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Generalized displacement correlation method for estimating stress intensity factors

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ABSTRACT

Conventional displacement-based methods for estimating stress intensity factors require special quarter-point finite elements in the first layer of elements around the fracture tip and substantial near-tip region mesh refinement. This paper presents a generalized form of the displacement correlation method (the GDC method), which can use any linear or quadratic finite element type with homogeneous meshing without local refinement. These two features are critical for modeling dynamic fracture propagation problems where locations of fractures are not known a priori. Because regular finite elements' shape functions do not include the square-root terms, which are required for accurately representing the near-tip displacement field, the GDC method is enriched via a correction multiplier term. This paper develops the formulation of the GDC method and includes a number of numerical examples, especially those consisting of multiple interacting fractures. We find that the proposed method using quadratic elements is accurate for mode-I and mode-II fracturing, including for very coarse meshes. An alternative formulation using linear elements is also demonstrated to be accurate for mode-I fracturing, and acceptable mode-II results for most engineering applications can be obtained with appropriate mesh resolution, which remains considerably less than that required by most other methods for estimating stress intensities.

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1. Introduction

The stress intensity factor (SIF) is an important concept in fracture mechanics for relating stress and energy release rate at the fracture tip to loading and crack geometry. Although closed-form analytical solutions are available for a number of special fracture-load configurations (many of which have been compiled in [1]), SIF's are often calculated in the context of numerical methods, especially the finite element method (FEM) for arbitrary fracture-load configurations. Several methods are available for calculating or estimating SIF's with the FEM, such as the *J*-integral [2] and its variants, the stiffness derivative technique [3], and a suite of methods based on the interpretation of near-tip nodal displacements. In the last category, there are at least three variants, including the displacement extrapolation method [4–7], the quarter-point displacement method [8], and the displacement correlation method [9,10]. These methods and others have been compared in a number of studies [e.g. 5,11–14]. One of the most significant advantages of the displacement-based methods is the simple formulation. Although the displacement-based methods were often found to be less accurate than the *J*-integral or the stiffness derivative method under certain conditions, the accuracy remains acceptable for most engineering applications [e.g. 5,14]. Many of the displacement-based methods were developed in the 1970s and 1980s in tandem with various special "quarter-point" finite element types [15–19] and transition elements [20] used in these methods. Though few new developments have been reported on the displacement-based methods in the intervening decades [21], they continue to be widely used.

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Nomenclature

a, b, c, h variables denoting key geometrical characteristics of fracture-solid systems
$C_{I}^{A}, C_{I}^{B}, C_{I}^{B}, C_{I}^{B}$ correction multipliers for the GDC method. The superscript denotes whether it is for Method A or Method B;
the subscript indicates the mode of fracture
$f_r^a, f_r^b, f_r^c, f_\theta^a, f_\theta^b, f_\theta^c$ functions of the angular coordinate of a point to simplify the expression of certain equations
<i>F_I</i> , <i>F_{II}</i> fracture-configuration correction factors related to geometrical characteristics of the fracture-solid system
<i>G</i> shear modulus of the solid
<i>K_I</i> , <i>K_{II}</i> mode-I and mode-II stress intensity factors
l_E characteristic length of a finite element
<i>P</i> point load applied on a bending beam with a mid-span notch
r, θ polar coordinates of a point
s span width of a bending beam with a mid-span notch
$u_r^a, u_r^b, u_r^c, u_{\theta}^a, u_{\theta}^b, u_{\theta}^c$ radial and angular displacements of a point due to the three boundary condition modes (<i>a</i> , <i>b</i> , and <i>c</i> de
noted in Fig. 1)
u_r , u_θ radial and angular displacements of a reference point. A superscript might be used to denote the reference point
β material constant depending on whether this is a plain-stress or plain-strain problem
δ mesh perturbation factor
<i>v</i> Poisson's ratio of the solid
$\sigma_{f_1} \tau_{f_2} \sigma_x, \sigma_y$ far-field stress components of the fracture-solid system
σ_c, τ_c fracture surface stress components
GDC generalized displacement correlation (method)
SIF stress intensity factor
-

The J-integral as well as its variants (e.g. M-integral) and the displacement-based methods all require accurate resolution of near-tip displacement/stress/strain fields. Therefore special element types (e.g. quarter-point elements) and/or near-tip element refinement are usually used as accuracy enhancement/assurance measures. The present study is motivated by engineering applications where such enhancement is impractical in terms of computational cost but a moderate error margin (e.g. 10%) is acceptable. The simulation of hydraulic fracturing in natural fracture systems represents this class of applications and is the direct motivation for the current study [22]. Simulating a hydraulic fracturing process usually involves modeling multiple cracks propagating along arbitrary paths, so the locations of the crack tips are not specified *a priori*. Employing any near-tip enhancement measures necessitates heterogeneous meshing, which is theoretically possible to handle but requires complex variable mapping between meshes every time a fracture advances. This is extremely costly for hydraulic fracturing simulations because very small time steps have to be adopted to handle the wide spectrum of length-scales that have to be resolved, spanning from tens of micrometers (the aperture width of typical rock joints) to hundreds of meters (dimension of the reservoir). Therefore, homogeneous and relatively coarse meshing without local refinement or frequent remeshing and variable mapping is the only viable option. Additionally, because we have to frequently handle the situation where two fractures are close to each other before they intersect, it is desired to only use information in the first layer of element surrounding a fracture tip. On the other hand, owing to the inherently high variability in rock properties and the high uncertainty in the determination of rock properties, an error of 10-20% in the estimation of SIF is considered well acceptable.

The goal of this study is to develop and verify a displacement-based method, termed the generalized displacement correlation (GDC) method for estimating SIF, which uses regular finite element types and does not require local mesh refinement. In the currently paper, we first review the mechanical and mathematical principles behind the original displacement-based methods in a generalized context in Section 2. Compared with the original derivation of these methods, the loading condition is generalized by including crack surface traction and the meshing scheme is generalized by circumventing the dependency on the specific shape functions of quarter-point elements. This new GDC formulation encompasses the original formulation based on quarter-point elements as a special case. Subsequently, we develop the new generalized formulation in Section 3 and further enhance its accuracy in Section 4 by introducing an empirical correction multiplier term. In Section 5, we test the new method against a number of fracture-load configurations with an emphasis on cases with inter-crack inter-actions, a situation critical to our hydraulic fracturing simulator development effort. The numerical examples in Sections 4 and 5 use the same Poisson's ratio and tip-region mesh configuration and use meshes based on a regular grid. The sensitivity of the results to the Poisson's ratio, near-tip mesh configurations, and mesh perturbation are evaluated in Sections 6 and 7.

2. Review of displacement-based methods in a generalized framework

Consider the two-dimensional (2D) continuum (linearly elastic and isotropic) around a crack tip as shown in Fig. 1, with far-field normal (σ_f) and shear (τ_f) stress existing along with crack surface traction (σ_c and τ_c). Note that "traction" in this



Fig. 1. The near-tip region of a 2D medium and the decomposition of fracture modes according to the superposition principle. The polar coordinate system used in this study is denoted in the figure. Fracture openings in this and other examples are exaggerated for illustration purposes.

paper refers to stress distributed along fracture surface while the same term is often used in cohesive zone models for a different meaning. Stresses σ_f , τ_f , σ_c , and τ_c are independent of each other, but the spatial variation of each of them is not considered. Their values can be either positive or negative, with the arrows in Fig. 1 indicating the positive stress directions. According to the superposition principle, the mechanical response of this system is the sum of the responses of the three cases [(a)-(c)] to the right of the equal sign in the figure. Case (a) and case (b) respectively correspond to the classical boundary/loading conditions for mode-I and mode-II fracturing, whereas in case (c) the crack surface traction balances the far-field stress. Only the stress conditions in the two former cases [(a) and (b)] induce stress/strain singularities in the near-tip region, while the latter case (c) generates homogeneous stress and displacement fields which contribute to the overall mechanical response but not the near-tip stress singularity. The loading conditions in case (a) and case (b) are the symmetric and skew-symmetric (antisymmetric) parts of the load that induce a near-tip stress singularity, respectively. Much of the development of *fracture mechanics* disregards the tractions along the crack surface, so case (a) and case (b) have been the focus of previous studies. In certain applications such as hydraulic fracturing, the pressure inside the fractures is the main mechanism for driving fracture extension with $\sigma_c < \sigma_f < 0$. Under such conditions, the stress condition in case (c) significantly contributes to the mechanical responses of the system and cannot be overlooked.

With higher-order terms omitted, the displacement field (relative to the crack tip displacement) induced by loading case (a) is

$$\begin{cases} u_r^a \\ u_\theta^a \end{cases} = \frac{K_I}{G} \sqrt{\frac{r}{2\pi}} \begin{cases} \cos\frac{\theta}{2} \\ -\sin\frac{\theta}{2} \end{cases} \left(\beta - \cos^2\frac{\theta}{2}\right)$$
(1)

where K_l is the mode-I stress intensity factor; G is the shear modulus of the medium; β is a constant depending on whether this is a plane strain ($\beta = 2[1 - v]$ with v being the Poisson's ratio) or a plane stress ($\beta = 2/[1 + v]$) problem. It we assume that the elasticity parameters (G and β) are constants for a given problem, the equation can be simplified as

$$\begin{cases} u_r^a \\ u_\theta^a \end{cases} = K_I \sqrt{r} \begin{cases} f_r^a(\theta) \\ f_\theta^a(\theta) \end{cases}$$
 (2)

where $f_r^a(\theta)$ and $f_{\theta}^a(\theta)$ are functions of the angular coordinate (θ) of the point where the displacement is measured. The effects of the elasticity parameters are incorporated into these two functions and they are considered constants for the purpose of this section. We can also write the corresponding equations for case (b), namely mode-II fracturing as

$$\begin{cases} u_r^b \\ u_{\theta}^b \end{cases} = \frac{K_{II}}{G} \sqrt{\frac{r}{2\pi}} \begin{cases} -\sin\frac{\theta}{2}(\beta - 3\cos^2\frac{\theta}{2}) \\ -\cos\frac{\theta}{2}(\beta + 2 - 3\cos^2\frac{\theta}{2}) \end{cases} = K_{II}\sqrt{r} \begin{cases} f_r^b(\theta) \\ f_{\theta}^b(\theta) \end{cases}$$
(3)

Loading in Fig. 1c induces a homogeneous stress field quantified by σ_c , σ_x , and τ_c . σ_x is the normal stress component (not denoted in Fig. 1) in the direction along the fracture tip, and is typically not concerned in fracture mechanics. The displacement induced by this homogeneous stress field is

$$\begin{cases} u_r^c \\ u_\theta^c \end{cases} = r \begin{cases} f_r^c(\theta, \sigma_c, \sigma_x, \tau_c) \\ f_\theta^c(\theta, \sigma_c, \sigma_x, \tau_c) \end{cases}$$

$$(4)$$

or

$$\begin{cases} u_r^c \\ u_\theta^c \end{cases} = r \begin{cases} f_r^c(\theta) \\ f_\theta^c(\theta) \end{cases}$$
 (5)

for any known stress state (σ_c , σ_x , τ_c). The explicit expression of functions f_r^c and f_{θ}^c can be derived based on Hooke's law, but it requires knowledge of the stress state and is not pursued here. Note that the f^c terms also encompass the effects of small rigid-body rotation of the system, but this is not explicitly discussed in the following development. The most important implication of Eq. (5) for the scope of this paper is that along any "ray" direction originating from the fracture tip, the displacement of any point relative to that of the tip is linearly proportional to its distance to the crack tip under the homogeneous stress condition.

Combining Eqs. (2), (3), and (5), we can write the overall displacement field for the arbitrary loading condition in Fig. 1 as

$$\begin{cases} u_r \\ u_\theta \end{cases} = \sqrt{r} \begin{cases} f_r^a K_I + f_r^b K_{II} \\ f_\theta^a K_I + f_\theta^b K_{II} \end{cases} + r \begin{cases} f_r^c \\ f_\theta^c \end{cases}$$
 (6)

with K_I and K_{II} being the unknowns while u_r and u_θ can be obtained from FEM solutions.

In order to apply any displacement-based stress intensity calculation method, the medium containing the fracture needs to be modeled using a finite element mesh. Quarter-point elements, with the inverse square root singularity embedded by shifting the mid-edge nodes on the ray edges to the quarter-points, are usually employed as the first layer of elements around the tip as shown in Fig. 2. Displacements along the crack face ($\theta = \pi$) at nodes *A* and *B* are obtained by solving the finite element model. Noticing that $f_{\theta}^{a}(\pi) = 0$ and $f_{\theta}^{b}(\pi) = 0$, we have

$$u_r^A = \frac{1}{2}\sqrt{l_E}f_r^b(\pi)K_{II} + \frac{1}{4}l_Ef_r^c(\pi)$$
(7)

$$u_r^B = \sqrt{l_E} f_r^b(\pi) K_{II} + l_E f_r^c(\pi)$$
(8)

$$u_{\theta}^{A} = \frac{1}{2}\sqrt{I_{E}}f_{\theta}^{a}(\pi)K_{I} + \frac{1}{4}I_{E}f_{\theta}^{c}(\pi)$$

$$\tag{9}$$

$$u_{\theta}^{B} = \sqrt{l_{E}} f_{\theta}^{a}(\pi) K_{I} + l_{E} f_{\theta}^{c}(\pi)$$
⁽¹⁰⁾

where l_E is the length of the element edge ($l_E = |TB| = 4|TA|$ in this particular case). By applying basic linear equation manipulation/solving techniques, we can eliminate the terms involving f_r^c or f_{θ}^c and obtain

$$K_I = \frac{4u_{\theta}^A - u_{\theta}^B}{\sqrt{l_E f_{\theta}^a(\pi)}} \quad \text{and} \quad K_{II} = \frac{4u_r^A - u_r^B}{\sqrt{l_E f_r^b(\pi)}}$$
(11a)

which is the core formulation for the displacement correlation method. The symmetry of the system can be exploited to improve the accuracy of the results with

$$K_{I} = \frac{4(u_{\theta}^{A} - u_{\theta}^{A'}) - (u_{\theta}^{B} - u_{\theta}^{B'})}{2\sqrt{I_{E}}f_{\theta}^{a}(\pi)} \quad \text{and} \quad K_{II} = \frac{4(u_{r}^{A} - u_{r}^{A'}) - (u_{r}^{B} - u_{r}^{B'})}{2\sqrt{I_{E}}f_{r}^{b}(\pi)}$$
(11b)



Fig. 2. Quarter-point element configurations near a crack tip.

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The formulation for the so-called quarter-point displacement method

$$K_I = \frac{u_{\theta}^A - u_{\theta}^{A'}}{\sqrt{l_E f_{\theta}^a(\pi)}} \quad \text{and} \quad K_{II} = \frac{u_r^A - u_r^{A'}}{\sqrt{l_E f_r^b(\pi)}}$$
(12)

is valid only if the terms involving f_r^c and f_{θ}^c in Eq. (6) vanish, implying the loading of the system is the sum of case (a) and case (b) excluding case (c) in Fig. 1, i.e. there is no traction along the crack faces. This limitation of the quarter-point displacement method was described by Tracey [10] but has largely been neglected, as it does not apply to the typical loading conditions in mechanical engineering, where crack surface tractions are absent. Although this limitation of the quarter-point displacement method does not lead to inaccuracies in many studies comparing these two methods in the context of mechanical engineering [12,13,19,23], it is highly deleterious if the method is to be used for hydraulic fracturing modeling or similar problems. The displacement extrapolation method suffers similarly since the loading scenario shown in case (c) of Fig. 1 is not supported in the assumptions underlying that method. Based on this, we select the displacement correlation method as the basis for further development.

The original development of the displacement correlation method and the quarter-point displacement method derive the same equations as Eqs. (11) and (12), respectively, through a different procedure. The purpose of the above development is to provide the necessary basis for the development of the new generalized method in the next section.

3. Formulation of the generalized method

From the procedure in Section 2, we see that the core of the displacement correlation method is to solve equations of the following form

$$u_i = f_i^a \sqrt{r_i} K_I + f_i^b \sqrt{r_i} K_{II} + r_i f_i^c$$
(13)

where u_i, f_i^a , and f_i^b are known from FEM solutions of the specific fracture-load configuration and near-tip region closed-form solutions; K_I and K_{II} are the two unknowns to solve; f_i^c can be removed by the following procedure. Because f_i^c is a function of the angular coordinate θ but not the radial coordinate r, we can use known displacements (either u_r or u_θ) and other information (r_i, f_i^a , and f_i^b) at two points with the same angular coordinate θ to eliminate the f_i^c term. The symmetry and/ or skew-symmetry of f_i^a and f_i^b can also be used to directly eliminate K_I or K_{II} when solving for the other. The choice of the four displacement components in obtaining Eqs. (7)–(10), namely $u_r^A = u_r(l_E/4, \pi)$, $u_r^B = u_r(l_E, \pi)$, $u_{\theta}^A = u_{\theta}(l_E/4, \pi)$, and $u_{\theta}^B = u_{\theta}(l_E/4, \pi)$ allows this approach. $r_i = l_E/4$ and $r_i = l_E$ are used for convenience to exploit nodal displacements in the quarter-point elements. However, displacements at other points (not necessarily nodes) can be used instead to solve Eq. (13).

Through this generalization of the original displacement correlation method, the special quarter-point element and neartip region mesh refinement can be eliminated, and we can substitute the displacements at appropriate reference points and other necessary information into Eq. (13) to solve for SIF's. In the selection of the reference points, we first consider points with $\theta = \pm \pi$, consistent with the original displacement correlation method, where the features of $f_r^a(\pi) = 0$ and $f_{\theta}^b(\pi) = 0$ simplifies solution. If quadratic elements (i.e. shape functions are second-degree polynomials) are used, we can use $r = l_E/2$ and $r = l_E$, which are both within the first layer of elements about the crack tip. Appealing to symmetry, we have

$$u_r(l_E/2,\pi) - u_r(l_E/2,-\pi) = \sqrt{2l_E} f_r^b(\pi) K_{II} + \frac{l_E}{2} \left[f_r^c(\pi) - f_r^c(-\pi) \right]$$
(14a)

$$u_r(l_E,\pi) - u_r(l_E,-\pi) = 2\sqrt{l_E} f_r^b(\pi) K_{II} + l_E [f_r^c(\pi) - f_r^c(-\pi)]$$
(14b)

$$u_{\theta}(l_{E}/2,\pi) - u_{\theta}(l_{E}/2,-\pi) = \sqrt{2l_{E}}f_{\theta}^{a}(\pi)K_{I} + \frac{l_{E}}{2}\left[f_{\theta}^{c}(\pi) - f_{\theta}^{c}(-\pi)\right]$$
(14c)

$$u_{\theta}(l_{E},\pi) - u_{\theta}(l_{E},-\pi) = 2\sqrt{l_{E}}f_{\theta}^{a}(\pi)K_{I} + l_{E}[f_{\theta}^{c}(\pi) - f_{\theta}^{c}(-\pi)]$$
(14d)

Solving the above equations yield the formulation for the generalized displacement correlation (GDC) method as

$$K_{I} = \frac{2u_{\theta}(l_{E}/2,\pi) - 2u_{\theta}(l_{E}/2,-\pi) - u_{\theta}(l_{E},\pi) + u_{\theta}(l_{E},-\pi)}{(2\sqrt{2}-2)\sqrt{l_{E}}f_{\theta}^{a}(\pi)}$$
(15)

$$K_{II} = \frac{2u_r(l_E/2,\pi) - 2u_r(l_E/2,-\pi) - u_r(l_E,\pi) + u_r(l_E,-\pi)}{(2\sqrt{2}-2)\sqrt{l_E}f_r^b(\pi)}$$
(16)

where the constants $f_{\theta}^{a}(\pi) = f_{r}^{b}(\pi) = -\beta/\sqrt{2\pi}G$ follow from Eqs. (1)–(3). This set of equations does not require quarter-point elements around the crack tip, but does require quadratic elements (6-node triangle or 8-node quadrilateral in 2D). Since the objective of this paper is to generalize the displacement correlation method, we further consider finite element models where linear elements (3-node triangle or 4-node quadrilateral) are used. Under this condition, Eqs. (15) and (16) result in zero SIF's owing to the linear shape functions. Using displacements across two layers of elements around the tip (i.e. at $r = l_E$ and $r = 2l_E$) and replacing $l_E/2$ in the above equations with l_E and l_E with $2l_E$ solve this problem, but renders the method impractical for modeling fractures with arbitrary paths. Fig. 3 shows two problematic scenarios commonly addressed through FEM modeling of fractures: (a) sawtooth-shaped fractures typical in perturbed meshes where minor perturbation

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Fig. 3. Two common scenarios where the locations of points $(2l_{e}, \pi)$ and $(2l_{e}, -\pi)$ are ambiguous.

to node locations in the undeformed mesh is adopted to introduce randomness into fracture paths, and (b) a fracture having changed the direction of propagation. In both scenarios, the locations of points $(2l_E, \pi)$ and $(2l_E, -\pi)$ are ambiguous, making the method above inapplicable. To address this, we use displacements of points with $\theta = -\pi/2$, 0, and $\pi/2$ and $r = l_E$ and $r = 2l_E$, and also exploit the symmetry of f_{θ}^a and f_{θ}^b and skew-symmetry of f_{θ}^a and f_{r}^b to obtain

$$u_r(l_E, \pi/2) + u_r(l_E, -\pi/2) = 2\sqrt{l_E} f_r^a(\pi/2) K_I + l_E[f_r^c(\pi/2) + f_r^c(-\pi/2)]$$
(17a)

$$u_r(2l_E, \pi/2) + u_r(2l_E, -\pi/2) = 2\sqrt{2l_E f_r^a(\pi/2)} K_I + 2l_E [f_r^c(\pi/2) + f_r^c(-\pi/2)]$$
(17b)

$$u_{\theta}(l_E, \mathbf{0}) = \sqrt{l_E f_{\theta}^b(\mathbf{0}) K_{II} + l_E f_{\theta}^c(\mathbf{0})}$$
(17c)

$$u_{\theta}(2l_{E},\mathbf{0}) = \sqrt{2l_{E}}f_{\theta}^{b}(\mathbf{0})K_{II} + 2l_{E}f_{\theta}^{c}(\mathbf{0}) \tag{17d}$$

which yield

$$K_{I} = \frac{2u_{r}(l_{E}, \pi/2) + 2u_{r}(l_{E}, -\pi/2) - u_{r}(2l_{E}, \pi/2) - u_{r}(2l_{E}, -\pi/2)}{(4 - 2\sqrt{2})\sqrt{l_{E}}f_{r}^{a}(\pi/2)}$$
(18)

$$K_{II} = \frac{2u_{\theta}(l_{E}, 0) - u_{\theta}(2l_{E}, 0)}{(2 - \sqrt{2})\sqrt{l_{E}}f_{\theta}^{b}(0)}$$
(19)

where the constants $f_r^a(\pi/2) = (2\beta - 1)/4\sqrt{\pi}G$ and $f_{\theta}^b(0) = (1 - \beta)/\sqrt{2\pi}G$. We term the GDC method based on Eqs. (15) and (16) "Method A", and that based on Eqs. (18) and (19) "Method B". Method B can be applied to any finite element types, and is therefore "more general" than Method A. Method A only requires displacements across one layer of elements around the tip while Method B requires two layers. Neither Method A nor Method B requires a special meshing scheme at the near-tip region, such as a mesh type or mesh resolution different from that of the remainder of the computation domain. Both methods are easy to implement in existing FEM packages. Note that the points where displacements are used in the calculation need not to be nodes of the finite element mesh.

4. Enhancement of the generalized method

Error in the calculated stress intensity factors using the GDC method can be attributed to at least two sources:

- (1) The inability of the adopted finite element's shape functions to accurately represent the near-tip displacement field. The quarter-point element family was originally formulated for the very purpose of better representing the near-tip field by including a square-root term in the shape functions in the ray directions.
- (2) Omission of higher-order terms in Eqs. (1) and (3). These equations are accurate at the near-tip region, where the distances to the fracture tip and other sources inducing high displacement gradient are much smaller than the length of the fracture itself. In the GDC method, displacements at distances l_E and $2l_E$ (or $l_E/2$ and l_E) are used. Therefore, error increases with the ratio of element size to the fracture length.

In order to demonstrate the accuracy of the GDC method, we use the proposed method on the simplest fracture system, i.e. a finite-length fracture in an infinite domain as shown in Fig. 4. The fracture system considered here is straight crack of length 2*a* in a 2D infinite medium. Since most FEM models can accurately represent the linear displacement field induced by the loading condition in Fig. 1c, only the loading conditions in Fig. 1a and b are combined and modeled. However, the effects of homogeneous stress fields are appropriately handled in the formulations of the GDC method, and the superposition of

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Fig. 4. A finite-length crack in an infinite medium.



Fig. 5. Four mesh configurations considered in this study. The conventional six-node triangle element is used in all the numerical examples of the present study but the mid-edge node is not shown in this figure.

such a field would not affect the calculated SIF's. The near-tip mesh configuration can have a considerable effect on the accuracy of the original displacement-based methods (e.g. [23]); in all the numerical examples in the current and next section, the mesh configuration shown in Fig. 5a is used, and fracture tips are located at nodes shared by eight triangular elements. The other mesh configurations shown in Fig. 5 will be investigated in Section 6. In linearly elastic problems, the shear modulus of the medium, *G* does not affect the calculated stress intensity factors and thus can be arbitrarily selected. The model is assumed to be a plain-stress problem with a Poisson's ratio of 0.2. The effects of the Poisson's ratio will also be discussed in Section 6. The finite element mesh is sufficiently large (with each dimension longer than 100*a*) such that the effects of the finite boundaries are minimal and the domain can be considered infinite. We use quadratic (6-node) triangle elements with full-integration (three Gaussian points) for both Methods A and B in this study, although Method B is not restricted to quadratic elements.

The theoretical solutions for the stress intensity factors in this crack configuration are $K_I = \sqrt{\pi a \sigma_y}$ and $K_{II} = \sqrt{\pi a \tau}$. Numerical solutions of the SIF's, denoted by K'_I and K'_{II} are obtained by solving finite element models with various levels of mesh resolutions (quantified by a/l_E , the ratio of the half crack length to element length) and substituting the obtained displacement values into Eqs. (15) and (16) or (18) and (19). We then seek an enhancement measure in the form of a "correction multiplier" to be added to Eqs. (15), (16), (18), and (19). We will test the performance of the corrected/enhanced formulation on a number of more complex crack systems in next section for Methods A and B. The values of $C_I = K_I/K'_I$ and $C_{II} = K_{II}/K'_{II}$, which are the multipliers that need to be applied to Eqs. (15) and (16) or (18) and (19), respectively to correct the numerical solutions are shown in Fig. 6 as functions of a/l_E . The correction factors are significantly larger than unity, since the 6-node triangular finite element cannot accurately represent the near-tip displacement field. C_I and C_{II} both converge to constant values as the element size becomes smaller relative to the crack length. We can fit the discrete data points with the following empirical relationship

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Fig. 6. The effects of the mesh resolution on the correction multipliers. (a) Results for Method A; (b) results for Method B.

$$C = \frac{\alpha_1}{\sqrt{1 - \alpha_2 l_E/a}} \tag{20}$$

which has a similar format as the correction term used in [24]. The regression results are

$$C_I^A = \frac{1.555}{\sqrt{1 + 0.640 l_E/a}} \tag{21a}$$

$$C_{II}^{A} = \frac{2.831}{\sqrt{1+1.163l_{e}/g}}$$
(21b)

$$C_{I}^{B} = \frac{1.260}{\sqrt{1 - 0.138l_{F}/a}}$$
(21c)

$$C_{II}^{B} = \frac{1.727}{\sqrt{1 - 0.845 l_{E}/a}}$$
(21d)

where the superscripts A and B of C_I and C_{II} indicate whether the correction multipliers are for Method A or Method B. The coefficients of determination (R^2) for all regressions are greater than 0.99.

The correction multipliers calculated using Eq. (21) converge but not to unity. This appears counterintuitive because even though the shape functions (quadratic for the above calculations and linear if linear elements were used) of a single element does not accommodate the square root terms in Eq. (6), refining the mesh (with smaller l_E) should result in piecewise quadratic shape functions for the mesh as a whole better representing the displacement field. However, regardless of the refinement level, only displacements within the first one (Method A) or two (Method B) layers of elements around the fracture tip are used. As the mesh is refined, the reference points where displacement information is used in the calculation are closer to the fracture tip. For infinitesimal elements, this mechanism can eliminate the error induced by the second source of error, but not the first. A similar phenomenon exist for the original displacement-based methods: Numerous studies have observed that errors of these methods do not converge to zero as the near-tip mesh is refined [12,13,18,19,23] and an explanation was offered by Harrop [25].

5. Accuracy of the generalized method for different fracture configurations

The values as well as the regression formula of the correction multipliers in Section 4 are obtained for a specific fractureload configuration. Considering that the main purpose of this correction term is to correct errors caused by the finite elements' inability to accurately represent the near-tip displacement field described by Eqs. (1)-(3), we hypothesize that the same multipliers can be applied to all other crack-load configurations and obtain reasonable SIF results. In this section, we apply the correction multipliers obtained from the special case in Section 4 to a spectrum of fracture configurations to test this hypothesis. Special attention is paid to coarse meshes and effects of interference between neighboring fractures and between fractures and free surfaces. Achieving acceptable accuracy under these conditions is crucial for managing the computational cost of the simulation of dynamic fracture propagation in complex fracture systems. Four test cases for which closed-form solutions of SIF's exist are carefully selected: The first case embodies the interference between fracture tip and free surface boundary; the second deals with heterogeneous stress field; the last two cases represent interactions between neighboring fractures. Both mode-I and mode-II SIF's are considered whenever applicable. Both Method A and Method B are

evaluated for the first case in Section 5.1. Since the mathematical and mechanical principles behind these two methods are similar, only the more general Method B is considered for the other three fracture-load configurations.

5.1. Center-cracked infinite strip with a finite width

Consider a center-cracked strip with an infinite length but finite width 2*b*. The crack is 2*a* long and perpendicular to the longitudinal direction of the strip as shown in Fig. 7a. The strip is subjected to a tensile stress σ in the longitudinal direction and a uniformly distributed shear stress τ along the fracture faces, inducing mode-I and mode-II stress concentration, respectively. The stress intensity factors are

$$K_I = \sigma \sqrt{\pi a} F_I(a/b)$$
 and $K_{II} = \tau \sqrt{\pi a} F_{II}(a/b)$ (22)

where F_I and F_{II} are the fracture-configuration correction factors that can be estimated using the modified Koiter's formula [1]:

$$F_{I}(a/b) = F_{II}(a/b) = [1 - 0.025(a/b)^{2} + 0.06(a/b)] \left(\cos\frac{\pi a}{2b}\right)^{-1/2}$$
(23)

with a relative error of less than 0.1% for any a/b value. In this and other examples, if F_I and F_{II} are close to unity, it means this fracture-load configuration is similar to the reference configuration of a single fracture in an infinite plane.

To apply the GDC method, the strip is discretized into a finite element mesh of a length that is more than 12 times longer than its width, which is found to sufficiently approximate the infinite length according to a sensitivity analysis. Different levels of mesh refinement with b/l_E ranging from 4 to 64 as well as various crack length-to-strip width ratios, i.e., a/b = 0.125, 0.25, 0.50, 0.75, and 0.875 are adopted to investigate the effects of these two factors. Due to the symmetry of the crack and mesh configuration, the tensile stress σ does not contribute to the calculated K_{II} and τ does not contribute to K_I . In all the numerical examples in Section 4, a Poisson's ratio of 0.2 and the crack tip mesh configuration shown in Fig. 5a (eight triangle elements connected to the tip) are used. The effects of the Poisson's ratio and crack tip mesh configuration will be studied in Section 6. To allow precise comparison, the calculation results of the GDC method (both Method A and Method B) with the correction multipliers computed using Eq. (21) applied, as well as the theoretical solution based on Eq. (23) are shown in Tables 1A–2B. Note that the values of F_I and F_{II} , instead of the stress intensity factors K_I and K_{II} are shown. F_I and F_{II} can be considered normalized values of the SIF's. Due to the relationships described in Eq. (22), the relatively errors for K_I and K_{II} are the same as those for F_I and F_{II} , respectively.

The results show that Method B for mode-I fracturing and Method A for both mode-I and -II are fairly accurate for all the scenarios considered, including those with very coarse meshes. The relative errors are generally smaller than 2% with few exceptions. The accuracy of Method-B for mode-II fracturing seems to be dependent on the fracture geometry and mesh resolution. For $b/l_E = 4$ with a/b = 0.5, $b/l_E = 8$ with a/b = 0.75, and $b/l_E = 16$ with a/b = 0.875, erroneous results are obtained.



Fig. 7. Center-cracked infinite strip with a finite width. (a) The crack configuration; (b) the mesh for the case where $b = 8l_E$ and a/b = 0.75 (with opening of the fracture exaggerated). The reference points used by Method A and Method B are indicated in the figure.

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able 1A
Calculated F_I values using the GDC method (Method A) for the center-cracked infinite strip.

a/b	F _I , numeri	Relative e	Relative error (%)								
	$b/l_E = 4$	8	16	32	64	$b/l_E = 4$	8	16	32	64	
0.125	N/A ^a	N/A ^a	1.011	1.004	1.007	N/A ^a	N/A ^a	0.1	-0.6	-0.2	1.009
0.25	N/A ^a	1.038	1.032	1.036	1.040	N/A ^a	-0.1	-0.7	-0.3	0.1	1.039
0.50	1.168	1.171	1.179	1.186	1.189	-1.5	-1.3	-0.6	0.0	0.3	1.186
0.75	N/A ^a	1.595	1.612	1.622	1.628	N/A ^a	-1.8	-0.8	-0.1	0.2	1.624
0.875	N/A ^a	N/A ^a	2.271	2.288	2.300	N/A ^a	N/A ^a	-1.3	-0.5	0.0	2.300

^a N/A, numerical results unavailable due to the incompatibility between the a/b value and the mesh configuration.

Table 1B

Calculated F₁ values using the GDC method (Method B) for the center-cracked infinite strip.

a/b	F _I , numeri	Relative e	$F_{I}(a/b)$ Eq. (23)								
	$b/l_E = 4$	8	16	32	64	$b/l_E = 4$	8	16	32	64	
0.125	N/A	N/A	1.008	1.011	1.009	N/A	N/A	-0.1	0.2	0.0	1.009
0.25	N/A	1.036	1.040	1.038	1.037	N/A	-0.3	0.1	-0.1	-0.1	1.039
0.50	1.196	1.186	1.182	1.183	1.184	0.8	0.0	-0.4	-0.3	-0.2	1.186
0.75	N/A	1.640	1.618	1.617	1.619	N/A	1.0	-0.4	-0.5	-0.3	1.624
0.875	N/A	N/A	2.325	2.295	2.291	N/A	N/A	1.1	-0.2	-0.4	2.300

Table 2A Calculated F_{II} values using the GDC method (Method A) for the center-cracked infinite strip.

a/b	F _{II} , numer	Relative e	Relative error (%)								
	$b/l_E = 4$	8	16	32	64	$b/l_E = 4$	8	16	32	64	
0.125	N/A	N/A	1.013	1.000	1.006	N/A	N/A	0.3	-0.9	-0.3	1.009
0.25	N/A	1.040	1.030	1.038	1.045	N/A	0.1	-0.8	-0.1	0.6	1.039
0.50	1.165	1.172	1.188	1.201	1.208	-1.8	-1.2	0.2	1.2	1.8	1.186
0.75	N/A	1.579	1.621	1.645	1.658	N/A	-2.8	-0.2	1.3	2.1	1.624
0.875	N/A	N/A ^a	2.241	2.294	2.323	N/A	N/A	-2.6	-0.3	1.0	2.300

Table 2B Calculated F_{II} values using the GDC method (Method B) for the center-cracked infinite strip.

a/b	F _{II} , numerical result						Relative error (%)					
	$b/l_E = 4$	8	16	32	64	$b/l_E = 4$	8	16	32	64		
0.125 0.25 0.50 0.75 0.875	N/A N/A 0.014 ^b N/A N/A	N/A 1.027 0.972 - 0.841 ^b N/A ^a	1.021 1.018 1.132 1.124 - 1.710 ^b	0.994 1.031 1.181 1.502 1.432	1.001 1.041 1.200 1.610 2.070	N/A N/A — 98.8 N/A N/A	N/A -1.2 -18.1 - 152 N/A	1.1 -2.0 -4.5 -30.8 - 174	-1.6 -0.8 -0.4 -7.6 -37.8	-0.8 0.2 1.2 -0.9 -10.0	1.009 1.039 1.186 1.624 2.300	

^b Degenerate results; see discussion below. The Bold typeface used in other tables highlights degenerate results owing to similar reasons.

In these three situations, the fracture tip is two elements away (i.e. $(b - a)/l_E = 2$) from the lateral boundary. One of the displacement components used in Eq. (19), $u_{\theta}(2l_E, 0)$ happens to be at the lateral boundary. The mechanical response at this point is substantially affected by the free-surface boundary condition and violate an assumption of the GDC method. This is not an issue for Method A or the calculation of K_I using Method B because none of the displacement components used in Eqs. 15, 16, and (18) is at the boundary. At the same mesh refinement level, if the distance between the crack tip and the lateral free-surface boundary is $4l_E$ instead of $2l_E$, the relative error for K_{II} (Method B) is approximately between 20% and 40%, which though suboptimal for typical mechanical engineering applications is often acceptable for geo-science or geo-engineering scenarios due to the high aleatoric uncertainty in geo-systems. Nevertheless, if the crack tip is $6l_E$ or farther away from the free surface, the error drops below 10% for K_{II} by Method B.

5.2. Three-point bending beam with a notch at mid-span

Consider a beam specimen with a span-to-height ratio of s/b = 4 with a notch of length *a* cut at the mid-span as shown in Fig. 8. The beam is subjected to a mid-span force *P*. Due to the symmetry of the configuration, the mode-II stress intensity factor is zero, and for mode-I



Fig. 8. Three-point bending beam with a mid-span notch.

Table 3Calculated *F_I* values using the GDC method for the three-point bend beam (Method B only).

a/b	F _I , numeri	Relative e	Relative error (%)								
	$b/l_E = 4$	8	16	32	64	$b/l_E = 4$	8	16	32	64	
0.125	N/A	N/A	0.944	0.965	0.972	N/A	N/A	-5.1	-3.0	-2.3	0.995
0.25	N/A	1.013	1.005	1.003	1.001	N/A	0.5	-0.2	-0.4	-0.6	1.007
0.50	1.581	1.468	1.422	1.409	1.406	11.7	3.7	0.4	-0.5	-0.7	1.416
0.75	N/A	3.623	3.439	3.369	3.352	N/A	8.2	2.7	0.6	0.1	3.349
0.875	N/A	N/A	9.469	9.075	8.929	N/A	N/A	7.1	2.6	1.0	8.843

$$K_I = \frac{3Ps}{2b^2} \sqrt{\pi a} F_I(a/b) \tag{24}$$

where $F_l(a/b)$ is the fracture-configuration correction factor, with similar meaning to its counterpart in Eq. (22) but different values. Its value can be calculated using the following dimensionless regression equation proposed by Srawley [26] with a relative error smaller than 0.5%

$$F(a/b) = \frac{1.99 - a/b(1 - a/b)[2.15 - 3.93a/b + 2.7(a/b)^2]}{(1 + 2a/b)(1 - a/b)^{3/2}\sqrt{\pi}}$$
(25)

To test the accuracy of the GDC method on this configuration, we perform FEM analysis with different levels of mesh refinement and different notch lengths. The results of Method-B are summarized in Table 3 in a manner similar to that of Tables 1 and 2. The results are generally accurate. In the worst case scenario, where the height direction of the beam is discretized into four element, the relative error is 11.7%, which remains acceptable for many engineering applications. As the mesh is refined, the numerical results for each geometrical configuration generally converge to the closed-form solution with some minor fluctuation (a few tenths of a percent), which is within the 0.5% error inherent in the closed-form solution. The accuracy is compromised when the notch is short or long compared with the beam height (e.g. a/b = 0.125 or 0.875). In both cases, the points where the displacements are used in the GDC method have similar distances to the notch tip and to the lower or upper free surface of the beam and are not within the near-tip region.

5.3. Two finite-length fractures along a single line

In Sections 5.3 and 5.4, we investigate the accuracy of the GDC method for scenarios with multiple fractures interacting with each other. We first consider the configuration shown in Fig. 9, where two finite-length fractures along a single line existing in an infinite plane. This configuration tends to strengthen the stress intensity at the two tips *A* and *B*, compared with the configurations whether each crack exists alone in an infinite plane. For any tip under a given far-field stress condition (σ and τ), the stress intensity factors (mode-I and mode-II) are dependent on certain geometrical features of the system, and the following closed-form solutions are available [1]

$$K_{I}^{A} = \sigma \sqrt{\pi b} F_{I}^{A}(a/b, c/b)$$

$$K_{II}^{A} = \tau \sqrt{\pi b} F_{II}^{A}(a/b, c/b)$$
(26b)

$$K_I^B = \sigma \sqrt{\pi a} F_I^B(a/b, c/b) \tag{26c}$$

and
$$K_{II}^B = \tau \sqrt{\pi a} F_{II}^B(a/b, c/b)$$
 (26d)

where

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Fig. 9. Two finite-length fractures along a single line in an infinite plane.

 Table 4

 Calculated stress intensity for the two-fracture case at crack tip A (Method B only).

b/c	F_{l} , numerical result F_{l} , relative error (%)		<i>F_{II}</i> , numerical result			F _{II} , relati	ve error (%	F_{I} , F_{II} anly. solu.					
	$b/l_E = 4$	8	16	$b/l_E = 4$	8	16	$b/l_E = 4$	8	16	$b/l_E = 4$	8	16	
2	1.027	1.036	1.041	-1.5	-0.6	-0.2	0.975	1.015	1.035	-6.5	-2.7	-0.7	1.043
4	1.044	1.071	1.090	-5.1	-2.7	-0.9	0.318	1.004	1.068	-71.1	-8.7	-2.9	1.100
8	1.137 ^c	1.113	1.163	-5.6 ^c	-7.7	-3.5	N/A	0.246	1.076	N/A	-79.6	-10.7	1.206
16	N/A	1.249 ^c	1.248	N/A	-9.3 ^c	-9.3	N/A	N/A	0.185	N/A	N/A	-86.5	1.377
32	N/A	N/A	1.445 ^c	N/A	N/A	-11.4 ^c	N/A	N/A	N/A	N/A	N/A	N/A	1.632

$$F_{I}^{A} = F_{II}^{A} = \frac{1}{\sqrt{1 - \alpha_{A}}} \left[1 - \frac{1}{\alpha_{B}} \left\{ 1 - \frac{E(k)}{K(k)} \right\} \right]$$
(27a)
$$F_{I}^{B} = F_{II}^{B} = \frac{1}{\sqrt{2 - \alpha_{A}}} \left[1 - \frac{1}{\alpha_{B}} \left\{ 1 - \frac{E(k)}{K(k)} \right\} \right]$$
(27b)

$$F_{II}^{B} = F_{II}^{B} = \frac{1}{\sqrt{1 - \alpha_{B}}} \left[1 - \frac{1}{\alpha_{A}} \left\{ 1 - \frac{L(\kappa)}{K(k)} \right\} \right]$$
(27b)

with $\alpha_A = a/(a+c)$, $\alpha_B = b/(b+c)$, and $k = \sqrt{\alpha_A \alpha_B}$ and

$$K(k) = \int_0^{\pi/2} (1 - k^2 \sin^2 \phi)^{-1/2} d\phi$$

$$E(k) = \int_0^{\pi/2} (1 - k^2 \sin^2 \phi)^{1/2} d\phi$$
(28a)
(28b)

In the numerical solutions, we fix the length ratio of the two fractures to be a/b = 0.5 and investigate the effects of the mesh refinement levels ($b/l_E = 4$, 8, and 16) and the distance between the two fracture tips (c/b = 1/2, 1/4, 1/8, and 1/16 whenever applicable). The finite element model is more than 50*b* long in each dimension to minimize the boundary effects. The numerical results for the two crack tips *A* and *B* are summarized in Tables 4 and 5, respectively.

The trends observed in this series of results are similar to those from Sections 5.1 and 5.2. Method B of the GDC method is more accurate for mode-I stress intensity than for mode-II. Even under pathological conditions, i.e. mesh coarseness limit reached and strong numerical coupling between the two tips, the error is of the order of 10%. The accuracy for mode-II is non-ideal but still acceptable for many applications. The only exceptions are when the two tips are only two elements away from each other. In this situation, $u_{\theta}(2l_{E}, 0)$ used in Eq. (19) for a tip is the displacement of the other tip, resulting in strong numerical coupling between the two fractures. In these situations, Method A is more appropriate since it uses displacements "behind" fracture tips, where less numerical coupling between the two fractures is expected.

5.4. An infinite array of parallel fractures in an infinite plane

Consider the fracture configuration shown in Fig. 10 where an infinite array of parallel finite-length cracks are periodically arranged on an infinite plane subjected to far-field stress. The interaction between fractures tends to reduce mode-I stress intensity but enhance mode-II stress intensity. The stress intensity factors are $K_I = \sigma \sqrt{\pi a} F_I(a/h)$ and $K_{II} = \tau \sqrt{\pi a} F_{II}(a/h)$ where F_I and F_{II} are the crack configuration correction factors as functions of the crack length and the interval between neighboring cracks. The analytical solutions for F_I and F_{II} are unavailable but well-accepted numerical solutions are presented in [1] and are plotted as continuous curves in Fig. 11. In the FEM solution of this study, we investigate the effects of mesh refinement level ($a/l_E = 16$, 8, 4, and 2) and distance between adjacent fractures (a/h). Due to the periodicity of the

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1	02	
	02	

Table 5
Calculated stress intensity for the two-fracture case at crack tip <i>B</i> (Method B only).

b/c	F _I , nume	rical resul	t	<i>F_I</i> , relativ	e error (%)		F _{II} , nume	rical result		F _{II} , relativ	ve error (S	%)	F_I , F_{II} anly. solu.
	$b/l_E = 4$	8	16	$b/l_E = 4$	8	16	$b/l_E = 4$	8	16	$b/l_E = 4$	8	16	
2	1.078	1.113	1.122	-4.2	-1.1	-0.3	0.978	1.058	1.098	-13.1	-6.0	-2.5	1.126
4	1.117	1.197	1.238	-11.2	-4.8	-1.5	-0.526	1.038	1.177	-142	-17.4	-6.3	1.257
8	1.304 ^c	1.287	1.387	-10.9 ^c	-12.1	-5.2	N/A	-0.370	1.204	N/A	-125	-17.7	1.464
16	N/A	1.533 ^c	1.541	N/A	−12.9 ^c	-12.5	N/A	N/A	-0.270	N/A	N/A	-115	1.761
32	N/A	N/A	1.878 ^c	N/A	N/A	−13.4 ^c	N/A	N/A	N/A	N/A	N/A	N/A	2.169

^c Limit of the mesh coarseness reached where only one element exist between tip *A* and tip *B*. *K*_{II} cannot be calculated at this level of mesh refinement using Method B.



Fig. 10. Parallel finite-length fractures in an infinite plane.



Fig. 11. Comparison of the GDC method results and well-accepted reference numerical solutions [1]. The latter are shown as continuous curves and they have an estimated error of less than 1%. a/(a + h) is used as the horizontal axis to be consistent with the notation in [1]. Note that a/(a + h) = 1/(1 + h/a).

configuration, only one crack and the surrounding medium need to be included in the mesh with appropriate periodic boundary conditions applied. The width of the mesh is more than 50 times the crack length to minimize the effects of the far-field lateral boundaries. As shown in Fig. 11, the results of the GDC methods (Method B only) are fairly accurate for mode-I with relative errors below 10%. The results for mode-II are less accurate and the most significant factor affecting

the accuracy is h/l_E . When $h/l_E = 4$ (i.e. eight elements between adjacent cracks), the relative error can be as high as 30% for large a/h values, but the ascending trend of the $F_{II} - a/(a + h)$ curve can still be reproduced. When $h/l_E = 2$, the relative error becomes unacceptably large and fails to represent the general trend of the $F_{II} - a/(a + h)$ curve. Among all the numerical cases, the shortest distance between neighboring cracks is $4l_E$ (i.e. $h/l_E = 2$). If the neighboring cracks are only $2l_E$ apart, Method B for mode-I will fail because all the displacement components used in Eq. (18) would be zero due to the symmetry of the problem, yielding zero stress intensity. This condition dictates the largest element size that can be used for mode-I.

6. The effects of mesh configurations and the Poisson's ratio

In all the numerical examples in Sections 4 and 5, the Poisson's ratio is assumed to be 0.2. As shown in Eq. (1), the Poisson's ratio is related to the value of β thereby affecting the near-tip displacement field. As mentioned in Section 3, the accuracy of the GDC method (without enhancement through the correction multipliers) depends on the ability of the finite element in representing the near-field displacement field. Therefore, it is expected that the values of C_I and C_{II} are dependent on the Poisson's ratio. We repeat the numerical examples on a single fracture in an infinite plane in Section 4 with Poisson's ratios ranging from 0 to 0.4, and the correction multipliers required for obtaining accurate SIF's for different mesh refinement levels are shown in Fig. 12. A unified regression model is established by assuming the two constants in Eq. (20) to vary linearly with respect to the Poisson's ratio, and the regression results are

$$C_{I}^{B} = \frac{1.226 + 0.206\nu}{\sqrt{1 - (0.349 - 1.125\nu)l_{E}/a}}$$
(29a)

$$C_{II}^{B} = \frac{1.757 - 0.048V}{\sqrt{1 - (0.874 - 0.179V)l_{E}/a}}$$
(29b)

The effects of the Poisson's ratio are more significant for mode-I than for mode-II. Even for mode-I, ignoring these effects by using the correction multipliers for v = 0.2 introduces less than 4% incremental error to the calculated SIF's for arbitrary Poisson's ratio.

The correction multipliers are also dependent on the near-tip mesh configuration. All the previous numerical examples are based on the mesh configuration shown in Fig. 5a where eight triangular elements are connected to the tip node. The other thee configurations in Fig. 5 are also common in FEM analysis. We repeat the numerical analysis in Section 4 with the additional mesh configurations to determine the correction multipliers for different configurations and the results for a Poisson's ratio of 0.2 are shown in Fig. 13. Note that mesh-i, mesh-ii, and mesh-iii use the same space discretization scheme with the only difference among them being in the location of the crack tip and the crack orientation. For a given mesh, the l_E value of mesh-iii is $\sqrt{2}$ times larger than that for mesh-i and mesh-ii. To use mesh configuration iv, l_E in Eq. (18) is replaced with $l'_F \sqrt{3} l_E/2$. This constrains the solution to only use the displacements of points within two element layers of the tip.

The trend of the variation of the correction multipliers with respect to the mesh refinement level is the same for all the mesh configurations. The curves become relatively flat when $a/l_E > 8$. In configurations i and iii, the near-tip region is discretized into eight elements in the angular direction while it is discretized into four elements for mesh-ii. Better refinement in the angular direction improves the displacement field representation, yielding correction multipliers closer to unity. In the region with a radius of $2l_E$ around the tip, more elements are involved in mesh-iii than in mesh-i (the mesh is the same for these two configurations but l_E for mesh-iii is longer), enabling a better displacement field representation. Despite these observations, the effects of the mesh configurations on the correction multipliers are moderate. If we used the correction multipliers for mesh-i on mesh configuration ii, it would induce an error of 4%.



Fig. 12. The effects of the Poisson's on the correction multipliers for (a) mode-I and (b) mode-II at different mesh refinement levels. The effects on C_{II} are small and the regression curves are not plotted. Only the results for Method B are shown.



Fig. 13. The effects of near-tip mesh configurations on the correction multipliers for (a) mode-I and (b) mode-II at different mesh refinement levels.

Additionally, though all the examples in this paper are for plane-stress conditions using Method B, application of the generalized Method B to plane-strain conditions or Method A to plane-strain and plane-stress conditions is straightforward.

7. The effects of mesh perturbation

In the previous numerical examples, the finite meshes are all based on regular grids and the fractures align with the grids. In this section, we investigate the effects of mesh perturbation on the accuracy of the GDC method. Only the mesh pattern shown in Fig. 5a is tested but the qualitative observations should apply to all mesh patterns. A mesh perturbation factor δ is introduced to quantify the degree of perturbation. The *x*- and *y*-coordinates of each end-edge node is moved from its original location in the regular mesh by a distance that follows a uniform distribution between $-\delta l_E$ and δl_E . The mid-edge nodes are moved accordingly. The nodes along external boundaries and existing fractures are not perturbed in order to maintain the geometrical configurations of the system. Three levels of perturbation with $\delta = 0.1, 0.2, and 0.3$ are considered. For a given level of perturbation, different mesh patterns can be obtained by altering the seed value for the random number generator used in the meshing routine. Eight individual and independent random realizations are analyzed as a simple random sample for each perturbation level. Both Method A and Method B are evaluated in this section wherever appropriate. For the perturbed mesh, we still use the characteristic element size parameter l_E of the parent regular mesh are applied.

The two fracture-loading configurations investigated in Sections 5.1 and 5.3, representing fracture-boundary and fracture-fracture interactions, respectively, are assessed. These two configurations are termed the "finite strip" case and "dual fracture" case in the following description. One relatively coarse mesh resolution is used for each case. For the finite strip scenario, $b/l_E = 16$ and $a/l_E = 12$, so the distance between the fracture tip and the lateral boundary is approximately four times the element size. One of the reference points used by Method B for mode-II is at the middle point between the fracture tip and the free surface boundary, not strictly speaking in the near-tip region. The error for this particular case is approximately 31% as shown in Table 2-B, and Method B for mode-II is inappropriate for this particular case. We present the results with mesh perturbation for this case anyway for the sake of completeness, but this limitation should be born in mind. For the dual fracture scenario, $b/l_E = 8$, $a/l_E = 4$, and $c = l_E$, so there is only two elements between the two fracture tips, which made Method B for mode-II inapplicable because one of the reference points for a fracture tip is at the tip of the other tip, as illustrated by the very poor results corresponding to this particular scenario in Table 4 (bold font). Therefore, K_{II} using Method B for the second fracture-loading configuration is not pursued here. The mesh patterns and locations of reference points are illustrated in Fig. 14.



Fig. 14. Perturbed mesh and location of reference points. (a–c) are all based on the finite strip configuration in Section 5.1 and they have perturbation factor δ = 0.1, 0.2, and 0.3, respectively; (d) is based on the dual fracture configuration investigated in Section 5.3 and only the mesh for δ = 0.3 is shown. Note that reference points for Method B on perturbed mesh generally do not coincide with nodes.

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Fig. 15. Results of the GDC method for the finite strip case with perturbed meshes. (a) Results for Method A; (b) results for Method B. Each data point represents one random realization of a mesh perturbation level. The horizontal coordinates of some data points are slightly offset to enable visually separating mode-I and mode-II data points. The margins for 10% relative error are shown in this figure.



Fig. 16. Results of the GDC method for the dual fracture case with perturbed meshes. (a) Results for Method A; (b) results for method B. Results for mode-II with method B are not presented for reasons described above.

The calculation results of the GDC method for individual random realizations of different mesh perturbation levels are shown in Figs. 15 and 16, for the finite strip case and the dual fracture case, respectively. Considering the stochastic nature of mesh perturbation, we also present some statistical analysis results in Tables 6A and 6B. As expected, mesh perturbation affects the GDC calculation results in a random manner, with a greater degree of perturbation causing greater variation of SIF results. Mode-II results seem to be more sensitive to mesh perturbation that mode-I results. The mean of SIF results by Method A appears to be unaffected by mesh perturbation, whereas mesh perturbation slightly increases the mean of mode-I SIF by Method B. Note that the results of mode-II SIF by Method B for both cases tested are inherently inaccurate due to inadequate mesh resolution for these particular cases. The results for the finite strip case are shown here only to illustrate the additional errors induced by mesh perturbation. In most cases, even relatively severe mesh perturbation induces less than 10% additional error, whereas practical needs for mesh perturbation more severe than $\delta = 0.3$ are very rare. More importantly, these numerical examples demonstrate that the GDC method is reasonably robust and its general accuracy does not rely on the symmetry or regularity of meshing.

8. Concluding remarks

Compared with the original displacement-based methods for calculating stress intensity factors, the generalized displacement correlation (GDC) method proposed in this paper has two advantages: (1) It is designed to work with conventional finite element types, and (2) it uses a homogeneous mesh without local refinement around fracture tips. The former feature makes it convenient to implement the new method in existing finite element packages. The latter is important for modeling dynamic fracture propagation problems where the locations of fractures are not known *a priori*. These two features are critical to engineering applications where adopting special element types and local refinement are impractical, such as in the simulation of hydraulic fracturing in complex natural fracture systems.

We propose two suites of formulations, termed Method A and Method B, for the GDC method. The former utilizes displacement information within one layer of elements around the fracture tip, and requires quadratic or higher-order finite elements. The latter can work with any element types, but requires displacements within two layers of elements. To enhance

1	0	6

		Theoretical	GDC results							
			$\delta = 0$	$\delta = 0.1$		$\delta = 0.2$		$\delta = 0.3$		
				Mean	St. dev.	Mean	St. dev.	Mean	St. dev.	
Method A	F _I F _{II}	1.624 1.624	1.612 1.621	1.616 1.622	0.029 0.050	1.611 1.617	0.054 0.094	1.626 1.651	0.111 0.203	
Method B	F_I F_{II}	1.624 1.624	1.618 1.124 ^d	1.638 1.114	0.016 0.035	1.654 1.095	0.032 0.063	1.667 1.040	0.046 0.102	

Table 6AGDC results for the finite strip case with various levels of mesh perturbation.

^d Poor results due to inappropriate mesh resolution, as explained in Section 5.1.

Table 6B

GDC results for the dual fracture case with various levels of mesh perturbation.

		Theoretical	GDC results							
			$\delta = 0$	$\delta = 0.1$		$\delta = 0.2$		$\delta = 0.3$		
				Mean	St. dev.	Mean	St. dev.	Mean	St. dev.	
Method A	F _I F _{II}	1.206 1.206	1.209 1.205	1.215 1.194	0.012 0.036	1.211 1.169	0.017 0.081	1.196 1.134	0.035 0.095	
Method B	F_I F_{II}	1.206 1.206	1.116 N/A	1.133 N/A	0.014 N/A	1.145 N/A	0.024 N/A	1.149 N/A	0.036 N/A	

accuracy of both methods, a correction multiplier is also proposed. Without this correction term, the accuracy of the GDC method is limited due to the inability of regular finite element types to accurately represent the near-tip displacement field. Through a series of numerical examples with a variety of crack configurations, we find that that the new GDC method is acceptably accurate for calculating mode-I stress intensity factors. Even in the limit of mesh coarseness when there is only one element between the two tips of the adjacent fractures, the error is of the order of 10%. The accuracy of Method B for mode-II is less than for mode-I, but acceptable results for most engineering applications, especially for geo-engineering applications, can be obtained even with coarse meshes. Severe errors are inevitable if the points where displacements are used for the calculation are very close to other fracture tips or boundaries of the computation domain. However, this is not unique to the GDC method, and other comparable methods suffer under the same conditions because the near-tip region is inadequately resolved. To correctly model these problems (e.g. tips close to each other or to the boundaries), sufficiently fine meshes must be adopted.

We found that the correction factor is a function of a number of variables for a give near-tip mesh configuration, including fracture length relative to element size, the Poisson's ratio, and random mesh perturbation. However, if we ignore these effects, by using correction factors derived for infinitely long fractures with a nominal Poisson's ratio of 0.2 on a regular mesh, the error is still likely to be within 10%.

Only the correction multipliers for quadratic six-node triangle elements are presented in this paper. Correction multipliers for any combination of element type and mesh configuration can be easily determined through a small number of FEM simulations following the procedure in Section 4. Only one crack-loading configuration needs be considered, and the resultant correction multipliers can be used in arbitrary fracture-load configurations with the same mesh.

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