# TWO FINITE ELEMENT TECHNIQUES FOR COMPUTING MODE I STRESS INTENSITY FACTORS IN TWO- OR THREE-DIMENSIONAL PROBLEMS 

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Sponsor: R. D. Cheverton Manuscript Completed: November 1979

Date Published: February 1981

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Prepared for the
Office of Nuclear Regulatory Research
U. S. Nuclear Regulatory Commission Washington, DC 20555
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NRC FIN No. B0119

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Union Carbide Corporation, Nuclear Division operating the
Oak Ridge Gaseous Diffusion Plant . Oak Ridge National Laboratory Oak Ridge Y-12 Plant . Paducah Gaseous Diffusion Plant
under Contract No. W-7405-eng-26
for the
Department of Energy
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## FOREWORD

The work reported here was performed at Oak Ridge Nationai Laboratory (ORNL) under sponsorship of the U.S. Nuclear Regulatory Commission's (NRC's) Heavy-Se son Steel Technolngy Program, which is directed by ORNL. The program is conducted as part of the C NL Pressure Vessel Technology Program, of which G. D. Whitman is manager. The manager for the 'RC is Milton Vagins.

This report is designated Heavy-Section Steel Technology Program Technical Report 58. Prior reports in this series are listed below:

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## ACKNOWLEDGMENTS


#### Abstract

The work described in this report was performed with funds provided by the Heavy-Section Steel Technolcgy (HSST) program under sponsorship of the U. S. Nuclear Regulatory Commission (NRC). The author gratefully acknowledges the encouragement provided by many individuals: R. D. Cheverton, G. D. Whitman, R. H. Bryan, and J. G. Merkle. The author gratefully acknowledges the help and support that W. G. Dodge has provided in connection with the ADINA code. Acknowledgment is also due R. S. Wallace for his assistance with many of the computations and P. G. Fowler for generating the 3-D plots.


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## NOMENCLATURE

## Lat in <br> Characters

| a | Crack length or crack radius |
| :---: | :---: |
| A | Crack area |
| B | Strain-displacement transformation matrix |
| $\mathrm{d}_{\text {ii }}$ | A term on the diagonal of the stress-strain transformation matrix |
| D | Diagonal matrix or stress-strain transformation matrix |
| E | Young's modulus of elasticity |
| $\mathrm{f}_{\text {i }}$ | Modified right-hand side of the upper triangularized matrix |
| F | Total force |
| $\mathrm{F}_{\mathrm{B}}$ | Body force per unit volume |
| $\mathrm{F}_{\mathrm{e}}$ | Nodal forces statically equivalent to the body forces plus the surface tractions |
| $\mathrm{F}_{0}$ | Nodal forces statically equivalent to the thermal plus initial strains |
| G | Potential energy released per unit extension or crack area (the so-called energy release "rate") |
| G | Modulus of rigidity |
| $\mathrm{K}_{\mathrm{I}}, \mathrm{K}_{12 \mathrm{D}}, \mathrm{K}_{\text {I3D }}$ | Stress intensity factor, in 2-D and 3-D, respectively |
| $\ell$ | Length parameter |
| L | Lower triangular matrix |
| N | Field to nodal values transformation matrix (interpolating functions) |
| P | Potential energy of the structure |
| Q | Thermal strain energy of the structure |
| r | Distance to crack root in polar coordinate system |


| Latin <br> Characters (Contd.) |  |
| :---: | :---: |
| T | Surface traction per unit area |
| u | Generalized nodal displacements |
| $u_{n}, u_{z}$ | Noalal displacements in the $n-$ and $z$-directions, respectively |
| $u^{\prime}$ | Generalized displacement field |
| U | Upper triangular matrix |
| V | Volume |
| w | Width parameter |
| W | Strain energy density function |
| Greek |  |
| Characters |  |
| $\alpha$ | Coefficient of linear expansion |
| $\varepsilon$ | Generalized strain matrix |
| $\varepsilon_{0}$ | Initial or thermal strain matrix |
| $v$ | Poisson's ratio |
| $\sigma$ | Gencralized stress matrix |
| $\sigma_{n z}$ | Stresses in the n - and z -directions, respectively |
| ${ }^{\tau} \mathrm{nz}$ | Shear stress in the $n-z$ direction |
| $\theta$ | Polar angle between $n-$ and r -directions |
| $\phi$ | Angular location of a point on the crack front |

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NOMENCLATURE (Contd.)
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| $2-D, 3-D$ | Two- and three-dimensional |
| :---: | :---: |
| ADINA | An FE computer code for Automatic Dynamic Incremental Nonlinear Analysis |
| ASTM | American Society for Testing and Materials |
| CPU | Central Processor Unit |
| DOF | Degrees of freedom |
| FE | Finite element(s) |
| FMECH | A 2-D FE based computer code to calculate $K_{I}$ using energy release methods |
| GEN8 | An FE mesh generation computer code |
| HEATINGS | A finite difference based heat conduction computer code |
| HSST | Heavy-Section Steel Technology |
| IO | Input or output computer operation |
| LEFM | Linear elastic fracture mechanics |
| MERGE | An FE mesh generation computer code |
| NRC | Nuclear Regulatory Commission |
| ORNL | Oak Ridge National Laboratory, Oak Ridge, Tennessee |
| PE | Potential energy |
| QMESH | An FE mesh generation |
| REDUCE | An FE mesh node-numbering optimizer |
| STRIP | An FE mesh generation computer code |
| TSE-4 | The fourth thermal shock experiment |
| V8 | An FE mesh generation computer code |
| V-8 | One of the experimental test cylinders of the HSST progran |

# STRESS INTENSITY FACTORS IN TWO- OR 

THREE-DIMENSIONAL PROBLEMS
S. K. Iskander


#### Abstract

Two finite element (FE) approaches were used to calculate opening mode I stress intensity factors ( $\mathrm{K}_{\mathrm{I}}$ ) in two- and threedimensional (2-D and 3-D) problems for the Heavy-Section Steel Technology (HSST) program. For problems that can be modeled in two dimensions, two techniques were used. One of these may be termed an "energy release rate" technique, and the other is based on the classical near-tip displacement and stress field equations. For three-dimensional problems, only the latter technique was used.

In the energy release technique, $K_{I}$ is calculated as the change in potential energy of the structure due to a small change in crack length. The potential energy is calculated by the FE method but without completely solving the system of linear equations for the displacements. Furthermore, the system of linear equations is only slightly perturbed by the change in crack length and, therefore, many computations need not be repeated for the second structure with the slight change in crack length. Implementation of these last two items has resulted in considerable savings in the calculation of $K_{I}$ as $c$ mpared to two complete FE analyses. These ideas are incor,orated in the FMECH code.

The accuracy of the methods has been checked by comparing the results of the two approaches with each other and with closed form solutions. It is estimated that the accuracy of the results is about $\pm 5 \%$.


## 1. INTRODUCTION

Flaws in structural components may occur during the fabrication processes or after the component has been put into service. Such flaws may escape detection in inspection. An important consideration in the design and analysis of such structural components is the stability of such flaws not only under normal operating conditions, but under hypo chetical accident conditions.

Among the significant parameters in judging the propensity of flaws to propagate is the stress intensity factor, which depends upon the geometry and loading of the structure. These factors may be classified into three basic types, each associated with a local mode of deformation near the crack tip as shown in Figure 1 [1]. Conceptually, the methods described it this report can be used for the determination of the stress intensity factor associated with any of the three modes. This report, however, deals with the opening mode I type of deformation, which implies a loading sym.etric with respect to the plane of the crack. All the methods described are based upon the basic assumptions of linear-elastic fracture mechanics (LEFM), viz., that all nonlinear effects are confined to a "small" region surrounding the crack tip and, consequently, may be neglected. A linear elastic analysis then approximates the essential features of such cases.

The number of "closed" form solutions in LEFM are limited. For the case of an arbitrarily shaped structure subjected to a general system of loads, numerical methods are used. There are many such methods, among which the influence function [2], the boundary integral [3], and the FE [4] approaches are noteworthy. Of these approaches, the ones based


Figure 1. The three basic modes of deformation associated with cracks
upon FE have been particularly popular due to the widespread availability of general-purpose computer codes, and this report discusses the implementation of two FE procerures used to calculate $K_{I}$. One of these is based upon the so-called energy release concept of LEFM and the other on the classical relationship between the neur-tip stress (or displacement) fields and the stress intensity factor $\mathrm{K}_{\mathrm{I}} .^{1}$ The energy release rate concept has been used with our 2-D models only, whereas the stress/displacement method has been used with both our 2-D and 3-D models. Verification of the methods has been carried out by comparison with closed form solutions and by comparing the results of the two different FE approaches for the 2-D problems.

The objective of this report is to document the energy release and stress/displacement methods as they have been implemented for the HSST program. Thus, in Chapter 2 two approaches based upon the energy release method are presented. The first approach is labeled the potential energy method, whereas the second one is termed the strain energy method. It is shown that, for special cases of either external loads only or thermal gradients only, both methods give the correct stress intensity factor. In cases where both types of loads are present, the strain energy approach will lead to erroneous results. Also presented is the complete derivation of a simple procedure to extract the potential energy from the upper triangularized system of equations arising in FE methods without solving explicitly for the displacements. The stress/displacement method is also documented.

[^0]In Chapter 3 , results from several $2-$ and $3-$ D examples are presented. First, as parial verification of the methods given in Chapter 2, a problem for which a closed form solution exists was solved in both $2-D$ and $3-D$ cases. Second, in $2-D$ cases, both the potential energy and stress/displacement methods were used to solve the same problem. In all cases, the results were within a few percentage points of the closed form solution or each other. Also presented are the results of two 3-D analyses performed for the HSST program. The first analysis explores, for several different cylinder lengths, the variation of $K_{I}$ along the length of a uniform depth axial crack in a cylinder subjected to a thermal shock. The second deals with the variation of $K_{I}$ along the crack front for several different load cases of the $V-8$ cylinder.

One important task in the FE modeling--that of mesh generation--is also one of the most time-consuming. Many codes have been developed by the author to facilitate this task and they will be mentioned briefly. The method used in optimizing the profile of the system of linear equations occurring in our 3-D FE analysis is also mentioned. Optimization is cricical since, in many cases, analysis would not be possible due to limitations imposed by the computer environment.

## 2. METHODS OF ANALYSIS

### 2.1. Introduction

In this chapter, both the energy release rate and the stress/displacement methods will be discussed.

### 2.2. Energy Release Rate Methods

This section presents the basis of the energy release rate method for calculation $K_{I}$ as it is implemented in the FMECH code. This method relies upon the calculation of the PE of the structure, and this is accomplished by reans of the FE method. A brief derivation of the governing equations of the FE method based upon PE considerations is first given. These equations will be referred to in a discussion of the errors that arise if the strain energy is used in place of the potential energy in situations where both thermal and mechanical loads act on the structure (since some authors have used the strain energy release rate to calculate $K_{I}$ ). The governing equations of the FE method are also referred to in reviewing an economical computational method for calculating the PE. It is shown that the PE of the structure is available at the end of the forward elimination process and, hence, if the displacements or stresses are not required, the FE solution can be terminated at this point, thereby saving the cost of back substitution and the cost of computing the displacements and stresses.

One of the earliest papers in fracture mechanics was that of Griffith [5]. It introduced the concept of energy release duc. the propagation of cracks in brittle materials. Later, Irwin [6] and

Orowan [7] extended Griffith's work to account for the "small" plastic zone ahead of the crack tip; so this concept may also be termed the Griffith-Irwin-Orowan theory.

Let $P$ be the total potential energy (PE) of deformation of a structure, and $G$ dA be the total elastic energy made available when the crack extends by an amount dA , where A is the crack surface area (one side of the crack); then

$$
\begin{equation*}
G=-\frac{\mathrm{dP}}{\mathrm{dA}} \tag{1}
\end{equation*}
$$

Furthermore, Irwin [8] showed that

$$
\begin{array}{rlr}
G & =\frac{1}{E} K_{I}^{2} & \text { for plane stress } \\
\text { and } G & =\frac{1-v^{2}}{E} K_{I}^{2} & \text { for plane strain } \tag{2}
\end{array}
$$

where

$$
\begin{aligned}
& \mathrm{K}_{\mathrm{I}}=\text { opening mode stress intensity factor, } \\
& \mathrm{E}=\text { modulus of elasticity } \\
& \mathrm{V}=\text { Poisson's ratio. }
\end{aligned}
$$

Thus, it is possible to est imate $K_{I}$ by evaluating the change in the PE of a structure $(\triangle P)$ due to a small extension of the crack surface $(\triangle A)$.

There are various approaches to implementing energy methods using FE analysis. One approach would be to perform an FE analysis on the cracked structure, then extend the crack by a small amount and perform a second FE analysis, thus determining the PE corresponding to each crack length [9]. The disadvantage of this method is the necessity of having to perform two complete FE analyses.

Rather than perform twc complete and separate FE analyses to obtain $K_{I}$, it is possible to optimize the process in a manner that reduces the costs to an amount equal to or less than that for a single complete FE analysis. In describing this optimization, many of the basic equations of the FE method will be referred to and, therefore, thry will now be reviewed from a PE viewpoint although such a derivation is common in the PE 1iterature.

The total PE of a structure is given by [11, p. 384]

$$
\begin{equation*}
P=\int_{\text {vol }} W d V-\int_{v o l} u^{\prime} F_{B} d V-\int_{\text {area }} u^{\prime} T d A \tag{3}
\end{equation*}
$$

where

$$
\begin{aligned}
& \mathrm{W}=\text { strain energy density }, \\
& \mathrm{u}^{\prime}=\text { displacement field }, \\
& \mathrm{F}_{\mathrm{B}}=\text { body forces per unit volume, } \\
& \mathrm{T}=\text { surface tractions per unit area. }
\end{aligned}
$$

The first term is the internal strain energy and the remaining two terms are the work done by the external forces.

Equation (3) will be rewritten in a form suitable for FE analysis.

The strain energy density can be expressed as ${ }^{2}$

$$
W=\frac{1}{2}\left(\varepsilon-\varepsilon_{0}\right)^{T} D\left(\varepsilon-\varepsilon_{0}\right)
$$

[^1]where

```
\(\varepsilon=\) total strain vector,
\(\varepsilon_{0}=\) thermal and/or initial strain vector, and superscript \(T\) denotes
        transpose,
\(\mathrm{D}=\) matrix \(0_{i}^{-}: \overbrace{\text { e }}\) stress-strain transformation matrix, i.e.,
    \(\sigma=\mathrm{D}\left(\varepsilon-\varepsilon_{0}\right), \sigma=\) stress vector.
```

Let $\varepsilon=B u$, where
$u=$ nodal displacement vector,
$B=$ strain-displacement transformation matrix.
Then for the structure

$$
\begin{aligned}
\int W d V= & \frac{1}{2} u^{T} \int_{v o l} B^{T} D B d V u-u^{T} \int_{v o l} B D \varepsilon_{0} d V \\
& +\frac{1}{2} \int_{v o l} \varepsilon_{0}^{T} D \varepsilon_{0} d V \\
= & \frac{1}{2} u^{T} K u-u^{T} F_{0}+Q \quad,
\end{aligned}
$$

where

$$
\begin{aligned}
\mathrm{K} & =\text { stiffness matrix for the structure } \\
& =\int_{\text {vol }} B^{T} D B d V \\
F_{0} & =\begin{array}{l}
\text { nodal loads statically equivalent to the thermal or initial } \\
\text { strains }
\end{array} \\
& =\int_{\text {vol }} B^{T} D \varepsilon_{0} d V \\
Q & =\text { thermal energy }=\frac{1}{2} \int \varepsilon_{0}^{T} D \varepsilon_{0} d V .
\end{aligned}
$$

Thus, the PE can be written as

$$
\mathrm{P}=\frac{1}{2} u^{\mathrm{T}} \mathrm{~K} u-u^{\mathrm{T}} \mathrm{~F}_{0}-u^{\mathrm{T}} \mathrm{~F}_{\mathrm{e}}+\mathrm{Q},
$$

where $F_{e}$ are the nodal forces statically equivalent to the body forces and the surface tractions, i.e.,

$$
\mathrm{F}_{\mathrm{e}}=\int_{\text {area }} \mathrm{N}^{\mathrm{T}} \mathrm{~T} d \mathrm{~A}+\int_{\text {vol }} \mathrm{N}^{\mathrm{T}} \mathrm{~F}_{\mathrm{B}} d V
$$

where $\mathrm{N}=$ the interpolating functions relating the displacement field within an element to fts nodal displacements, viz., $u^{\prime}=N u$.

Thus, the PE van be written in the form

$$
\begin{equation*}
P=\frac{1}{2} u^{T} K u-u^{T} F+Q \tag{4}
\end{equation*}
$$

where $F=F_{e}+F_{0}$ (the sum of the external loads $F_{e}$ and the initial or thermal loads $\mathrm{F}_{0}$ ).

The use of the theorem of miaimum potential energy, viz., the first variation on $P$ with respect to the displacement $\delta u$ [11, p. 382]

$$
\text { viz., } \begin{align*}
\delta \mathrm{P} & =0 \quad \text { leads to } \\
K u & =F \tag{5}
\end{align*}
$$

Equation (5) forms the basis for the FE method. Substituting for F from Eq. (5) into Eq. (4)

$$
P=\frac{1}{2} u^{T} K u-u^{T} K u+Q
$$

or

$$
\begin{equation*}
P=-\frac{1}{2} u^{T} K u+Q \tag{6}
\end{equation*}
$$

Reca!1 that the term $Q$ depends upon the thermal strains $\varepsilon_{0}$ and the material properties $D$. The $\varepsilon_{0}$ are functions of temperature and the coefficient of thermal expansion (which is not contained in D) only. The term $Q$ is therefore iadependent of the boundary conditions. Thus, the
change in PE with respect to a change in crack area can be written as

$$
\begin{equation*}
G=-\frac{d P}{d A}=\frac{d}{d A}\left(\frac{1}{2} u^{T} K u\right) \tag{7}
\end{equation*}
$$

In the absence of external mechanical loads, the PE of a structure, Eq. (3), is given by $\int w d v$, the internal strain energy of the structure. On the other hand, if there are no thermal stresses, the strain is equal to the work done by the external forces ${ }^{3}$ (Clapeyron's Theorem [11, p. 86]). In both these cases, the use of the strain energy in place of the PE in Eq. (1) is valid, except for the sign. This has lead many authors $[1,12]$ to use the term "strain energy release rate" in place of the "potential enersy release rate." Where both external and thermal loads act on a structure, the use of strain energy can lead to errors, since this does not account for the work done by thermal expansion on the external loads.

For example, consider a bar supported at one end in such a manner that a uniform temperature change only causes a stress-free thermal expansion. If the bar is subjected to external mechanical loads, a unfform temperature change will cause no change in the strain energy, but will alter the PE of the structure.

As mentioned above, one obvious and straightforward method of calculating $K_{I}$ by the energy release rate method would be to perform two complete and separate FE analyses. However, two simplifications are possible. First, the element stiffnesses calculated for the first FE analysis are saved and, since they are a function of the dimensions and material properties only, they may be utilized in the second FE analysis

$$
{ }^{3} \text { Mathematically expressed as } \int_{\text {vol }} u^{\prime} F_{B} d V+\int_{\text {area }} u^{\prime} T d A=2 \int W d V \text {. }
$$

and only the boundary conditions are modified to account for the new crack root. Since the cost of the calculation of the element stiffness comprises a major portion of the cost of an FE analysis, the savings are significant.

The second simplification concerns the solution of Eq. (5) and the computation of the displacements and stresses. About half the cost associated with these operations can be saved by means of a scheme due to Hellen [10]. It is shown that the PE of the structure is available at the end of the forward elimination process and thus, if the displacements or the stresses are not required, their costs can be eliminated by cerminating the FE computation at this point. Basically, this scheme enables the computation of the term ( $u^{T} \mathrm{~K} \mathrm{u}$ ) or, equivalently, the work done by the total forces acting through their corresponding displacements $\left(F^{T} u\right)$. The method relies upon the symmetry of the stiffness matrix $K$ and upon the ability to decompose the matrix into the following form [13]

$$
\begin{equation*}
K=L D V, \tag{8}
\end{equation*}
$$

where L and U are lower and upper triangular matrices with units along their principal diagonals, and $D$ is a diagonal matrix whose elements are the "pivots" of Gaussian elimination. Moreover, since $K$ is symmetric, $\mathrm{U}=\mathrm{L}^{\mathrm{T}}$. Then

$$
\begin{equation*}
\mathrm{K}=\mathrm{LD} \mathrm{~L}^{\mathrm{T}} \tag{9}
\end{equation*}
$$

> The solution of Eq. (5) then proceeds as follows:

$$
K u=F
$$

or
L. D $L^{T} \quad u=F$

Forward elimination is symbolically represented as:

$$
\begin{equation*}
\text { D } L^{T} u=L^{-1} F \tag{10a}
\end{equation*}
$$

or, if the pivots are reduced to unity, then

$$
\begin{equation*}
\mathrm{L}^{\mathrm{T}} \mathrm{u}=\mathrm{D}^{-1} \mathrm{~L}^{-1} \mathrm{~F} \tag{10b}
\end{equation*}
$$

For brevity, let

$$
\mathrm{F}^{\prime \prime}=\mathrm{L}^{-1} \mathrm{~F}
$$

and

$$
F^{\prime}=D^{-1} L^{-1} F
$$

then

$$
\mathrm{L}^{\mathrm{T}} \mathrm{u}=\mathrm{F}^{\prime}
$$

In other words, $\mathrm{F}^{\prime}$ (or $\mathrm{F}^{\prime \prime}$ ) is the state of the total force vector at the end of the forvard elimination process. The remaining back substitution process is symbolized as

$$
\begin{equation*}
\mathrm{u}=\mathrm{L}^{\mathrm{T}^{-1}} \mathrm{~F} \tag{11}
\end{equation*}
$$

then

$$
\begin{align*}
u^{T} K u & =\left(F^{, T} L^{-1}\right)\left(L D L^{T}\right)\left(L^{T^{-1}} F^{\prime}\right) \\
& =F^{\prime}{ }^{T} D F^{\prime} \tag{12a}
\end{align*}
$$

or

$$
\begin{equation*}
u^{T} K u=F^{\prime \prime}{ }^{T} D^{-1} F^{\prime \prime} \tag{12~b}
\end{equation*}
$$

Here the form (12b) would be used if, at the end of the forward elimination, the pivots had been normalized; otherwise, (12a) would be used.

Thus,

$$
\begin{equation*}
u^{T} k u=\sum_{i=1}^{n}\left(f_{i}^{\prime}\right)^{2} d_{i i} \tag{13}
\end{equation*}
$$

In other words, the above sum can be accumulated during the forward elimination process and there is no need to carry out the back substitution stage or its attendant cost in $\mathrm{CPU}^{4}$ and $\mathrm{IO},{ }^{5}$ nor is there a need to calculate either the displacements or the stıasses.

There are many advantages to the energy method and these have been documented in detail in [9]. For a given mesh, energy methods give better accuracy than stress/displacement methods and they are insensitive to a large extent to the details of the mesh layout. This may be due to errors that cancel out since the meshes used in both analyses (the first with a crack length $A$ and the second wich a crack length of $(A+\triangle A)$ are identical except for the boundary conditions on the crack front. A significant adzantage is also the high degree of automation which energy methods lend themseives to, as contrasted with the stress or displacement methods mentioned later.

[^2]
### 2.3. Stress and Displacement Methods

These methods rely upon the classical relationships between the near field solution and the stress intensity factor $K_{I}[14]$ (see Figure 2 for coordinate system notation). The stresses are given by

$$
\begin{align*}
\sigma_{n} & =\frac{K_{I}}{(2 \pi r)^{1 / 2}} \cos \frac{\theta}{2}\left|1-\sin \frac{\theta}{2} \sin \frac{3 \theta}{2}\right|+\ldots \\
\sigma_{z} & =\frac{K_{I}}{(2 \pi r)^{1 / 2}} \cos \frac{\theta}{2}\left|1+\sin \frac{\theta}{2} \sin \frac{3 \theta}{2}\right|+\ldots \\
\tau_{n z} & =\frac{K_{I}}{(2 \pi r)^{1 / 2}} \sin \frac{\theta}{2} \cos \frac{\theta}{2} \cos \frac{3 \theta}{2}+\ldots \tag{14a}
\end{align*}
$$

The displacements are (assuming plane strain conditions)

$$
\begin{align*}
& u_{n}=\frac{K_{I}}{G}\left(\frac{r}{2 \pi}\right)^{1 /-} \cos \frac{\theta}{2}\left(1-2 v+\sin ^{2} \frac{\theta}{2}\right)+\ldots \\
& u_{z}=\frac{K_{I}}{G}\left(\frac{r}{2 \pi}\right)^{1 / 2} \sin \frac{\theta}{2}\left(2-2 v-\cos ^{2} \frac{\theta}{2}\right)+\ldots \tag{14b}
\end{align*}
$$

where $\ldots$ represent terms of order $r$ and $r^{(2 n+1) / 2}(n=0,1, \ldots, \alpha)$ that are small with respect to the first term when $r$ becomes small,
$G=$ modulus of rigidity $=E / 2(1+\nu)$ for isotropic materials,
$\nu=$ Poisson's ratio,
$\mathrm{E}=$ Young's modulus of elasticity.
In the region close enough to the crack, $K_{I}$ is calculated from
Eq. (14a) or (14b). If the displacements are used, by taking $\theta=\pi$, we get

$$
\begin{equation*}
K_{I}=\frac{E \sqrt{2 \pi}}{4\left(1-v^{2}\right)}\left(\frac{u_{z}}{\sqrt{r}}\right) \tag{15}
\end{equation*}
$$



Figure 2. The two coordinate systems of axes used. The orientation of the axes is determined by the tangent ( $t$ ) to the crack root edge and the normal ( $n$ ) to the crack edge, with both ( t ) and ( n ) in the plane of the crack. The z -axis is then normal to both $t$ - and $n$-axes. The polar system $(r-\theta)$ is in the $n-z$ plane.

Similarly, if the stresses are used, by taking $\theta=0$, we get

$$
\begin{equation*}
\mathrm{K}_{\mathrm{I}}=(2 \pi)^{1 / 2}\left(\sigma_{z} \sqrt{r}\right) \tag{16}
\end{equation*}
$$

The displacements are the primary solution in most FE computer codes. The stresses are obtained by numerical differentiation with some loss in accuracy. In the interests of accuracy, the stresses are computed at the so-called gaussian integration points, which are interior to the element and not exactly at $\theta=0$. Thus, it may be preferable to use the displacements whenever possible. However, the stresses would be used to calculate $K_{I}$ where the displacements are not available, for example, in 3-D problems in which the crack front does not intersect the free surface at a right angle.

The procedure used to calculate $K_{I}$ is to plot either the term $\mathrm{ur}^{-1 / 2}$ or $\sigma \mathrm{r}^{1 / 2}$ as a function of $r$. The stresses or displacements are taken from the FE solution along a ray normal to the crack front. By extrapolating the curve back to $r=0$, an estimate of $K_{I}$ may be obtained. It has been found by solving problems with closed form solutions that a straight line extrapolation gives reasonably good results [31]. (Also see Sections 3.2.1. and 3.3.1. of this report.)

The total number of degrees of freedom (DOF) and, hence, the computational effort can be decreased considerably by the use of so-called "singularity elements" around the crack tip. These elements reproduce the $\frac{1}{\sqrt{r}}$ singularily in the stresses, thereby allowing the use of relatively few elements in the crack tip region; otherwise, the high stress gradients in that recton necessitate a highly refined mesh so that these gradients can be modeled with reasonable accuracy by either constant or linear
stress elements. One type of singularity element can be produced from any element with mid-side nodes by locating the mid-side nodes adjacent to the crack front at the $\frac{1}{4}$ point (Figure 3). ${ }^{6}$ The stresses in the element will possess the correct singularity in the stresses as well as the correct stiffness, but the displacements at the $\frac{1}{4}$ nodes will be inaccurate [16]. Such an element is often referred to as the $\frac{1}{4}$ point element.

One advantage of the energy methods mentioned is the immediate computation of $K_{I}$ as compared to stress or displacement methods which require a certain amount of manual plotting and computation, thereby lending themselves to a high degree of automation.

The energy methods described above hare been implemented in the finite element code FMECH (Fracture Mechanics). This code is designed to compute $\mathrm{K}_{\mathrm{I}}$ for problems with any form of loading, mechanical or thermal. However, it is particularly useful and relatively efficient in LEFM problems with transient thermal loadings. In such cases, the aim is to calculate $\mathrm{K}_{\mathrm{I}}$ for a large number of different temperature distributions as well as different values of the crack depth.

[^3]DWG. NO. K/G-80-217,
(u)


Figure 3. Quarter point singularity element. Number adjacent to the nodes refer to the $E E$ displacements at these nodes in
Table 7

## 3. EXAMPLES OF ANALYSIS

### 3.1. Introduction

This chapter presents some problems that have been solved using the techniques previously described in this report. The object of this is twofold. First, by solving problems that have either well-established or "closed form" solutions, a partial validation of the method and some assessment of the accuracy of the solution can be made. Second, it presents examples of the type of problems arising in the HSST program that have been solved with these techniques.

The first section deals with $2-$ D problems, the second one with 3-D problems. Both sections start with a problem for which a closed form solution exists then deal with the problems arising in the HSST program.

### 3.2. Two-Dimensional Examples

Two problems will be discussed in this section. The first has a closed form solution and thus gives a rough assessment of the accuracy of the stress/displacement method. The second problem, one that arises in the HSST program, was solved using all the different approaches described in this report and the results have been compared.

### 3.2.1. Finite Length Strip With a Central Crack

The closed form solution of the centrally cracked strip has been given by Isida [17] and has been used extensively to verify numerical solutions. Such a strip with a crack-to-width ratio $(a / w)$ of 0.10 and a length-to-width $(\ell / w)$ of 3 has been modeled with 14 elements and is shown in Figure 4. The

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Figure 4. Finite element idealization for a centrally cracked strip
displacenent method was used. The strass intensity factor $K_{I}$ was normalized by dividing it by $T \sqrt{\pi a},{ }^{7}$ where $T$ is the externally applied traction and 2 a is the crack length (Figure 4). The theoretical vaiue given by Isida is 1.006 . The normalized stress intensity factor calculated from the displacements ${ }^{8}$ is plotted against $r$, the distance from the crack tip, and is shown in Figure 5. The two straight-1ine fits give the values obtained when using either two or three integrition points ${ }^{9}$ in the FE code. The results for the two integration points are better ( $2 \%$ higher than theoretical) than three integration points ( $7 \%$ lower). Hellen $[29,30]$ has reported results showing similar trends. Two integration points were used in all subsequent work of this type.
3.2.2. Long Axial Flaw in a Cylinder Under Thermal Shock

The cylinder analyzed is the one used in the TSE-4 experiment [19]. In this experiment, the test specimen was a cylinder of 0.533 m ( 21 in. ) outside diameter, 0.241 m ( 9.5 in. ) inside diameter, and 0.914 m ( 36 in. ) long. The material was A508, class 2 with a quench only heat treatment (from $871^{\circ} \mathrm{C}$ ). The long axial flaw was 11 mm ( 0.42 in. ) deep. Initially, t e whole cylinder was heated to $291^{\circ} \mathrm{C}\left(555^{\circ} \mathrm{F}\right)$ then the inside surface was suddenly cooled by a methyl alcohol/water mixture.

[^4]

Figure 5. Estimates of $K_{I}$ as calculated from nodal displacements

The coolant temperature was initially $-25^{\circ} \mathrm{C}\left(-13^{\circ} \mathrm{F}\right)$. The actual experimental temperatures obtained by the thermocouples on the cylinder during the transient were used in the calculation of $K_{I}$. The analysis was performed using two different codes, each with different techniques. The FMECH code utilizes energy-based techniques, but two slightly differing approaches were used. The first approach with this code, which is documented in Reference [9], may be termed the "strain" energy approach and is based upon the fact that, in the absence of external forces, the PE and the internal strain energy are identical (see Eq. (3)). The displacements and stresses are first computed by the FMECH code and then the internal strain energy term is calculated from a quadratic expression in the stresses [9]. The other approach uses a more economical method described in Section 2.2. of this report and has been termed the "potential energy" approach. A typical inesh for both approaches is shown in Figure 6.

The mesh for the displacement method using the ADINA code is shown in Figure 7. The results of the analysis using all of these different approaches are given in Table 1. The maximum difference between the $K_{I}$ values for any particular time interval is $7 \%$. The results calculated by the potential energy approach are more accurate than those obtained by the strain energy method since fewer computations are performed and, therefore, round-off errors are less.

Considering the accuracy demonstrated by the stress/displacement method in Section 3.2., the close agreement between the

(a)


Figure 6. A typical FE mesh used with energy release methods


Figure 7. A typical model for use with displacement or stress methods

Table 1
Comparison of $\mathrm{K}_{\mathrm{I}}$ obtained by energy release rate methods to those obtained by the displacement method

| Time in Transient (Minutes) | $\mathrm{K}_{\mathrm{I}}{\mathrm{MPa} \cdot \mathrm{m}^{1 / 2}\left(\mathrm{ksi}^{1} \cdot \mathrm{~m}^{1 / 2}\right)}^{\text {a }}$ |  |  |
| :---: | :---: | :---: | :---: |
|  | 0.4 | 5.03 | 15.03 |
| Energy Release Rate Methods (FMECH Code) |  |  |  |
| Strain Energy | $\begin{gathered} 26.9 \\ (24.5) \end{gathered}$ | $\begin{gathered} 94.3 \\ (85.8) \end{gathered}$ | $\begin{gathered} 69.0 \\ (62.8) \end{gathered}$ |
| Potential Energy | $\begin{gathered} 27.8 \\ (25.3) \end{gathered}$ | $\begin{gathered} 97.1 \\ (88.4) \end{gathered}$ | $\begin{gathered} 71.1 \\ (64.7) \end{gathered}$ |
| Displacement Method (ADINA Code) |  |  |  |
| 251 Nodes | $\begin{gathered} 25.1 \\ (22.8) \end{gathered}$ | $\begin{gathered} 90.7 \\ (82.5) \end{gathered}$ | $\begin{gathered} 66.5 \\ (60.5) \end{gathered}$ |
| 625 Nodes | $\begin{gathered} 25.5 \\ (23.2) \end{gathered}$ | $\begin{gathered} 91.2 \\ (83.0) \end{gathered}$ | $\begin{gathered} 66.9 \\ (60.9) \end{gathered}$ |


#### Abstract

stress/displacement and the energy methods also indicates that the energy method gives good accuracy.

The main advantage of the energy method over the stress/displacement method is the high degree of automation that the former one lends itself to. Stress/displacement methods as described in this report require some manual effort: extracting the stresses or displacements at the proper locations, calculating the $K_{I}$ values, plotting and extrapolating the values to $r=0$. Some of this effort may be eliminated by autonating some of it. However, an element of judgment will always still be required in fitting the straight line through the data points. The energy method on the other hand yields $\mathrm{K}_{\mathrm{I}}$ values immediately.


### 3.3. Three-Dimensional Examples

In this section, analyses of some 3-D problems will be presented. Although the stress/displacement method is easily applied to 3-D problems, the usual difficulties of generating 3-D FE meshes become more acute in 3-D crack problems in which it is difficult to obtain the number of nodes necessary for a reasonable degree of accuracy and still remain within the available computer capacity. Some of the problems described here were close to the maximum size problems that could be handled with the software and hardware existing at the time.

A brief description of the mesh generation process is given in Appendix $A$, and the details of the FE modeling and analysis are given in Appendix B.

The questions of validating the model and the mesh for 3-D problems are the same as for $2-D$ problems. Partial verification of the overall method was obtained by solving the 3-D problem of an imbedded plane circular crack located centrally in a cylinder which has a closcd form solution. The other examples presented are problems arising in the HSST program.

One of these problems, that of the variation of $K_{I}$ axially along a longitudinal crack of uniform depth in cylinders of differing lengths, implicitly provides some verification of the results since, as the length increases, the stress intensity factors at mid-length should approach those of the corresponding 2-D plane strain problem.

Another problem presented arose during the pretest analysis of the V-8 cylinder. A part-through crack was to be produced on the outside surface of a cylinder in an HSST program test whose purpose was the determination of the influence of a residual stress field on the behavior of a flaw in a pressurized vessel.
3.3.1. Plane Circular Crack Located Centrally in a Cylinder

Th. closed form solution for the case of a long cylinder subjected to a uniform axial tensile load applied at the ends and stress-free cylindrical sides and a centrally located plane circular crack was obtained by Sneddon and Welch [20]. Values of $K_{I}$ from that solution for Poisson's ratio of 0.25 have been reproduced in a graphical form by Rooke [21]. A particular case of such a cylinder was modeled and analyzed. Table 2 gives the dimensions, material properties, and loading used in the analysis. The axisymmetry of the problem allows a sector of any angular

Table 2

Dimensions and material properties used in the analysis of centrally cracked cylinder

## Dimensions

Radius of cylinder Cracked radius Distance from crack plane to ends

254 mm (10 in.)
127 mm ( 5 in.)
812.8 mm (32 in.)

## Macerial Properties

Young's modulus E
$200 \mathrm{GPa}\left(29 \times 10^{6} \mathrm{psi}\right)$
Poisson's rav $\quad v$
0.25

## Loading

End traction
68.9 MPa ( 10 ksi )
dimension to be analyzed, but the restrictions imposed by the code ${ }^{10}$ lead to the choice of a quarter of the cylinder to be modeled. Figure 8 shows the $2-D$ mesh used to generate the $3-1)$ model shown in Figure 9. Details of the FE modeling and analysis are given in Appendix $B$.

The nodal displacements from the FE analysis are given in Table 3. As an illustration of the process of estimating $K_{I}$ from these displacements, the term $u_{z} r^{-1 / 2}$ has been plotted in Figure 10. A best fit straight line extrapolated to $r=0$ gives a value of

$$
\left.u_{z} r^{-1 / 2}\right|_{r=0, \theta=\pi}=7.14 \times 10^{-3} \mathrm{~mm}^{1 / 2} 1.42 \times 10^{-3} \mathrm{in}^{1 / 2}
$$

and the use of expression (15) yields a value of

$$
\mathrm{K}_{\mathrm{I}}=30.2 \mathrm{MPa} \cdot \mathrm{~m}^{1 / 2}\left(27.5 \mathrm{ksi} \cdot \mathrm{in}^{1 / 2}\right)
$$

When this value is normalized by dividing by $2 T \sqrt{a / \pi}$, the stress intensity factor for a plane circular crack of radius a in an infinite medium subjected to a uniform stress of $T$ gives a value of 1.088 . Rooke [19] gives a corresponding value of 1.074 . Thus, the method gives values within $1.5 \%$ of the clnsed form solution. No convergence studies with differing mesh sizes were performed to insure that this result is not fortuitous. However, in the V-8 analysis described later in this report, convergence studies with a mesh similar to this one and a mesh with $50 \%$ more nodes

[^5]

Figure 8. Two-dimensional mesh used to generate the 3-D FE mesh by propagating it normal to the plane of the paper. The resulting 3-D mesh, shown in Figure 9, was used for the analysis of the circular central crack


Figure 9. Three-dimensional FE mesh used with the circular crack problem

## Table 3

Nodal displacements from FE analysis for a central circular crack in the cylindrical bar with $\nu=0.25$ (Ratio of crack-to-cylinder radii $=0.5$ )



Figure 10. Plot of the $u_{z} r^{-1 / 2}$ results from the circular crack problem

## were performed. The results of the finer mesh were almost identical to those from the coarser one. <br> 3.3.2. Variation of $K_{I}$ in an Axially Cracked Cylinder Subjected to Thermal Shock (the End Effect Problem) ${ }^{11}$

The object of this series of analyses was the determination of the effect of the free ends on the values of $K_{I}$ for a specific thermal shock problem. The problem is that of a thick cylinder with an axial crack extending along the whole length, Figure 11 (a). As the length of the cylinder increases, the $K_{I}$ values along $a$ central portion of the crack will approack the $2-$ D plane strain value. The problem of concern is to determine the minimum lengths that will satisfy a required relationship between the $2-D$ and $3-D$ $K_{I}$ values over a specified central portion of the cylinder.

The cylinder analyzed has an inner radius of 343 mm and a wall thickness of 152 mm . It is initially at a uniform temperature of $129^{\circ} \mathrm{C}$. The inside surface is quenched with coolant at $-196^{\circ} \mathrm{C}$. The thermal analysis was performed on a one-dimensional axisymmetrjc model using the HEATING5 [23] code with the data shown in Table 4 and some of the resulting temperature distributions through the wall thickness is given in Figure 12. The raaial temperature distributions were used in the FE analysis of the cracked cylinder. Figure $11(\mathrm{~b})$ shows the models ${ }^{12}$ analyzed while Table 5 gives the cylinder lengths, times in the transient,

[^6]

Figure 11. Notation used in the end effects study: (a) geometry
of the problem, (b) nomenclature

Table 4
Input data for HEATING5 analysis, case designation 13C
(One-dimensional axisymmetric solution)

## Cylinder Radii

| $\quad$Inside <br> Outside | $343 \mathrm{~mm}(13.5 \mathrm{in})$. <br> $\quad 495 \mathrm{~mm}(19.5 \mathrm{in})$ |
| :--- | :--- |
| Initial Temperature |  |$\quad 129{ }^{\circ} \mathrm{C}\left(264{ }^{\circ} \mathrm{F}\right)$.

Specific Heat ${ }^{\text {a }}$

|  | SI |
| :---: | :---: |
| ${ }^{\circ} \mathrm{C}$ | $\mathrm{J} \cdot \mathrm{Kg}^{-1} \mathrm{~K}^{-1}$ |
| -184 | 184 |
| -157 | 251 |
| -129 | 312 |
| -101 | 356 |
| - 73 | 385 |
| - 46 | 406 |
| - 18 | 423 |
| 38 | 452 |
| 93 | 479 |
| 149 | 503 |
| 204 | 528 |
| 260 | 551 |
| 316 | 574 |


|  | US Units |
| :---: | :---: |
| $\frac{{ }^{\circ} \mathrm{F}}{}$ | $\frac{\text { Btu } \cdot 1 \mathrm{~b}^{-1}}{}{ }^{\circ}{ }^{\circ} \mathrm{F}^{-1}$ |
| -300 | 0.044 |
| -250 | 0.060 |
| -200 | 0.0745 |
| -150 | 0.085 |
| -100 | 0.092 |
| -50 | 0.097 |
| 0 | 0.101 |
| 100 | 0.108 |
| 200 | 0.1145 |
| 300 | 0.1202 |
| 400 | 0.1260 |
| 500 | 0.1315 |
| 600 | 0.1370 |

Table 4 (Contd.)

Surface Heat Transfer Coefficient ${ }^{\text {a }}$

| ${ }^{0} \mathrm{C}^{\mathrm{b}}$ | $\mathrm{kW} \cdot \mathrm{m}^{-2} \mathrm{~K}^{-1}$ | ${ }^{{ }^{\circ} \mathrm{F}^{\mathrm{b}}}$ | $\mathrm{Btu} \cdot \mathrm{hr}^{-1} \mathrm{ft}^{-2}{ }^{\circ} \mathrm{F}^{-1}$ |
| :---: | :---: | :---: | :---: |
| -176 | 5.31 | -285 | 936 |
| -162 | 4.05 | -260 | 714 |
| -148 | 3.01 | -235 | 530 |
| -134 | 2.22 | -210 | 391.2 |
| -121 | 1.64 | -185 | 288 |
| -107 | 1.21 | -160 | 213.6 |
| -92.8 | 0.886 | -135 | 156 |
| -78.9 | 0.668 | -110 | 117.6 |
| -71.9 | 0.586 | -97.5 | 103.2 |
| -65 | 0.531 | -85 | 93.6 |
| -58.1 | 0.511 | -72.5 | 90 |

${ }^{\text {a }}$ For temperatures outside this range, the values at the end of the table are used.
$b$ These are the film temperatures used by the HEATING5 code, and are the average of the surface temperature and the coolant temperature.


## Table 5

Data used in the free-end-effects study on cylinders

Temperature distribution designation 13C

Nominal Axial Lengths 2L

| rmm |  | $(\mathrm{ln})$. |
| ---: | ---: | ---: |
|  |  | $(16)$ |
| 406 |  | $(27)$ |
| 686 |  | $(36)$ |
| 914 |  | $(48)$ |
| 1219 |  |  |

Time in Transient
minutes 5
5
1, 5, 10
5

## Radii

Inside
343 mm (13.5 in.)
Outside

495 mm (19.5 in.)

## Material Properties


the material properties, and the $2-D$ plane strain $K_{I}$ values. These $\mathrm{K}_{\mathrm{I}}$ values were calculated using both the energy and the displacement methods. The results are very close to each other, an indication of the accuracy of either method.

The results of calculations for 914 mm (36 in.) length and three different times in the transient (1, 5, and 10 minutes) are shown in Figure 13. The 3-D $K_{I}$ values ( $\mathrm{K}_{\mathrm{I} 3 \mathrm{D}}$ ) were normalized by dividing them by the $2-D K_{I}$ value ( $K_{I 2 D}$ ).

It appears that for this particular problem the end effects are largely independent of the time in the transient. Thus, calculations for other cylinder lengths were made for a single time in the transient.

The results of analyzing different cylinder lengths are shown in Figure 14. The $K_{I}$ values near the mid-length are all greater than the corresponding $2-D$ value, except for the very short cylinder. In the shortest cylinders, the $\mathrm{K}_{\text {I3D }}$ values never attained the $K_{I 2 D}$ values. For the longer cylinders, the $K_{I 3 D}$ values exceeded $K_{I 2 D}$ for some portion of the cylinder length and then dropped rapidly to zero. The percentage of the cylinder length with $K_{I 3 D}$ values within $\pm 15 \%$ of $K_{I 2 D}$ was $23 \%, 47 \%, 67 \%$, and $73 \%$ for the $406,686,914$, and 1219 mm cylinders.

It may be worth noting that the free ends rotate inwards so that the ends appear to be concave, with the deepest point towards the inside radius (Figure 15). This is, of course, due to the inside being cooler than the outside, which tends to result in closing of the crack near the ends. In the 3-D
(u)

NORMALIZED AXIAL DISTANCE FROM MID-LENGTH

Figure 13. Variation of $K_{I}$ with distance from mid-length for different times in the transient for the 914 mm (36 in.) cylinder


[^7]OWG. NO.K/G-80-2174
(u)

CYLINDER LENGTH


TIME IN TRANSIENT
TEMPERATURE DESIGNATION CRACK DEPTH-TO.WALL THICKNESS


Figure 15. Axial displacements st the free end of radial lines on the plane of symmetry
analysis, the nodes on the crack front were free to move and the two nodes at the end of the cylinder and closest to the crack root moved in such a way as to have "crossed" over the line of symmetry (see Figure 16). Such an anomalous behavior, of course, has no physical reality since the other face of the crack will prevent such an occurrence. This effect can be prevented by the use of crack elements, gap elements, etc. However, a simple stratagem was to fix a priomi the nodes that had exhibited such a behavior in a previous analysis. This was tried out and had a negligible effect on the $K_{I 3 D}$ values. The tendency of the nodes at the free ends and closest to the crack tip to close took place in all the cylinder lengths analyzed and, hence, the $K_{\text {I3D }}$ values at the free ends were assigned zero values.
3.3.3. Analysis of the $V-8$ Cylinder

Another example of the 3-D fracture mechanics analysis is that performed on a cylindrical vessel designated as $\mathrm{V}-8$. This vessel was tested in 1979 as a part of the ongoing HSST program, and full details may be found in Reference [24] with a summaty in Reference [25]. It consists of an internally pressurized cylinder, closed at the ends, and with an external flaw.

Figure 17 gives the dimensions of the $\mathrm{v}-8$ cylinder and the associated crack as modeled. Symmetry allowed the modeling of only one quarter of the cylinder. As explained in Appendix A, a 2-D mesh is produced from which the $3-D$ one is generated. Furthermore, the 2-D mesh itself is composed of two separate portions (coarse and fine) that are later merged together, as illustrated in Figure 18.

DWG.NO. K/G-80-2172


Figure 16. Anomaly in crack opening displacements near the free end

(b)

Figure 17. FE model used in the analysis of the $V-8$ cylinder: (a) quarter cylinder analyzed, (b) crack geometry


Figure 18. Two-dimensional , nerating plane for the $V-8$ cylinder analysis showing the "fine" and "coarse" mesh regions

For the purpose of a convergence study, two FE models were analyzed in which the basic 2-D meshes of the cylinder wall differed in the "fine mesh" region (Figure 18) close to the flaw. The "coarse mesh" portion was the same for both meshes. The 2-D models of the fine mesh region are shown in Figures 19 and 20, and these are designated Mesh 1 and Mesh 2, respectively. Mesh 2 produced about $50 \%$ nore degrees of freedom (LOF) in the $3-\mathrm{D}$ mesh. The 3-D meshes were generated from these $2-D$ meshes as described in the Appendix A. Figure 21 illustrates such a 3-D mesh generated from Mesh 1. The angles subtended by the elements in the circumferential direction and wed to generate the $3-\mathrm{D}$ meshes, the number of nodes, elements, DOF, etc., together with the computer resources required in the analyses are given in Appendix $B$. Under an internal pressure of $69 \mathrm{MPa}(10 \mathrm{ksi})$ the nodal displacement at the outside surface of the cylinder at the center of the crack for Mesh 1 and Mesh 2 were $99.5 \mu \mathrm{~m}\left(3.92 \times 10^{-3} \mathrm{in}.\right)$ and $99.8 \mu \mathrm{~m}\left(3.93 \times 10^{-3} \mathrm{in}.\right)$, respectively, a change of about $0.25 \%$. It was concluded that no significant benefit would be realized in analyzing the finer mesh.

The four analyses performed are summarized in Table 6; the resulting displacements normal to both the crack face and the crack front at various angles $\theta$ (Figure $17(b)$ ) are given in Table 7. These displacements were used to calculate the $\mathrm{K}_{\mathrm{I}}$ values given in Table 8 by means of the displacement method described in Chapter 2. These $K_{I}$ values were compared to those calculated by other methods [24, Table 4.6]. It appears that the $\mathrm{K}_{\mathrm{I}}$ values calculated by the method described above are accurate.

CRACK FRONT


Figure 19. Two-dimensional "fine" mesh portion of the generating plane (Mesh 1) for the V-8 cylinder

## CRACK FRONT



Figure 20. Two-dimensional "fine" mesh portion of the generating plane (Mesh 2) for the V-8 cylinder with more nodes than Mesh 1


Figure 21. Three-dimensional mesh generated from that shown in Figure 18 for the V-8 cylinder analysis

$$
\begin{aligned}
& \text { Table } 6 \\
& \text { Cases analyzed for } V-8 \text { test }^{\text {a }} \\
& \text { Young's Modulus } \\
& \text { Poisson's Ratio }
\end{aligned} \quad \mathrm{E}=200 \mathrm{GPa}=0.29 \mathrm{GP}
$$

| Case Number | Loading |
| :---: | :--- |
| 1 | Internal pressure 68.95 MPa |
| 2 | Ditto, using finer Mesh No. 2 |
| 4 | Uniform Pressure on Crack Surface 58.95 MPa <br> from 127.6 MPa at the deepest point in the <br> flaw corresponding to a radius of 431.8 mm <br> to 255 MPa at the outside radius of 495.3 |

${ }^{\text {a }}$ Except for Case 2, these were all performed using Mesh No. 1.

Table 7
Nodal displacements obtained from FE analysis of the $V-8$ cylinder for the four load cases

Load Case 1
Internal Pressure 68.95 MPa with Mesh No. 1

Node Designation ${ }^{a}$
r mm
Angle $\phi$
0
10
20
30
40
50

## .

$\qquad$
14.10
$\qquad$
21.15
$\square$ 4 28.19

| Nodal Displacements $\mu \mathrm{m}$ |  |  |
| :--- | :---: | :---: |
| 47.22 | 57.58 | 65.58 |
| 47.09 | 57.45 | 65.46 |
| 46.74 | 57.05 | 65.13 |
| 46.08 | 56.34 | 64.59 |
| 45.01 | 55.22 | 63.86 |
| 42.47 | 52.93 | 62.89 |

Load Case 2
Internal Pressure 68.95 MPa and "Finer" Mesh (Mesh No. 2)
Node Designation ${ }^{\text {a }}$
r mm

| 2 | $\frac{3}{2}$ | $\frac{4}{18.10}$ |  | $\frac{5}{18.80}$ |
| :--- | :--- | :--- | :--- | :--- |

Angle $\phi$
0
10
20
30
40
50

Nodal Displacements $\mu \mathrm{m}$
$\begin{array}{lllll}38.40 & 47.63 & 54.58 & 60.50 & 65.8\end{array}$
$\begin{array}{lllll}38.30 & 47.47 & 54.43 & 60.38 & 65.74\end{array}$
$\begin{array}{lllll}37.95 & 47.09 & 54.03 & 59.99 & 65.41\end{array}$
$\begin{array}{lllll}37.31 & 46.38 & 53.31 & 59.33 & 64.90\end{array}$
36.32
45.24
52.22
58.42
64.24
33.99
43.13
50.67
57.48
64.29

Table 7 (Contd.)

| Crack Surface Pr | $\begin{array}{r} \text { Load } \\ \text { d with } \end{array}$ | ressur | Pa |
| :---: | :---: | :---: | :---: |
| Node Designation ${ }^{\text {a }}$ | 2 | 3 | 4 |
| r mm | 14.10 | 21.15 | 28.19 |
| Angle $\Phi$ | Nodal Displacements $\mu \mathrm{m}$ |  |  |
| 0 | 23.61 | 28.96 | 33.20 |
| 10 | 23.59 | 28.93 | 33.17 |
| 20 | 23.54 | 28.88 | 33.15 |
| 30 | 23.40 | 28.73 | 33.07 |
| 40 | 23.11 | 28.42 | 32.94 |
| 50 | 21.98 | 27.38 | 32.54 |

Load Case 4
Crack Surface Pressurized with a Linearly Varying Pressure, 127.6 MPa at Crack Root and 255.1 MPa at Outer Surface

$\mathrm{a}_{\text {These }}$ node designations are indicated in Figure 3. Displacements at node 1 are inaccurate as mentioned in text.

Table 8

$$
K_{I} \text { in } \mathrm{MNm}^{-3 / 2} \text { as a function of } \phi^{a}
$$

|  | Angle $\phi$ |  |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| Case Load | $0^{\circ}$ | $10^{\circ}$ | $20^{\circ}$ | $30^{\circ}$ | $40^{\circ}$ | $50^{\circ}$ |
| 1 | 55.2 | 55.5 | 54.9 | 53.5 | 52.0 | 46.4 |
| 2 | 55.7 | 55.9 | 55.7 | 53.7 | 52.2 | 47.4 |
| 3 | 48.3 | 48.3 | 47.9 | 47.7 | 47.5 | 43.2 |
| 4 | 71.4 | 71.9 | 73.8 | 76.7 | 80.1 | 78.2 |

${ }^{a} 60^{\circ}$ is the point at the free surface.

## 4. DISCUSSIONS AND CONCLUSIONS

The energy and stress/displacement methods used to calculate $\mathrm{K}_{\mathrm{I}}$ in 2- and 3-D problems have been documented in this report. These methods have been used for solving problems of interest to the HSST program. The stress/displacement methods are based upon the near field equations relating the stresses and displacements to the stress intensity factors. Some approximation is involved in the use of these equations, namely, the higher order terms are neglected since these terms are, in general, unknown. The term used is valid only in the "vicinity" of the crack. It was therefore important to validate the method through the solution of test problems with closed form solutions. The resulting methodology was used successfully for the solution of both 2- and 3-D problens, some of which have been presented in this report.

Energy methods described in this report do not involve such an approximation at the theoretical level. The approximations made are those associated with the numerical methods used and the resulting errors apply to both energy and stress/displacement methods. As such, energy methods are more reliable and accurate than stress/displacement methods. Their main limitation is the practicability of their application to 3-D problems.

This report also discussed the errors that may arise if the strain energy is used in place of the potential energy in problems where both mechanical and thermal loads occur.

Examples of 2- and 3-D problems solved have been presented. For 2-D problems, these have been solved by both energy and displacement methods, and the results of either method compare well with each other.

This indicates that the methods as applied to the type of problems described in this report are accurate and reliable.

One of the erro-s that may arise in the FE method is that associated with a sufficiently fine mesh, especially in the vicinity of st sep stress gradients such as those associated with cracks. Some assurance that an adequately fine mesh was used has been obtained by meins of limited convergence studies in which the same problem was solved with two meshes, one of which had more degrees of freedom than the other No significant change in the results was observed.

The energy methods as implemented in the FMECH zode use a fournoded quadrilateral, and the mesh in the crack region is fine enough to allow the depth of the crack to be changed by the small amount $\triangle A$ by means of releasing the node at the crack root. This method was preferred to altering the coordinates of crack root node by a small amount to achieve the same purpose because altering the coordinates will distort the two elements sharing that node with a possible loss in accuracy in that region. The stress/displacement method uses an eight-noded $1 / 4$-point element. In terms of computer resources, the eight-noded element is more efficient than a four-noded element. However, the total effort required to compute $K_{I}$ using the stress/displacement method is more than that required for the energy method since the former requires some manual effort. However, for the 3-D problems the stress/displacement method has proven useful since it is easily implemented if a suitable FE code is available and once the required input data is prepared.

In conclusion, each of the methods discussed in this report have their advantages and limitations, but within these limitations both methods are capable of yielding $K_{I}$ values within about $\pm 5 \%$ accuracy.

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## APPENDIX A

## FINITE ELEMENT MESH GENERATION AND PROFILE REDUCTION

## 1. Mesh Generation

One important task (in fact, one of the most time-consuming ones in FE modeling) is that of mesh generation. A large number of codes were written, others acquired and implemented, to facilitate this task.

For the cases described in this report, the geometry of the problem allows 3-D meshes to be generated from $2-D$ ones. By rotating the $2-\mathrm{D}$ mesh, such as that shown in Figure 18 about an axis lying in its plane, a 3-D one may be obtained (Figure 21). Similarly, by propagating a 2-D one, such as that shown in Figure 8, in a direction normal to its glane, the $3-\mathrm{D}$ mesh shown in Figure 9 is produced.

The mesh generating effort may be conveniently divided into two tasks. The first will be termed the " $2-\mathrm{D}$ level" and consists of generating the $2-D$ mesh. The second task, the " $3-\mathrm{D}$ level" one, comprises taking the $2-\mathrm{D}$ mesh and producing from ic the $3-\mathrm{D}$ mesh.

At the 2-D level, the most important task was the coding of a mesh plotting program and implementing it on the time-sharing PDP-10 system [26]. All the $2-\mathrm{D}$ plots shown in this report were produced by this package. It is capable of plotting the whole structure or, by means of user-defined coordinaze limits, it can plot certain areas of the mesh thereby "zooming" on detailed areas. The node and element numbers may be displayed or suppressed.

This :lows the analyst to generate and view the mesh. If it is deemed unsatisfactory, the whole process may be repeated. A satisfactory
mesh may be produced in less than an hour. A similar process, when attempted in a batch mode, would take several days to complete.

Generation of the $2-D$ models of cylinders is accomplished by means of the GEN8 code [26], producing semicircular annuli similar to that shown in Figure A.1(a). The number of elements in the radial and circumferential directions, geometric progression factors for both these directions, and the inside and outside radii are the input parameters.

Strips such as that shown in Figure A. 2 have been useful in the modeling process. The centrally cracked strip and the coarse mesh portion of the $V-8$ cylinder model (both mentioned earlier in this report) as well as many others have been generated by the "STRIP" code [26]. This allows the user to define the number of elements horizontally and vertically as well as their dimensions. The number of horizontal elements may be doubled in number as the mesh is gezerated vertically down, with a onetime transition from four to five elements possible. This allows a mesh to be produced that can be made finer in regions of high stress gradients.

A third code "MERCE" [26j swaps any two elements designated by the user with a so-called singularity block (Figure A.1(b) and A.2(b)) to produce automatically the $1 / 4$-point singularity elements surrounding the crack tip (Figures A.1(c) and A.2(c)).

The "QMESH" system [27], a very useful tool that generates more complex 2-D meshes, was implemented and used to generate the meshes used, for example, in the centrally cracked circular bar and the $V-8$ cylinder. Being a general-purpose package, it is, however, too unwieldy to use for simple meshes such as those mentioned above and local mesh refinements are difficult to accomplish.

(c)

Figure A.1. (a) Typical finite element mesh generated by the "GEN8" code. The singularity block (b) is then "merged" into a pair of elements, e.g., those shown shaded in (a) to produce a model with a "crack" located at $a / w=0.5$, a portion of which is shown in (c)

(b)

Figure A.2. (a) Example of type of mesh produced by STRIP code, and the incorporation of the singularity block (b) by the MERGE code to produce the mesh (c) used in numerical experiments on the center cracked or edge notched bars. Note that the singularity block can be relocated at different positions giving in this case nine "specimens" with $a / w=0.1$ through $a / w=0.9$

The generation of $3-D$ meshes from any of the above $2-D$ meshes is accomplished by means of the $3-D$ and $V-8$ codes [26]. These codes atso produce output that allows the user to relate the $3-\mathrm{D}$ models to the $2-\mathrm{D}$ ones. It also produces output identifying element faces in a particular area, if desired. Such output would be useful in the FE analysis of pressure loaded structures, e.g., the $V-8$ cylinder analysis mentioned in Chapter 3.

The concept of the subdivision of a major task into smaller ones has allowed a considerable amount of modularity in coding the various programs. Furchermore, it permits the user to monitor the satisfactory completion of one task before proceeding to the next one. A major drawback of the $3-D$ mesh generation codes described above is that the whole mesh is penalisea by the presence, at the $2-D$ level, of the fine mesh in the crack region which is then propagated into the third dimension. It is more efficient $\& \rightarrow$ implement a merging scheme at the $3-\mathrm{D}$ level similar to that shown in Figures A .1 and A .2 for $2-\mathrm{D}$ meshes. For 3-D meshes, the process is accomplished in three steps. The first consists of generating a relitively coarse mesh for the $3-\mathrm{D}$ structure to be analyzed. Then, a "block" containing the finer mesh required to adequately model the crack region is generated. The final stage would be co merge this block into the mesh for the $3-D$ structure.

## 2. Profile Reduction

The bandwidth of automatically generated meshes is, generally speaking, larger than necessary. This condition not only leads to uneconomical FE analyses but, in some instances, to cases impossible
to run on the computer. This is causcd primarily by the very high number of input/output ( $1 / 0$ ) operations necessary to store and solve out-of-core the large system of linear equations occurring in the FE method of analysis. This, in turn, causes large wall clock times (WCT) during which large segments of memory (1750K) are tied up for say, 10 hours, and unacceptable amounts of scratch disk space are required. Thus, it is imperative to optimize the node numbering scheme. This is accomplished by means of the computer code "REDUCE" [28].

As an illustration of sone of the computer resources saved by optimizing the node numbering scheme of the FE models, some pertinent numbers from two separate analyses are given in Tables A.1 and A.2. The problem analyzed is the circular crack described earlier in this report. The "before" refers to values associated with the meshes as generated (i.e., with unoptimized meshes) while the "after" are values associated with the optimized meshes. The figures were compiled from the data generated by the ADINA and the REDUCE codes. These figures show a reduction of $20 \%$ to $47 \%$ in the computer resources required. The example problem selected was a small one. In cases like that of Mesh 2 of the $V-8$ analysis, the savings were significant enough to have made the difference between the feasibility or nonfeasibility of the FE analysis.

```
                    Table A.1
Effect of node renumbering on the IBM 360/195
    computer resources required by
                the ADINA computer code
        (Approx. 1700K memory used)}\mp@subsup{}{}{\textrm{a}
            Number of Nodes 1286
            Number of Elements }23
```

|  | Before | After | \% Reduction |
| :---: | :---: | :---: | :---: |
| Number of Matrix Elements (approx.) | $1.897 \times 10^{6}$ | $1.498 \times 10^{6}$ | 21 |
| Maximum Half-Bandwidth | 1415 | 962 | 32 |
| Mean Half-Bandwidth | 552 | 436 | 21 |
| Number of Blocks | 29 | 23 | 21 |
| CPU Minutes | 14 | 9 | 36 |
| 1/0 in Thousands | 39 | 28 | 28 |
| Wall Clock Time ${ }^{\text {b }}$ (Minutes) | 70 | 37 | 47 |
| Approximate Cost ${ }^{\text {c }}$ | \$87 | \$58 | 33 |

${ }^{a^{P}}$ Plane Circular Crack problem.
${ }^{\text {b }}$ On the IBM $360 / 195$ MVT environment, this can vary widely.
${ }^{\mathrm{c}}$ Printed output to tape.

Table A. 2
REDUCE code data

|  | Half-Bandwidth |  |  | Profile (approx.) ${ }^{\text {a }}$ |  |  | $\begin{aligned} & \mathrm{CPU} \\ & \mathrm{Min} \\ & \hline \end{aligned}$ | $\begin{aligned} & \cos ^{\mathrm{b}} \\ & \hline \end{aligned}$ |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  | Before | After | \% | Before | After | \% |  |  |
| Circular Crack | 506 | 331 | 35 | 256,000 | 193,000 | 25 | 6 | 28 |
| Mesh No. $1^{\text {c }}$ | 1270 | 626 | 51 | 653,000 | 392,000 | 40 | 6 | 44 |
| Mesh No. 2 | 1656 | 593 | 64 | 1,278,000 | 777,000 | 39 | 24 | 115 |

${ }^{a_{\text {Profile }}}$ is the sum of the number of terms in the so-called "sky-line", based on the element connectivity, not upon the actual number of DOF.
${ }^{\mathrm{b}}$ Includes print cost.
${ }^{c}$ IBM FORTRAN EXTENDED PLUS compiler generated object code.


#### Abstract

APPENDIX B

\section*{details of the fe fudeling and analysis}

In this appendix, details of the FE modeling and analysis are given. Apart from documenting the data used in the 3-D analyses described in Ch nter 3, this information has proven useful as a basis for estimating the computer resources required for other similar 3-D computations. This information is important since an upper-1imit estimate of the computer resources required in a particuiar run must be declared before the run is made.


Table B. 1
Data used to propagate meshes from $2-D$ into $3-D$


## Table B. 2

## Miscellaneous data of interest related to the 3-D FE analysis using the ADINA code ${ }^{\text {a }}$ <br> (Fast Memory Used 1700 K Bytes)

|  | $\begin{gathered} \text { Circular } \\ \text { Crack } \\ \text { Problum } \\ \hline \end{gathered}$ | End Effects Study Cylinder Lengths, mm |  |  | V-8 Cylinder |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  |  | $406^{b}$ | $686^{\text {c }}$ | $914^{\text {d }}$ | $\begin{aligned} & \text { Mesh } \\ & \text { No. } 1 \\ & \hline \end{aligned}$ | $\begin{gathered} \text { Mesh } \\ \text { No. } 2 \\ \hline \end{gathered}$ |
| Nodes | 1286 | 1045 | 2573 | 2573 | 2328 | 3456 |
| Elements | 234 | 160 | 480 | 480 | 434 | 666 |
| Equations (Degrees of Freedom) | 3439 | 2726 | 7179 | 7179 | 6389 | 9603 |
| $\begin{aligned} & \text { Matrix Elements } \\ & \text { in millions (approx.) } \end{aligned}$ | 1.5 | 0.64 | 3.2 | ?. 2 | 3.1 | 6.2 |
| Max. Half Bandwidth | 962 | 509 | 900 | 900 | 1789 | 1543 |
| Mean Half Bandwidth | 436 | 231 | 943 | 943 | 485 | 645 |
| Max. Block Size | 67565 | 60127 | 39386 | 39386 | 44354 | 29226 |
| Number of Blocks | 29 | 11 | 82 | 82 | 71 | 216 |
| CPU Time (Minutes) <br> IBM 360/195 | 9 | 3.1 | 27 | 28 | 22 | 63 |
| I/C (Thousands) ${ }^{\text {e }}$ | 28 | 46 | 50 | 175 | 105 | 492 |
| $\begin{aligned} & \text { Wal1 Clock Time } \\ & \text { (Minutes) } \end{aligned}$ | 37 | 90 | 175 | 250 | 125 | 585 |
| Approximate Cost (\$) | 58 | 57 | 178 | $235^{\mathrm{g}}$ | $165^{g}$ | $600^{g}$ |

${ }^{\text {a }}$ Object code obtained with the IBM FORTRAN H compiler. This version still had the sequential I/O routines.
${ }^{b} 1500 \mathrm{~K}$ region and 3 K blocking factor for scratch units, unoptimized profile.
${ }^{c}$ One time step of the transient (also applies to 1219 mm cylinder).
$\mathrm{d}_{\text {Three }}$ time steps of the transient.
e Depends upon blocking factor of the scratch units.
${ }^{\mathrm{f}}$ In an "Multiprogramming with a Variable Number of Tasks" (MVT) environment, this can vary considerably for identical runs.
${ }^{g}$ Printed output to tape.

Conversion Factors for Units Used in This Report
(Reference: Metric Practice ASTM E380-76)

| To Convert From | To | Multiply By |  |
| :--- | :--- | :--- | :--- |
| Stress/Pressure | $\mathrm{lbf} \cdot \mathrm{in}^{-2}(\mathrm{psi})$ | $\mathrm{kPa}\left(\mathrm{kN} \cdot \mathrm{m}^{-2}\right)$ | 6.894757 |
| Stress Intensity <br> Factor | $\mathrm{ksi} \cdot \mathrm{in}^{-1 / 2}$ | $\mathrm{MPa} \cdot \mathrm{m}^{1 / 2}\left(\mathrm{MN}^{2} \cdot \mathrm{~m}^{-3 / 2}\right)$ | 1.098843 |
| Coefficient of <br> Heat Transfer | $\mathrm{Btu} \cdot \mathrm{h}^{-1} \cdot \mathrm{ft}^{-2} \cdot{ }^{\circ} \mathrm{F}^{-1}$ | $\mathrm{~W} \cdot \mathrm{~m}^{-2} \cdot \mathrm{~K}^{-1}$ | 5.678263 |
| Heat Capacity | $\mathrm{Btu} \cdot \mathrm{lb}^{-1} \cdot{ }^{\circ} \mathrm{F}^{-1}$ | $\mathrm{~kJ} \cdot \mathrm{~kg}^{-1} \cdot \mathrm{~K}^{-1}$ | 4.186800 |
| Thermal <br> Conductivity | $\mathrm{Btu} \cdot \mathrm{h}^{-1} \cdot \mathrm{ft}^{-2} \cdot{ }^{\circ} \mathrm{F}^{-1}$ | $\mathrm{~W} \cdot \mathrm{~m}^{-1} \cdot \mathrm{~K}^{-1}$ | 1.730735 |
| Density | $\mathrm{lb} \cdot \mathrm{ft}^{-3}$ | $\mathrm{~kg} \cdot \mathrm{~m}^{-3}$ | 16.01846 |

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[^0]:    ${ }^{1}$ This method will be referred to as the stress/displacement method for brevity.

[^1]:    ${ }^{2}$ Vectors and matrices will not be given any special symbol (e.g., $\left\{\varepsilon_{0}\right\}$ or $\left.[D]\right)$ since their definition and context of their use make their meaning unambiguous.

[^2]:    ${ }^{4} \mathrm{CPU}=$ Central Processor Unit--used synonomously for computations.
    ${ }^{5}$ IO $=$ Input or Output operations to disks or tapes.

[^3]:    ${ }^{6}$ Such elements were introduced almost simultaneously by Barsoum [15] and Henshell [16].

[^4]:    ${ }^{7}$ The stress intensity factor for an infinite plate.
    ${ }^{8}$ All the FE analyses with the stress and displacement method were performed using the ADINA Code [18].
    ${ }^{9}$ Implying a $2 \times 2$ or $3 \times 3$ integration order in $2-D$ and $2 \times 2 \times 2$ or $3 \times 3 \times 3$ in 3-D integration order in 3-D.

[^5]:    ${ }^{10}$ In the ADINA code, boundary conditions at the planes of symmetry can only be applied in directions parallel to the $x-y-z$ coordinate axes.

[^6]:    ${ }^{11}$ If the ratio cylinder-length-tomoutside-radius is of the same order of magnitude as unity, neither the plane strain nor the plane stress assumptions are useful and complete 3-D theory must be employed [22, p. 289].
    ${ }^{12}$ Details of the method of mesh generation and profile reduction are given in Appendix $A$. The $2-D$ mesh used to generate the $3-D$ meshes is similar to that shown in Figure 7.

[^7]:    Figure 14. Variation of $\mathrm{K}_{\mathrm{I}}$ with axial distance from fid-length for different cylinder lengths

