

BOUNDARY INTEGRAL AND FINITE ELEMENT SIMULATION OF ELECTROMAGNETIC NDE PHENOMENA

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INTRODUCTION

Finite element (FE) studies of energy/material interactions associated with the nondestructive evaluation (NDE) of materials have not only yielded useful information concerning the physics of new NDE phenomena [1] but also provided "test-beds" for the simulation of NDE situations too difficult to replicate in a laboratory environment [2]. FE code has been developed for the analysis of those NDE processes governed by elliptic [3], parabolic [4] and hyperbolic [5] partial differential equation (PDE) types taking advantage of axisymmetry wherever possible in order to conserve computer capacity. In those situations requiring fine spatial and/or temporal discretization, it has been found that the FE code makes excessive demands on even the best computer resources. Examples of this situation include the finite element modeling of the remote field effect in large diameter pipelines [6] and the simulation of ultrasonic wave propagation through large structures [7].

The boundary element (BE) method, on the other hand, is ideally suited to infinite domain problems and generally requires a smaller set of equations to be solved, particularly when a solution is needed at only discrete points on the boundary of the material, as is often the case in NDE studies.

Rizzo [8] describes the common theoretical background of both FE and BE methods. Zienkiewicz et al [9], Li et al [10], and Brebbia and Georgiou [11] discuss the application of both methods to structural mechanics problems and also show how FE and BE discretized regions can be coupled to give a hybrid solution procedure.

This research is concerned with the relative merits and limitations of FE and BE methods as applied to the simulation of electromagnetic NDE phenomena. To illustrate the procedure, a comparative study is described in this paper of the direct current potential drop (DCPD) technique as used in fatigue crack monitoring.

POTENTIAL DROP METHOD

The DCPD method has been accepted in fracture mechanics as one of the most accurate and efficient methods for monitoring crack growth [12]. It has been applied to monitor fatigue, stress corrosion and creep cracks. It is also useful for measuring velocities of fast running cleavage cracks and evaluating the extent of crack closure in fatigue crack propagation studies [12,13].

In the DCPD method (Fig. 1), a constant current is passed through a cracked specimen, and the electric potential drop across the crack is monitored. As the crack increases, the uncracked cross-sectional area of the test piece decreases, the current path resistance increases and thus the potential increases. In practice for a particular geometry, calibration curves are given in the form of U/U_0 vs A/W where U_0 is the reference potential drop across the initial cracked specimen (Fig. 2), U is the potential drop as the crack length increases, and A/W is the crack length to width ratio. These calibration curves have been determined experimentally [16-18], analytically (Johnson's formula) [14,15], and numerically [12,13], and they are accepted for use in fracture mechanics.

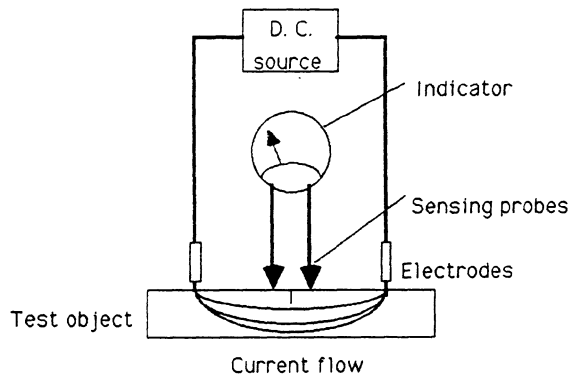


Fig. 1. DCPD experimental setup.

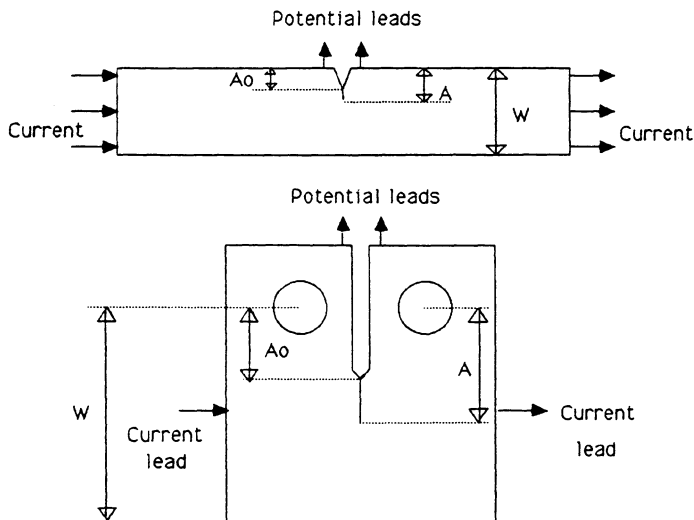


Fig. 2. SEN and CT geometry.

The underlying elliptic PDE describing the DCPD technique is Laplace's equation

$$\nabla^2 V = 0 \quad (1)$$

where V is the steady state voltage in the specimen geometry, with a constant current I in the plane of the geometry. The next section describes briefly the procedure used in FE and BE analysis to solve Laplace's equation.

FINITE ELEMENT FORMULATION

Laplace's equation is solved by the FE method through a variational formulation. The functional for Eq. (1) becomes [12,13,20]

$$F = \int_v \frac{1}{2} \sigma \left[\left(\frac{\partial V}{\partial x} \right)^2 + \left(\frac{\partial V}{\partial y} \right)^2 \right] dv - \int_s V_s i_s ds \quad (2)$$

where v is the volume enclosed, S is the surface that bounds the volume v , V_s and i_s are the voltage and current specified at the boundary. By taking the first variation of F and setting it to zero, that is

$$\delta F = 0 \quad (3)$$

the potential V is obtained by solving the matrix equation

$$[S][V] = [Q] \quad (4)$$

where $[S]$ is the conductivity matrix and $[Q]$ is the forcing vector due to the boundary conditions. The FE meshes for the two geometries are shown in Figs. 3a and 3b.

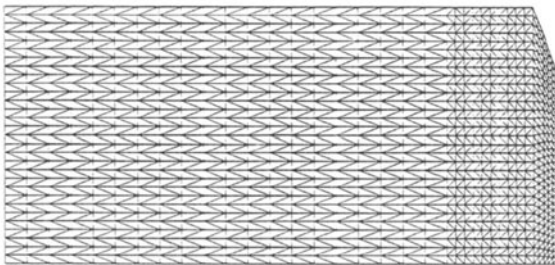


Fig. 3a. SEN finite element mesh.

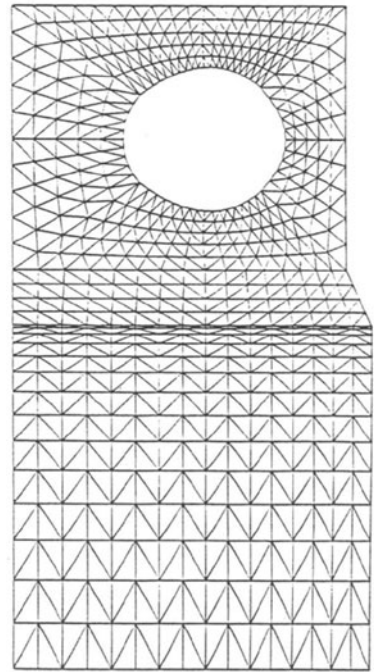


Fig. 3b. CT finite element mesh.

BOUNDARY ELEMENT FORMULATION

The boundary integral equation corresponding to Eq. (1) can be written as

$$c(\xi)V(\xi) = \int_{ds} \left[F(x, \xi) \frac{\partial V}{\partial n}(x) - \frac{\partial F}{\partial n}(x, \xi)V(x) \right] ds(x) \quad (5)$$

where ds is the boundary, F denotes a fundamental solution to the elliptic pde, and the coefficient $c(\xi)$ is the interior angle 2π or π [21]. The fundamental solution is the Green's function, which is

$$F(x, \xi) = -\frac{1}{2\pi\sigma} \ln(r) \quad (6)$$

where r is the vector from ξ to x . This integral equation is solved by standard numerical techniques for the potential V on the boundary.

RESULTS

The geometries considered for comparison are the single edge notched (SEN) specimen and the compact tension (CT) specimen (Fig. 2). Fig. 4 compares the FE and BE results, and Johnson's formula for the SEN specimen. Calibration curves for the CT test piece are shown in Fig. 5. The FE and BE results indicate less than 10 percent deviation from Johnson's formula. The deviation is more pronounced for the CT specimen as the input current is applied only at a point. The memory requirements and the execution times are outlined in Table I. The results in Table I are rough estimates as an initial study.

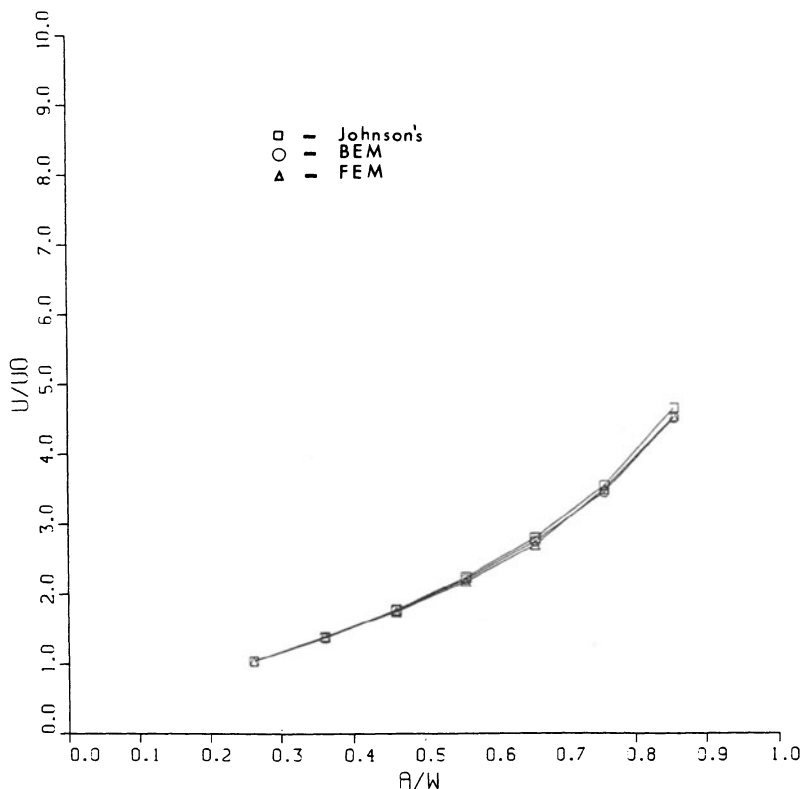


Fig. 4. Calibration curves for the SEN specimen.

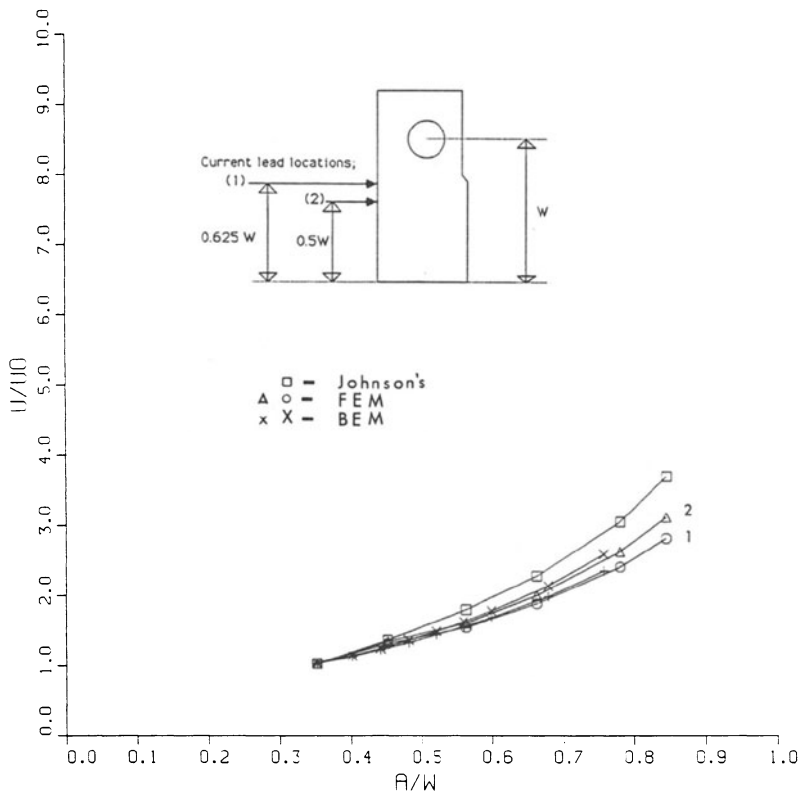


Fig. 5. Calibration curves for the CT specimen.

Table I. Comparison of computer resources for FE and BE.

| | SINGLE EDGE NOTCH | | COMPACT TENSION | |
|--------------------|-------------------|-------|-----------------|------|
| | FE | BE | FE | BE |
| Number of Nodes | 1116 | 38 | 972 | 68 |
| Number of Elements | 2100 | 19 | 1720 | 34 |
| Bandwidth | 33 | 39 | 54 | 69 |
| CPU Time (seconds) | 3.05 | 0.566 | 4.06 | 1.07 |
| Total Memory | 46K | 3K | 53K | 9K |

Figs. 6a and 6b are the equi-potential contours obtained by post-processing of the FE method results. These contour plots help to visualize the effect of the current input and the behavior of the voltage contours as the crack propagates in the material. (The current contours, if plotted, would be orthogonal to the voltage

contours). The equi-potential contours were used to derive the optimum condition for the current input and voltage measurement positions [18]. To generate the calibration curves, the voltage is monitored at only one point near the crack edge, as the crack length increases. Thus, the BEM is more convenient in the generation of the calibration curves. The computer resources in Table I are self explanatory and indicate the tradeoffs required for the two numerical techniques.

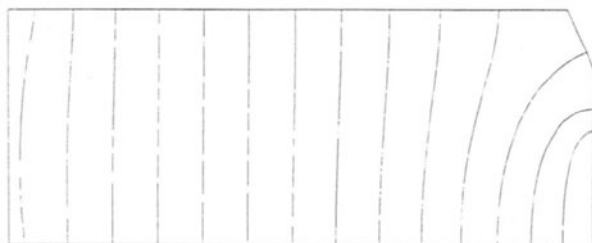


Fig. 6a. Equipotential contours for the SEN specimen using FE.

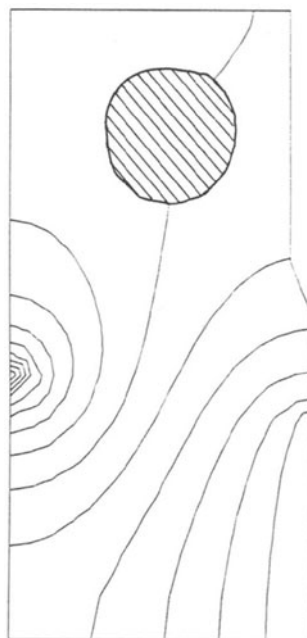


Fig. 6b. Equipotential contours for the CT specimen using FE.

CONCLUSIONS

Despite the preliminary nature of this study, a number of tentative observations can be drawn from the BE/FE results.

1. Where specific output data is needed at discrete points on a test geometry (as is often the case in NDT), the BE approach requires much less computational effort.
2. Where a complete qualitative and quantitative picture of the energy/material interaction physics is required, then the FE technique is most efficient.

These comments apply to the solution for Laplace's equation for the DCPD testing of isotropic materials described in this paper. A more detailed study is already underway to extend the comparison to ACPD and eddy current NDT methods.

With regard to the DCPD method for measuring fatigue crack growth, it is interesting to note that for the compact tension specimen, Johnson's formula appears to be unduly pessimistic in its estimation of crack length.

ACKNOWLEDGEMENT

The authors would like to thank Mr. Zhongqing You for his helpful suggestions.

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