Strain



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Procedure for accurate calculation of the *J*-integral from Digital Volume Correlation displacement data

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Abstract

The application of Digital Image and Volume Correlation techniques to obtain displacement fields from images has become ubiquitous in experimental fracture mechanics. In this paper, a procedure to extract the J-integral (J) from three-dimensional displacement fields obtained using digital volume correlation is presented. The procedure has been specifically adapted to allow for experimental noise and errors, such as poorly defined crack front displacements, smearing of the displacement field across the crack faces and knowledge of the imprecise crack front location. The implementation is verified using analytical crack tip fields perturbed with synthetic image correlation errors to characterise the response of J. The method is then applied to experimental results using a Magnesium alloy WE43 loaded elastically in mixed-mode. The steps outlined are intended as a guideline for application of the volume integral from displacement fields to allow for accurate calculation of J along a crack front embedded within the volume.

Keywords

digital volume correlation, fracture mechanics, J-integral, X-ray computed tomography

Symbols	
A_c, \hat{A}_c	Crack extension in the global and reference state respectively
σ_{ii}	Cauchy stress tensor
\mathcal{E}_{ij}	Infinitesimal strain tensor
	Displacement vector in the global, interpolated and reference state respectively
0 [] u[] u[W	Strain energy density
V, \overline{V}	Integration volume in the global and reference state respectively
q	Crack extension function
v	Poisson's ratio
E	Young's modulus
J	J-integral
S	Crack front
n	Crack front segment
ΔS	Integration volume breadth (x_3)
x_i , \hat{x}_i	Co-ordinate system in the global and reference state respectively
Δx	Integration volume width (x_1) or height (x_2)
<i>O</i> _{<i>i</i>}	Crack tip location
L_i	Crack front direction
M _i	Crack normal direction
N _i	Crack propagation direction
P_{i}^{1}, P_{i}^{2}	Start and end points of crack path.
R	Rotation matrix
\hat{r}_m	mask height in reference state
t	displacement vector
Τ	Transformation matrix
	mask height in reference state displacement vector Transformation matrix

1 Introduction

The application of Digital Image Correlation (DIC) and Digital Volume Correlation (DVC) techniques to obtain displacement data fields from images has become ubiquitous in experimental fracture mechanics. Many examples now exist that demonstrate the successful extraction of fracture parameters, such as stress-intensity factors obtained by McNeill *et al.* [1], Lim *et al.* [2] and Roux *et al.* [3], mixed-mode problems addressed by Rethore and Gravouil [4], Yoneyama *et al.* [5,6] and Sistaninia *et al.* [7] and energy approaches such as the J-integral by Yoneyama et al. [8] and the authors of this work [9,10], to name but a few.

A significant limitation of image correlation approaches (compared to others such as Moire interferometry) is the reduced accuracy experienced near the strain concentration at the crack tip. State-of-the art subset based correlation techniques suffer from smearing of information across the crack faces resulting in spurious displacement and strain readings. It is therefore imperative for DIC and DVC based fracture property extraction methods to account for these sources of error by: i) using significantly more complex (and generally in-house developed) correlation techniques (e.g. subset splitting [11]); ii) replacing poorly defined or missing data in these regions (e.g. with numerical approximations of the material after the crack front has been found [12,13]); or iii) developing methods that allow for the selection of data remote from the crack tip thereby eliminating the need for accurate displacement data near the crack faces and crack front [10].

Similarly, the knowledge of the crack tip location may be critical when extracting fracture parameters from experimentally obtained displacement fields. Often, it is not possible to define an exact crack tip or front location and complex segmentation and optimisation techniques have been developed that aim to best identify the crack tip location [14,15].

Random permutations in the signal (light or x-ray) also result in noise in the individual vectors of the displacement fields. DIC noise is typically reported at 0.01 and 0.02 pixel for in-plane and out-of-plane respectively [16], while DVC tends to achieve much lower accuracies, which are typically reported as 0.05 to 0.1 voxels [17]. This is mainly due to the reliance of DVC on there being a suitable distribution (shape, size, brightness) of natural features within the material volume [18], whereas DIC can utilise artificially applied and optimal speckle patterns with high contrast.

Some of these problems can be overcome when using field fitting approaches, whereby optimal analytical solutions to crack tip displacement fields are fitted to the measured displacement data [5]. However, field fitting approaches may be ill-suited to cases where the crack tip region is poorly defined, as in large-scale plasticity, crack branching, or micro-cracking (e.g. graphite or concrete) [19]. Furthermore, displacement fields in regions near the crack tip are a challenge to measure [20]. This is because both DIC and DVC usually assume a small set of deformation parameters (i.e. translation, dilatation and shear strains and in some cases rotation) for efficiency purposes. As a result, sharp discontinuities and crack tip singularities tend to be inconsistent with the assumed deformations within a subset (region of analysis), causing poor correlation in crack tip regions [11,21]. As such, the J-integral provides better suited tool for extracting, from experimentally determined displacement fields, the fracture parameters that describe the action on the crack tip. In the DIC form, as previously presented by the authors [9,10], high strain regions near the crack can be omitted under the assumption of a plane stress state, provided that the crack path does not leave the J-integral contour. In the volumetric form of the J-integral [22], the displacement fields around the crack front are required since the

integral encapsulates a predefined region of interest that necessitates data along the crack front [22].

This presents a problem when using DVC computed displacement fields to obtain J-integral based fracture parameters. The poorly defined displacement fields, in combination with the need of accurate displacement data around the crack front, creates a need for an unique implementation of the J-integral. One method to overcome this is the replace the missing, or censored, data using the Finite Element (FE) framework, as presented by Bahrli *et al.* [12]. However, this requires extrapolation from the measured data in order to define a displacement field in the cracked region, which in turn requires assumptions on the material properties (plasticity, microcracking) in this region, and the method cannot consider the effects of complex fracture mechanics such as crack branching or bridging. This extrapolation also requires knowledge of the precise crack front location. Another method, which the authors are aware of, is to adapt the volumetric formulation to allow for poorly defined crack front displacement data. However, such work has not been reported in the open literature.

We present an alternative procedure for computing the J-integral directly from DVC displacement data. The work is therefore similar to the analysis performed on surface J-integrals that has previously been presented by the authors [9,10]. This paper presents a procedure and formulation of the J-integral that is better suited for DVC displacement data. Typical sources of errors are studied using synthetically generated displacement fields and errors. Subsequently, the procedure is applied to an experimental case study. The experimental results are obtained at sub-critical loads, with elastic deformation, and validation is performed against simulations using the finite element software Abaqus that uses the volume integral as detailed in [22,23].

2 Three-dimensional framework

In this work, the premise for the three-dimensional (volumetric) framework of the J-integral (J) is that global DVC computed displacement data is available in the vicinity of the crack front in a Cartesian co-ordinate system X_i given by U_i , and that the mechanical (deformation) properties for the material are known (i.e. the relationship between the strain and stress tensors). To subsequently compute J values along the crack front, the volumetric form presented by Shih *et al.* [22] is adopted. It is noted that the volumetric J formulation differs from the classical two-dimensional formulation in that the contour is not represented by a line or area, but as a surface encapsulating a specified region, i.e. volume, V.

J is calculated for *V*, which is centred at O_i along the crack front S_i as show in **Error! Reference source not found.** A convention is used where the global coordinates system X_i is designated by an uppercase and a local Cartesian co-ordinate system x_i is designated by a lowercase at the crack front. The local co-ordinate system is centred at O_i with x_1 in the crack plane and perpendicular to the crack front, x_2 normal to the crack plane and x_3 tangential to the crack front.

Similar to the adaptation of the line integral to an equivalent domain integral in the twodimensional formulation, the volumetric formulation takes on an equivalent volumetric domain integral given by Eq. (1). The volumetric integral allows for more data points to be considered, compared to alternative formulations, and is therefore better suited for full-field data [9].

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$$J = \frac{1}{A_c} \int_{V} \left(\sigma_{ij} \frac{\partial u_i}{\partial x_1} - w \delta_{1j} \right) \frac{\partial q}{\partial x_j} dV$$
(1)

Eq. (1) contains a displacement derivative, $\partial u_i/\partial x_1$, in the x_1 direction, a Cauchy stress tensor, σ_{ii} , a strain energy density, w and $\partial q/\partial x_i$, which is the derivative of a virtual crack extension function, q, defined by Eq. (5). A_c is obtained using Eq. (6) and is directly dependent on the choice in q. δ_{1i} is the Kronecker delta.

 σ_{ij} is calculated using Eqs. 2 and 3. In this work, Eq. (2) is given under the assumptions of a linear-elastic isotropic material, however, more complex mechanical behaviour models may be used such as presented by [12].

$$\sigma_{ij} = \frac{E}{1+v} \Big[\varepsilon_{ij} + \frac{v}{1-2v} \varepsilon_{kk} \delta_{ij} \Big]$$
⁽²⁾

E and v are the Young's modulus and Poisson's ratio respectively. ε_{ii} is the symmetric strain tensor calculated using the displacement gradients, $\partial u_i / \partial x_i$, in Eq. (3).

$$\varepsilon_{ij} = \frac{1}{2} \left(\frac{\partial u_i}{\partial x_j} + \frac{\partial u_j}{\partial x_i} \right)$$
(3)

The strain energy density, w, is calculated using Eq. (4).

$$w = \int \sigma_{ij} d\varepsilon_{ij} \tag{4}$$

The crack extension function, q, requires a function that is sufficiently smooth and that is constrained at its boundaries by Eq. (5)

$$q = \begin{cases} 1 & \text{on } O_i \\ 0 & \text{on the outer edge of } V \end{cases}$$
(5)

 A_c is obtained by integrating q along crack front segment ΔS as defined by Shih et al. [22] using Eq. (6).

$$A_c = \int_{\Delta s} q \, dS \tag{6}$$

Eqs. 2 and 3 assume material specific mechanical behaviour properties within V and so I is computed on the basis that the mechanical (deformation) behaviour is known. In the presented form, J is a function of U_i , E and v. No load or crack length measurements are required to evaluate J.

The distinct feature of the *I*-integral is its contour independence. In the two-dimensional formulation, with an absence of out-of-plane (x_3) stresses under a plane stress assumption (or strains, under a plane strain assumption), contours can be arbitrarily chosen so that the crack tip region is excluded. This significantly reduces error in the computation of J and is especially Strain

important when considering experimental data that have a significantly higher error than, for example, finite element computed data [9]. In the three-dimensional formulation, the contour will inherently include crack tip data as out-of-place stresses are present [22]. Therefore, the contours cannot be taken in such a way that the crack tip region is excluded. However, as shown in Section 3, careful consideration of q can allow for a computation of J from experimental data with significantly reduced error without affecting contour independence.

Numerical implementation

Consider the crack front depicted in **Error! Reference source not found.**a. In the presented approach, three aspects are required for the computation of J from the respective DVC displacement data, namely crack-tip segmentation, integration volume mapping, masking and summation of the integrand. The procedure is outlined below and the implementation is done using Matlab ver. 9.6 software package, where the DVC computed displacement field data with the respective position data are imported into Matlab's workspace.

3.1 Crack tip segmentation

The first step requires V to be aligned with the crack front and therefore necessitates knowledge of the crack tip location and orientation. Since J is computed at specific locations along the crack front, the number of J values obtained is dependent on the choice in n number of segments. Using the procedure outlined by [24], S_i is divided in n segments, between P_i^1 and P_i^2 with each segment located at O_i^n with known orientation L^n (crack path), M^n (crack direction) and N^n (crack normal) as depicted in **Error! Reference source not found.**a. Each segment has a width of ΔS . In this work, the crack tip location and direction is visually, however, other methods are available [12,15,25].

3.2 Integration Volume mapping

The second step involves mapping V to the respective crack tip location and orientation. For this, additional steps are required that are outlined below, which involve: i) setting up a reference volume \hat{V} ; ii) mapping of the reference volume in the global domain; iii) extracting nodal displacements by means of interpolation; and iv) mapping these back to the reference state.

i) \hat{V} is of unit size, has an orientation that is aligned with the horizontal, vertical and normal coordinate system for \hat{x}_1 , \hat{x}_2 and \hat{x}_3 respectively, the crack plane is aligned with \hat{x}_1 , the crack tip is located the origin, and the crack is assumed to propagating in the \hat{x}_1 direction, i.e. the crack normal aligned with \hat{x}_2 as shown in in **Error! Reference source not found.**a. A convention is adopted, whereby the reference volume's symbols use a top hat.

The reference volume has regularly spaced position data in \hat{x}_i to allow for simple differentiation and integration, as shown in **Error! Reference source not found.**b (cross-section of the mid-plane). No advantage is seen in numerical differentiation and integration approaches that require point wise evaluation (i.e. such as in FE implementation). This requires the knowledge of shape function derivatives and their respective Jacobian at each data point. The relatively noisy nature of DVC coupled with the inherent subset smearing does not, in the authors' opinion, warrant such implementation complexities. It is shown

later, that the choice in q has the greatest influence when dealing with smeared and noisy crack tip data and crack tip position accuracy.

The crack extension function, q, is defined as per Eq. (5) and is a function of \hat{x}_i . J is written with respect to the local co-ordinate system of \hat{V} , with q defined as vector components in which J is the projection onto the q vector field ($J = J_i q_i$ (i = 1,2,3)). The form given in Eq. (5) is equivalent to computing J in vector form if q_i is pointing in the crack growth direction (i.e. $q_2 = q_3 = 0$). The reader is referred to Kuna [26] for more details on the vector form of the volume integral.

Three functions are considered for q. It is noted that for the evaluation of J, Eq. (1) requires the derivates of q with respect to \hat{x}_i . Thus, the choice in q is such to either allow for simple numerical implementation (i.e. constant gradients) or to inherently remove poorly defined displacement data near the crack tip (i.e. by setting a near zero gradient at the crack tip). A representation of the q function derivatives, $\partial q/\partial \hat{x}_j$, is shown in **Error! Reference source not found.**

The choice of a 'linear' function results in constant gradient with a minimal gradient value. A polynomial (in this work a 4th order polynomial was chosen) results in steeper gradient at the outer edges in \hat{V} (away from the crack tip) and near zero gradient at the crack tip region. A piecewise linear function (termed 'contour') results in a constant gradient near the outer edges in \hat{v} and zero gradients at the crack tip. As mentioned above, the inherent nature of J in the volume, which encapsulates a segment along a crack front, is that crack tip data will always be included. As such the contour function does not consider a piecewise linear function in \hat{x}_3 . This ensures that the gradients in \hat{x}_3 are at a minimum near the crack tip. \hat{A}_c is based on the choice in q, and is calculated using Eq. (6).

A mask is defined over \hat{V} that is centred along the crack (Fig. 2b). The mask defines those displacement data points that are judged to have been significantly smeared near the crack and therefore are not considered to be accurate enough for the evaluation of J. In the present analysis of a planar crack, the mask is cuboid in shape, has unit thickness, a length of 0.5 and defined by height \hat{r}_m , as shown in **Error! Reference source not found.**b. Data points included in the mask are replaced by means of extrapolation (detailed in Section 3.3).

ii) Mapping of U_i to \hat{V} is achieved by mapping the position co-ordinates of the reference volume domain \hat{x}_i to the global domain x_i . This operation requires the knowledge of a transformation matrix T at each integral segment computation n. T^n comprises of a scaling matrix α (taken as constant across all segments), a rotation matrix R^n and a translation vector t^n , which are assembled as given in Eq. (7).

$$\boldsymbol{T}^{n} = \begin{bmatrix} \boldsymbol{\alpha}^{n} \boldsymbol{R}^{n} & 0\\ \boldsymbol{t}^{n} & 1 \end{bmatrix}$$
(7)

 α is a diagonal matrix and determines the size of V. In this work $\alpha_{11} = \alpha_{22} = c$ so that the volume length and height are defined by c and are equal in x_1 and x_2 . $\alpha_{33} = \Delta S$ represents a volume width that is equal to the segment length. \mathbf{R}^n is dependent on the crack tip orientation and is obtained using Eq. (8). \mathbf{t}^n is dependent on the crack tip position and is obtained using Eq. (9).

$$\boldsymbol{R}^{n} = \left[\boldsymbol{M}^{n}, \boldsymbol{N}^{n}, \boldsymbol{L}^{n}\right]$$

$$\tag{8}$$

$$\boldsymbol{t}^n = \boldsymbol{O}^n \tag{9}$$

The mapping of the \hat{V} to V is undertaken by using Eq. (10).

$$\begin{bmatrix} \boldsymbol{x} \\ 1 \end{bmatrix} = \boldsymbol{T}^n \begin{bmatrix} \hat{\boldsymbol{x}} \\ 1 \end{bmatrix}$$
(10)

- iii)Interpolation of the displacement field is undertaken using linear interpolation to obtain displacement values at x. The interpolation method uses a Delaunay triangulation method as outlined by [27]. Similar to the argument presented in i) it is in the authors' opinion that higher order interpolation schemes do not improve the accuracy in the computation of J. In fact, it is argued that due to the noise nature of DVC data, higher interpolation schemes may result in an increase of error as a result of Runge's phenomenon [28].
- iv)Subsequently, the interpolated displacement field is mapped back to the reference state using Eq. (11).

$$\hat{u}_i = \left[\boldsymbol{R}^n \right]^T u_i \tag{11}$$

where $[\mathbf{R}^n]^T$ is the transpose of \mathbf{R}^n and u_i is the interpolated displacement data. A_c is scaled by ΔS to compute its equivalent value in the reference state, i.e. $\hat{A}_c = A_c \cdot \Delta S$. dV is scaled by $2\Delta S\alpha$ to compute its equivalent value in the reference state, i.e. $d\hat{V} = c^2 \Delta S$.

3.3 Masked data

Before the computation of *J*, it is important to address the masked data close to the crack faces, particularly if a mixed-mode type loading case is expected. Similar to the method presented by the author's for computing mixed-mode J-integral solutions from surface displacement data [9], masked field data is replaced using an iterative least-squares polynomial fitting method developed by Garcia *et al.* [29] and Wang *et al.* [30]. The number of iterations used is 100. It is noted that this approach is limited to low displacement gradient regions and smaller masked areas. This is due to the inherent inability of polynomial functions to represent crack tip fields that represent discontinuous and singular values.

3.4 Summation of the integrand

The numerical integration of \hat{V} requires the computation of its displacement gradients $d\hat{u}_i/d\hat{x}_j$, which in turn are used to compute the strain tensor using Eq. (4), the stress tensor using Eq. (3) and the strain energy density using Eq. (5) at each data point in \hat{x}_i . The choice of q determines $dq/d\hat{x}_i$ and \hat{A}_c as illustrated in Fig. 3. The computation of J then becomes a simple summation of the terms listed in Eq. (1).

4 Errors of the volume J-integral

There several sources of error to consider when computing *J* from DVC displacement data. Firstly, as with any experimental measurement technique, the DVC computed displacement field contains random noise. Furthermore, regions near the crack faces and tip are prone to errors. These can arise from the effects of image artefacts introduced in the reconstruction process [31], or may be due to DVC smearing at the crack faces [32]. Secondly, knowledge of the crack alignment and location is required in order to position the integration volume, and while sophisticated crack segmentation algorithms exist to aid segmentation, these always include some measure of uncertainty [32]. Thirdly, a reasonably accurate knowledge of material properties is required in order to compute stress values from the strain field [10].

A pragmatic approach was adopted to quantify the influence of the above errors in the computation of J. Synthetic displacement data were generated using the finite element (FE) method, whereby a known error was imposed, and its effect in the computation of J was therefore isolated. Furthermore, J's contour independence served as an indicator to assess valid computations. Contour independence was inferred when fluctuations of less than 5% over five contours were achieved. Three contour type functions were considered, as shown in Fig. 3; the contour (piecewise linear) function used a contour width of three.

The error analysis utilised a parallel approach of an Abaqus and an in-house developed linearelastic FE model. A unit size cuboid geometry was generated, with the crack tip orientated at the origin (Fig. 4). The Abaqus model (ver. 6.13) comprised 65800 quadratic 20-node brick elements (C3D20) with collapsed quarter points along the crack front. J-integral values were computed over ten contours using Abaqus' built-in functionality. The in-house developed FE model comprised of 127500 linear eight-node brick elements. The grid of element nodes were arranged in such a way that the computation of J does not require any interpolation (Section 3.2.iii) and thereby errors due to interpolation are excluded (no crack tip mesh refinement was undertaken). J values were computed in this model, using the above-presented technique, over ten contours. Both models used Young's modulus of 10 GPa and a Poisson's ratio of 0.3. Displacement boundary conditions were applied at the top and bottom surface nodes of the unit size cuboid, such that the sample is loaded either in a crack opening mode (i.e. in the U_2 direction, designated as BC_1), in-plane shear (i.e. in the U_1 direction, designated as BC_2), outof-plane shear (i.e. in the U_3 direction, designated as BC_3) or a combination of the three (designated as $BC_{1,2,3}$). Fig. 4a illustrates the application of BC_1 .

The error analysis presented in Fig. 5 and 6 considered a Monte Carlo analysis of 1000 runs applied on the in-house FE generated displacement data, providing the influence on the computation of *J* of normally distributed random displacement noise and rigid body (crack tip) misalignment. The respective displacement boundary conditions are applied in such a way that an average *J* along the crack path of unity is obtained. For BC_1 , BC_2 , BC_3 and $BC_{1,2,3}$ this equated to displacement magnitudes of $7.45 \cdot 10^{-2}$, $2.96 \cdot 10^{-1}$, $2.98 \cdot 10^{-1}$ and $5 \cdot 10^{-2}$ units respectively. The error in *J* was subsequently calculated on the basis of the difference between the unit value (i.e. no error) and error analysis for each crack segment along the crack path, expressed as a percentage.

It is worth noting, that the above presented error analysis did not consider the influence of reconstruction and crack smearing errors on the computation of J. Initial analysis showed that

these sources of error do not result in contour independence and therefore do not obtain valid J computations. Errors in the knowledge of material properties are simply a scaling error when a linear-elastic material is assumed, and so have not been considered.

Fig. 5 shows the mean, upper and lower bounds of displacement noise versus the error in the computation of *J*. Considering a typical limit of DVC measurement accuracy of approximately 0.1 voxels (LaVision product manual, DaVis 8.4) and the unit size sample geometry, an assumed (typical for experimental data) image volume of 1000x1000x1000 voxels would equate to a displacement accuracy of $1 \cdot 10^{-4}$ units. For BC_1 , BC_2 and BC_3 this equate to a 1.34%, 0.0338% and 0.0336% displacement noise respectively. Similarly, Fig. 6 shows the mean, upper and lower bounds of crack tip misalignment (i.e. rigid body movement of *V* in X_1 and X_2 directions) to establish the sensitivity of *J* with respect to crack tip alignment. The percentage misalignment reported is with respect to the integration volume size, which at its limit (10 contours) represents 200 voxels or 0.2 units (i.e. a 1% crack tip location error is equivalent to 2 voxels or $2 \cdot 10^{-3}$ units). Since the fracture mechanics framework uniquely represents the displacement field around any crack tip, an equivalent error estimate in *J* may be calculated by scaling the calculated *J* by the error estimates shown in Fig. 5 and 6 with the *J* value obtained in other experiments (for an approximately similar displacement accuracy).

Fig. 5 and 6 illustrate two features: Firstly, a shearing type loading results in a significantly higher error. This is consistent with findings reported for J values computed from DIC (surface) data [9] and is attributed to the higher strain energy density values along the crack faces. It highlights the importance of adequate masking and data replacement when mixed-mode loading is expected. Further testing, not shown, included a mask of varying width \hat{r}_m . The results show that the data replaced using the iterative least-squares polynomial fitting (Section 3.3) was robust in maintaining contour independence and accurate J computations. Secondly, the type of q function (Fig. 3) has a notable effect on the J error. A compromise exists - a linear q function shows less sensitivity to random displacement noise, whereas a contour type function shows an order of magnitude reduction in error for crack tip location. The polynomial function does not seem to offer any significant improvement for either of these considerations. This compromise exists due to the gradients of the q function. As mentioned above, a linear q results in the lowest possible gradients and thereby, the displacement noise error is minimised. However, data along the crack face and tip are considered, thereby making a linear contour function significantly more susceptible to alignment errors; which is less of the case for the contour function as only gradients in the x_3 are considered.

Considering the noisy nature of DVC displacement data, reconstruction errors, crack smearing and crack segmentation difficulties, it is recommended that a contour function is chosen with a low gradient, i.e. with a large contour width. A 1 to 2 % displacement error and a 1 to 2 % crack tip location uncertainty equate to uncertainty in the computation of *J* between 0.8 to 3.2 % for a crack opening mode. This error significantly increases for shear type loading to approximately 5 to 18.5 %.

5 Verification on experimentally obtained DVC displacement fields

The previous section highlighted two prominent sources of error that are related explicitly to DVC computed displacement data. It was shown that in the volumetric form, J is sensitive to mixed-mode data, for which displacement noise and crack alignment can play an important

role. Careful consideration in q allows for less sensitivity concerning crack tip alignment in the crack propagation (x_1) and crack plane (x_2) directions. To test the robustness of the presented method, a material with natural contrast pattern was studied under mixed-mode loading. Magnesium alloy (WE43), which has high commercial value due to its good castability, high strength, high creep resistance, high corrosion resistance and high flame (ignition) resistance, was chosen. WE43 is widely used for structural components in aircraft for weight savings, and orthopaedic implants due to its biocompatibility [33].

5.1 Materials

Magnesium alloys, such as WE43 that contains 4wt% Yt and 3wt% Nd, produce good speckle contrast in X-CT reconstructions due to the attenuation variations between the Magnesium bulk material and precipitates that are richer in rare earth elements [34,35] and is therefore well suited for DVC [36]. Precipitates with strong X-ray attenuation are quite visible in the reconstructed image (shown by a central slice in Fig. 7b). Brightening effects close to the edges of the sample and the notch should not be mistaken for variations in precipitate density. Rather, this is due to beam-hardening, which is a common (and persistent in this case) X-CT image artefact. The beam hardening is a constant artefact in this dataset, and so did had little effect the DVC analysis [37].

The specimen geometry is shown in Fig. 7. The specimen was rectangular over the gauge region of 50 mm in length (X_3) . The gauge cross-section had a width (X_1) and thickness (X_2) of 10 x 20 mm respectively. The notch was machined using wire-Electrode Discharge Machining (wire-EDM with a 0.3 mm diameter wire) to minimize residual stresses and work hardening that would result from mechanical notch cutting methods. The notch was machined at an angle of 30° to each of the global axes X_i . The loading of the sample was such that the bottom end was loaded in shear in the X_1 direction under displacement control using a translation stage with an accuracy of 0.01 mm.

5.2 Tomography

Tomographs were obtained using Stellenbosch University's X-CT scanner facility using a General Electric Phoenix VTomeX L240 microCT scanner with a 2048×2048 pixel detector with a 16-bit depth [38]. Optimized parameters were selected according to the guidelines set out in [39]. X-ray settings included 220 kV and 200 μ A and copper beam filtration of 1.5 mm, with a voxel size set to 15 μ m. Image acquisition time was 500 ms per image and images were recorded in 2000 rotation steps during a full 360-degree rotation of the sample. At each step, the first image was discarded, and the subsequent three images averaged to provide high image quality. Detector shift was activated to minimize ring artefacts, and automatic scan optimizer was activated to eliminate artefacts due to possible sample movement or X-ray spot drift. Reconstruction was performed in system-supplied Datos reconstruction software.

Three X-CT images were taken to obtain two DVC datasets: one loaded reference image, one unloaded image with an applied displacement of 1 mm in X_1 at the bottom grip region while the top grip remained fixed, and one unloaded image with a rigid body displacement of 1 mm. All displacements were assumed to be elastic as correlation was taken with reference to the loaded image.

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A centralized region consisting of 1260 x 1860 x 2024 voxels (X_1 , X_2 and X_3) was extracted and reduced to 8 bits for improved computational efficiency when computing DVC displacements (Fig. 7a). The crack geometry was inferred from the known notch geometry and the tomography images. To improve its estimate, a best-fitting linear plane was fitted by using a least-squares approach, whereby the crack front shape was assumed linear between P_i^1 and P_i^2 with constant N_i and M_i vectors.

5.3 DVC

The DVC analysis was performed with DaVis (ver. 8.3) volume correlation software using a fast Fourier transform (FFT). Image correlation used an initial subset size of 200 and a final subset size of 80 voxels and a step size of 20 voxels (75% overlap). The rigid body displacement dataset revealed an experimental accuracy of 1.4 μ m (i.e. ~ 0.1 voxels), which is comparable to precision reported in the literature on a similar subset size [21,40]. Background noise was calculated as 0.36 μ m by correlating the two identically loaded datasets. The U_1 displacement field over the ROI is shown in Fig. 8a at $X_1 = 0$ and $X_2 = 0$.

The DVC datasets resulted in 44 x 74 x 60 displacement data points (X_1 , X_2 and X_3). The computed DVC displacement data and crack geometry were imported into the Matlab environment, following rigid body rotation removal using the Euler angle extraction methodology of Mostafavi *et al.* [36].

5.4 Finite element simulation

Three-dimensional finite element (FE) simulation of the experimental setup was carried out using Abaqus ver. 6.13. The elastic properties of Mg WE43 were taken as provided by the manufacturer with Young's modulus of 44 GPa and a Poisson's ratio of 0.35. The linear-elastic model comprised of 105200 quadratic 20-node brick elements (C3D20) as shown in Fig. 9. The model used quadratic elements with quarter-point elements along the crack front, an initial radial element width of 0.3 mm at the notch mouth, converging to 0.01 mm (after 10 rings of elements) at the crack front, as shown in the magnified window in Fig. 9. Displacement boundary conditions were applied using the displacements applied at the BC. This way, the specimen loading could be simulated as close as possible to the experimental setup. Subsequently, *J* values were computed using Abaqus' built-in functionality over ten contours, termed as J_{Abaqus} . In parallel, the FE computed nodal displacements and the crack front geometry were imported into the Matlab environment for processing using the procedure presented in Section 3.

5.5 J-integral computation

The crack segment between P_i^1 and P_i^2 and integration volume V are shown in Fig. 8b. The integration volume thickness was chosen as $\Delta S = 0.6$ mm (i.e. approximately three DVC data points in x_3), and width and height was $\Delta x = 4.8$ mm thus considering nine J computation points along the crack path (i.e. in x_1 and x_2). It is estimated that the crack tip alignment concerning the integration volume was < 1% (to 5 voxels or 75 µm), as the notch tip of radius 150 µm was easily visible in the Tomographs. A crack mask $\hat{r}_m = 1.2$ mm (i.e. four subsets) was used. J values were computed using the contour function.

Fig.10 shows the DVC computed displacement data over the integration volume V at $\hat{x}_3 = 0$. The data shows the mixed-mode nature of the crack loading, whereby significant mode III type loading is present. This is to be expected as the sample loaded in shear perpendicular to the notch tip.

Both DVC and FE displacement data were used to compute J along the defined crack (termed J_{DVC} and J_{FE} respectively) using the procedure as outlined in Section 3. Both DVC and FE displacement data was chosen to compare best: i) the course mesh of the presented approach with an ideally meshed notch tip (FE) and; ii) estimate the influence of experimental error (displacement noise and crack tip position).

5.6 Results and discussion

The mean normalised J computations are shown in Fig. 11a for the equivalent DVC, FE, and Abaqus computations. Fig. 11b illustrates the contour independence for the first four computations along the crack path for the DVC displacement data. The difference when computing J from FE and DVC data is 6.2 % indicating the combined error of displacement noise and crack tip position uncertainty. This is in agreement with the estimated error, presented in Section 4, of 7 %. The experimental error in the computing in a maximum deviation of 5.2 %, indicating the strong influence of DVC displacement noise when computing J from DVC displacement data.

Considering the differences between the FE and Abaqus J computations a smoothing effect near the edges of the specimen is highlighted. The relatively large integration volume of 4.8 mm, when considering that the Abaqus contours are taken 0.1 mm from the crack tip, results in less sensitivity in capturing variations in J along the crack front. This is attributed to the larger integration volume being less sensitive to localised strain gradients, which are more apparent near the edge of the specimen where the stress-strain conditions change from plane strain to plane stress.

The agreement between the measured and predicted J values demonstrates the effectiveness of using DVC data to obtain fracture parameters of the loaded sample. This study shows that high-resolution X-CT can be used to reliably quantify the energy release rate, measured employing the J-integral. The study furthermore highlights that J is highly effective for the computations of fracture properties in the volume when using DVC data. Whiel sensitive to displacement noise, J does not require the exact knowledge of the crack tip position. When a mode I type loading is expected, an error less than 1 % can be expected in J. This is especially advantageous for materials where the location of the crack tip is difficult to identify, such as nuclear graphite [19].

6 Conclusions

A simple procedure for evaluating the J-integral in its volumetric form from DVC computed displacement data has been presented and verified against experimental data for a mixed-mode loaded crack. DVC computed displacement data present unique challenges due to noise, erroneous data near crack faces and imprecise knowledge of the crack tip position.

- The procedure for the robust extraction of *J*-integral values uses a specifically defined *q* function, along a three-dimension crack front.
- Common errors associated with DVC computed displacement data were quantified using analytical displacement data to assign confidence intervals. In the presented form, the J-integral shows sensitivity to displacement noise; however, it does not require the exact knowledge of the crack tip.
- Analysis of the J-integral obtained from an experimental study of a mixed-mode problem, using a Magnesium WE43 alloy with an inclined single edge rounded notch to represent a crack, found a good agreement with the analytical solution from an FE model. *J*-integral values along the crack front could be determined to within 6.2 % of the finite element solution.

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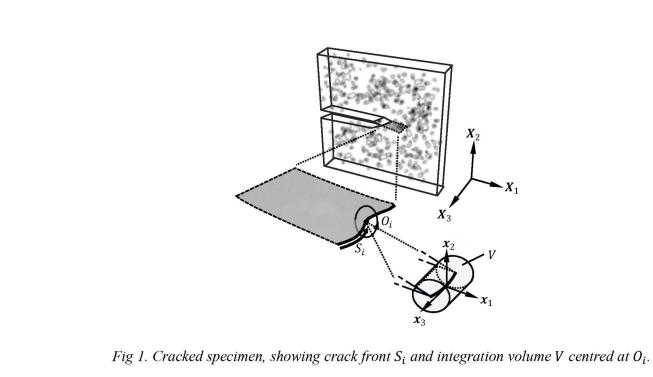
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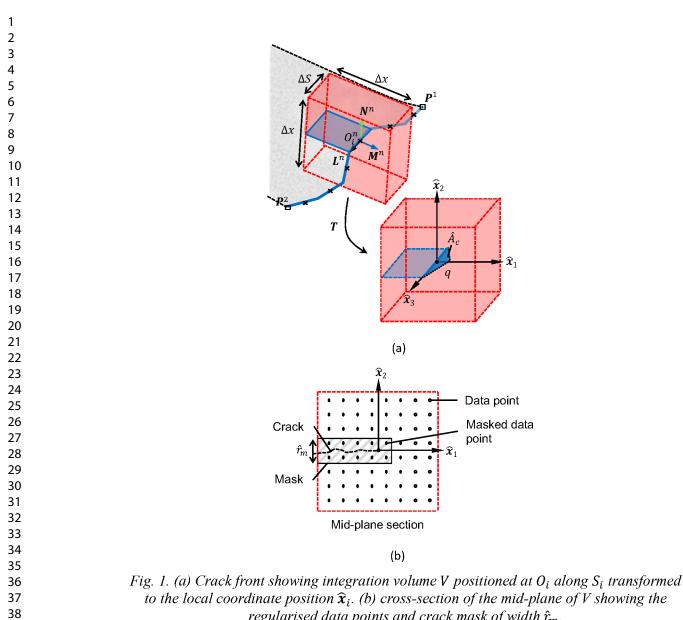
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Strain





regularised data points and crack mask of width \hat{r}_m .

Strain

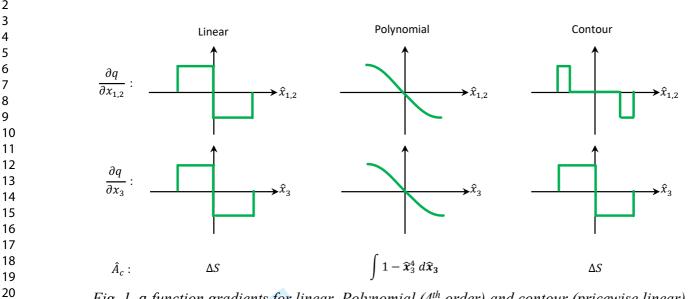
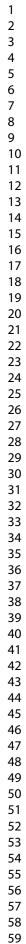


Fig. 1. q function gradients for linear, Polynomial (4th order) and contour (pricewise linear) functions.

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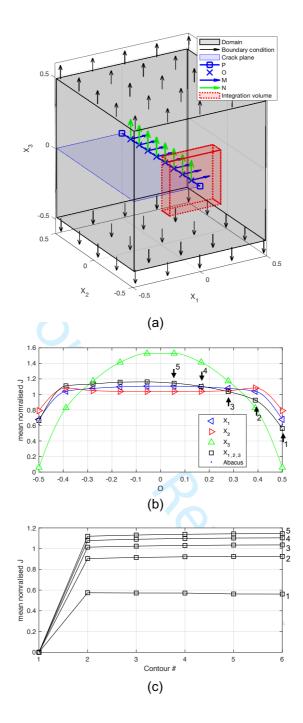


Fig. 4. (a) Unit volume showing boundary conditions (arrow), crack path P, crack segmentation and integration volume V. (b) showing the respective J values commuted along the crack path P. (c) J contours for points 1 – 5 shown in (b) indicating J convergence.

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Strain

BC

1 2 3 4 U noise [%]

4 5 0

J error [%]

 3.9 ± 0.59

 15.3 ± 2.3

 95 ± 14

0

1

5

BC

1 2 3 4 U noise [%]

J error [%]

 4.4 ± 0.24

 16.8 ± 1.1

 $103\pm\ 6.9$

5

BC₁

1 2 3 4 U noise [%]

J error [%]

 0.72 ± 0.079

 2.9 ± 0.32

 18 ± 2.1

1 %

2 %

5 %

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Fig. 5. Influence of random displacement noise on the computation of J. The table below
provides values for the contour q function.

ιcemen. values for .

BC.

2 3 X₁ [%]

J error [%]

0.27 - 0.10

0.42 - 0.25

4 5 0

1

BC.

Linear

Polynomial

Contour

X, [%] 3

3

[%] ×

X, [%] 3

1 %

2 %

0

0 1 2 3 X₁ [%]

J error [%]

0.20 - 0.10

0.72 - 0.27

5 0

4

BC,

3 10

2 3 X₁ [%]

J error [%]

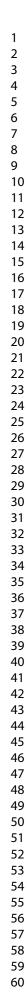
5.2 - 1.3

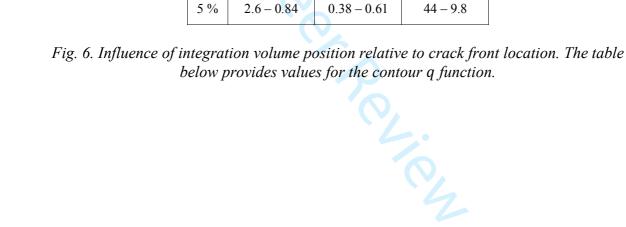
13 - 3.3

44 - 9.8

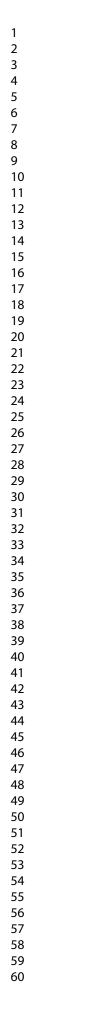
4

R





Strain



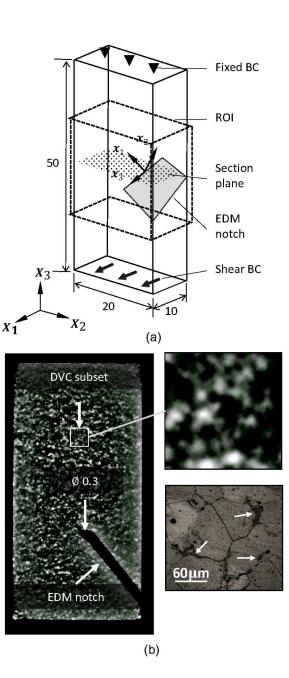


Fig. 7.(a) Schematic of mixed-mode specimen, highlighting BCs, ROI over which displacements were computed using DVC, and the mid-plane section. (b) X-CT scan image of mid-plane showing the EDM notch of 0.3 mm, DVC subset and micrograph (highlighting contrasting features).

Strain

[mm]

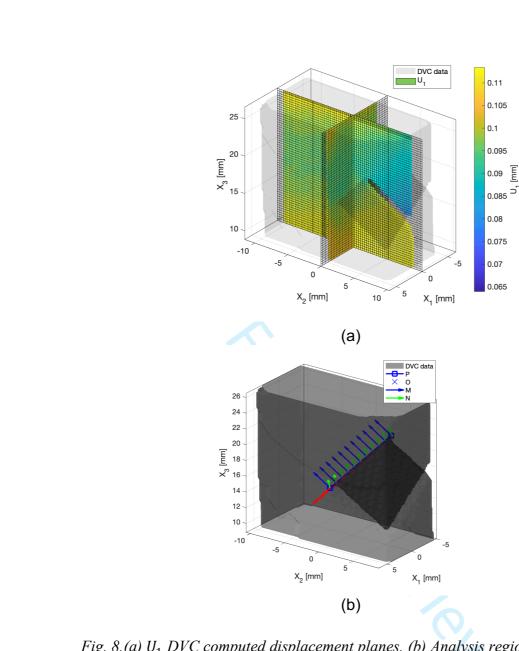


Fig. 8.(a) U_1 DVC computed displacement planes. (b) Analysis region showing the crack path P, crack segmentation centres O_i , and crack orientation defined by M_i and N_i .

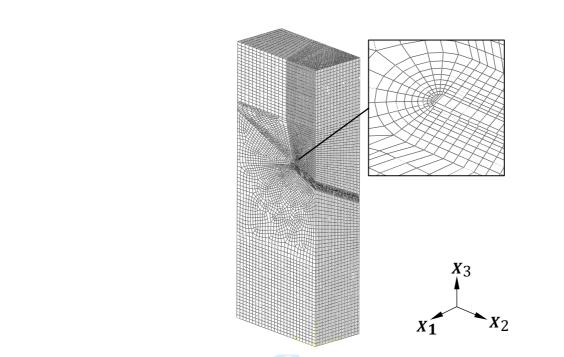
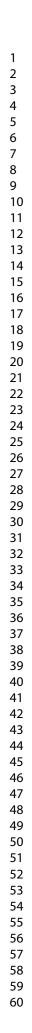


Fig. 9. Equivalent FE model of experimental setup. Insert showing notch tip mesh detail allowing for 10 J contours.



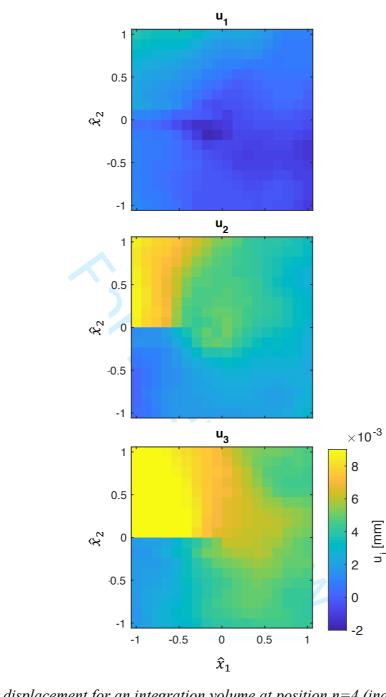


Fig. 10. DVC displacement for an integration volume at position n=4 (indicated in Fig. 11). Displacements are shown for at $x_3 = 0$.

Strain

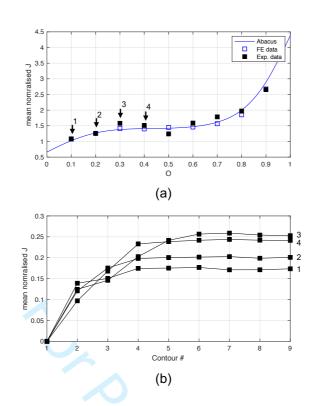


Fig. 11. Mean normalised J values along the crack path from the experiment and equivalent FE model. (b) J contours for points 1 - 4 shown in (a) indicating J convergence.

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