int_{\Omega} g'(c) \Psi_{0}^{+}(\boldsymbol{\epsilon}(\boldsymbol{u})) w \, dV + G_{c} \int_{\Omega} \left(\frac{1}{2\ell_{0}} (c - 1) w + 2\ell_{0} \boldsymbol{\nabla} c \cdot \boldsymbol{\nabla} w \right) \, dV$$
(3)

with g'(c) = 2(1 - k)c, and

$$E'_{d}(\boldsymbol{u}, d; \boldsymbol{v}, w) := \frac{\partial E_{d}}{\partial \boldsymbol{u}} \cdot \boldsymbol{v} + \frac{\partial E_{d}}{\partial d} w$$

$$= \int_{\Omega} \left[g(d) \frac{\partial \Psi_{0}^{+}}{\partial \boldsymbol{\epsilon}} (\boldsymbol{\epsilon}(\boldsymbol{u})) + \frac{\partial \Psi_{0}^{-}}{\partial \boldsymbol{\epsilon}} (\boldsymbol{\epsilon}(\boldsymbol{u})) \right] : \boldsymbol{\epsilon}(\boldsymbol{v}) \, dV$$

$$+ \int_{\Omega} g'(d) \Psi_{0}^{+}(\boldsymbol{\epsilon}(\boldsymbol{u})) w \, dV + G_{c} \int_{\Omega} \left(\frac{1}{\ell} dw + \ell \boldsymbol{\nabla} d \cdot \boldsymbol{\nabla} w \right) \, dV + \frac{1}{\gamma} \int_{CR_{l-1}} (d-1) w \, dV$$

$$(4)$$

with g'(d) = 2(1-k)(d-1). The problem is then solved by seeking a solution $(\boldsymbol{u}, c) \in \boldsymbol{V}_1 \times H^1(\Omega, [0, 1])$, such that

$$E'_c(\boldsymbol{u}, c; \boldsymbol{v}, w) = 0 \quad \forall \ (\boldsymbol{v}, w) \in \boldsymbol{V}_0 \times H^1(\Omega, [0, 1])$$
(5)

or alternatively, a solution $(\boldsymbol{u},d) \in \boldsymbol{V}_1 \times H^1(\Omega,[0,1])$, such that

$$E'_d(\boldsymbol{u}, d; \boldsymbol{v}, w) = 0 \quad \forall (\boldsymbol{v}, w) \in \boldsymbol{V}_0 \times H^1(\Omega, [0, 1])$$
(6)

3 Residual and tangent for the c-formulation

Equation (5) is rewritten into a system of coupled equations:

$$Q_{1}(\boldsymbol{u}, c; \boldsymbol{v}) = \int_{\Omega} \left[g(c) \frac{\partial \Psi_{0}^{+}}{\partial \boldsymbol{\epsilon}} (\boldsymbol{\epsilon}(\boldsymbol{u})) + \frac{\partial \Psi_{0}^{-}}{\partial \boldsymbol{\epsilon}} (\boldsymbol{\epsilon}(\boldsymbol{u})) \right] : \boldsymbol{\epsilon}(\boldsymbol{v}) \, dV = 0$$
 (7)

$$Q_2(\boldsymbol{u}, c; w) = \int_{\Omega} \left[g'(c) \Psi_0^+(\boldsymbol{\epsilon}(\boldsymbol{u})) w + G_c \left(\frac{1}{2\ell_0} (c - 1) w + 2\ell_0 \boldsymbol{\nabla} c \cdot \boldsymbol{\nabla} w \right) \right] dV = 0$$
 (8)

To ensure crack irreversibility, the tensile strain energy density Ψ_0^+ in Equation (8) is replaced by a history field $\mathcal{H}(\boldsymbol{x},t)$, with the property

$$\mathcal{H}(\boldsymbol{x}, t_0) = \frac{G_c}{4\ell_0} \left(\frac{1}{C} - 1 \right) \left(1 - \max \left\{ \frac{\delta(\boldsymbol{x}, \Gamma)}{\ell_0}, 1 \right\} \right)$$
(9)

$$\mathcal{H}(\boldsymbol{x}, t_n) = \max \left\{ \mathcal{H}(\boldsymbol{x}, t_{n-1}), \Psi_0^+(\boldsymbol{\epsilon}(\boldsymbol{u}(\boldsymbol{x}, t_n))) \right\}, n > 0$$
 (10)

where C is a constant and $\delta(\boldsymbol{x}, \Gamma)$ denotes the shortest distance from the point \boldsymbol{x} to the initial crack Γ . Moreover, Equation (8) is scaled by $\frac{\ell}{G_c} = \frac{2\ell_0}{G_c}$ to obtain

$$Q_2(\boldsymbol{u}, c; w) = \int_{\Omega} \left[\frac{g'(c)\ell}{G_c} \mathcal{H}(\boldsymbol{x}, t) w + (c - 1)w + \ell^2 \nabla c \cdot \nabla w \right] dV = 0$$
 (11)

The Equations (7) and (11) are solved in a staggered manner, where in the former the phase field c is assumed known and in the latter the displacement field u is assumed known. We then do a linearization of the functionals Q_1 and Q_2 about a certain known configuration (u_0, c_0) for the unknown variables, i.e.

$$Q_1(\boldsymbol{u}_0 + \Delta \boldsymbol{u}, c_0; \boldsymbol{v}) \approx Q_1(\boldsymbol{u}_0, c_0; \boldsymbol{v}) + \frac{\partial Q_1}{\partial \boldsymbol{u}} \cdot \Delta \boldsymbol{u} = 0$$
 (12)

$$Q_2(\boldsymbol{u}_0, c_0 + \Delta c; \boldsymbol{w}) \approx Q_2(\boldsymbol{u}_0, c_0; \boldsymbol{v}) + \frac{\partial Q_2}{\partial c} \Delta c = 0$$
(13)

from which we obtain the tangent operators

$$\frac{\partial Q_1}{\partial \boldsymbol{u}} \cdot \Delta \boldsymbol{u} = \int_{\Omega} \boldsymbol{\epsilon}(\Delta \boldsymbol{u}) : \left[g(c) \frac{\partial^2 \Psi_0^+}{\partial \boldsymbol{\epsilon}^2} (\boldsymbol{\epsilon}(\boldsymbol{u})) + \frac{\partial^2 \Psi_0^-}{\partial \boldsymbol{\epsilon}^2} (\boldsymbol{\epsilon}(\boldsymbol{u})) \right] : \boldsymbol{\epsilon}(\boldsymbol{v}) \, dV$$
(14)

$$\frac{\partial Q_2}{\partial c} \Delta c = \int_{\Omega} \left[\Delta c \left(\frac{g''(c)\ell}{G_c} \mathcal{H}(\boldsymbol{x}, t) + 1 \right) w + \ell^2 \boldsymbol{\nabla} (\Delta c) \cdot \boldsymbol{\nabla} w \right] dV$$
 (15)

Since g''(c) = 2(1-k) is a constant, the latter is linear in Δc and w, and independent of c, hence a linear solve is sufficient for Equation (13). In the *IFEM* implementation, we introduce the following variables

$$s_1 = 1 + 2(1 - k)\mathcal{H}(\mathbf{x}, t)\frac{\ell}{G_c}$$
 and $s_2 = \ell^2$ (16)

with which Equation (11) can be written

$$Q_2(\boldsymbol{u}, c; w) = \int_{\Omega} \left[(s_1 c - 1)w + s_2 \boldsymbol{\nabla} c \cdot \boldsymbol{\nabla} w \right] dV = 0$$
(17)

Alternatively to using the history field, we may include a penalty term to enforce the crack irreversibility as in Equation (2), i.e.:

$$E_{c}(\boldsymbol{u},c) = \int_{\Omega} \left[g(c) \Psi_{0}^{+}(\boldsymbol{\epsilon}(\boldsymbol{u})) + \Psi_{0}^{-}(\boldsymbol{\epsilon}(\boldsymbol{u})) \right] dV$$

$$+ G_{c} \int_{\Omega} \left(\frac{1}{4\ell_{0}} (1-c)^{2} + \ell_{0} \boldsymbol{\nabla} c \cdot \boldsymbol{\nabla} c \right) dV + \frac{1}{2\gamma} \int_{CR_{l-1}} c^{2} dV$$
(18)

such that the alternatives to Equations (11) and (15) become, respectively

$$Q_{2}(\boldsymbol{u}, c; w) = \int_{\Omega} \left[\frac{g'(c)\ell}{G_{c}} \Psi_{0}^{+}(\boldsymbol{\epsilon}(\boldsymbol{u})) w + (c-1)w + \ell^{2} \boldsymbol{\nabla} c \cdot \boldsymbol{\nabla} w \right] dV + \frac{1}{\gamma} \int_{CR_{c}} \frac{\ell}{G_{c}} cw dV = 0 \quad (19)$$

and

$$\frac{\partial Q_2}{\partial c} \Delta c = \int_{\Omega} \left[\Delta c \left(\frac{g''(c)\ell}{G_c} \Psi_0^+(\boldsymbol{\epsilon}(\boldsymbol{u})) + 1 \right) w + \ell^2 \boldsymbol{\nabla} (\Delta c) \cdot \boldsymbol{\nabla} w \right] dV + \frac{1}{\gamma} \int_{CR_{l-1}} \Delta c \frac{\ell}{G_c} w dV \qquad (20)$$

Equation (19) can be rewritten as Equation (17) where the parameter s_1 now reads

$$s_1 = 1 + 2(1 - k)\Psi_0^+(\boldsymbol{\epsilon}(\boldsymbol{u}))\frac{\ell}{G_c} + \begin{cases} \frac{\ell}{\gamma G_c} & \forall \boldsymbol{x} \in CR_{l-1} \\ 0 & \forall \boldsymbol{x} \notin CR_{l-1} \end{cases}$$
(21)

4 Residual and tangent for the d-formulation

Equation (6) is rewritten into a system of coupled equations:

$$Q_{1}(\boldsymbol{u}, d; \boldsymbol{v}) = \int_{\Omega} \left[g(d) \frac{\partial \Psi_{0}^{+}}{\partial \boldsymbol{\epsilon}} (\boldsymbol{\epsilon}(\boldsymbol{u})) + \frac{\partial \Psi_{0}^{-}}{\partial \boldsymbol{\epsilon}} (\boldsymbol{\epsilon}(\boldsymbol{u})) \right] : \boldsymbol{\epsilon}(\boldsymbol{v}) \, dV = 0$$
(22)

$$Q_{2}(\boldsymbol{u},d;w) = \int_{\Omega} \left[g'(d)\Psi_{0}^{+}(\boldsymbol{\epsilon}(\boldsymbol{u}))w + G_{c}\left(\frac{1}{\ell}dw + \ell\nabla d \cdot \nabla w\right) \right] dV + \frac{1}{\gamma} \int_{CR_{l-1}} (d-1)w dV = 0$$
(23)

These two equations are solved in a staggered manner, where in Equation (22) the phase field d is assumed known and in Equation (23) the displacement field u is assumed known. We then do a linearization of the functionals Q_1 and Q_2 about a certain known configuration (u_0, d_0) for the unknown variables, i.e.

$$Q_1(\boldsymbol{u}_0 + \Delta \boldsymbol{u}, d_0; \boldsymbol{v}) \approx Q_1(\boldsymbol{u}_0, d_0; \boldsymbol{v}) + \frac{\partial Q_1}{\partial \boldsymbol{u}} \cdot \Delta \boldsymbol{u} = 0$$
 (24)

$$Q_2(\boldsymbol{u}_0, d_0 + \Delta d; w) \approx Q_2(\boldsymbol{u}_0, d_0; \boldsymbol{v}) + \frac{\partial Q_2}{\partial d} \Delta d = 0$$
 (25)

from which we obtain the tangent operators

$$\frac{\partial Q_1}{\partial \boldsymbol{u}} \cdot \Delta \boldsymbol{u} = \int_{\Omega} \boldsymbol{\epsilon}(\Delta \boldsymbol{u}) : \left[g(d) \frac{\partial^2 \Psi_0^+}{\partial \boldsymbol{\epsilon}^2} (\boldsymbol{\epsilon}(\boldsymbol{u})) + \frac{\partial^2 \Psi_0^-}{\partial \boldsymbol{\epsilon}^2} (\boldsymbol{\epsilon}(\boldsymbol{u})) \right] : \boldsymbol{\epsilon}(\boldsymbol{v}) \, dV$$
 (26)

$$\frac{\partial Q_2}{\partial d} \Delta d = \int_{\Omega} \left[\Delta d \left(g''(d) \Psi_0^+(\boldsymbol{\epsilon}(\boldsymbol{u})) + \frac{G_c}{\ell} \right) w + G_c \ell \boldsymbol{\nabla} (\Delta d) \cdot \boldsymbol{\nabla} w \right] dV + \frac{1}{\gamma} \int_{\Omega R_c} \Delta dw \, dV (27)$$

Since g''(d) = 2(1-k) is a constant, the latter is linear in Δd and w, and independent of d, hence a linear solve is sufficient for Equation (25).

References

- [1] M. J. Borden, C. V. Verhoosel, M. A. Scott, T. J. R. Hughes, and C. M. Landis. A Phase-field Description of Dynamic Brittle Fracture. *Computer Methods in Applied Mechanics and Engineering*, 217–220:77–95, 2012.
- [2] T. Gerasimov and L. De Lorenzis. A Line Search Assisted Monolithic Approach for Phase-field Computing of Brittle Fracture. *Computer Methods in Applied Mechanics and Engineering*, 312:276–303, 2016.