

## IV. CONCLUSION

The calculation of eigenmodes in uniform wave guides or in resonant cavities by integral equation and moment method sometimes generates non physical solutions. In this paper a practical criterion for a correct choice of the weighting functions is demonstrated. Its rigorous implementation would be calculus intensive but we show that an approximate implementation efficiently eliminates spurious solutions generated by the conventional Galerkin's method without lengthening computations.

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## A New FEM Approach for Open Boundary Laplace's Problem

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**Abstract**—An efficient improved finite element method (FEM) is presented for electromagnetic Laplace's problems with open boundary. The whole infinite domain is divided into a set of infinite elements instead of ordinary finite elements. Since a special FEM discretization and FEM solving procedure are used, it can not only take much less computer memory than that the conventional FEM needs, but also avoid the calculation error introduced by the truncated boundary or absorbing boundary condition used in conventional FEM.

## I. INTRODUCTION

Recently, FEM has been more and more widely used for many electromagnetic problems with open boundary. Because in FEM the solution domain is discretized with finite elements, only finite domain problems can be handled directly. So, the solution domain with open boundary must be truncated. A relatively simple technique to implement is to select an external boundary with a zero potential to truncate the solution domain. Another alternative is to use an appropriate absorbing boundary condition or infinite element [1]–[3].

This paper presents a new FEM technique for static electromagnetic problems with open boundary. It utilizes a special discretization form to divide the whole infinite domain into infinite triangular elements. So no truncated boundary or absorbing boundary is needed and the calculation error produced by appropriate boundary condition is avoided. As an example, the capacitance matrix for the two coupled microstrips with open-boundary is calculated and the numerical results are compared with those obtained by other methods.

## II. NEW FEM PROCEDURE

We begin with our discussion of a two-dimensional electrostatic open boundary problem. Assuming that the solution domain is  $\Omega_s$  and its open boundary is  $\Gamma_s$ . We use a regular polygon  $\Gamma_0$ , which contains the solution domain  $\Omega_s$ , to divide the whole infinite domain  $\Omega$  into two parts (see Fig. 1). Defining that the region within  $\Gamma_0$  is  $\Omega_{in}$  and the one out of  $\Gamma_0$  is  $\Omega_{out}$ .  $\Gamma_0$  may be placed very close to  $\Gamma_s$ . If  $\Gamma_s$  itself is a regular polygon, then it is selected as  $\Gamma_0$ .

A. Analysis of  $\Omega_{in}$ 

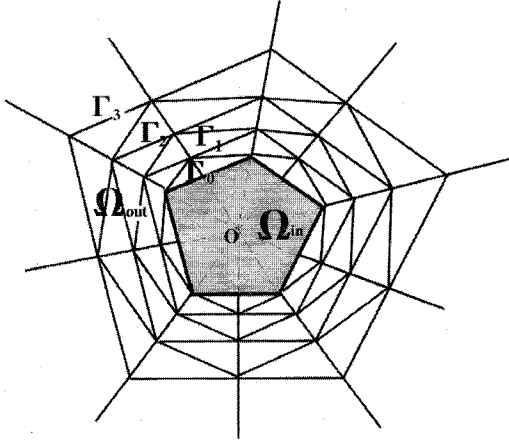
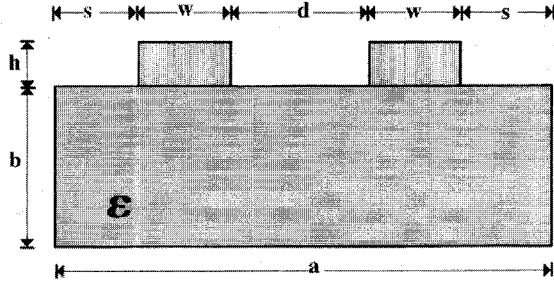
Because the region  $\Omega_{in}$  is a finite domain, it can be treated by conventional FEM. The electromagnetic field distribution can be obtained from the scalar potential  $\Phi(x, y)$  satisfying the Laplace's equation with associated Dirichlet or Neumann boundary condition. The region  $\Omega_{in}$  is subdivided to triangular finite elements and the stiffness matrix  $K_{in}$  can be obtained by assembling each element coefficient matrix. Assuming that the number of nodes on  $\Gamma_0$  is  $M_0$  and the scalar potentials on  $\Gamma_0$  form a column vector  $\Phi_0$  of order  $M_0$ , the number of remaining nodes is  $M_{in}$  and the corresponding potentials form a column vector  $\Phi_{in}$  of order  $M_{in}$ . Thus, the functional within the region  $\Omega_{in}$  can be obtained by

$$F_{in} = \frac{1}{2} (\Phi_{in}^T \quad \Phi_0^T) \begin{pmatrix} K_{in}^{11} & K_{in}^{12} \\ K_{in}^{21} & K_{in}^{22} \end{pmatrix} \begin{pmatrix} \Phi_{in} \\ \Phi_0 \end{pmatrix}. \quad (1)$$

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Fig. 1. Similar discretization of infinite region  $\Omega_{out}$ .Fig. 2. Coupled microstrips ( $\epsilon = 2\epsilon_0$ ,  $a = 11b$ ,  $h = b$ ,  $w = 3b$ ,  $d = 2b$ ,  $s = 1.5b$ ).

### B. Analysis of $\Omega_{out}$

The region  $\Omega_{out}$  is a infinite domain. The FEM discretization is taken as following: first, a coordinate center O within  $\Gamma_0$  is selected. Next, a set of infinite similar polygons of  $\Gamma_0$  are taken as shown in Fig. 2. The point O is the similar center of polygons and the constant  $\xi$  is similar ratio. The set of similar polygons are denoted as  $\Gamma_1, \Gamma_2, \dots, \Gamma_k, \dots$  and the region between  $\Gamma_{k-1}$  and  $\Gamma_k$  is denoted as  $k$ th layer. Each layer is divided into quadrilaterals of number  $M_0 - 1$  by the rays out from similar center point O to nodes on  $\Gamma_0$ . (Note that the polygon apexes must be nodes.) Then each quadrilateral is cut into two triangles. It is necessary that the discretization of every layer is similar as shown in Fig. 1. Assuming that scalar potentials on  $\Gamma_k$  is  $\Phi_k$ . After the element stiffness matrix is obtained by means of conventional FEM, the layer stiffness matrix can be formed by assembling all finite elements on one layer. It can be written as

$$\begin{pmatrix} K_0 & -A^T \\ -A & K'_0 \end{pmatrix} \quad (2)$$

where  $K_0$ ,  $K'_0$ , and  $A$  are all matrices of order  $M_0$ . Because every layer is discretized similarly, all layer stiffnesses are identical. So, if the functional of  $k$ th layer is

$$F_k = \frac{1}{2} (\Phi_{k-1}^T \quad \Phi_k^T) \begin{pmatrix} K_0 & -A^T \\ -A & K'_0 \end{pmatrix} \begin{pmatrix} \Phi_{k-1} \\ \Phi_k \end{pmatrix} \quad (3)$$

then, the functional of the whole infinite domain  $\Omega_{out}$  is a infinite sum as

$$F_{out} = \sum_{k=1}^{\infty} F_k. \quad (4)$$

### C. Synthesis of $\Omega_{in}$ and $\Omega_{out}$

Considering both  $\Omega_{in}$  and  $\Omega_{out}$ , the functional of the whole  $\Omega$  can be written as

$$F = F_{in} + F_{out}. \quad (5)$$

Because the first derivatives of the functional (5) with respect to the node variables should be zero, we can have a set of equations of infinite order

$$\begin{cases} K_{in}^{11} \Phi_{in} + K_{in}^{12} \Phi_0 = 0 \\ K_{in}^{21} \Phi_{in} + K_{in}^{22} \Phi_0 + K_0 \Phi_0 - A^T \Phi_1 = 0 \\ -A \Phi_0 + K \Phi_1 - A^T \Phi_2 = 0 \\ \dots \\ -A \Phi_{k-1} + K \Phi_k - A^T \Phi_{k+1} = 0 \\ \dots \end{cases} \quad (6)$$

where  $K = K_0 + K'_0$ . It is impossible to solve the (6) directly. It is easy to prove that there exists a matrix  $X$  of order  $M_0$  which satisfies

$$\Phi_{k+1} = X \Phi_k \quad (7)$$

where  $X$  is called transfer matrix. Letting  $k = 0$ , from (6) and (7) we have

$$K_{in}^{21} \Phi_{in} + K_{in}^{22} \Phi_0 + K_0 \Phi_0 - A^T X \Phi_0 = 0 \quad (8)$$

$$K_{out} = K_0 - A^T X \quad (9)$$

$$K_{in}^{21} \Phi_{in} + K_{in}^{22} \Phi_0 + K_{out} \Phi_0 = 0. \quad (10)$$

Combining (6) and (10), we have

$$\begin{pmatrix} K_{in}^{11} & K_{in}^{12} \\ K_{in}^{21} & K_{in}^{22} + K_{out} \end{pmatrix} \begin{pmatrix} \Phi_{in} \\ \Phi_0 \end{pmatrix} = 0. \quad (11)$$

Imposing boundary condition within the solution domain  $\Omega_0$ , the scalar potentials  $\Phi_{in}$  and  $\Phi_0$  can be solved from (11). Then from (7), the scalar potentials  $\Phi_k$  of the region  $\Omega_{out}$  can be obtained.

### D. Calculation of Transfer Matrix $X$

Combine (7) with (6), then

$$-A + KX - A^T X^2 = 0. \quad (12)$$

Let

$$\begin{aligned} R_1 &= \begin{pmatrix} K & -A \\ I & 0 \end{pmatrix} \\ R_2 &= \begin{pmatrix} A^T & 0 \\ 0 & I \end{pmatrix} \end{aligned} \quad (13)$$

where  $I$  is unit matrix. From (6) we have:

$$R_1 \begin{pmatrix} \Phi_k \\ \Phi_{k-1} \end{pmatrix} = R_2 \begin{pmatrix} \Phi_{k+1} \\ \Phi_k \end{pmatrix}. \quad (14)$$

Assuming that  $\lambda$  and  $g$  are the eigenvalues and eigenvectors of  $X$ , then

$$Xg = \lambda g. \quad (15)$$

Letting  $k = 1$  and  $\Phi_0 = g$  in (14), (14) together with (7) and (15) will take the form

$$R_1 \begin{pmatrix} \lambda g \\ g \end{pmatrix} = \lambda R_2 \begin{pmatrix} \lambda g \\ g \end{pmatrix}. \quad (16)$$

Equation (16) can now be solved by any standard generalized eigenvalue analysis computer package for the eigenvalues of number

TABLE I  
CAPACITANCE MATRIX FOR THE COUPLED MICROSTRIPS OF FIG. 2

	Reference	Shield	ABC	Higher Order	Infinite FEM	the paper
	[6]	[4]	[4]	ABC[4]	[5]	
$C_{11} \cdot 10^{-10}$	0.9224	1.091	0.9249	0.9230	0.9311	0.9228
$C_{12} \cdot 10^{-11}$	-0.8504	-0.4712	-0.8061	-0.8377	-0.8015	-0.8401

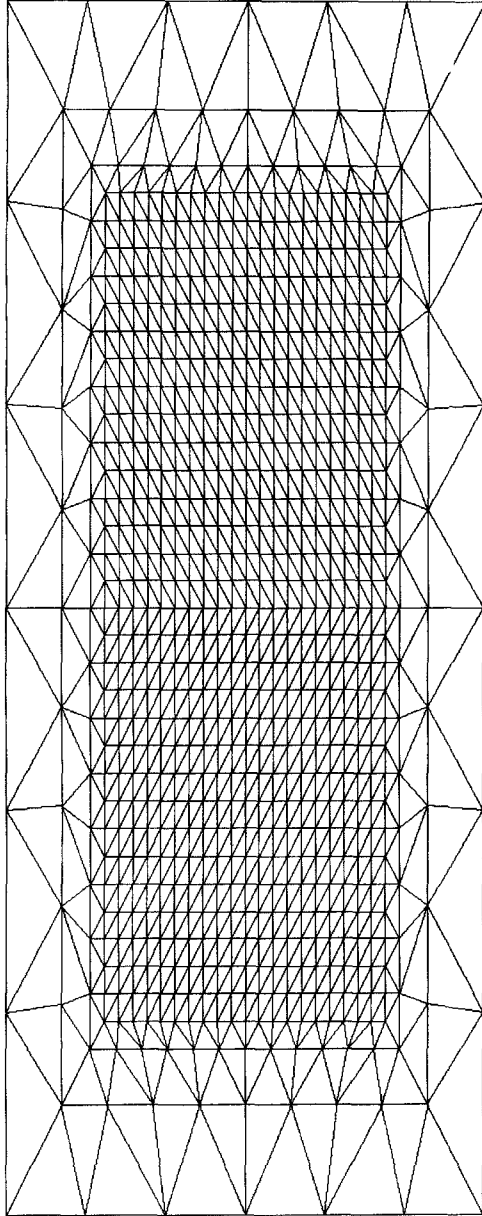


Fig. 3. Finite element mesh with transition layers.

$2M_0$  and the associated eigenvectors, which include the eigenvalues and the eigenvectors of transfer matrix  $X$ . Then we calculate

$$\begin{aligned} \det(R_1 - \lambda R_2) &= \det \begin{pmatrix} K - \lambda A^T & -A \\ I & -\lambda I \end{pmatrix} \\ &= (-1)^n \det(-A + \lambda K - \lambda^2 A^T) \end{aligned} \quad (17)$$

where  $\det$  denotes the determinant of matrix. Because the matrix  $K$  is symmetric, (17) is a symmetric polynomial with respect to  $\lambda$ . If  $\lambda \neq 0$  is a eigenvalue then  $1/\lambda$  will be another one. It can be proved from (7) that all eigenvalues of  $X$  must be not greater than one, so it is easy to obtain the eigenvalues and the associated eigenvalues of  $X$ . Then  $X$  can be obtained from (15).

#### E. Proof of (7)

The scalar potential within the infinite region  $\Omega_{out}$  satisfies the Laplace's equation. Assuming that the scalar potentials  $\Phi_0$  on  $\Gamma_0$  is known, all scalar potentials within  $\Omega_{out}$  can be uniquely determined. That is to say,  $\Phi_1$  can be uniquely determined by  $\Phi_0$ . Because the Laplace's equation and the associated functional are linear,  $\Phi_1$  must be determined linearly by  $\Phi_0$ . Then, there must exist a real matrix which satisfies

$$\Phi_1 = X\Phi_0. \quad (18)$$

Next, we take the linear transform  $x' = \xi x$  and  $y' = \xi y$ , where  $\xi$  is the similar ratio between the layers of  $\Omega_{out}$ . The positions of  $\Gamma_0, \Gamma_1, \dots, \Gamma_k$  now have been shifted to the position of  $\Gamma_1, \Gamma_2, \dots, \Gamma_{k+1}$ , correspondingly. Since the scalar potentials on  $\Omega_{out}$  have the identical values in different coordinate systems, the region within boundary  $\Gamma_0$  and the one within  $\Gamma_1$  have the same solutions if the initial condition on  $\Gamma_0$  and  $\Gamma_1$  are identical. Therefore, if the scalar potentials  $\Phi_1$  is known,  $\Phi_2$  can also be obtained

$$\Phi_2 = X\Phi_1. \quad (19)$$

Continuing to make the same linear transform, we can obtain the (7).

### III. APPLICATION

Consider the two coupled microstrips shown in Fig. 2. Its finite element mesh is shown in Fig. 3. Because the number  $M_0$  of boundary nodes is related to the stiffness matrix bandwidth and the computation time of the transfer matrix, some transition layers has been applied in Fig. 3. The capacitance matrix is calculated by the above method. Table I shows the numerical results obtained by the present method, together with those obtained by using a p.e.c shield, the ABC, the higher order ABC [4], infinite FEM [5], and other method [6]. Table I indicates that the present method can yield more accurate results.

### IV. CONCLUSION

The above mentioned new FEM procedure is useful for two-dimension and three-dimension electromagnetic Laplace's problems with open boundary. Because the whole region is discretized with infinite elements and no appropriate boundary is needed, the calculation error produced by the truncated boundary or absorbing boundary condition is avoided. The numerical results show that the calculation precision is very high.

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## Thru Characteristics of a Coaxial Gap (FDTD Model and Measurements)

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**Abstract**—Thru characteristics of a coaxial cable interrupted by a small gap are modeled and measured. Finite-difference time-domain (FDTD) modeling is applied in cylindrical coordinates to semirigid coaxial cable and to the intervening gap material. Both dispersive and nondispersive gap materials are investigated. Gap loss and phase shift are accurately predicted by this two-dimensional model which accounts for TEM and TM modes in the gap and coaxial apertures. An application of the model is to establish reference data for thin sample permittivity or moisture measurements.

### I. INTRODUCTION

A FDTD model in cylindrical coordinates is presented for a coaxial cable and intervening small gap. That is the cable is severed and a gap of up to one probe diameter (3.58 mm) is opened between the ends. Modeled results are verified with measurements to demonstrate accuracy for several gap sizes and materials. The coaxial gap is proposed as a sensor for very localized permittivity measurements of thin samples as an alternative to the coaxial reflection method which requires stacking of thin samples [1], [2] or the semiempirical approach used in [2] which requires two measurements of the same sample. Neither method is suitable for continuous thin sample moisture measurement. Since FDTD results are not readily invertible the procedure for determination of permittivity would be to characterize the gap response for a number of materials or moisture levels as part of calibration and to then use an interpolation procedure to determine the actual permittivity from the measured values. With sufficient computational speed an iterative solver would be feasible. This paper presents the FDTD numerical model along with measurements of two thin samples for verification. This presentation deals with the transmission response of the gap while in [3] the reflection properties

of the coaxial cable have been considered and found to be accurately predicted by this approach. Now with the reflection and transmission properties well modeled and through the use of [4] one can either iteratively or through a look-up procedure determine the material properties.

### II. NUMERICAL ANALYSIS

A lossless and nondispersive cable dielectric is assumed while the gap material may have loss and dispersive characteristics. Dispersion is accounted for through recursive relations similar to those in [5]. The model is two-dimensional with the center of the center conductor forming a line of symmetry as in Fig. 1. This two-dimensional model accounts for TEM and TM modes at the apertures and throughout the gap but is not capable of modeling TE modes. This is not expected to be a limiting factor since the actual problem has circular symmetry and thus there are no mechanisms to excite TE modes [6]. Due to the small cable size chosen for this study and the need to precisely model gap details a cell size of 0.2 mm is chosen. This yields 52 cells from the inner conductor to the outer conductor in 3.58 mm (0.141") cable or 392 cells per wavelength at 26.5 GHz. With this small element size the Courant condition requires a correspondingly small time step, in the tens of picosecond range, depending upon the dielectric material selected. Field components used in this solution are the axial and radial electric fields and the circumferential magnetic field. Absorbing boundaries are of the first order Mur [7] type applied to the axial electric field while the line of symmetry is accounted for through symmetry of the magnetic field. Governing field equations in cylindrical coordinates are given in pseudo code form below for dispersive materials where the Debye equation and an additional conductivity term are used to describe their frequency dependent behavior as follows

$$\epsilon^*(\omega) = \epsilon_\infty + \frac{\epsilon_S - \epsilon_\infty}{1 + j\omega\tau_0} - j \frac{\sigma}{\omega\epsilon_0}. \quad (1)$$

Where the terms are defined as follows with their corresponding values for water at 25° shown in brackets,  $\epsilon_\infty$  = permittivity at infinite frequency (4.9),  $\epsilon_S$  = static permittivity (78.52),  $\tau_0$  = relaxation time ( $8.38 \times 10^{-12}$ ),  $\sigma$  = conductivity (0.0), and  $\omega$  = angular frequency

$$HC^{n+1/2}(I, J) = HC^{n-1/2}(I, J) + F^*((EA^n(I+1, J) - EA^n(I, J)) - (ER^n(I, J+1) - ER^n(I, J))) \quad (2)$$

$$ER^{n+1}(I, J) = B^*ER^n(I, J) + C^*(HC^{n+1/2}(I, J-1) - HC^{n+1/2}(I, J)) + D^*SR^n(I, J) \quad (3)$$

$$SR^{n+1}(I, J) = A^*SR^n(I, J) + 0.5^*(A^*ER^n(I, J) + ER^{n+1}(I, J)) \quad (4)$$

$$EA^{n+1}(I, J) = B^*EA^n(I, J) + E^*(HC^{n+1/2}(I, J) - HC^{n+1/2}(I-1, J)) + C^*(HC(I, J) - HC(I-1, J)) + D^*SA^n(I, J) \quad (5)$$

$$SA^{n+1}(I, J) = A^*SA^n(I, J) + 0.5^*(A^*EA^n(I, J) + EA^{n+1}(I, J)) \quad (6)$$

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