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**USER'S MANUAL FOR  
THREE DIMENSIONAL FDTD VERSION A  
CODE FOR SCATTERING FROM FREQUENCY-INDEPENDENT  
DIELECTRIC MATERIALS**

by

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## TABLE OF CONTENTS

I.	INTRODUCTION . . . . .	4
II.	FDTD METHOD . . . . .	4
III.	OPERATION . . . . .	5
IV.	RESOURCE REQUIREMENTS . . . . .	7
V.	VERSION A CODE CAPABILITIES . . . . .	8
VI.	DEFAULT SCATTERING GEOMETRY . . . . .	8
VII.	SUBROUTINE DESCRIPTION . . . . .	8
	MAIN ROUTINE . . . . .	9
	SUBROUTINE SETFZ . . . . .	9
	SUBROUTINE SAVFZ . . . . .	9
	SUBROUTINE FAROUT . . . . .	10
	SUBROUTINE BUILD . . . . .	10
	SUBROUTINE DCUBE . . . . .	10
	SUBROUTINE SETUP . . . . .	11
	SUBROUTINE EXSFLD . . . . .	11
	SUBROUTINE EYSFLD . . . . .	11
	SUBROUTINE EZSFLD . . . . .	11
	SUBROUTINES RADEYX, RADEZX, RADEZY, RADEXY, RADEXZ and RADEYZ . . . . .	11
	SUBROUTINE HXSFLD . . . . .	12
	SUBROUTINE HYSFLD . . . . .	12
	SUBROUTINE HZSFLD . . . . .	12
	SUBROUTINE DATSAV . . . . .	12
	FUNCTIONS EXI, EYI and EZI . . . . .	12
	FUNCTION SOURCE . . . . .	12
	FUNCTIONS DEXI, DEYI and DEZI . . . . .	13
	FUNCTION DSRCE . . . . .	13
	SUBROUTINE ZERO . . . . .	13
VIII.	INCLUDE FILE DESCRIPTION (COMMONA.FOR) . . . . .	13
IX.	RCS COMPUTATIONS . . . . .	14
X.	RESULTS . . . . .	14
XI.	SAMPLE PROBLEM SETUP . . . . .	14
XII.	NEW PROBLEM CHECKLIST . . . . .	15
	COMMONA.FOR . . . . .	16
	SUBROUTINE BUILD . . . . .	16
	SUBROUTINE SETUP . . . . .	16
	FUNCTIONS SOURCE and DSRCE . . . . .	17
	SUBROUTINE DATSAV . . . . .	17

## TABLE OF CONTENTS (cont.)

XIII. REFERENCES . . . . .	17
XIV. FIGURE TITLES . . . . .	17

## I. INTRODUCTION

The Penn State Finite Difference Time Domain Electromagnetic Scattering Code Version A is a three dimensional numerical electromagnetic scattering code based upon the Finite Difference Time Domain Technique (FDTD). The supplied version of the code is one version of our current three dimensional FDTD code set. This manual provides a description of the code and corresponding results for the default scattering problem. The manual is organized into fourteen sections: introduction, description of the FDTD method, Operation, resource requirements, Version A code capabilities, a brief description of the default scattering geometry, a brief description of each subroutine, a description of the include file (COMMONA.FOR), a section briefly discussing Radar Cross Section (RCS) computations, a section discussing the scattering results, a sample problem setup section, a new problem checklist, references and figure titles.

## II. FDTD METHOD

The Finite Difference Time Domain (FDTD) technique models transient electromagnetic scattering and interactions with objects of arbitrary shape and/or material composition. The technique was first proposed by Yee [1] for isotropic, non-dispersive materials in 1966; and has matured within the past twenty years into a robust and efficient computational method. The present FDTD technique is capable of transient electromagnetic interactions with objects of arbitrary and complicated geometrical shape and material composition over a large band of frequencies. This technique has recently been extended to include dispersive dielectric materials, chiral materials and plasmas.

In the FDTD method, Maxwell's curl equations are discretized in time and space and all derivatives (temporal and spatial) are approximated by central differences. The electric and magnetic fields are interleaved in space and time and are updated in a second-order accurate leapfrog scheme. The computational space is divided into cells with the electric fields located on the edges and the magnetic fields on the faces (see Figure 1). FDTD objects are defined by specifying dielectric and/or magnetic material parameters at electric and/or magnetic field locations.

Two basic implementations of the FDTD method are widely used for electromagnetic analysis: total field formalism and scattered field formalism. In the total field formalism, the electric and magnetic field are updated based upon the material type present at each spatial location. In the scattered field formalism, the incident waveform is defined analytically and the scattered field is coupled to the incident field through the different material types. For the incident field, any waveform, angle of incidence and polarization is possible. The separation of the incident and

scattered fields conveniently allows an absorbing boundary to be employed at the extremities of the discretized problem space to absorb the scattered fields.

This code is a scattered field code, and the total E and H fields may be found by combining the incident and scattered fields. Any type of field quantity (incident, scattered, or total), Poynting vector or current are available anywhere within the computational space. These fields, incident, scattered and total, may be found within, on or about the interaction object placed in the problem space. By using a near to far field transformation, far fields can be determined from the near fields within the problem space thereby affording radiation patterns and RCS values. The accuracy of these calculations is typically within a dB of analytic solutions for dielectric and magnetic sphere scattering. Further improvements are expected as better absorbing boundary conditions are developed and incorporated.

### III. OPERATION

Typically, a truncated Gaussian incident waveform is used to excite the system being modeled, however certain code versions also provide a smooth cosine waveform for convenience in modeling dispersive materials. The interaction object is defined in the discretized problem space with arrays at each cell location created by the discretization. All three dielectric material types for E field components within a cell can be individually specified by the arrays IDONE(I,J,K), IDTWO(I,J,K), IDTHRE(I,J,K). This models arbitrary dielectric materials with  $\mu = \mu_0$ . By an obvious extension to six arrays, magnetic materials with  $\mu \neq \mu_0$  can be modeled.

Scattering occurs when the incident wave, marched forward in time in small steps set by the Courant stability condition, reaches the interaction object. Here a scattered wave must appear along with the incident wave so that the Maxwell equations are satisfied. If the material is a perfectly conductive metal then only the well known boundary condition

$$E_{\tan}^{\text{scat}} = -E_{\tan}^{\text{inc}} \quad (1)$$

must be satisfied. For a nondispersive dielectric the requirement is that the total field must satisfy the Maxwell equations in the material:

$$\nabla \times E^{\text{tot}} = \nabla \times (E^{\text{inc}} + E^{\text{scat}}) = -\frac{1}{\mu_0} \frac{\partial H^{\text{tot}}}{\partial t} = -\frac{1}{\mu_0} \frac{\partial (H^{\text{inc}} + H^{\text{scat}})}{\partial t} \quad (2)$$

$$\nabla \times \mathbf{H}^{\text{tot}} = \nabla \times (\mathbf{H}^{\text{inc}} + \mathbf{H}^{\text{scat}}) = \epsilon \frac{\partial \mathbf{E}^{\text{tot}}}{\partial t} + \sigma \mathbf{E}^{\text{tot}} \quad (3)$$

$$= \epsilon \frac{\partial (\mathbf{E}^{\text{inc}} + \mathbf{E}^{\text{scat}})}{\partial t} + \sigma (\mathbf{E}^{\text{inc}} + \mathbf{E}^{\text{scat}}) \quad (4)$$

Additionally the incident wave, defined as moving unimpeded through a vacuum in the problem space, satisfies everywhere in the problem the Maxwell equations for free space

$$\nabla \times \mathbf{E}^{\text{inc}} = -\frac{1}{\mu_0} \frac{\partial \mathbf{H}^{\text{inc}}}{\partial t} \quad (5)$$

$$\nabla \times \mathbf{H}^{\text{inc}} = \epsilon_0 \frac{\partial \mathbf{E}^{\text{inc}}}{\partial t} \quad (6)$$

Subtracting the second set of equations from the first yields the Maxwell equations governing the scattered fields in the material:

$$\nabla \times \mathbf{E}^{\text{scat}} = -\frac{1}{\mu_0} \frac{\partial \mathbf{H}^{\text{scat}}}{\partial t} \quad (7)$$

$$\nabla \times \mathbf{H}^{\text{scat}} = (\epsilon - \epsilon_0) \frac{\partial \mathbf{E}^{\text{inc}}}{\partial t} + \sigma \mathbf{E}^{\text{inc}} + \epsilon \frac{\partial \mathbf{E}^{\text{scat}}}{\partial t} + \sigma \mathbf{E}^{\text{scat}} \quad (8)$$

Outside the material this simplifies to:

$$\nabla \times \mathbf{E}^{\text{scat}} = -\frac{1}{\mu_0} \frac{\partial \mathbf{H}^{\text{scat}}}{\partial t} \quad (9)$$

$$\nabla \times \mathbf{H}^{\text{scat}} = \epsilon_0 \frac{\partial \mathbf{E}^{\text{scat}}}{\partial t} \quad (10)$$

Magnetic materials, dispersive effects, non-linearities, etc., are further generalizations of the above approach. Based on the value of the material type, the subroutines for calculating scattered E and H field components branch to the appropriate expression for that scattered field component and

that component is advanced in time according to the selected algorithm. As many materials can be modeled as desired, the number equals the dimension selected for the flags. If materials with behavior different from those described above must be modeled, then after the appropriate algorithm is found, the code's branching structure allows easy incorporation of the new behavior.

#### IV. RESOURCE REQUIREMENTS

The number of cells the problem space is divided into times the six components per cell set the problem space storage requirements

$$\text{Storage} = NC \times 6 \text{ components/cell} \times 4 \text{ bytes/component} \quad (11)$$

and the computational cost

$$\text{Operations} = NC \times 6 \text{ comp/cell} \times 10 \text{ ops/component} \times N \quad (12)$$

where  $N$  is the number of time steps desired.

$N$  typically is on the order of ten times the number of cells on one side of the problem space. More precisely for cubical cells it takes  $\sqrt{3}$  time steps to traverse a single cell when the time step is set by the Courant stability condition

$$\Delta t = \frac{\Delta x}{\sqrt{3}c} \quad \Delta x = \text{cell size dimension} \quad (13)$$

The condition on  $N$  is then that

$$N = 10 \times (\sqrt{3}NC)^{\frac{1}{3}} \quad NC^{\frac{1}{3}} \sim \begin{array}{l} \text{number cells on a side} \\ \text{of the problem space} \end{array} \quad (14)$$

The earliest aircraft modeling using FDTD with approximately 30 cells on a side required approximately 500 time steps. For more recent modeling with approximately 100 cells on a side, 2000 or more time steps are used.

For  $(100 \text{ cell})^3$  problem spaces, 24 MBytes of memory are required to store the fields. Problems on the order of this size have been run on a Silicon Graphics 4D 220 with 32 MBytes of memory, IBM RISC 6000, an Intel 486 based machine, and VAX 11/785. Storage is only a problem as in the case of the 486 where only 16 MBytes of memory was available. This limited the problem space size to approximately  $(80 \text{ cells})^3$ .

For  $(100 \text{ cell})^3$  problems with approximately 2000 time steps, there is a total of  $120 \times 10^9$  operations to perform. The speeds of the previously mentioned machines are 24 MFLOPs (4 processor upgraded version), 10 MFLOPs, 1.5 MFLOPs, and 0.2 MFLOPs. The run times are then  $5 \times 10^3$  seconds,  $12 \times 10^3$  seconds,  $80 \times 10^3$  seconds and  $600 \times 10^3$  seconds, respectively. In hours the times are 1.4, 3.3, 22.2 and 167 hours. Problems of this size are possible on all but the last machine and can in fact be performed on a personal computer (486) if one day turnarounds are permissible.

## V. VERSION A CODE CAPABILITIES

The Penn State University FDTD Electromagnetic Scattering Code Version A has the following capabilities:

- 1) Ability to model lossy dielectric and perfectly conducting scatterers.
- 2) First and second order outer radiation boundary condition (ORBC) operating on the electric fields for dielectric or perfectly conducting scatterers.
- 3) Near to far zone transformation capability to obtain far zone scattered fields.
- 4) Gaussian and smooth cosine incident waveforms with arbitrary incidence angles.
- 5) Near zone field, current or power sampling capability.
- 6) Companion code for computing Radar Cross Section (RCS).

## VI. DEFAULT SCATTERING GEOMETRY

The code as delivered is set up to calculate the far zone backscatter fields for an infinitely thin, 29 cm square, perfectly conducting plate. The problem space size is 60 by 60 by 49 cells in the x, y and z directions, the cells are 1 cm cubes, and the incident waveform is a  $\phi$ -polarized Gaussian pulse with incidence angles of  $\theta=45$  and  $\phi=30$  degrees. The output data files are included as a reference along with a code (RCS3D.FOR) for computing the frequency domain RCS using these output data files. The ORBC is the second order absorbing boundary condition set forth by Mur [2].

## VII. SUBROUTINE DESCRIPTION

In the description for each subroutine, an asterisk (\*) will be placed by the subroutine name if that particular subroutine is normally modified when defining a scattering problem.



## MAIN ROUTINE

The main routine in the program contains the calls for all necessary subroutines to initialize the problem space and scattering object(s) and for the incident waveform, far zone transformation, field update subroutines, outer radiation boundary conditions and field sampling.

The main routine begins with the include statement and then appropriate data files are opened, and subroutines ZERO, BUILD and SETUP are called to initialize variables and/or arrays, build the object(s) and initialize the incident waveform and miscellaneous parameters, respectively. Subroutine SETFZ is called to initialize parameters for the near to far zone transformation if far zone fields are desired.

The main loop is entered next, where all of the primary field computations and data saving takes place. During each time step cycle, the EXSFLD, EYSFLD, and EZSFLD subroutines are called to update the x, y, and z components of the scattered electric field. The six electric field outer radiation boundary conditions (RADE??) are called next to absorb any outgoing scattered fields. Time is then advanced 1/2 time step according to the Yee algorithm and then the HXSFLD, HYSFLD, AND HZSFLD subroutines are called to update the x, y, and z components of scattered magnetic field. Time is then advanced another 1/2 step and then either near zone fields are sampled and written to disk in DATSAV, and/or the near zone to far zone vector potentials are updated in SAVFZ. The parameter NZFZ (described later) in the common file defines the type of output fields desired.

After execution of all time steps in the main field update loop, subroutine FAROUT is called if far zone fields are desired to compute the far zone fields and write them to disk. At this point, the execution is complete.

## SUBROUTINE SETFZ

This subroutine initializes the necessary parameters required for far zone field computations. The code as furnished computes backscatter far zone fields and can compute bistatic far zone fields for one scattering angle (i.e. one  $\theta$  and  $\phi$  angle). Refer to reference [3] for a complete description of the near to far zone transformation. Other versions of this subroutine provide for multiple bistatic angles.

## SUBROUTINE SAVFZ

This subroutine updates the near zone to far zone vector potentials.

## SUBROUTINE FAROUT

This subroutine changes the near zone to far zone vector potentials to far zone electric field  $\theta$  and  $\phi$  components and writes them to disk.

## SUBROUTINE BUILD \*

This subroutine "builds" the scattering object(s) by initializing the IDONE, IDTWO, and IDTHRE arrays. The IDONE-IDTHRE arrays are for specifying perfectly conducting and lossy dielectric materials. Refer to Figure 1 for a diagram of the basic Yee cell. For example, setting an element of the IDONE array at some I,J,K location is actually locating dielectric material at a cell edge whose center location is I+0.5,J,K. Thus, materials with diagonal permittivity tensors can be modeled. The default material type for all ID??? arrays is 0, or free space. By initializing these arrays to values other than 0, the user is defining an object by determining what material types are present at each spatial location. Other material types available for IDONE-IDTHRE are 1 for perfectly conducting objects and 2-9 for lossy non-magnetic dielectrics. It is assumed throughout the code that all dielectric materials are non-magnetic (i.e. the materials have a permeability of  $\mu_0$ ). This subroutine also has a section that checks the ID??? arrays to determine if legal material types have been defined throughout the problem space. The actual material parameters ( $\epsilon$  and  $\sigma$ ) are defined in subroutine SETUP. The default geometry is a 29 cm square perfectly conducting plate.

The user must be careful that his/her object created in the BUILD subroutine is properly formed.

When it is important to place the object in the center of the problem space (to have lowest possible cross-pol scattering for symmetric objects), NX etc. should be odd. This is due to the field locations in the Yee cell and also the placement of the E field absorbing boundary condition surfaces.

If the object being modeled has curved surfaces, edges, etc. that are at an angle to one or more of the coordinate axes, then that shape must be approximately modeled by lines and faces in a "stair-stepped" (or stair-cased) fashion. This stair-cased approximation introduces errors into computations at higher frequencies. Intuitively, the error becomes smaller as more cells are used to stair-case a particular object.

## SUBROUTINE DCUBE

This subroutine builds cubes of dielectric material by defining four each of IDONE, IDTWO and IDTHRE components corresponding to one spatial cube of dielectric material. It can

also be used to define thin (i.e. up to one cell thick) dielectric or perfectly conducting plates. Refer to comments within DCUBE for a description of the arguments and usage of the subroutine.

#### SUBROUTINE SETUP \*

This subroutine initializes many of the constants required for incident field definition, field update equations, outer radiation boundary conditions and material parameters. The material parameters  $\epsilon$  and  $\sigma$  are defined for each material type using the material arrays EPS and SIGMA respectively. The array EPS is used for the total permittivity and SIGMA is used for the electric conductivity. These arrays are initialized in SETUP to free space material parameters for all material types and then the user is required to modify these arrays for his/her scattering materials. Thus, for the lossy dielectric material type 2, the user must define EPS(2) and SIGMA(2). The remainder of the subroutine computes constants used in field update equations and boundary conditions and writes the diagnostics file.

#### SUBROUTINE EXSFLD

This subroutine updates all x components of scattered electric field at each time step except those on the outer boundaries of the problem space. IF statements based upon the IDONE array are used to determine the type of material present and the corresponding update equation to be used. These scattered field equations are based upon the development given in [4].

#### SUBROUTINE EYSFLD

This subroutine updates all y components of scattered electric field at each time step except those on the outer boundaries of the problem space. IF statements based upon the IDTWO array are used to determine the type of material present and the corresponding update equation to be used.

#### SUBROUTINE EZSFLD

This subroutine updates all z components of scattered electric field at each time step except those on the outer boundaries of the problem space. IF statements based upon the IDTHRE array are used to determine the type of material present and the corresponding update equation to be used.

#### SUBROUTINES RADEYX, RADEZX, RADEZY, RADEXY, RADEXZ and RADEYZ

These subroutines apply the outer radiation boundary conditions to the scattered electric field on the outer

boundaries of the problem space.

#### SUBROUTINE HXSFLD

This subroutine updates all x components of scattered magnetic field at each time step. The standard non-magnetic update equation is used.

#### SUBROUTINE HYSFLD

This subroutine updates all y components of scattered magnetic field at each time step. The standard non-magnetic update equation is used.

#### SUBROUTINE HZSFLD

This subroutine updates all z components of scattered magnetic field at each time step. The standard non-magnetic update equation is used.

#### SUBROUTINE DATSAV \*

This subroutine samples near zone scattered field quantities and saves them to disk. This subroutine is where the quantities to be sampled and their spatial locations are to be specified and is only called if near zone fields only are desired or if both near and far zone fields are desired. Total field quantities can also be sampled. See comments within the subroutine for specifying sampled scattered and/or total field quantities. When sampling magnetic fields, remember the  $\delta t/2$  time difference between E and H when writing the fields to disk. Sections of code within this subroutine determine if the sampled quantities and the spatial locations have been properly defined.

#### FUNCTIONS EXI, EYI and EZI

These functions are called to compute the x, y and z components of incident electric field. The functional form of the incident field is contained in a separate function SOURCE.

#### FUNCTION SOURCE \*

This function contains the functional form of the incident field. The code as furnished uses the Gaussian form of the incident field. An incident smooth cosine pulse is also available by uncommenting the required lines and commenting out the Gaussian pulse. Thus, this function need only be modified if the user changes the incident pulse from Gaussian to smooth cosine. A slight improvement in computing speed and vectorization may be achieved by moving this function inside each of the incident field functions EXI, EYI and so on.

## FUNCTIONS DEXI, DEYI and DEZI

These functions are called to compute the x, y and z components of the time derivative of incident electric field. The functional form of the incident field is contained in a separate function DSRCE.

## FUNCTION DSRCE \*

This function contains the functional form of the time derivative of the incident field. The code as furnished uses the time derivative of the Gaussian form of the incident field. A smooth cosine pulse time derivative is also available by uncommenting the required lines and commenting out the Gaussian pulse. Thus, the function need only be modified if the user changes from the Gaussian to smooth cosine pulse. Again, a slight improvement in computing speed and vectorization may be achieved by moving this function inside each of the time derivative incident field functions DEXI, DEYI and so on.

## SUBROUTINE ZERO

This subroutine initializes various arrays and variables to zero.

## VIII. INCLUDE FILE DESCRIPTION (COMMONA.FOR) \*

The include file, COMMONA.FOR, contains all of the arrays and variables that are shared among the different subroutines. This file will require the most modifications when defining scattering problems. A description of the parameters that are normally modified follows.

The parameters NX, NY and NZ specify the size of the problem space in cells in the x, y and z directions respectively. For problems where it is crucial to center the object within the problem space, then NX, NY and NZ should be odd. The parameter NTEST defines the number of near zone quantities to be sampled and NZFZ defines the field output format. Set NZFZ=0 for near zone fields only, NZFZ=1 for far zone fields only and NZFZ=2 for both near and far zone fields. Parameter NSTOP defines the maximum number of time steps. DELX, DELY, and DELZ (in meters) define the cell size in the x, y and z directions respectively. The  $\theta$  and  $\phi$  incidence angles (in degrees) are defined by THINC and PHINC respectively and the polarization is defined by ETHINC and EPHINC. ETHINC=1.0, EPHINC=0.0 for  $\theta$ -polarized incident field and ETHINC=0.0, EPHINC=1.0 for  $\phi$ -polarized incident fields. Parameters AMP and BETA define the maximum amplitude and the  $e^2$  temporal width of the incident pulse respectively. BETA automatically adjusts when the cell size is changed and normally should not be changed by the user. The far zone scattering angles are defined by THETFZ and PHIFZ. The code as furnished

performs backscatter computations, but these parameters could be modified for a bistatic computation.

## IX. RCS COMPUTATIONS

A companion code, RCS3D.FOR, has been included to compute RCS versus frequency. It uses the file name of the FDTD far zone output data (FZOUT3D.DAT) and writes a data file of far zone electric fields versus time (FZTIME.DAT) and RCS versus frequency (3DRCS.DAT). The RCS computations are performed up to the  $10 \text{ cell}/\lambda_0$  frequency limit. Refer to comments within this code for further details.

## X. RESULTS

As previously mentioned, the code as furnished models an infinitely thin, 29 cm square, perfectly conducting plate and computes backscatter far zone scattered fields at angles of  $\theta=45$  and  $\phi=30$  degrees.

Figures 2-3 shows the co-polarized far zone electric field versus time and the co-polarized RCS for the 29 cm square perfectly conducting plate.

Figures 4-5 shows the cross-polarized far zone electric field versus time and the cross-polarized RCS for the 29 cm square perfectly conducting plate.

## XI. SAMPLE PROBLEM SETUP

The code as furnished models an infinitely thin, 29 cm square, perfectly conducting plate and computes backscatter far zone scattered fields at angles of  $\theta=45$  and  $\phi=30$  degrees. The corresponding output data files are also provided, along with a code to compute RCS using these data files. In order to change the code to a new problem, many different parameters need to be modified. A sample problem setup will now be discussed.

Suppose that the problem to be studied is RCS backscatter versus frequency from a 28 cm by 31 cm perfectly conducting plate with a 3 cm dielectric coating with a dielectric constant of  $4\epsilon_0$  using a  $\theta$ -polarized field. The backscatter angles are  $\theta=30.0$  and  $\phi=60.0$  degrees and the frequency range is up to 3 GHz.

Since the frequency range is up to 3 GHz, the cell size must be chosen appropriately to resolve the field IN ANY MATERIAL at the highest frequency of interest. A general rule is that the cell size should be  $1/10$  of the wavelength at the highest frequency of interest. For difficult geometries,  $1/20$  of a wavelength may be necessary. The free space wavelength at 3 GHz is  $\lambda_0=10$  cm and the wavelength in the dielectric coating at 3 GHz is 5 cm. The cell size is chosen as 1 cm, which provides a

resolution of 5 cells/ $\lambda$  in the dielectric coating and 10 cells/ $\lambda_0$  in free space. Numerical studies have shown that choosing the cell size  $\leq 1/4$  of the shortest wavelength in any material is the practical lower limit. Thus the cell size of 1 cm is barely adequate. The cell size in the x, y and z directions is set in the common file through variables DELX, DELY and DELZ. Next the problem space size must be large enough to accommodate the scattering object, plus at least a five cell boundary (10 cells is more appropriate) on every side of the object to allow for the far zone field integration surface. It is advisable for plate scattering to have the plate centered in the x and y directions of the problem space in order to reduce the cross-polarized backscatter and to position the plate low in the z direction to allow strong specular reflections multiple encounters with the ORBC. A 10 cell border is chosen, and the problem space size is chosen as 49 by 52 by 49 cells in the x, y and z directions respectively. As an initial estimate, allow 2048 time steps so that energy trapped within the dielectric layer will radiate. Thus parameters NX, NY and NZ in COMMONA.FOR would be changed to reflect the new problem space size, and parameter NSTOP is changed to 2048. If all transients have not been dissipated after 2048 time steps, then NSTOP will have to be increased. Truncating the time record before all transients have dissipated will corrupt frequency domain results. Parameter NZFZ must be equal to 1 since we are interested in far zone fields only. To build the object, the following lines are inserted into the BUILD subroutine:

```

C
C      BUILD THE DIELECTRIC SLAB FIRST
C
      ISTART=11
      JSTART=11
      KSTART=11
      NXWIDE=28
      NYWIDE=31
      NZWIDE=3
      MTYPE=2
      CALL DCUBE(ISTART,JSTART,KSTART,NXWIDE,NYWIDE,NZWIDE,MTYPE)
C
C      BUILD PEC PLATE NEXT
C
      ISTART=11
      JSTART=11
      KSTART=11
      NXWIDE=28
      NYWIDE=31
      NZWIDE=0
      MTYPE=1
      CALL DCUBE(ISTART,JSTART,KSTART,NXWIDE,NYWIDE,NZWIDE,MTYPE)

```

The PEC plate is built last on the bottom of the dielectric slab to avoid any air gaps between the dielectric material and the PEC plate. In the common file, the incidence angles THINC and PHINC have to be changed to 30.0 and 60.0 respectively, the cell sizes (DELX, DELY, DELZ) are set to 0.01, and the polarization is set to ETHINC=1.0 and EPHINC=0.0 for  $\theta$ -polarized fields. Since dielectric material 2 is being used for the dielectric coating, the constitutive parameters EPS(2) and SIGMA(2) are set to  $4\epsilon_0$  and 0.0 respectively, in subroutine SETUP. This completes the code modifications for the sample problem.

## XII. NEW PROBLEM CHECKLIST

This checklist provides a quick reference to determine if all parameters have been defined properly for a given scattering problem. A reminder when defining quantities within the code: use MKS units and specify all angles in degrees.

### COMMONA.FOR:

- 1) Is the problem space sized correctly? (NX, NY, NZ)
- 2) For near zone fields, is the number of sample points correct? (NTEST)
- 3) Is parameter NZFZ defined correctly for desired field outputs?
- 4) Is the number of time steps correct? (NSTOP)
- 5) Are the cell dimensions (DELX, DELY, DELZ) defined correctly?
- 6) Are the incidence angles (THINC, PHINC) defined correctly?
- 7) Is the polarization of the incident wave defined correctly (ETHINC, EPHINC)?
- 8) For other than backscatter far zone field computations, are the scattering angles set correctly? (THETFZ, PHIFZ)

### SUBROUTINE BUILD:

- 1) Is the object completely and correctly specified?

### SUBROUTINE SETUP:

- 1) Are the constitutive parameters for each material specified correctly? (EPS and SIGMA)



**FUNCTIONS SOURCE and DSRCE:**

- 1) If the Gaussian pulse is not desired, is it commented out and the smooth cosine pulse uncommented?

**SUBROUTINE DATSAV:**

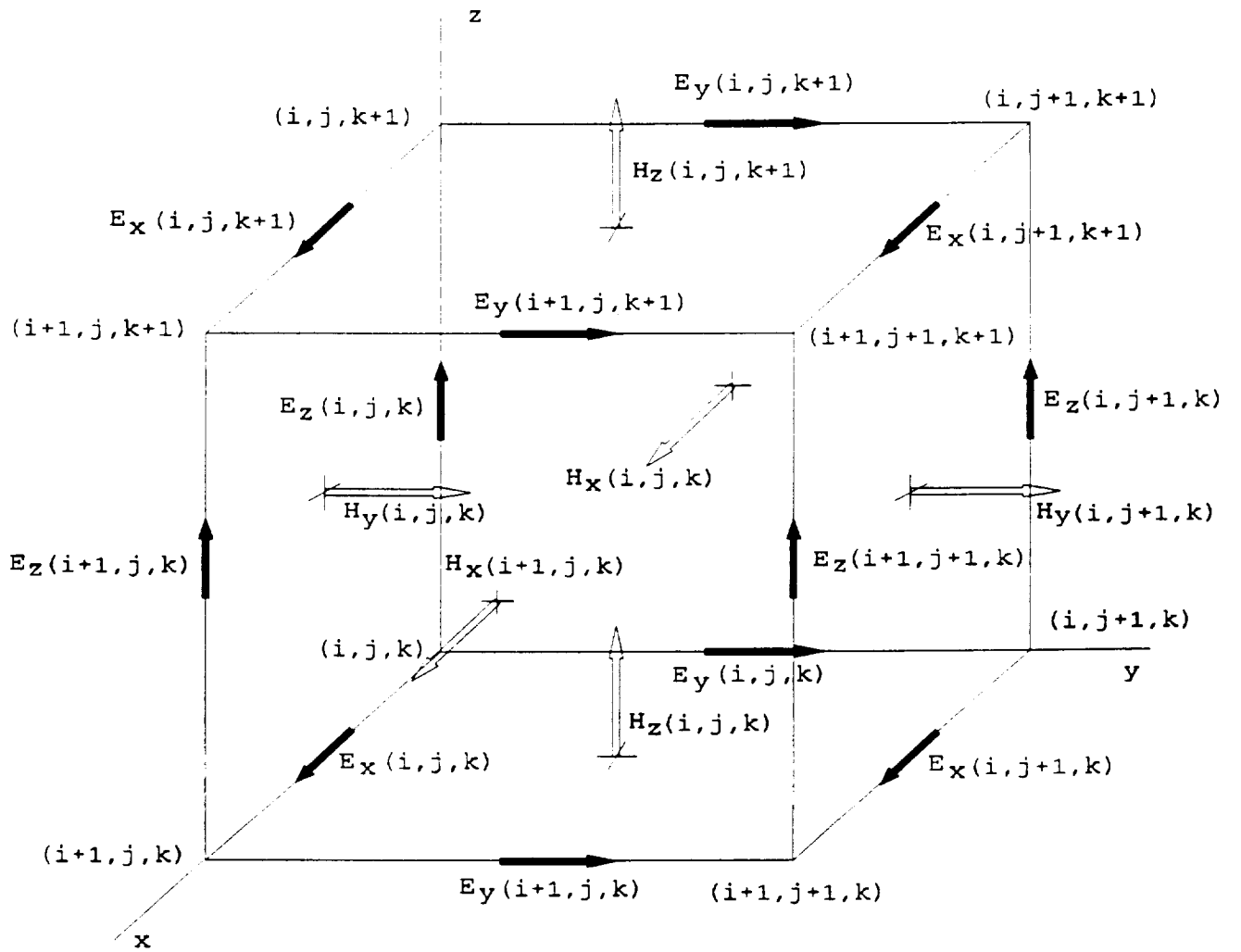
- 1) For near zone fields, are the sampled field types and spatial locations correct for each sampling point? (NTYPE, IOBS, JOBS, KOBS)

**XIII. REFERENCES**

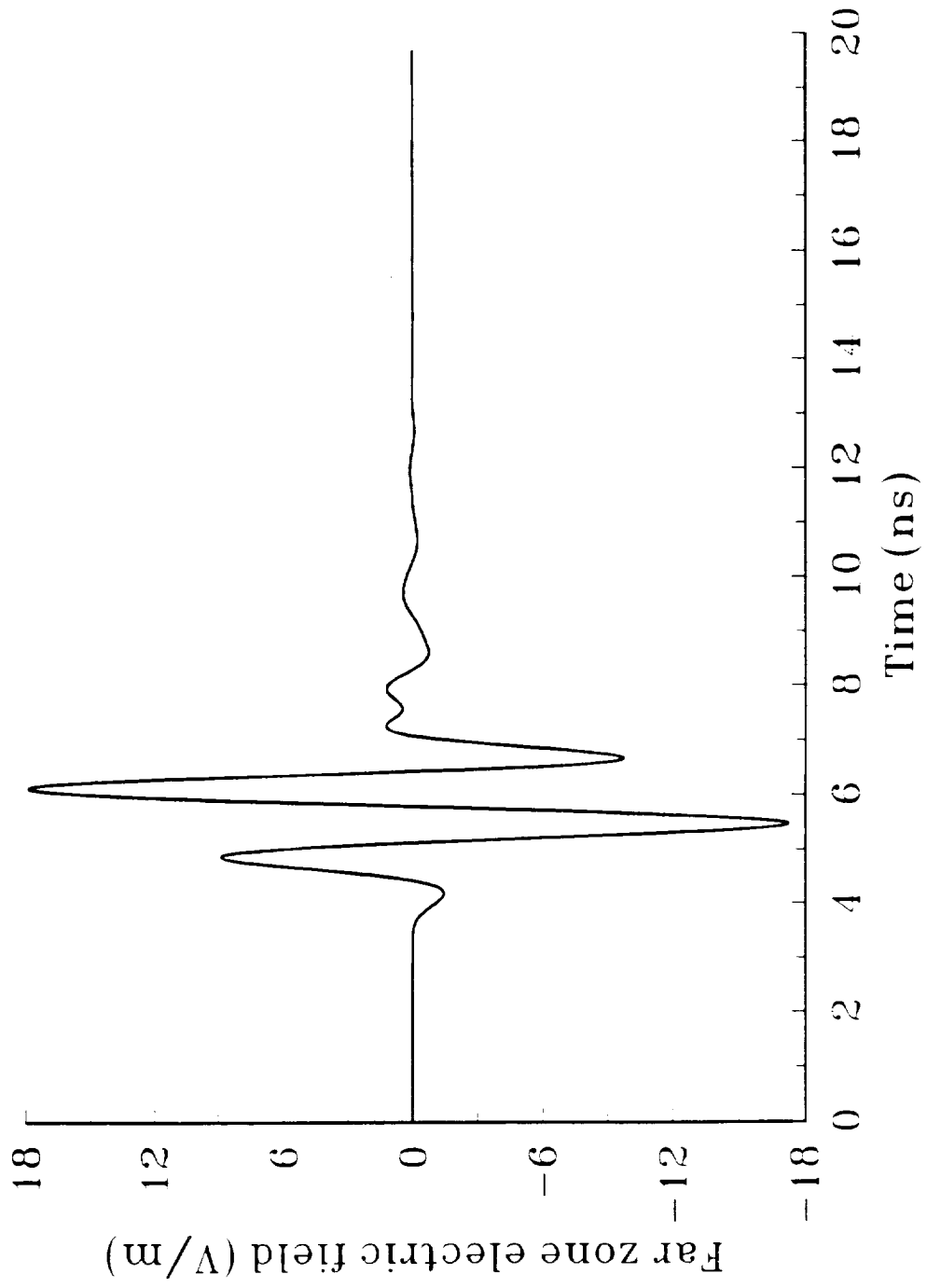
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**XIV. FIGURE TITLES**

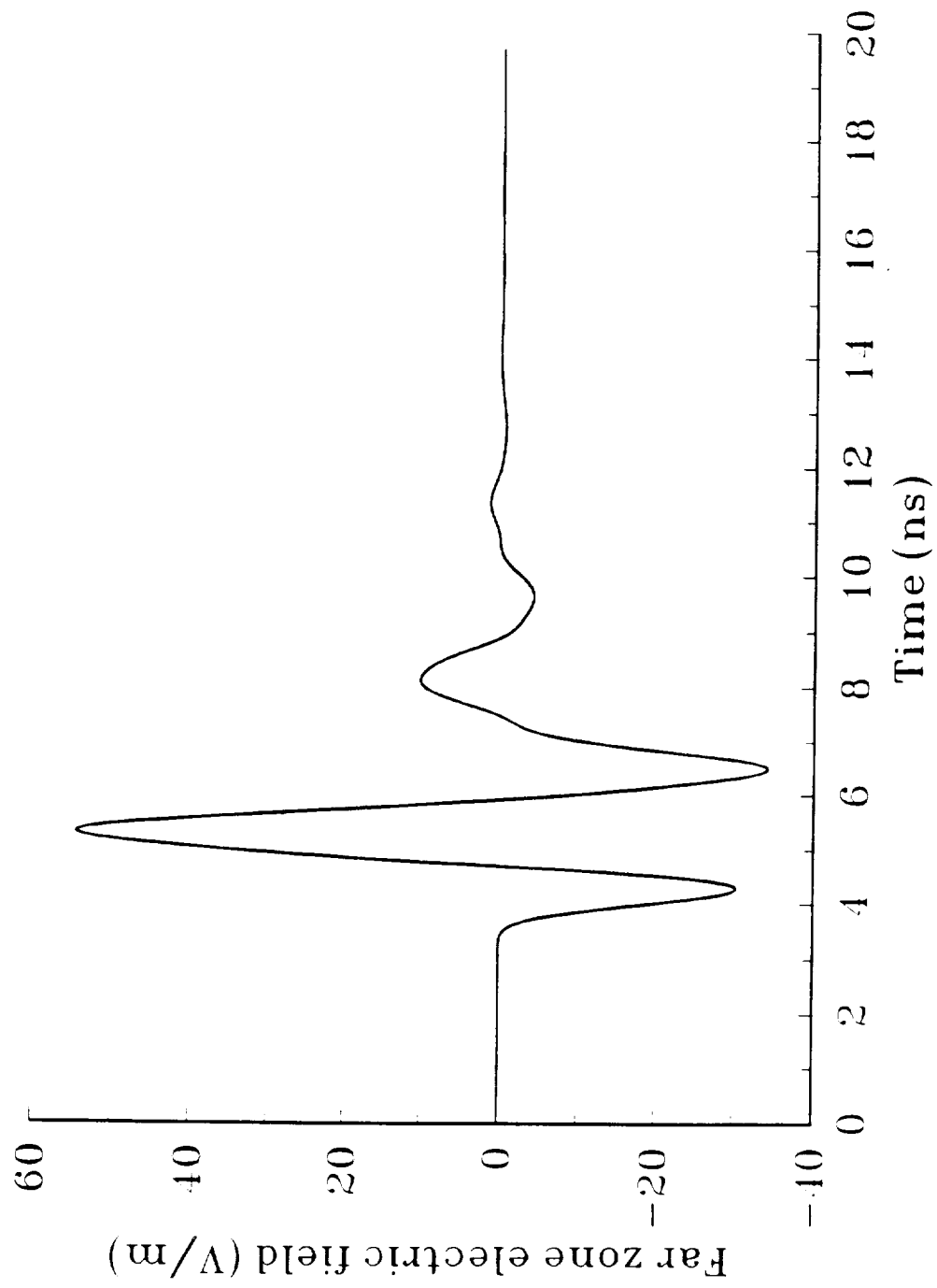
- Fig. 1 Standard three dimensional Yee cell showing placement of electric and magnetic fields.
- Fig. 2 Co-polarized far zone scattered field versus time for 29 cm square perfectly conducting plate.
- Fig. 3 Co-polarized RCS versus frequency for 29 cm square perfectly conducting plate.
- Fig. 4 Cross-polarized far zone scattered field versus time for 29 cm square perfectly conducting plate.
- Fig. 5 Cross-polarized RCS versus frequency for 29 cm square perfectly conducting plate.



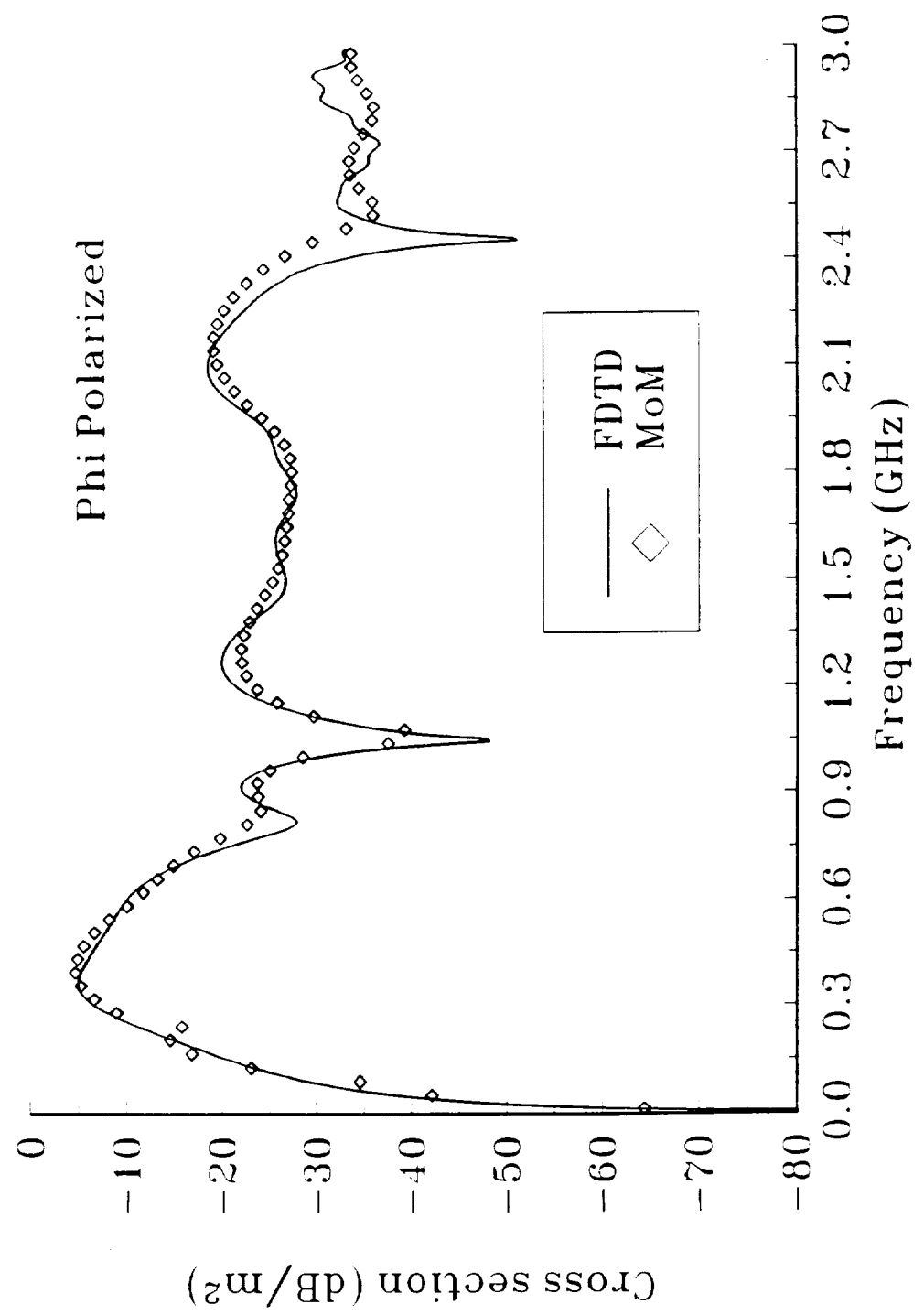
Cross-polarized far zone scattered field  
29 cm square PEC plate;  $\vartheta=45^\circ$ ,  $\varphi=30^\circ$



Co-polarized far zone scattered field  
29 cm square PEC plate;  $\vartheta=45^\circ$ ,  $\varphi=30^\circ$  degrees



Co-polarized Radar Cross Section  
29 cm square PEC plate;  $\Theta=45^\circ$ ,  $\varphi=30^\circ$  degrees



Cross-polarized Radar Cross Section  
29 cm square PEC plate;  $\theta=45^\circ$ ,  $\varphi=30^\circ$  degrees

