

A Multi Phase-Field-Cohesive Zone Model for Laminated Composites: Application to Delamination Migration

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Abstract

Failure processes in Laminated Fiber-Reinforced Composites (LFRCs) entail the development and progression of different physical mechanisms and, in particular, the interaction between inter-laminar and intra-laminar cracking. Reliable modeling of such complex scenarios can be achieved by developing robust numerical predictive tools that allow for the interaction of both failure modes. In this study, a novel Multi Phase-Field (MPF) model relying on the Puck theory of failure for intra-laminar failure at ply level is coupled with a Cohesive Zone Model (CZM) for inter-laminar cracking. The current computational method is numerically implemented as a system of non-linear partially coupled equations using the finite element method via user-defined UMAT and UEL subroutines in ABAQUS. The computational tool is applied to qualitatively predict delamination migration in long laminated fiber-reinforced polymers composites comprising 44 cross-ply laminates. The reliability of the current approach is examined via the correlation with experimental results. Finally, the present study is complemented with additional representative examples with the aim of providing further insight into the potential role of different aspects of the system in the delamination migration, including (i) the variation of the ply angle in the migration zone, (ii) the load application point, and (iii) initial crack length.

Keywords: A. Fiber Reinforced Composites; B. Fracture Mechanics; C. Finite Element Method (FEM); D. Phase-Field Fracture; E. Cohesive Zone Model

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1. Introduction

The widespread use of Laminated Fiber Reinforced Composites (LFRCs) due to their high strength to weight ratios has paved their way into many practical applications in different industrial sectors, with remarkable impact in aerospace and aeronautics, and more recently in the automotive sector [1, 2], to quote a few of them. In the presence of ever-increasing demands of new technological advancements, load-bearing capabilities and failure modes of these materials are not yet fully understood, posing notable restrictions on their use and leading to the introduction of high safety factors.

Within this context, delamination failure can be conceived as one of the most critical failure mechanisms in Laminated Fiber Reinforced Composites (LFRCs). Such cracking events are generally associated with low through-thickness stiffness in layered disposals and can emerge from manufacturing defects, the occurrence of post-buckling phenomena [3, 4], among many others. From a modeling standpoint, delamination events in layered composites structures have been analyzed using linear elastic fracture mechanics (LEFM), enforcing initiation and growth along the ply interface with the use of strain energy-based methods to define the corresponding fracture toughness [5, 6]. Delamination can be found in low-velocity impact [7], skin debonding

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[8], defects from notches [9], [10], among many other practical cases. These phenomena have been extensively investigated based on cohesive-like cracking methods in pre-notched coupon configurations. An ad-hoc inter-laminar failure model can be inserted into the most critical locations of the specimen for triggering such cracking events [11, 12, 13, 14]. From a mechanical perspective, delamination can be seen as a result of the coalescence of micro-cracks at the ply interface, perpendicular to the tensile stress [15], [16], [17].

However, for some specific loading cases and configurations, it has been reported that mostly due to the sign of the change of shearing stresses, a pre-existing crack along the ply interface can kink out from such location, propagating into the adjacent ply. Matrix-cracking induced by delamination can further progress through the ply thickness and then "propagates out" at another interface as reported in [15, 18, 19], leading to the so-called "delamination migration" failure mode. Thus, following the terminologies used in [20], the turning of delamination crack front into one of the adjacent plies is referred to as "kinking," whereas, when a crack propagates across the thickness of plies and propagates out in adjacent interface is called "migration" [21]. The main causes of such phenomena can be motivated by the analysis of the stress field around the interface crack tip, see the experimental studies conducted in [15, 22, 19] for a $0^\circ/90^\circ$ interface and [16] for a $0^\circ/\theta^\circ$ interface.

In the last decade, the advent of new numerical capabilities and modeling techniques has promoted the thorough analysis of the potential reasons for the driving mechanisms that provoke delamination migration failures [23]. Note also that the migration phenomenon is also observed in the models without the presence of pre-existing cracks [22]. Numerical methods such as eXtended FEM (XFEM) [24], augmented finite element methods [25], floating node methods [26], cohesive-crack element methods [27] have been widely applied to study delamination migration. Moreover, delamination migration can be explained using micro-mechanics as reported by Arteiro and coauthors [28].

Concerning the different modeling methods for triggering intra-laminar failure in LFRCs, the use of continuum damage mechanics (CDM)-based methods have been of notable importance; see the phenomenological models proposed in [29, 30] and the references given therein. Notwithstanding, the local versions of CDM models generally suffer from mesh pathological issues that can be remedied using alternative methodologies. Nonlocal damage models [31, 32] have been used extensively, which also share as part of phase-field formalism while providing a detailed description of the damage behavior. On the other hand, theories such as crack band theory have been extensively used in the literature to mitigate spurious mesh sensitivity in the modeling location. In this direction, substantial contributions from Bažant are noteworthy. See [33, 34, 35, 36, 37] and the reference therein. Within this context, Phase-field methods, originally proposed by Francfort and Marigo [38] and subsequently developed in [39], have become plausible nonlocal modeling alternatives that prevent most of the main limitations of alternative modeling tools for fracture in solids. With strong foundations on Griffith's approach to Fracture Mechanics, PF methods exploit a multi-field variational formalism, whereby fracture is accounted for via diffusive representation within a particular region of the domain characterized by the length-scale l . The Phase-field method for fracture integrates classical damage mechanics and Griffith's theory of failure. This generalized formalism can handle topologically complex fractures such as initiation, branching of the crack, intersection of crack, coalescence of damage, etc., or combining them in two and three-dimensional settings without any ad-hoc. This approach encompasses an energy minimization of the potential energy of the body (split into bulk and fracture counterparts, respectively), resulting in a multi-field FE formulation. Therefore, complex cracking phenomena can be naturally captured due to the implicit consideration of the corresponding crack phase-field variables' evolution equation. Posterior developments of PF techniques for fracture have been oriented towards the analysis of alternative formulations [40, 41], shells [30], ductile fracture [42, 43, 44], composite materials [45, 46, 47, 48, 49, 50, 51], heterogeneous media [48], hydrogen assisted cracking [52], among many others, and from a numerical point of view striving for different solution strategies [53]. The authors have recently proposed incorporating phenomenological failure criteria for composite materials widening the range of applicability of PF methods for different composite materials [54].

In this investigation, the two phase-field fracture model proposed by the authors in [55]) is combined with a cohesive zone model following a bi-linear traction separation law to predict the development of such failure events in LFRCs. Consequently, at the intra-ply level, the driving force for each of the failure mechanisms within the PF technique, i.e., fiber failure and inter-fiber failure, is exploited independently. Moreover,

this fracture model is readily applied to capture the kinking and migration events along with delamination. The proposed model is examined against experimental data available in the related literature to validate the current simulation technique's predictive capabilities. In addition to the previous aspects, the present study also encompasses the analysis of the variation of design parameters for which physical solution favors delamination-migration, such as load application point, angle of the ply at the interfaces, and initial crack length. Such effects are studied to understand each of the variation's load-bearing capacity and show the applicability of the proposed model to predict complex failure phenomena.

The manuscript is organized as follows. In Section 2, the Multi Phase-Field-Cohesive Zone (MPF-CZ) formulation based on the Puck failure theory is introduced. In Section 3, the model's variational formulation, along with the strong form, is outlined with a special focus on each failure mechanism's construction particular contribution. In Section 4, the finite element implementation of the model is derived. The assessment of the proposed framework for its validation is discussed in Section 5 in conjunction with a further investigation concerning the system's variation of numerous design parameters. Finally, Section 6 summarizes the main conclusions of the present study.

2. Multi Phase-Field-Cohesive Zone (MPF-CZ) formulation based on the Puck failure theory

In this section, the computational framework herein proposed for capturing delamination migration in layered composite structures is outlined. The current method relies on the combination of multiple phase-fields. Each phase-field based on different physically motivated failure mechanisms such as Puck failure criterion [55] driving the crack propagation in long fiber reinforced polymer composites along with the cohesive zone model.

Following the standard Phase-field formulation, consider an arbitrary body in the general n_{dim} Euclidean space, occupying the placement $\mathcal{B} \in \mathbb{R}^{n_{dim}}$, with its external delimiting boundary $\partial\mathcal{B} \in \mathbb{R}^{n_{dim}-1}$, see Figure 1. For any material point, the position vector is denoted by $\mathbf{x} \in \mathcal{B}$. The displacement field is identified by the vector $\mathbf{u} : \mathcal{B} \rightarrow \mathbb{R}^{n_{dim}}$, with infinitesimal strain tensor $\boldsymbol{\varepsilon} := \nabla^s \mathbf{u}$ for $\boldsymbol{\varepsilon} : \mathcal{B} \rightarrow \mathbb{R}^{n_{dim} \times n_{dim}}$. The displacement boundary conditions are prescribed as $\mathbf{u} = \bar{\mathbf{u}}$ on $\partial\mathcal{B}_u$ and traction conditions are given by $\bar{\mathbf{t}} = \boldsymbol{\sigma} \cdot \mathbf{n}$ on $\partial\mathcal{B}_t$ such that, kinematic and static boundary conditions satisfy: $\overline{\partial\mathcal{B}_t \cup \partial\mathcal{B}_u} = \partial\mathcal{B}$ and $\partial\mathcal{B}_t \cap \partial\mathcal{B}_u = \emptyset$, where \mathbf{n} is outward normal vector and $\boldsymbol{\sigma}$ is the Cauchy stress tensor.

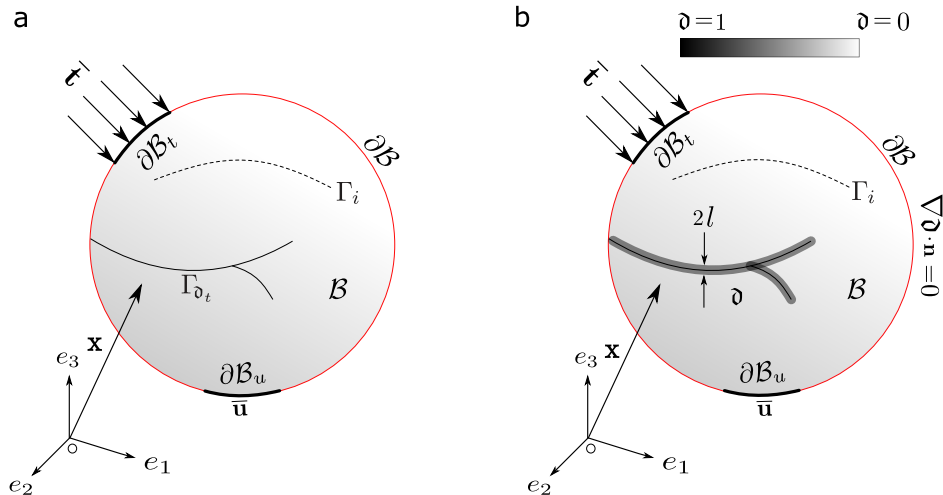


Figure 1: Body under consideration: (a) sharp crack representation and (b) Regularized crack topology.

In addition to the previous definitions, let Γ be a crack set incorporating interface cracks Γ_i arising from the cohesive interface and cracks in the bulk Γ_{δ_t} from the multi phase-field such that, $\Gamma_i \cup \Gamma_{\delta_t} = \Gamma$ and $\Gamma_i \cap \Gamma_{\delta_t} = \emptyset$ for each discrete $t \in [0, T]$ with $\Gamma_t \subseteq \Gamma_{t+1}$. The displacement jumps along the interface

as the relative displacement between two homogeneous points at the flanks are denoted by $\mathbf{g} = \mathbf{u}_i^+ - \mathbf{u}_i^-$, representing the difference between kinematic field along the interface Γ_i^+ and Γ_i^- in line with [56, 57]. On the other hand, $\Gamma_{\mathfrak{d}_i}$ is defined as the set of discontinuous points \mathbf{x} , where \mathbf{u} has one sided approximate limits $\mathbf{u}_{\mathfrak{d}}^+ \neq \mathbf{u}_{\mathfrak{d}}^-$ with respect to a suitable direction \mathbf{v}_u normal to $\Gamma_{\mathfrak{d}_i}$ inline with [58, 59].

To account for multiple energies in the system, the total free energy functional describing the mechanics of body \mathcal{B} , is given as the sum total of internal and external energies acting on the system as follows:

$$\Pi(\mathbf{u}, \Gamma) = \Pi_{\text{int}}(\mathbf{u}, \Gamma) + \Pi_{\text{ext}}(\mathbf{u}). \quad (1)$$

Phase-field approximation of the fracture works on the idea that the fracture propagates due to competition between elastic energy stemming from the state function of strain tensor $\boldsymbol{\varepsilon}$ and fracture energy (surface energy) stemming from scalar internal damage like variable \mathfrak{d} [55]. The consistent generalization of the isotropic damage formulation for the consideration of different failure mechanisms can be postulated by the additive decomposition of total internal energy into multiple contributions, in which each of them is associated with a certain failure mechanism. In such a postulation, a scalar damage variable \mathfrak{d}_i ($i = 1, \dots, n$) is associated with each one of the n failure mechanisms, such that $\mathfrak{d}_i = 0$ for intact material state and $\mathfrak{d}_i = 1$ for fully broken state and so that $\mathfrak{d}_i \in [0, 1]$ for each $i = 1, \dots, n$. Moreover, to account for non-local damage evolution, the respective gradients $\nabla_{\mathbf{x}} \mathfrak{d}_i$ are incorporated in the formulation. This additive decomposition postulation for intralaminar failure was successfully applied in [29] and later extended to incorporate failure criteria such as Puck failure theory so as to distinguish between the fiber and matrix dominated damage mechanisms in LFRPs [55, 30]. However, the consideration of dissipative energies stemming from debonding along the interface Γ_i , and crack propagation in the bulk $\Gamma_{\mathfrak{d}_i}$ is still a matter of investigation. The prediction of such crack topology becomes increasingly complex due to branching and coalescence phenomenons, as well as the interaction with diffusive cracks, which may induce the debonding process along the existing interface Γ_i . One possibility to achieve this is by employing a phase-field model incorporating multiple diffusive crack fields within the bulk and interface elements relying on cohesive zone methodologies at a prescribed interface inline with [57]. Hence, the total internal energy is now an amalgamation of (i) total elastic energy constituting from bulk (fiber and inter-fiber) energy, (ii) surface energy (crack energy) stemming from bulk (fiber and inter-fiber) failure, and (iii) cohesive interface energy obeying a bi-linear traction-separation law:

$$\Pi_{\text{int}}(\mathbf{u}, \Gamma) = \Pi_{\text{int},b}(\mathbf{u}, \Gamma_{\mathfrak{d}_i}) + \Pi_{\text{int},c}(\Gamma_i), \quad (2)$$

where $\Pi_{\text{int},b}(\mathbf{u}, \Gamma)$ is the internal energy stemming from the bulk (fiber and inter-fiber) and $\Pi_{\text{int},c}(\Gamma_i)$ is the dissipative energy associated with cohesive debonding which are addressed in detail in the sequel. The bulk energy $\Pi_{\text{int},b}(\mathbf{u}, \Gamma)$ is further decomposed into:

$$\Pi_{\text{int},b}(\mathbf{u}, \Gamma_{\mathfrak{d}_i}) = \Pi_{\text{int},FF}(\mathbf{u}, \Gamma_{\mathfrak{d}_{FF}}) + \Pi_{\text{int},IFF}(\mathbf{u}, \Gamma_{\mathfrak{d}_{IFF}}), \quad (3)$$

where $\Pi_{\text{int},FF}(\mathbf{u}, \Gamma_{\mathfrak{d}_{FF}})$ and $\Pi_{\text{int},IFF}(\mathbf{u}, \Gamma_{\mathfrak{d}_{IFF}})$ correspond to the energies associated with fiber failure and inter-fiber failure, respectively.

With such decomposition at hand, the scheme herein used recalls that the dissipated energy arising from each of the individual failure mechanisms only affects their corresponding counterparts in the elasticity tensor, therefore precluding the coupling between fiber inter-fiber failures and each of these failures with respect to the cohesive debonding. Note that, due to this preclusion of interaction between the energies, the elastic energy is strongly coupled with the surface energies stemming from fiber, inter-fiber, and cohesive debonding. In contrast, surface energies among themselves are only weakly coupled.

The effective Helmholtz free energy function $\hat{\Psi}$ is defined as:

$$\hat{\Psi}(\boldsymbol{\varepsilon}, \mathbf{A}) = \frac{1}{2} \boldsymbol{\varepsilon} : \mathbb{C}^e : \boldsymbol{\varepsilon}, \quad (4)$$

where \mathbb{C}^e is the undamaged elastic constitutive tensor defined as:

$$\mathbb{C}^e := \partial_{\boldsymbol{\varepsilon}\boldsymbol{\varepsilon}} \hat{\Psi} = \lambda \mathbf{1} \otimes \mathbf{1} + 2\mu_T \mathbb{I} + \alpha(\mathbf{1} \otimes \mathbf{A} + \mathbf{A} \otimes \mathbf{1}) + 2(\mu_L - \mu_T) \mathbb{I}_A + \beta \mathbf{A} \otimes \mathbf{A}, \quad (5)$$

where $\mathbf{1}$ is the second order identity tensor, $\mathbb{I}_{A,ijkl} = A_{im}\mathbb{I}_{jmkl} + A_{jm}\mathbb{I}_{mikl}$ represents the fourth-order identity matrix, and λ , α , β , μ_T and μ_L are to the elastic constants taking the form:

$$\lambda = E_{22}(\nu_{23} + \nu_{31}\nu_{13})/D, \quad (6)$$

$$\alpha = E_{22}[\nu_{31}(1 + \nu_{32} - \nu_{13}) - \nu_{32}]/D, \quad (7)$$

$$\beta = E_{11}(1 - \nu_{32}\nu_{23})/D - E_{22}[1 - \nu_{21}(\nu_{12} + 2(1 + \nu_{23}))]/D - 4G_{12}, \quad (8)$$

$$\mu_L = G_{12}; \quad \mu_T = G_{23}, \quad (9)$$

with $D = 1 - \nu_{32}^2 - 2\nu_{13}\nu_{31} - 2\nu_{32}\nu_{13}\nu_{31}$. The material direction is denoted by \mathbf{a} and $\mathbf{A} := \mathbf{a} \otimes \mathbf{a}$ is so called structural tensor.

Here, 1-direction corresponds to the fiber orientation, 2-direction is transverse in-plane orientation with respect to the fiber direction, and 3-direction stands for transverse out-of-plane orientation.

2.1. Bulk energies

In the light of previous developments, the total energy of the fiber can be established as the sum total of the elastic energy and fracture energy associated with the fiber:

$$\begin{aligned} \Pi_{\text{int},FF}(\mathbf{u}, \Gamma_{\mathfrak{d}_{FF}}) &\approx \Pi_{\text{int},FF}(\mathbf{u}, \mathfrak{d}_{FF}) = \int_{\mathcal{B}} (1 - \mathfrak{d}_{FF})^2 \hat{\Psi}_{FF}(\boldsymbol{\varepsilon}, \mathbf{A}) d\Omega \\ &+ \int_{\mathcal{B}} \mathcal{G}_{c,FF} \left[\frac{1}{2l_{FF}} \mathfrak{d}_{FF}^2 + \frac{l_{FF}}{2} |\nabla \mathfrak{d}_{FF}|^2 \right] d\Omega, \end{aligned} \quad (10)$$

where $\hat{\Psi}_{FF}$ is the elastic contribution associated with the fiber: $\hat{\Psi}_{FF} = \frac{1}{2} \boldsymbol{\varepsilon} : \mathbb{C}_{FF}^e : \boldsymbol{\varepsilon}$, where:

$$\mathbb{C}_{FF}^e = \begin{bmatrix} \mathbb{C}_{11}^e & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 \end{bmatrix}, \quad (11)$$

and $\mathcal{G}_{c,FF}$ is the fracture energy, and l_{FF} is the material characteristic length associated with fiber failure in phase-field related to the apparent material strength [60] as:

$$l_{FF} = \frac{27}{256} \frac{E_{11} \mathcal{G}_{c,FF}}{\sigma_{s,FF}^2}, \quad (12)$$

where $\sigma_{s,FF}$ is the apparent material strength associated with fiber failure.

Similarly, the inter-fiber contribution to the total internal energy can be expressed as:

$$\begin{aligned} \Pi_{\text{int},IFF}(\mathbf{u}, \Gamma_{\mathfrak{d}_{IFF}}) &\approx \Pi_{\text{int},IFF}(\mathbf{u}, \mathfrak{d}_{IFF}) = \int_{\mathcal{B}} (1 - \mathfrak{d}_{IFF})^2 \hat{\Psi}_{IFF}(\boldsymbol{\varepsilon}, \mathbf{A}) d\Omega \\ &+ \int_{\mathcal{B}} \mathcal{G}_{c,IFF} \left[\frac{1}{2l_{IFF}} \mathfrak{d}_{IFF}^2 + \frac{l_{IFF}}{2} |\nabla \mathfrak{d}_{IFF}|^2 \right] d\Omega, \end{aligned} \quad (13)$$

where $\hat{\Psi}_{IFF} = \frac{1}{2} \boldsymbol{\varepsilon} : \mathbb{C}_{IFF}^e : \boldsymbol{\varepsilon}$ is the elastic contribution associated with the inter-fiber failure, and:

$$\mathbb{C}_{IFF}^e = \begin{bmatrix} 0 & \mathbb{C}_{12}^e & \mathbb{C}_{13}^e & 0 & 0 & 0 \\ \mathbb{C}_{21}^e & \mathbb{C}_{22}^e & \mathbb{C}_{23}^e & 0 & 0 & 0 \\ \mathbb{C}_{31}^e & \mathbb{C}_{32}^e & \mathbb{C}_{33}^e & 0 & 0 & 0 \\ 0 & 0 & 0 & \mathbb{C}_{44}^e & 0 & 0 \\ 0 & 0 & 0 & 0 & \mathbb{C}_{55}^e & 0 \\ 0 & 0 & 0 & 0 & 0 & \mathbb{C}_{66}^e \end{bmatrix}. \quad (14)$$

where $\mathcal{G}_{c,IFF}$ and l_{IFF} are the fracture energy and corresponding length scale associated with inter-fiber failure, respectively. Similarly, the length scale parameter is estimated for the apparent material strength $\sigma_{s,IFF}$ of inter-fiber failure as:

$$l_{IFF} = \frac{27}{256} \frac{E_{22} \mathcal{G}_{c,IFF}}{\sigma_{s,IFF}^2}. \quad (15)$$

Note that with these definitions at hand, the damaged constitutive matrix \mathbb{C} renders:

$$\mathbb{C}(\mathfrak{d}_{FF}, \mathfrak{d}_{IFF}) = (1 - \mathfrak{d}_{FF})^2 \mathbb{C}_{FF}^e + (1 - \mathfrak{d}_{IFF})^2 \mathbb{C}_{IFF}^e, \quad (16)$$

$$\mathbb{C}(\mathfrak{d}_{FF}, \mathfrak{d}_{IFF}) = \begin{bmatrix} \mathcal{P}_1 \mathbb{C}_{11}^e & \mathcal{P}_2 \mathbb{C}_{12}^e & \mathcal{P}_2 \mathbb{C}_{13}^e & 0 & 0 & 0 \\ \mathcal{P}_2 \mathbb{C}_{21}^e & \mathcal{P}_2 \mathbb{C}_{22}^e & \mathcal{P}_2 \mathbb{C}_{23}^e & 0 & 0 & 0 \\ \mathcal{P}_2 \mathbb{C}_{31}^e & \mathcal{P}_2 \mathbb{C}_{32}^e & \mathcal{P}_2 \mathbb{C}_{33}^e & 0 & 0 & 0 \\ 0 & 0 & 0 & \mathcal{P}_{12} \mathbb{C}_{44}^e & 0 & 0 \\ 0 & 0 & 0 & 0 & \mathcal{P}_{12} \mathbb{C}_{55}^e & 0 \\ 0 & 0 & 0 & 0 & 0 & \mathcal{P}_2 \mathbb{C}_{66}^e \end{bmatrix}, \quad (17)$$

where $\mathcal{P}_1 = (1 - \mathfrak{d}_{FF})^2$, $\mathcal{P}_2 = (1 - \mathfrak{d}_{IFF})^2$, and $\mathcal{P}_{12} = \min(\mathcal{P}_1, \mathcal{P}_2)$. Notice that \mathcal{P}_{12} is not differentiable due to the existence of "min". Hence, we first take the minima so that, any minimum of \mathcal{P}_{12} (i.e \mathcal{P}_1 or \mathcal{P}_2) is differentiable.

2.2. Interface energies

The energy stemming from the interface is governed by a bi-linear traction separation law which constitutes of a linear elastic stage characterized by an initial stiffness of K_n , K_{t1} and K_{t2} corresponding to the normal and shear components, respectively, followed by a linear softening as in Figure 2. The irreversibility is accounted by introducing a damage variable \mathfrak{d}_c depending on the relative kinematic critical normal and tangential openings, \mathfrak{g}_n^f and $\mathfrak{g}_{t_j}^f$ ($j=1$ for 2D and $j=1,2$ for 3D), respectively [12, 13].

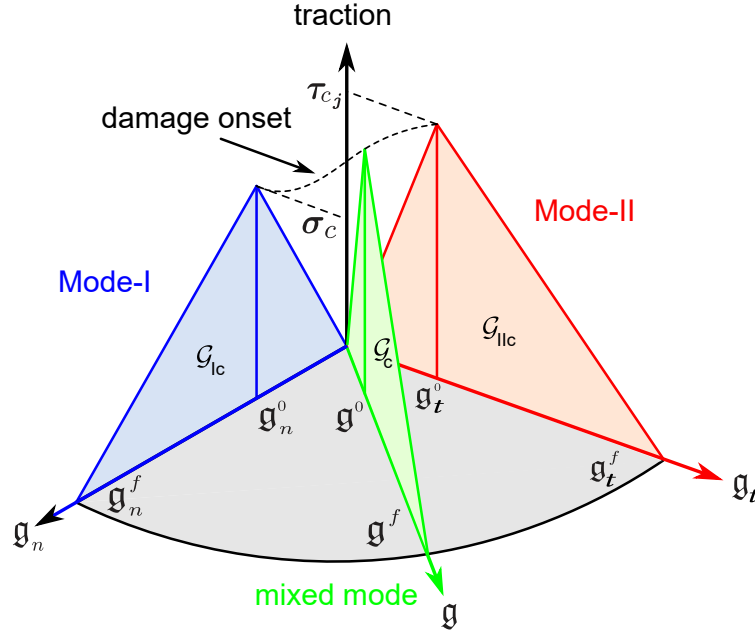


Figure 2: Schematic representation of the bi-linear cohesive zone model traction-separation law for mixed mode

The corresponding interface laws governing the normal and tangential tractions $\boldsymbol{\sigma}_n$ and $\boldsymbol{\tau}_{t_j}$, respectively, take the form:

$$\boldsymbol{\sigma}_n = \begin{cases} k_n \mathbf{g}_n & \text{if } \mathbf{g}_n \leq \mathbf{g}_n^0 \\ (1 - \mathfrak{d}_c) k_n \mathbf{g}_n & \text{if } \mathbf{g}_n^0 < \mathbf{g}_n < \mathbf{g}_n^f, \\ 0 & \text{Otherwise} \end{cases}, \quad (18)$$

$$\boldsymbol{\tau}_{t_j} = \begin{cases} k_{t_j} \mathbf{g}_{t_j} & \text{if } \mathbf{g}_{t_j} \leq \mathbf{g}_{t_j}^0 \\ (1 - \mathfrak{d}_c) k_{t_j} \mathbf{g}_{t_j} & \text{if } \mathbf{g}_{t_j}^0 < \mathbf{g}_{t_j} < \mathbf{g}_{t_j}^f, \\ 0 & \text{Otherwise} \end{cases}, \quad (19)$$

for each $j = 1, 2$ to account for the tangential traction, whereas \mathbf{g}_n and \mathbf{g}_{t_j} are the relative normal and tangential displacements, respectively. The Mixed-mode fracture energy of the interface reads:

$$\mathcal{G}_C^i = \mathcal{G}_{IC} + (\mathcal{G}_{IIC} - \mathcal{G}_{IC}) \left(\frac{\mathcal{G}_{II} + \mathcal{G}_{III}}{\mathcal{G}_I + \mathcal{G}_{II} + \mathcal{G}_{III}} \right)^\eta, \quad (20)$$

where \mathcal{G}_{IC} , \mathcal{G}_{IIC} , and \mathcal{G}_{IIIC} represent the corresponding fracture toughness associated with normal (Mode I) and shear (Mode II and III), respectively, and computed as the area under the traction separation curve, η identifies an experimental fitting parameter elucidating the effects of fracture mode mixities [4]. Finally, \mathcal{G}_I , \mathcal{G}_{II} , and \mathcal{G}_{III} are energy release rate associated with Mode I, Mode II, and III, respectively. Damage is initiated using the quadratic interaction function of nominal stress ratios. The damage is evolved using Benzeggagh-Kenane (BK) failure criterion based on cumulative energy such that mixed fracture toughness \mathcal{G}_C^i reaches the critical energy.

Based on the energy considerations, the evolution of the damage variable \mathfrak{d}_c is estimated based on the effective displacement of the cohesive element, where \mathfrak{d}_c is evolving linearly with:

$$\mathfrak{d}_c = \frac{\mathbf{g}_m^f (\mathbf{g}_m^{\max} - \mathbf{g}_m^0)}{\mathbf{g}_m^{\max} (\mathbf{g}_m^f - \mathbf{g}_m^0)}, \quad (21)$$

where \mathbf{g}_m^0 is effective displacement at initiation and $\mathbf{g}_m^f = \frac{2\mathcal{G}_C^i}{T_{eff}^0}$ is the effective displacement at failure. Here, T_{eff}^0 is the effective traction at damage initiation estimated using effective tractions $T_{eff} = \sqrt{(\boldsymbol{\sigma}_n)^2 + (\boldsymbol{\tau}_{t_1})^2 + (\boldsymbol{\tau}_{t_2})^2}$, as $T_{eff}^0 = \min \{T_{eff}(\mathfrak{d}_c > 0)\}$. And \mathbf{g}_m^{\max} refers to the maximum value of effective displacement during loading history estimated using $\mathbf{g}_m^{\max} = \max_{[0, T]} \mathbf{g}_m$, where $\mathbf{g}_m := \sqrt{(\mathbf{g}_n^2) + (\mathbf{g}_{t_1})^2 + (\mathbf{g}_{t_2})^2}$. Finally, with this at hand, the corresponding total energy generated by the cohesive interface takes the form:

$$\Pi_{\text{int},c}(\Gamma_i) \approx \Pi_{\text{int},c}(\mathfrak{d}_c) = \int_{\Gamma_i} \mathcal{G}_c^i(\mathbf{g}, \mathfrak{d}_c) d\mathcal{S} = \int_{\Gamma_i} \mathbf{g}^T \mathcal{T} d\mathcal{S}, \quad (22)$$

with $\mathbf{g} = [\mathbf{g}_n, \mathbf{g}_{t_1}]$ and $\mathcal{T} = [\boldsymbol{\sigma}_n, \boldsymbol{\tau}_{t_1}]$.

2.3. Fundamentals of Puck failure criterion

Damage evolution in bulk relies on Puck failure theory [61] whose corresponding failure criterion accounts for the independent assessment of fiber and inter-fiber failure surfaces. For the fiber failure, with usual notations, that is \parallel (subscript 1), \perp (subscript 2 and subscript 3) representing fiber direction, normal to the fiber direction in-plane and, normal to the fiber direction out-of-plane, respectively, for the ply co-ordinates in a local setting $0 - \mathbf{e}_1 - \mathbf{e}_2 - \mathbf{e}_3$ as in Figure 3.

According to the Puck theory, fiber failure is triggered based on energetic considerations that, the exposure factor (denoted $f_{E,FF+}$, for tensile only) reaches the value 1, where $f_{E,FF+}$ is given by:

$$f_{E,FF+} = \frac{1}{R_{\parallel}^t} \left[\sigma_{11} - \left(\nu_{\perp\parallel} - \frac{E_{\parallel}}{E_{\parallel f}} \nu_{\perp\parallel f} \right) (\sigma_{22} + \sigma_{33}) \mathcal{P}_2 \right], \quad (23)$$

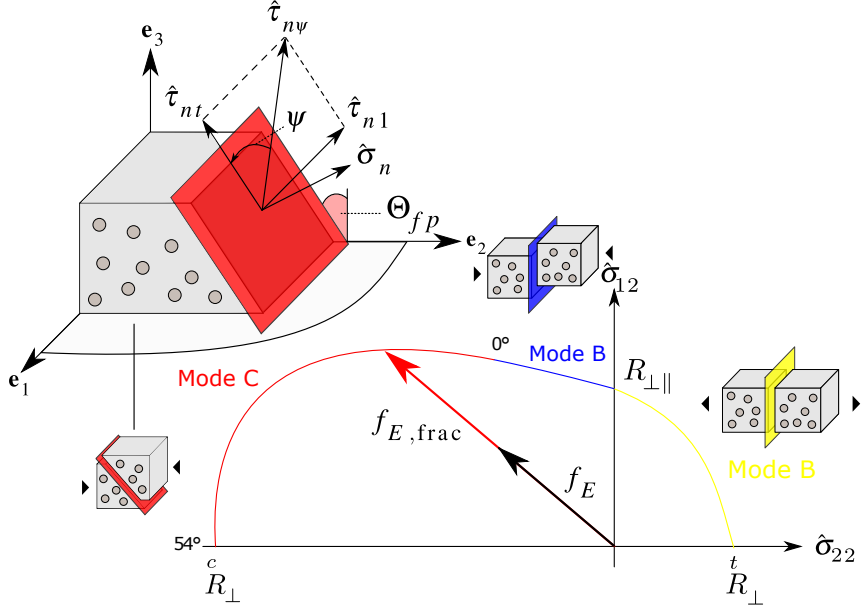


Figure 3: Puck failure theory: (a) definition of acting stresses on the fracture plane by θ_{fp} angle and (b) the exposure factor f_E , where $f_{E,(fr)}$ is the exposure factor at the failure point.

where R_{\parallel}^t stands for the tensile longitudinal strength in fiber direction. $\nu_{\perp\parallel}$ and $\nu_{\perp\parallel f}$ identify the major Poisson's ratios of the ply and the fibers, respectively, and $E_{\parallel f}$ is the elastic modulus of the fibers. $FF+$ denotes the fiber failure in tension, whereas the compression is omitted for the sake of brevity but can easily be incorporated as in [55]. The incorporation of \mathcal{P}_2 is to scale the influence of the transverse stress components on the longitudinal stress (lateral contraction) with respect to the state of matrix damage. It is assumed that the contraction in the longitudinal direction due to transverse stress will vanish in a case $\mathcal{P}_2 \rightarrow 1$, i.e., matrix rupture parallel to the fibers. In the case of compressive longitudinal stress, reduced compressive longitudinal fracture resistance of the plies is assumed in the case of increasing shear stress.

Puck theory also distinguishes the inter-fiber failure by introducing the so-called action plane [62, 61], which corresponds to identifying the potential fracture plane derived from the maximum stress states $\{\bar{\sigma}_n(\Theta), \bar{\tau}_{nt}(\Theta), \bar{\tau}_{n1}(\Theta)\}$, where $\bar{\sigma}_n(\Theta)$, $\bar{\tau}_{nt}(\Theta)$, and $\bar{\tau}_{n1}(\Theta)$ are the normal stress component, shear stress component transverse to the fiber, and the shear stress component plane parallel to the fiber, respectively, each on the action plane. The determination of the fracture plane is usually performed via the assessment of the most critical stress state in terms of the local components by calculating the value of inter-fiber exposure factor $F_{E,IFF}$ for all angles Θ within the interval $-90^\circ \leq \Theta \leq +90^\circ$, using an increment of one degree. The transformation from the local ply setting to the action plane system yields:

$$\begin{bmatrix} \sigma_n(\Theta) \\ \tau_{nt}(\Theta) \\ \tau_{n1}(\Theta) \end{bmatrix} = \begin{bmatrix} \cos^2 \Theta & \sin^2 \Theta & 2 \cos \Theta \sin \Theta & 0 & 0 \\ -\cos \Theta \sin \Theta & \cos \Theta \sin \Theta & \cos^2 \Theta - \sin^2 \Theta & 0 & 0 \\ 0 & 0 & 0 & \sin \Theta & \cos \Theta \end{bmatrix} \begin{bmatrix} \sigma_{22} \\ \sigma_{33} \\ \sigma_{23} \\ \sigma_{13} \\ \sigma_{12} \end{bmatrix}. \quad (24)$$

In particular, the expression for inter-fiber failure under tensile conditions on the action plane takes the form:

$$f_{E,IFF+}(\Theta) = \begin{cases} \sqrt{\left[\left(\frac{1}{R_{\perp}^{At}} - \frac{p_{\perp\psi}^t}{R_{\perp\psi}^A} \right) \sigma_n(\Theta) \right]^2 + \left(\frac{\tau_{nt}(\Theta)}{R_{\perp\perp}^A} \right)^2 + \left(\frac{\tau_{n1}(\Theta)}{R_{\parallel\perp}^A} \right)^2 + \frac{p_{\perp\psi}^t}{R_{\perp\psi}^A} \sigma_n(\Theta)} \\ \frac{1}{\eta_w} \text{ for } \sigma_n(\Theta) \geq 0, \end{cases} \quad (25)$$

where R_{\perp}^{At} and $R_{\perp\perp}^A$ are transverse tensile strength and fracture strength respectively, whereas $R_{\perp\psi}^A$ is the transverse tensile strength at any ply angle ψ , with $f_{E,IFF+}$ representing the exposure factor and the failure is triggered when $f_{E,IFF+}$ reaches the value 1.

In the previous expression, $\cos^2 \psi$ and $\sin^2 \psi$ are:

$$\cos^2 \psi = \frac{\tau_{nt}^2}{\tau_{nt}^2 + \tau_{n1}^2} \text{ and } \sin^2 \psi = \frac{\tau_{n1}^2}{\tau_{nt}^2 + \tau_{n1}^2}. \quad (26)$$

Finally, the definition of the inclination parameters $p_{\perp\psi}^t$ and $p_{\perp\psi}^c$ for any angle ψ renders:

$$\frac{p_{\perp\psi}^i}{R_{\perp\psi}^A} = \frac{p_{\perp\perp}^i}{R_{\perp\perp}^A} \cos^2 \psi + \frac{p_{\perp\parallel}^i}{R_{\perp\parallel}^A} \sin^2 \psi, \quad i = t, c; \quad (27)$$

with:

$$R_{\perp\psi}^A = \left[\left(\frac{\cos \psi}{R_{\perp\perp}^A} \right)^2 + \left(\frac{\sin \psi}{R_{\perp\parallel}^A} \right)^2 \right]. \quad (28)$$

Due to the modelling assumptions herein made, the proposed model requires the following fracture energy values: (i) $\mathcal{G}_{c,FF}$ fiber fracture energy, (ii) $\mathcal{G}_{c,IFF}$ matrix dominated fracture energy, and (iii) \mathcal{G}_c inter-laminar fracture energy. These properties can be determined via experimental procedures, see [63, 28, 64, 65] and the references therein.

3. Variational formulation and strong forms

Relying on the considerations given in Section 2, the total energy functional of the solid body \mathcal{B} , along with the cracks Γ_i and Γ_{∂_i} at any arbitrary instance $t \in [0, T]$ takes the form:

$$\Pi(\mathbf{u}, \Gamma) \approx \Pi(\mathbf{u}, \mathfrak{d}_i, \mathfrak{d}_c) = \Pi_{\text{int}}(\mathbf{u}, \mathfrak{d}_i, \mathfrak{d}_c) + \Pi_{\text{ext}}(\mathbf{u}), \quad (29)$$

where the internal and external contribution to the energy functional $\Pi(\mathbf{u}, \mathfrak{d}_i, \mathfrak{d}_c)$ read, respectively:

$$\begin{aligned} \Pi_{\text{int}}(\mathbf{u}, \mathfrak{d}_i, \mathfrak{d}_c) &= \int_{\mathcal{B}} (1 - \mathfrak{d}_{FF})^2 \hat{\Psi}_{FF}(\boldsymbol{\varepsilon}, \mathbf{A}) + (1 - \mathfrak{d}_{IFF})^2 \hat{\Psi}_{IFF}(\boldsymbol{\varepsilon}, \mathbf{A}) d\Omega \\ &+ \int_{\mathcal{B}} P_{FF} \mathcal{G}_{c,FF} \left[\frac{1}{2l_{FF}} \mathfrak{d}_{FF}^2 + \frac{l_{FF}}{2} |\nabla \mathfrak{d}_{FF}|^2 \right] + P_{IFF} \mathcal{G}_{c,IFF} \left[\frac{1}{2l_{IFF}} \mathfrak{d}_{IFF}^2 + \frac{l_{IFF}}{2} |\nabla \mathfrak{d}_{IFF}|^2 \right] d\Omega \\ &+ \int_{\Gamma_i} \mathfrak{g}^T \mathcal{T} d\mathcal{S}, \end{aligned} \quad (30)$$

$$\Pi_{\text{ext}}(\mathbf{u}) = - \int_{\mathcal{B}} \mathbf{f}_v d\Omega - \int_{\partial \mathcal{B}_t} \bar{\mathbf{t}} d\partial\Omega, \quad (31)$$

where \mathbf{f}_v is the deformation-independent volume-specific loads.

In the above expression, Eqn.(30), P_{FF} and P_{IFF} are the activation flags, which are activated when the corresponding pucks failure criterion is met. The approximate solution of the stated problem with MPF-CZ can be obtained as a quadruplet of the solution of the following minimization problem:

$$(\mathbf{u}^*, \mathfrak{d}_{FF}^*, \mathfrak{d}_{IFF}^*, \mathfrak{g}^*) = \underset{\mathcal{S}}{\operatorname{argmin}} \Pi(\mathbf{u}, \mathfrak{d}_i, \mathfrak{d}_c). \quad (32)$$

where, $\mathcal{S} = \{\dot{\mathfrak{d}}_{FF}, \dot{\mathfrak{d}}_{IFF} \geq 0 \text{ for all } \mathbf{x} \in \mathcal{B} \setminus \Gamma_i, \dot{d}_c \geq 0 \text{ for all } \mathbf{x} \in \Gamma_i\}$ enforcing irreversibility of the evolution parameters. One may also argue that \mathfrak{g}^* should be replaced by the \mathfrak{d}_c in the minimization problem, however, it is easy to see that the minimum value of \mathfrak{g}^* also leads to a minimum value of \mathfrak{d}_c from Eqn. (21).

The quadruplet set $(\mathbf{u}^*, \mathfrak{d}_{FF}^*, \mathfrak{d}_{IFF}^*, \mathfrak{g}^*)$ in Eqn.(32) is solved by taking a variational form of the total internal energy functional (Gateaux derivative) assuming enough regularity of the terms involved. In the infinitesimal setting, considering the irreversibility of the damage variables set \mathcal{S} for any admissible test function $(\delta \mathbf{u}, \delta \mathfrak{d}_{FF}, \delta \mathfrak{d}_{IFF}, \delta \mathfrak{g})$ in the appropriate space of distribution, a strong form of the field equations can be reduced to the following:

$$\operatorname{div} [(1 - \mathfrak{d}_{FF})^2 \mathbb{C}_{FF}^e : \boldsymbol{\varepsilon} + (1 - \mathfrak{d}_{IFF}) \mathbb{C}_{IFF}^e : \boldsymbol{\varepsilon}] + \mathbf{f}_v = \mathbf{0} \text{ in } \mathcal{B} \setminus \Gamma_i, \text{ and } \boldsymbol{\sigma} \cdot \mathbf{n} = 0 \text{ on } \partial \mathcal{B}_t, \quad (33)$$

$$2(1 - \mathfrak{d}_{FF}) P_{FF} \mathcal{H}_{FF}(\mathbf{x}, t) = \mathcal{G}_{c,FF} \delta \mathfrak{d}_{FF} \gamma(\mathfrak{d}_{FF}, \nabla_{\mathbf{x}} \mathfrak{d}_{FF}) \text{ in } \mathcal{B} \setminus \Gamma_i \text{ and } \nabla_{\mathbf{x}} \mathfrak{d}_{FF} \cdot \mathbf{n} = 0 \text{ in } \partial \mathcal{B}, \quad (34)$$

$$2(1 - \mathfrak{d}_{IFF}) P_{IFF} \mathcal{H}_{IFF}(\mathbf{x}, t) = \mathcal{G}_{c,IFF} \delta \mathfrak{d}_{IFF} \gamma(\mathfrak{d}_{IFF}, \nabla_{\mathbf{x}} \mathfrak{d}_{IFF}) \text{ in } \mathcal{B} \setminus \Gamma_i \text{ and } \nabla_{\mathbf{x}} \mathfrak{d}_{IFF} \cdot \mathbf{n} = 0 \text{ in } \partial \mathcal{B}, \quad (35)$$

$$\operatorname{div}_{\mathfrak{g}} [\boldsymbol{\sigma}_n(\mathfrak{d}_c) + \tau_n(\mathfrak{d}_c)] = 0 \text{ in } \Gamma_i, \quad (36)$$

wherein the previous expressions $\operatorname{div}_{\mathfrak{g}}[\bullet]$ represents the divergence operator taken with respect to \mathfrak{g} . The terms \mathcal{H}_{FF} and \mathcal{H}_{IFF} are the crack driving forces related to fiber and inter-fiber failure, respectively. In accordance with the Pucks failure criteria, the crack driving force of each $j = FF, IFF$ are given by:

$$\mathcal{H}_j(\mathbf{x}, t) = \xi_j \left[\left\langle \frac{\max_{\tau \in [0, t]} \hat{\Psi}_j(\mathbf{x}, \tau)}{\hat{\Psi}_{j,init}} - 1 \right\rangle_+ \right], \quad (37)$$

where ξ_j is a dimensionless fitting parameter that characterizes the damage activation and post peak behaviours from experimental results. $\hat{\Psi}_{j,init}$ is the effective elastic energy for damage initiation in each of $j = FF, IFF$.

The unilateral stationary condition of the total internal energy functional implies that $\delta \Pi_{int} = 0$ for all $(\delta \mathbf{u}, \delta \mathfrak{d}_{FF}, \delta \mathfrak{d}_{IFF}, \delta \mathfrak{g}) > 0$ and $\delta \Pi_{int} > 0$ for $(\delta \mathbf{u}, \delta \mathfrak{d}_{FF}, \delta \mathfrak{d}_{IFF}, \delta \mathfrak{g}) = 0$ along with the irreversibility and boundedness of $\mathfrak{d}_{FF}, \mathfrak{d}_{IFF}, \mathfrak{d}_c$ leads to the first-order optimality (KKT) conditions for the quasi-static evolution [39] [66]. Γ -convergence of the multi phase-field problem in the absence of the interface can be easily derived following the results obtained in [67].

It is worth noting that the irreversible character of the phase-fields and cohesive zone can be reduced to \mathcal{S} , this is fulfilled by the history variable embedded by the history variable Eqn.(37). The boundedness of the phase-field variables $\mathfrak{d}_{FF}, \mathfrak{d}_{IFF} \in [0, 1]$ is ensured due to the choice of degradation function $(1 - \mathfrak{d}_j)^2$ as in [68]. Also, it is important to note that, we have assumed $\nabla_{\mathbf{x}} \mathfrak{d}_j = 0$ due to the compactness property and $\delta \mathbf{u} = 0$ on $\partial \Omega_t$ and $\delta \mathfrak{d}_j = 0$ on $\partial \Omega$, from the variational form which are reflected in the choice of approximate spaces of test functions as in Section 4, see [58, 59].

4. Finite element implementation

In this section, details of the finite element implementation of the proposed model are outlined. A staggered solution scheme is used to solve the system of coupled Partial Differential Equations (PDEs) using an alternating minimization scheme [69]. Consider the discretization of the domain defined as $\mathcal{B} \rightarrow \mathcal{B}^e$,

$\Gamma_i \rightarrow \Gamma_i^e$ such that the functions $\mathbf{u}^e \in \mathcal{U}_h$, $\mathfrak{d}_j^e \in \mathcal{U}_0(\mathfrak{d}_j)$ for $j = FF, IFF$ are well defined along with the space of approximate functions:

$$\mathcal{U}_h(u) = \left\{ u \in H^1(\mathcal{B}) \left| \nabla u \in L^2(\mathcal{B}); u = u_d \text{ on } \partial\mathcal{B}_d \right. \right\}, \quad (38a)$$

$$\mathcal{U}_0(\mathfrak{d}_j) = \left\{ \mathfrak{d}_j \in H^1(\mathcal{B}) \left| \mathfrak{d}_j(\mathbf{x}) \in [0, 1], \mathfrak{d}_j \geq 0, \forall \mathbf{x} \in \mathcal{B} \right. \right\}. \quad (38b)$$

Similarly, the approximate space for the test functions (distributonal spaces) for $\delta \mathbf{u}^e \in \mathcal{V}_h$, $\delta \mathfrak{d}^e \in \mathcal{U}_{\delta 0}(\mathfrak{d}_j)$ for $j = FF, IFF$ takes the form:

$$\mathcal{V}_h(\delta u) = \left\{ \delta u \in H^1(\mathcal{B}) \left| \nabla \delta u \in L^2(\mathcal{B}); \delta u = 0 \text{ on } \partial\mathcal{B}_d \right. \right\}, \quad (39a)$$

$$\mathcal{U}_{\delta 0}(\delta \mathfrak{d}_j) = \left\{ \delta \mathfrak{d}_j \in H^1(\mathcal{B}) \left| \delta \mathfrak{d}_j \geq 0, \forall \mathbf{x} \in \mathcal{B} \right. \right\}. \quad (39b)$$

$$(39c)$$

At each element level, in the isoparametric space settings, the triplet of field variables $\{\mathbf{u}^e, \mathfrak{d}_{FF}^e, \mathfrak{d}_{IFF}^e\}$ as well as their variations $\{\delta \mathbf{u}^e, \delta \mathfrak{d}_{FF}^e, \delta \mathfrak{d}_{IFF}^e\}$ are approximated using linear first order Lagrangian triplet of shape functions $\{\mathbf{N}_i^u, \mathbf{N}_i^{\mathfrak{d}}, \mathbf{N}_i^{\mathfrak{d}}\}$ at i^{th} node of each element satisfying partition of unity is defined as:

$$\mathbf{u}^e = \sum_{i=1}^{\mathbf{N}_{node}} \mathbf{N}_i^u \mathbf{u}_i^e, \quad \mathfrak{d}_j^e = \sum_{i=1}^{\mathbf{N}_{node}} \mathbf{N}_i^{\mathfrak{d}} \mathfrak{d}_{j,i}^e, \quad (40)$$

$$\delta \mathbf{u}^e = \sum_{i=1}^{\mathbf{N}_{node}} \mathbf{N}_i^u \delta \mathbf{u}_i^e, \quad \delta \mathfrak{d}_j^e = \sum_{i=1}^{\mathbf{N}_{node}} \mathbf{N}_i^{\mathfrak{d}} \delta \mathfrak{d}_{j,i}^e, \quad \text{for each } j = FF, IFF. \quad (41)$$

The triplet of spatial derivatives $\{\nabla \mathbf{u}^e, \nabla \mathfrak{d}_{FF}^e, \nabla \mathfrak{d}_{IFF}^e\}$ are approximated using the gradients of the shape functions $\{\mathbf{B}_i^u, \mathbf{B}_i^{\mathfrak{d}}, \mathbf{B}_i^{\mathfrak{d}}\}$ at i^{th} node of each element takes the form:

$$\boldsymbol{\varepsilon}^e = \sum_{i=1}^{\mathbf{N}_{node}} \mathbf{B}_i^u \mathbf{u}_i^e, \quad \nabla \mathfrak{d}_j^e = \sum_{i=1}^{\mathbf{N}_{node}} \mathbf{B}_i^{\mathfrak{d}} \mathfrak{d}_{j,i}^e,$$

$$\delta \boldsymbol{\varepsilon}^e = \sum_{i=1}^{\mathbf{N}_{node}} \mathbf{B}_i^u \delta \mathbf{u}_i^e, \quad \nabla \delta \mathfrak{d}_j^e = \sum_{i=1}^{\mathbf{N}_{node}} \mathbf{B}_i^{\mathfrak{d}} \delta \mathfrak{d}_{j,i}^e, \quad \text{for each } j = FF, IFF.$$

Complying with the formulation of interface cohesive element, the displacement jump vector \mathbf{g} is represented in terms of local frames across the interface Γ_i [57]. Hence, the jump \mathbf{g} and its variation $\delta \mathbf{g}$ is approximated using the kinematic jump-displacement operator $\mathbf{B}^g = R\mathbf{N}^g L$ as:

$$\mathbf{g}^e = \mathbf{B}^g \mathbf{u}^e L, \quad \delta \mathbf{g}^e = \mathbf{B}^g \delta \mathbf{u}^e L, \quad (42)$$

where L matrix estimates the difference between the displacement of the upper and lower interface points and R is a rotation matrix that converts integration points from global to the local frame, and \mathbf{N}^g represents the standard cohesive shape function, see [13].

The discrete elemental residual vectors for the quadruplet $\{\mathbf{u}^e, \mathfrak{d}_{FF}^e, \mathfrak{d}_{IFF}^e, \mathbf{g}^e\}$ can be reduced to the following system of equations:

$$\mathbf{R}_e^{\mathbf{u}} = \int_{\mathcal{B}^e} [(1 - \mathfrak{d}_{FF})^2 (\mathbf{B}^u)^T \bar{\boldsymbol{\sigma}}_{FF} + (1 - \mathfrak{d}_{IFF})^2 (\mathbf{B}^u)^T \bar{\boldsymbol{\sigma}}_{IFF}] d\mathcal{B} - \int_{\mathcal{B}^e} (\mathbf{N}^u)^T \mathbf{f}_v d\mathcal{B} - \int_{\partial \mathcal{B}_u^e} (\mathbf{N}^u)^T \bar{\mathbf{u}} d\partial \mathcal{B} + \mathbf{R}_e^{\mathfrak{g}}, \quad (43)$$

$$\mathbf{R}_e^{\mathfrak{d}_{FF}} = \int_{\mathcal{B}^e} \left\{ \left[\frac{\mathcal{G}_{c,FF}}{l_{FF}} \mathfrak{d}_{FF} - 2(1 - \mathfrak{d}_{FF}) P_{FF} \mathcal{H}_{FF}(\mathbf{x}, t) \right] (\mathbf{N}^{\mathfrak{d}})^T + \mathcal{G}_{c,FF} l_{FF} (\mathbf{B}^{\mathfrak{d}})^T \nabla \mathfrak{d}_{FF} \right\} d\mathcal{B}, \quad (44)$$

$$\mathbf{R}_e^{\mathfrak{d}_{IFF}} = \int_{\mathcal{B}^e} \left\{ \left[\frac{\mathcal{G}_{c,IFF}}{l_{IFF}} \mathfrak{d}_{IFF} - 2(1 - \mathfrak{d}_{IFF}) P_{IFF} \mathcal{H}_{IFF}(\mathbf{x}, t) \right] (\mathbf{N}^{\mathfrak{d}})^T + \mathcal{G}_{c,IFF} l_{IFF} (\mathbf{B}^{\mathfrak{d}})^T \nabla \mathfrak{d}_{IFF} \right\} d\mathcal{B}. \quad (45)$$

where:

$$\mathbf{R}_e^{\mathfrak{g}} = \int_{\Gamma_i^e} (\mathbf{R}\mathbf{B}^{\mathfrak{g}})^T \mathcal{T}(\mathfrak{g}, \mathfrak{d}_c) d\mathcal{S}, \quad (46)$$

is the residual vector associated with the cohesive interface. It is clear that the displacement field \mathbf{u} is strongly coupled with the phase-fields \mathfrak{d}_{FF} , \mathfrak{d}_{IFF} and the displacement jump \mathfrak{g} . Whereas, the phase-fields are among themselves and with jump \mathfrak{g} are decoupled which are evident from Eq. (34), (35), (36) and from the assumption that $\Gamma_i \cap \Gamma_{\mathfrak{d}_i} = \emptyset$. Due to the existence of multiple phase-field and interface, the system of equations describing the fracture is non-linear. Hence, an iterative Newton-Raphson solver is used until the convergence in the sense of cauchy sequence $[(\mathbf{u}_n^{t+1} - \mathbf{u}_n^t)]$ is reached. Here, \mathbf{u}_n^t is the t^{th} iteration at n^{th} step. The corresponding Newton-Raphson iteration to estimate $(n + 1)$ time step takes the form:

$$\begin{bmatrix} \mathbf{u} \\ \mathfrak{d}_{FF} \\ \mathfrak{d}_{IFF} \end{bmatrix}_{n+1} = \begin{bmatrix} \mathbf{u} \\ \mathfrak{d}_{FF} \\ \mathfrak{d}_{IFF} \end{bmatrix}_n - \begin{bmatrix} K^{\mathbf{u}\mathbf{u}} + K^{\mathfrak{g}\mathfrak{g}} & 0 & 0 \\ 0 & K^{\mathfrak{d}_{FF}\mathfrak{d}_{FF}} & 0 \\ 0 & 0 & K^{\mathfrak{d}_{IFF}\mathfrak{d}_{IFF}} \end{bmatrix}_{n+1}^{-1} \begin{bmatrix} \mathbf{R}^{\mathbf{u}} \\ \mathbf{R}^{\mathfrak{d}_{FF}} \\ \mathbf{R}^{\mathfrak{d}_{IFF}} \end{bmatrix}_n. \quad (47)$$

where the corresponding element stiffness matrices read:

$$\mathcal{K}_e^{\mathbf{u}\mathbf{u}} := \frac{\partial \mathcal{R}_e^{\mathbf{u}}}{\partial \mathbf{u}^e} = \int_{\mathcal{B}^e} (\mathbf{B}^u)^T \mathbb{C}^{ep\mathfrak{d}} \mathbf{B}^u d\Omega, \quad (48)$$

$$\mathcal{K}_e^{\mathfrak{g}\mathfrak{g}} := \frac{\partial \mathcal{R}_e^{\mathfrak{g}}}{\partial \mathbf{u}^e} = \int_{\Gamma_i^e} (\mathbf{B}^{\mathfrak{g}})^T \partial_{\mathfrak{g}} \tau \mathbf{B}^{\mathfrak{g}} d\mathcal{S},$$

$$\mathcal{K}_e^{\mathfrak{d}_{FF}\mathfrak{d}_{FF}} := \frac{\partial \mathcal{R}_e^{\mathfrak{d}_{FF}}}{\partial \mathfrak{d}_{FF}^e} = \int_{\mathcal{B}^e} \left[\frac{\mathcal{G}_{c,FF}}{l_{FF}} + 2\mathfrak{P}_{FF} \mathcal{H}_{IFF} \right] \mathbf{N}^{\mathfrak{d}} (\mathbf{N}^{\mathfrak{d}})^T + \mathcal{G}_{c,FF} l_{FF} (\mathbf{B}^{\mathfrak{d}})^T \mathbf{B}^{\mathfrak{d}} d\Omega, \quad (49)$$

$$\mathcal{K}_e^{\mathfrak{d}_{IFF}\mathfrak{d}_{IFF}} := \frac{\partial \mathcal{R}_e^{\mathfrak{d}_{IFF}}}{\partial \mathfrak{d}_{IFF}^e} = \int_{\mathcal{B}^e} \left[\frac{\mathcal{G}_{c,IFF}}{l_{IFF}} + 2\mathfrak{P}_{IFF} \mathcal{H}_{IFF} \right] \mathbf{N}^{\mathfrak{d}} (\mathbf{N}^{\mathfrak{d}})^T + \mathcal{G}_{c,IFF} l_{IFF} (\mathbf{B}^{\mathfrak{d}})^T \mathbf{B}^{\mathfrak{d}} d\Omega. \quad (50)$$

In the Eqn (48), $\mathbb{C}^{ep\mathfrak{d}}$ represents the material consistent tangent estimated using the finite difference method as

$$\mathbb{C}^{ep\mathfrak{d}} = \frac{d\boldsymbol{\sigma}_{ij}}{d\boldsymbol{\varepsilon}_{kl}} \approx \frac{\boldsymbol{\sigma}_{ij}(\hat{\boldsymbol{\varepsilon}}^{kl}) - \boldsymbol{\sigma}_{ij}(\hat{\boldsymbol{\varepsilon}})}{\Delta \boldsymbol{\varepsilon}} \quad (51)$$

with $(\hat{\boldsymbol{\varepsilon}}^{kl}) = \boldsymbol{\varepsilon} + \frac{\Delta \boldsymbol{\varepsilon}}{2} (e_k \otimes e_l + e_l \otimes e_k) = \boldsymbol{\varepsilon} + \Delta \boldsymbol{\varepsilon}^{kl}$. Here e_k and e_l are the k^{th} and l^{th} unit vectors, $\Delta \boldsymbol{\varepsilon}^{kl}$ is the strain perturbation with respect to the components kl and $\Delta \boldsymbol{\varepsilon}$ is the scalar perturbation parameter. Moreover, $\boldsymbol{\sigma}$ is estimated from Eqn (17) as $\boldsymbol{\sigma} = \mathbb{C} : \boldsymbol{\varepsilon}$.

The previous non-linear system of equations has been implemented in the finite element software **ABAQUS**. For this purpose, a user-defined **UMAT** to define the material behaviour is written for the solution of equilibrium equations associated with the displacement field. Whereas, **UEL** is utilized to create residual and stiffness matrix for each of the two phase-fields using the material behaviour from **UMAT** for solving the fracture-associated problem.

5. Representative examples

In this section, modeling application to delamination migration is presented. A comprehensive numerical analysis is presented to validate the model against experimental results. In the sequel, a holistic sensitivity analysis is carried out utilizing the variation of the loading application point, ply angle, and initial pre-crack length to understand their effects on delamination migration.

5.1. Description of the numerical model: general aspects

Figure 4 depicts the baseline configuration under investigation herein. The corresponding numerical model consists of 44 cross-ply IM7/8552 laminates with the layup sequence

$[90_4^{\circ}/0_3^{\circ}/(90^{\circ}/0^{\circ})_{2s}/0_3^{\circ}/CL/90_4^{\circ}/CL/0^{\circ}/0^{\circ}/(90^{\circ}/0^{\circ})_{2s}/0^{\circ}/0^{\circ}/90_3^{\circ}/0^{\circ}/90^{\circ}]$, where CL refers to a cohesive layer. Each ply has a thickness of 0.125mm. Compared with the experimental sequence as in [15], a PTFE (Polytetrafluoroethylene) layer is replaced by a cohesive layer and in addition, another cohesive layer is added at the interface between the 90_4° and 0° sequence in order to account for delamination migration.

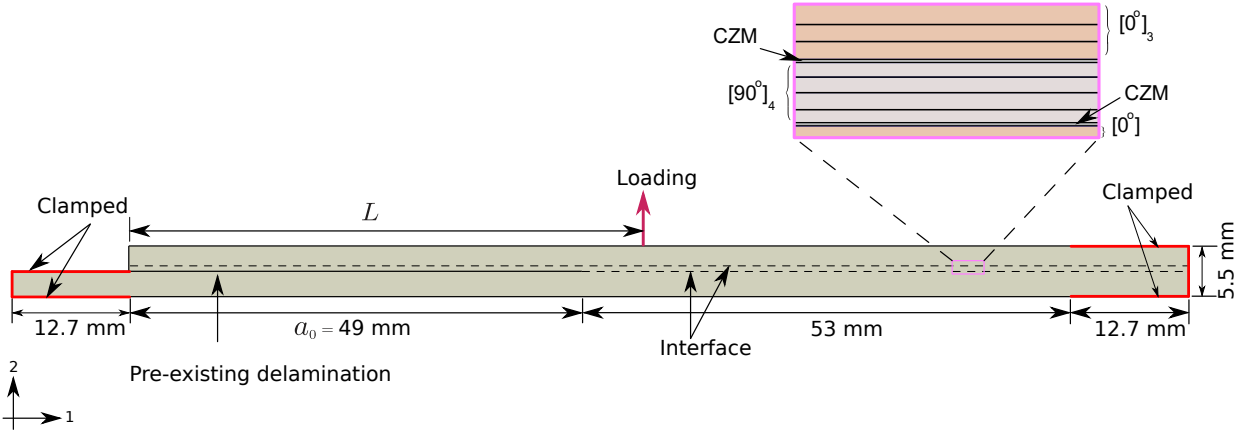


Figure 4: Schematic representation of the delamination migration model.

It is evident from the experimental results reported in [16] that, when shear stresses in the model change sign, migration/kinking occurs, and the crack propagates to the 90_4° layers facilitating the inter-fiber failure.

For each of the numerical simulations conducted in the sequel, a 2D analysis is carried with an out-of-plane thickness of 8.37mm. Since, in order to resolve the gradient of the phase field, the mesh size \bar{h} is restricted to $\bar{h} \approx l/4$. Hence, the domain is discretized by employing 960000 4-node quadrilateral plane stress elements with an average element size of 0.04mm. Each layer of the cross-ply contains at least four elements across its thickness.

The material properties of IM7/8552 ply are shown in Table 1 consistent with the experimental results in [15]. The fracture energy and length scale parameters for the phase-fields chosen according to the material fracture properties are shown in Table 2. The properties of the cohesive layer following [27] are listed in Table 3.

E_{11} (GPa)	E_{22} (GPa)	G_{12} (GPa)	ν_{12}	ν_{23}
161.0	11.38	5.17	0.03	0.43

Table 1: IM7/8552: Elastic properties.

$\mathcal{G}_{c,FF}$ (N/mm)	$\mathcal{G}_{c,IFF}$ (N/mm)	l_{FF} (mm)	l_{IFF} (mm)
81.5	0.2774	0.237	0.07

Table 2: IM7/8552: intra-laminar fracture properties and phase-field parameters.

Nominal stress (MPa)	Nominal stress in shear (MPa)	Nominal fracture energy (N/mm)	Fracture energy in shear (N/mm)	Power (BK law)
15	15	0.5	0.65	2.17

Table 3: Cohesive layer properties.

5.2. Numerical-experimental validation

The global failure response for the specimen under consideration is investigated for the load case $L = a_0 = 49\text{mm}$. The numerical-experimental correlation corresponding to the load-displacement curve is given in Figure 5. As shown in this figure, the failure response can reasonably be divided into three main zones: (i) delamination of the cohesive zone, (ii) kinking followed by migration, and (iii) delamination of the top cohesive zone. Overall, a satisfactory agreement between the numerical and the experimental data can be observed.

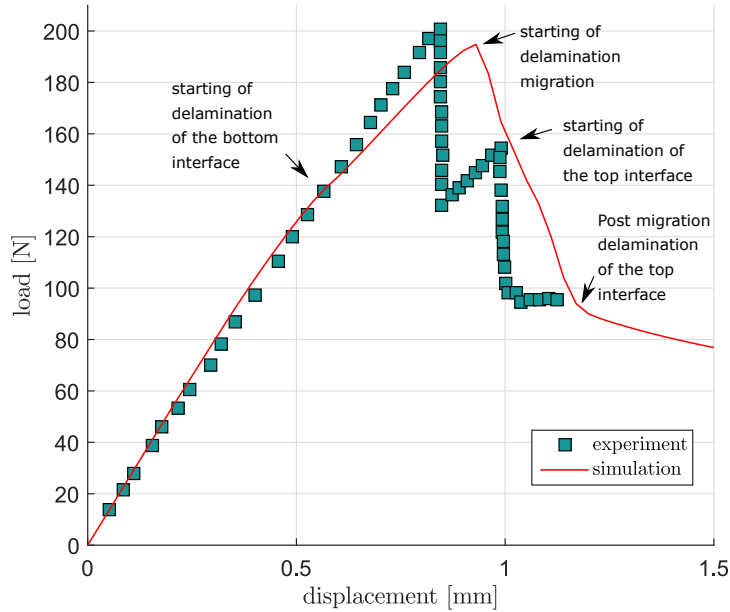


Figure 5: Numerical-experimental correlation corresponding to the load-displacement curve for $L = a_0 = 49\text{mm}$

As was previously discussed, based on postulations made in [15, 16] and the corresponding thorough discussion, delamination migration occurs due to a change of sign in the shear stress components. Negative shearing stresses promote delamination growth at the $0^\circ/90^\circ$ interface, and positive shearing stresses promote migration/kinking into 90° plies. The kinking happens at multiple sites across the specimen. Due to the diffusive nature of the bulk cracks, the shearing stress change can easily be noticed by the inter-fiber phase-field initiation as depicted in Figure 6. Notice that, due to the negative sign initially, delamination propagates until a certain point until shearing stresses are positive. Meanwhile, inter-fiber failure is already initiated, but from the opposite direction, i.e., $90^\circ/0^\circ$ interface, but is not nucleated. Whereas, when the

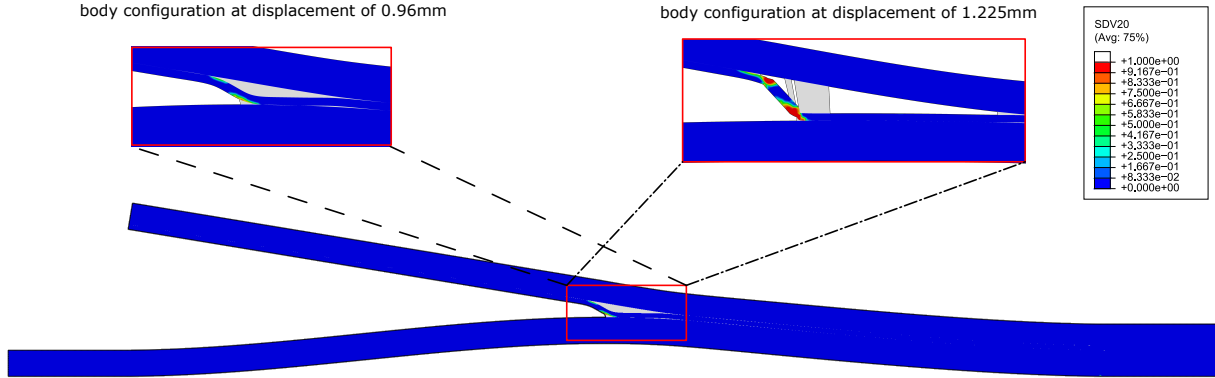


Figure 6: Numerical-experimental correlation corresponding to the load-displacement curve for $L = a_0 = 49\text{mm}$

shear stresses become positive in the adjacent increments, the migration starts developing, with a crack front now migrating into the $90^\circ/0^\circ$ interface.

For a phenomenology of embodiment, in the cohesive layers, once the failure criterion is met, the cohesive layer starts delaminating. Similarly, when the Puck criterion is violated, the inter-fiber failure phase-field is activated due to shearing stresses in the model. As long as the P_{IFF} is active, the inter-fiber failure phase-field crack \mathfrak{d}_{IFF} grows and migrates into the 90° plies until the $90^\circ/0^\circ$ interface. Simultaneously, the top cohesive layer at the $90^\circ/0^\circ$ interface starts to delaminate. When the migration crack front crosses the $90^\circ/0^\circ$ interface, the crack front is again propagating due to the negative shear stress leading to the delamination of the top cohesive layer. Here onward, the residual stresses dominate crack propagation in the model as in Figure 5. The discrepancies between the experimental and numerical results can be attributed to quasi-static load conditions and smooth initiation and propagation of phase-field approximations. However, a satisfactory agreement between the results can be observed.

5.3. Sensitivity analysis: role of different parameters

This section aims at proving a further understanding with regard to the potential role of different design parameters that can favor delamination-migration events.

5.3.1. Effect of position of loading application

The first aspect under analysis concerns the variation of load application point along the specimen. This parameter in the experimental setting might have a strong influence on the activation of migration phenomena by simply inducing a different local stress field at critical locations. Moreover, from a global standpoint, this can have notable effects on the load-displacement curve as the shear stresses acting on the specimen are significantly different from one another. Keeping $a_0 = 49\text{ mm}$, the variation of load L for $L = 0.7a_0, 0.8a_0, 0.9a_0, 1.0a_0, 1.1a_0$ are plotted in Fig. 7.

Based on the current results, it can be observed that for $a_0 > L$, all the cases exhibit delamination prior to migration and show a sudden drop in the load-carrying capacity when migration starts, see Figure 8. Whereas, for $a_0 < L$, the shearing stress sign is favorable for migration at the beginning, and hence there is smooth migration, with delamination spreading over the whole experiment, see Figure 8 which is consistent with [16]. It is also to notice that after delamination migration is finished, the residual stiffness for delamination converges to a single value for all the load variations.

5.3.2. Effect of the variation of ply angle

The second aspect under study has an inherent local effect since it focused on investigating the delamination migration at $0^\circ/\theta_4^\circ$ interface. For this purpose, the original stacking sequence is replaced by a new

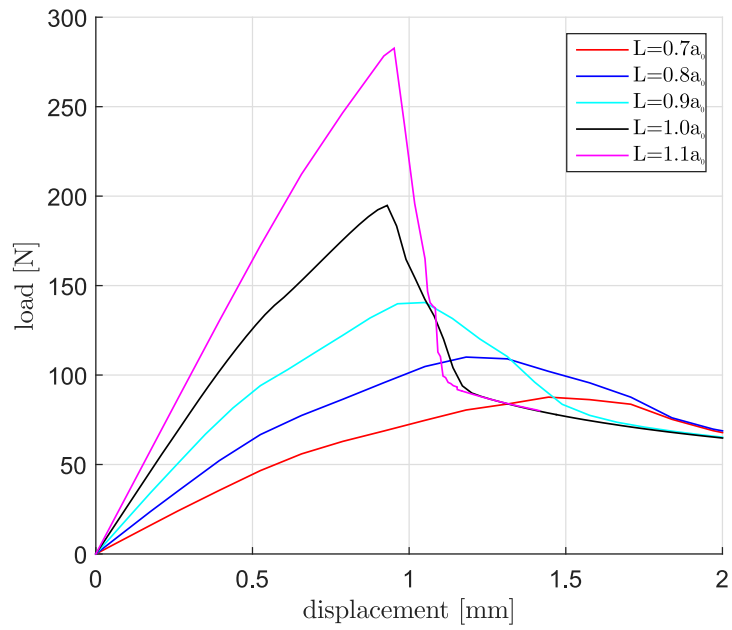


Figure 7: Variation of load across the specimen for $a_0 = 49\text{mm}$.

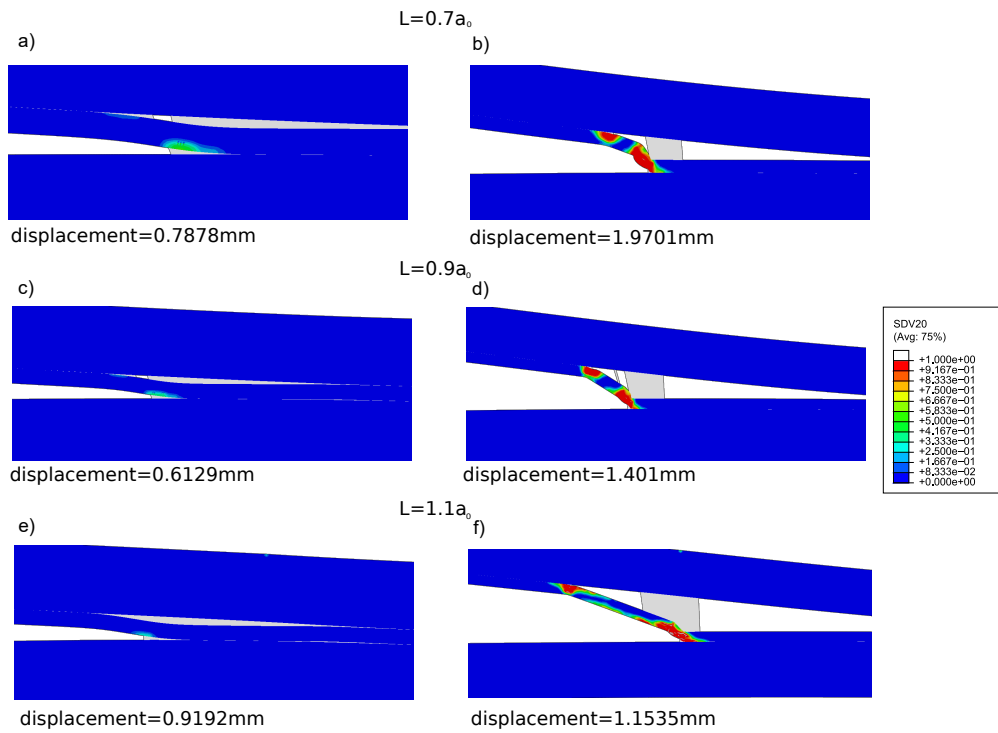


Figure 8: Inter-fiber crack field (SDV20) from 3 cases at $L = 0.7a_0$, $L = 0.9a_0$, $L = 1.1a_0$

layup sequence near the cohesive zones as $0_3^{\circ}/CL/\theta_4^{\circ}/CL/0^{\circ}/0^{\circ}$. The fiber orientation, θ° of 30° , 45° , and 60° is studied along with 90° , and the results are shown in Figure 9.

From this graph, it can be seen that the global pre-peak response is almost unaltered by the variation of the local orientation of the adjacent layers to the $0^{\circ}/\theta_4^{\circ}$ interface. However, this aspect has a notable influence on the post-peak response, delaying the delamination initiation and the subsequent delamination event of the top interface, according to the description given in Figure 6, but however, the cracking migration is predicted to occur at almost the same loading level (along the post-peak evolution).

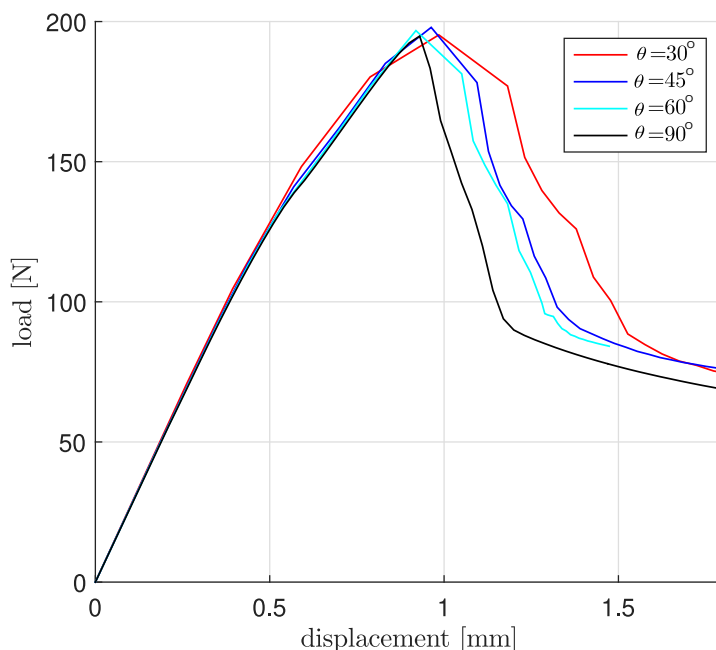


Figure 9: Variation of angle across the specimen for $L = a_0 = 49\text{mm}$.

5.3.3. Effect of the variation of initial crack length

The last effect under consideration corresponds to the initial crack length. Focusing on very specific cases, we vary initial crack size from $a_0 = 49\text{ mm}$ to $a_0 = 55\text{ mm}$.

The global load-displacement evolution curves for such cases are shown in Figure 10. According to these data, it can be stated that a simple variation of the initial crack length has a very remarkable role in the specimen response. Thus, observing the pre-peak evolution, before any inelastic process commences, the larger the initial crack length is set, the higher the maximum load is achieved. Moreover, concerning the post-peak evolution, it is observable that while the shorter initial crack-length case evidence similar evolution with respect to those previously described, i.e., with the occurrence of delamination events and the posterior cracking migration to the adjacent layer, the response of $a_0 = 55\text{mm}$ (and $a_0 > 55\text{mm}$) exhibited an entirely different evolution with no evidence of cracking migration, see Figure 10, and 11 where the matrix-failure maps for both configurations are depicted. These differences in the response are directly associated with the local stress state's discrepancies at the crack tips at the interface and the intermediate layer. This is again in line with previous studies as in [16, 15].

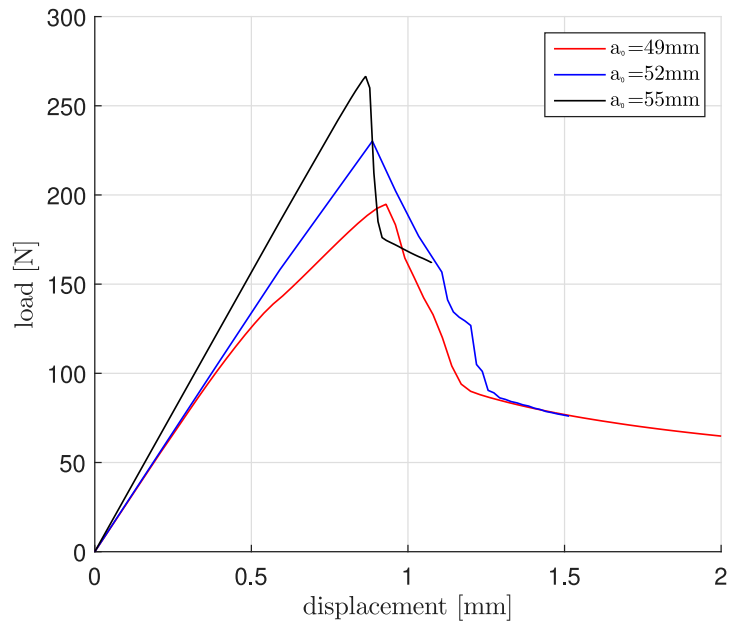


Figure 10: Variation of initial crack length across the specimen for $L = a_0$.

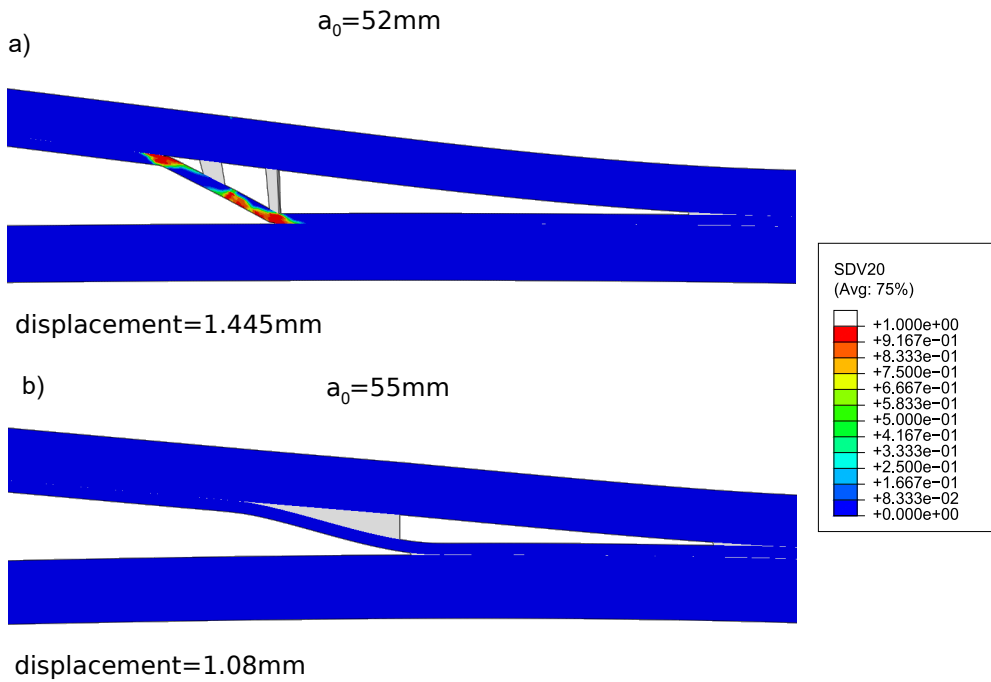


Figure 11: Inter-fiber crack field (SDV20) for initial crack length $a_0 = 52\text{mm}$, $a_0 = 55\text{mm}$ with $L = 1.0a_0$

6. Conclusion

A consistent multi-phase field-cohesive zone model relying on phenomenological failure criterion has been proposed for matrix-dominated cracking in the presence of interfaces. Due to fiber failure, inter-fiber failure,

and interface debonding, the failure mechanism is all accommodated in the model by considering a multi-phase field, each characterized by their failure mechanism with corresponding fracture energy and internal length scale, plus a cohesive zone model.

The computational framework has been carefully derived via multi-field variational formulation with multiple dissipative mechanisms within the spirit of phase-field and cohesive zone models such that thermodynamics consistency is preserved.

The model has been applied to the study of delamination migration in composite materials to illustrate the capabilities of the model. The variation of the design parameters, such as the load application point, the angle between the plies, and the initial crack length, have been presented to assess the model response's sensitivity to such design parameters.

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