

**173366**

MIDDLE EAST TECHNICAL UNIVERSITY  
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**173366**

NUMERICAL AND EXPERIMENTAL INVESTIGATION OF THE  
SCATTERING OF ELECTROMAGNETIC WAVES BY PERFECTLY  
CONDUCTING ENSEMBLES

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by

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Approved by :

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## ABSTRACT

In this thesis, the problem of electromagnetic scattering from an ensemble of perfectly conducting smooth bodies is investigated numerically and experimentally. Numerical part of the work done is based on a theory which employs multipole expansion techniques.

The scattering problem is considered as a problem of radiation from the induced surface sources. The re-radiated fields are expanded in terms of spherical vector wave functions. The unknown multipole coefficients are determined from the linear relations obtained from the properties of the spherical vector wave functions, expansion of the incident field and surface current relations.

Numerical techniques are developed for use of a general ensemble of arbitrary bodies. Due to difficulties arising from the limitations of the available computer, toroidally symmetric ensembles are investigated using a similar technique to the one used for the general problem.

Back-scattering cross-sections are measured by a method based on comparison of the scattering from standard scatterers (i.e. spheres) and the ensembles. A satisfactory agreement is obtained with numerical results.

LIST OF PRINCIPLE SYMBOLS

$\bar{a}, \bar{u}$	Base Vectors
$\alpha E, \alpha M, X, Y$	Multipole Coefficients
$\delta_{ij}$	Kronecker's Delta
$\Delta, \gamma, \xi, \eta$	Indical Functions
$G_0$	Free Space Green's Function
$\bar{H}$	Magnetic Field
$H_j^{(l)}(r)$	Hankel Functions (Spherical)
$h$	Scale Factors
$\bar{I}$	Identity Dyadic
$JE, IM, CE, CM$	Scattering Matrix Submatrices
$J_k(r)$	Spherical Bessel's Function
$k$	Propagation Constant
$L$	Lower Triangle Matrix
$P_k^m(\cos\theta)$	Associate Legendre Functions
$P$	Permutation Matrices
$Q_k^m(\theta)$	Associate Legendre Functions Divided by $\sin\theta$
$(r, \theta, \phi), (\rho, \beta, \varphi), (x, y, z)$	Coordinate Variables
$S$	Transformation Matrix
$S$	Scattering Matrix
$\sigma(\theta, \varphi)$	Re-radiation Diagram
$\sigma_b$	Back Scattering Cross-section
$U$	Upper Triangle Matrix
$U$	Excitation Matrix
$\bar{X}(\theta, \varphi)$	Angular Vector Function
$Z(r)$	Radial Function

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## 1. INTRODUCTION

### 1.1. DESCRIPTION OF THE PROBLEM

Electromagnetic fields, produced by time varying current and charge distributions localized in some region of an unbounded homogenous medium, can be calculated by the inhomogeneous wave equation. If perfectly conducting ensembles are present in the free space subject to illuminating electromagnetic waves, total wave distribution can be separated into two as the incident and the re-radiated fields. Re-radiated fields may be thought as if a radiation due to induced surface sources. The problem is, given the incident wave and the obstacle, to find the re-radiated fields. The incident wave may be thought as a collection of monochromatic waves, in particular plane waves.

### 1.2. A BRIEF SURVEY OF THE PREVIOUS TECHNIQUES AND THE METHOD USED

In classical scattering problems, generally, the shapes of the bodies of concern are taken to be coincident with some coordinate system, thus the method of separation of variables is easily applicable. By this method, only single bodies were investigated. The incident and the scattered fields are expanded in terms of the vector wave functions and the unknown coefficients are determined applying boundary conditions for the surfaces.

For instance, in the case of a sphere, the scattering coefficients can be found in closed form [1], [2].

An exact formulation by means of an integral equation was given by Maue and a derivation of this equation is given by Van Bladel [2]. For a three dimensional problem, integral equation solution by ordinary numerical methods is hard. Too many grid points are necessary for this and a convergent iterative scheme has to be developed. For every excitation this scheme must be repeated from the beginning.

Another method is to approximate the surface currents by the first N terms of an infinite series, thus it is possible to obtain N linear algebraic equations. This method was used by Mei and Bladel [3].

Waterman has developed a theory based on the Rayleigh [4] hypothesis, which is approximate for shapes other than a sphere.

For multibodies, the most significant research is by John H. Bruning and Yuen T. Lu [5]. They made theoretical, numerical and experimental investigations on the problem of two spheres illuminated by a plane wave. A method of shifting has been used for the vector wave functions. Their experiments are also worth mentioning. With the equipment of the Radiation Laboratory, Illinois, they have made perfect measurements on every case of two spheres.

The only significant research on toroidal bodies was made for thin wire loops, by Angelakos [6].

Victor Twersky [7] proposed a method for the general scattering problem which makes use of the isolated scatterer values to find multipole coefficients. This paper is purely theoretical, and in practice no numerical research is made.

The method proposed by Altunkan Hızal [8] in 1970 is employed in this research with a slight modification. The method and the modification is given in Chapter 2 . Though he has proposed the theory, the obtained results were for isolated rotational bodies illuminated with a plane wave in the direction of the axis of revolution.

In this thesis an introduction to numerical analysis of the arbitrary multibody problem is made and toroidal and rotational configurations are investigated completely.

## 2. FORMULATION OF THE PROBLEM

### 2.1. ORIGIN EXCLUSIVE TYPE OF FORMULATION

It is assumed that a finite number of smooth,conducting scatterers are existing within the limits  $r_o < r < r_i$  with respect to some coordinate frame. For  $r > r_i$  and  $r < r_o$  the incident and the scattered fields are investigated separately, but for the bulky region,  $r_o < r < r_i$ , the total field is considered(Fig.2.1.1).

Since the bodies are assumed to be perfect conductors , re-radiated field is given by the integral equation

$$\bar{H}_{sc}(\bar{r}) = \nabla \times \oint_S \bar{n} \times (\bar{H}_{in}(\bar{r}') + \bar{H}_{sc}(\bar{r}')) G_o(\bar{r}|\bar{r}') dS' \quad (2.1.1)$$

where S is the total surface of the bodies,  $\bar{H}_{sc}(\bar{r})$  is the re-radiated field at  $\bar{r}$ ,  $\bar{H}_{in}(\bar{r})$  is the incident field(its sources are assumed to be free from interactions),  $G_o(\bar{r},\bar{r}')$  is the free space Green's function which can be expressed as

$$G_o(\bar{r}|\bar{r}') = \frac{e^{ik|\bar{r}-\bar{r}'|}}{4\pi |\bar{r}-\bar{r}'|} \quad (2.1.2)$$

where  $k = \omega/c = 2\pi/\lambda$  , is the propagation coefficient.

Green's function can be treated as a dyadic since  $\bar{F} G_o = \bar{F} \cdot G_o \bar{I}$ . This dyadic can be expanded in terms of vector wave functions as follows [8]

For  $r > r'$

$$G_o(\bar{r}|\bar{r}') \bar{I} = \sum_{P,m} ik \left\{ \bar{M}_{P,m}^{1*}(\bar{r}') \bar{M}_{P,m}^s(\bar{r}) + \frac{1}{k^2} \bar{N}_{P,m}^{1*}(\bar{r}') \bar{N}_{P,m}^s(\bar{r}) \right\} \quad (2.1.3)$$

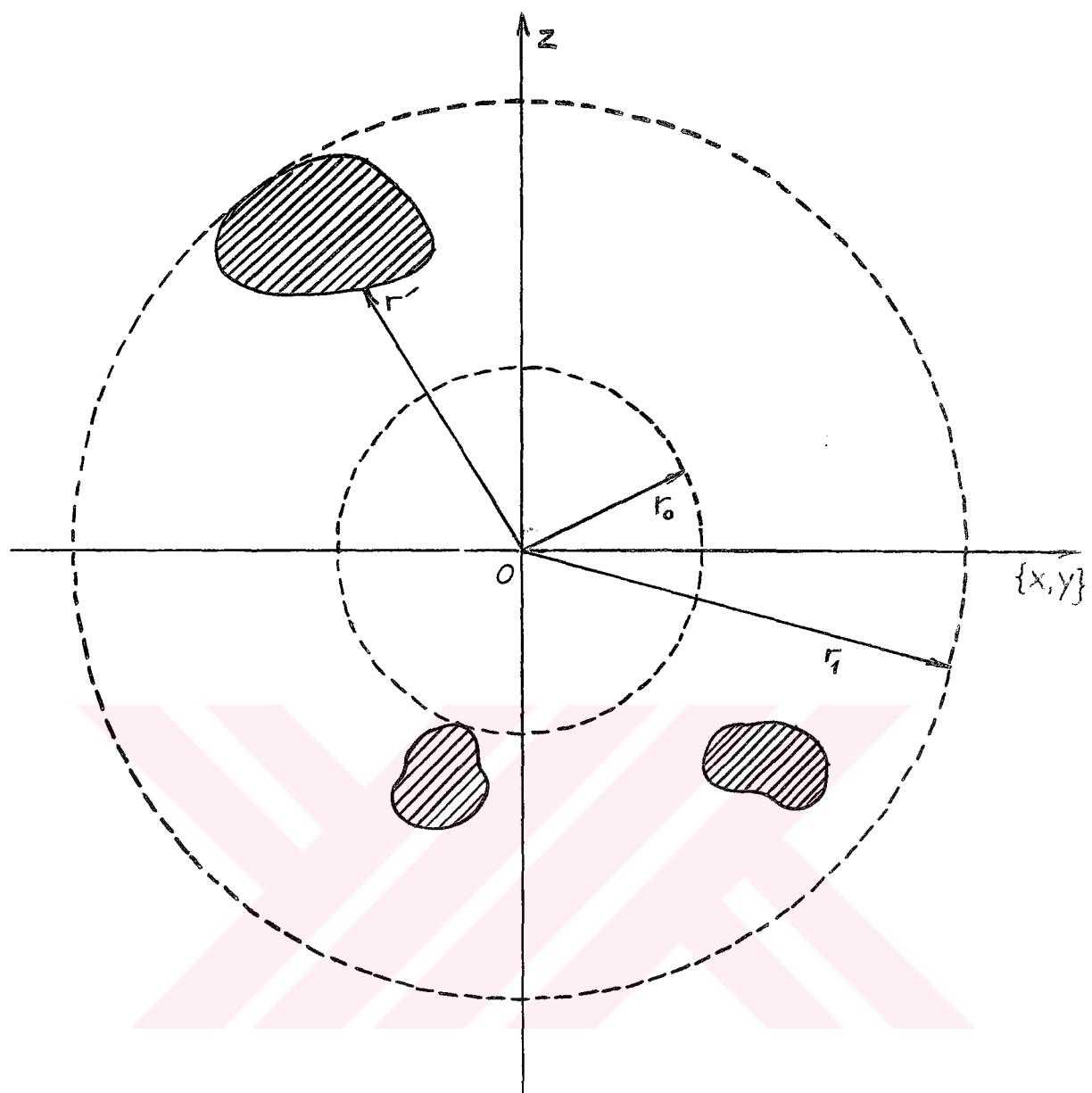


Fig 2.1.1.Scattering Geometry for Multibody Scattering.

For  $r < r'$ ,

$$G_o(\vec{F}|\vec{F}') \vec{I} = \sum_{p,m} ik \left\{ \bar{M}_{p,m}^1(\vec{F}) \bar{M}_{p,m}^{2*}(\vec{F}') + \frac{1}{k^2} \bar{N}_{p,m}^1(\vec{F}) \bar{N}_{p,m}^{2*}(\vec{F}') \right\} \quad (2.1.4)$$

where the vector wave functions  $\bar{M}_{pm}^i$  and  $\bar{N}_{pm}^i$  are given as

$$\bar{M}_{p,m}^i(\vec{F}) = Z_p^i(kr) \bar{X}_{p,m}(\theta, \varphi) = \frac{1}{k^2} \nabla \times \bar{N}_{p,m}^i(\vec{F}) \quad (2.1.5)$$

$$\bar{N}_{p,m}^i(\vec{F}) = \nabla \times \bar{M}_{p,m}^i(\vec{F}) \quad (2.1.6)$$

where the superscript takes the values 1, 2, s. The general radial function  $Z_p^i(kr)$  replaces the following spherical functions :

$$Z_p^1(kr) = J_p(kr) \quad (2.1.7)$$

$$Z_p^2(kr) = H_p^2(kr) = H_p^{1*}(kr) \quad (2.1.8)$$

$$Z_p^s(kr) = H_p^1(kr) = H_p^{2*}(kr) \quad (2.1.9)$$

where  $J_p$  is the spherical Bessel function and  $H_p^1, H_p^2$  are the spherical Hankel functions of the first and the second kind.

$X_{pm}$  is the vector spherical harmonic and it can be written as

$$\bar{X}_{p,m}(\theta, \varphi) = \frac{\vec{F} \times \nabla}{i\sqrt{p(p+1)}} \sqrt{\frac{(2p+1)(p-m)!}{4\pi (p+m)!}} P_p^m(\cos\theta) e^{im\varphi} \quad (2.1.10)$$

When Eqs (2.1.3) and (2.1.4) are substituted in (2.2.1) two different multipole expansions for the re-radiated fields are found, depending on  $r$  and  $r'$ . These are

$$\text{For } r < r_0, \quad \bar{H}_o = \sum_{p,m} (\alpha E_{p,m} \bar{M}_{p,m}^1 - \frac{i}{k} \alpha M_{p,m} \bar{N}_{p,m}^1) \quad (2.1.11)$$

$$\text{For } r > r_1, \quad \bar{H}_s = \sum_{p,m} (\alpha E_{p,m} \bar{M}_{p,m}^s - \frac{i}{k} \alpha M_{p,m} \bar{N}_{p,m}^s) \quad (2.1.12)$$

where the multipole coefficients are equal to :

$$\alpha EO_n = ik \oint_s \bar{n}' \times (\bar{H}_{in} + \bar{H}_{sc}) \cdot \bar{N}_{p,m}^{2*} ds' \quad (2.1.13)$$

$$\alpha MO_n = -k^2 \oint_s \bar{n}' \times (\bar{H}_{in} + \bar{H}_{sc}) \cdot \bar{M}_{p,m}^{2*} ds' \quad (2.1.14)$$

$$\alpha ES_n = ik \oint_s \bar{n}' \times (\bar{H}_{in} + \bar{H}_{sc}) \cdot \bar{N}_{p,m}^{1*} ds' \quad (2.1.15)$$

$$\alpha MS_n = -k^2 \oint_s \bar{n}' \times (\bar{H}_{in} + \bar{H}_{sc}) \cdot \bar{M}_{p,m}^{1*} ds' \quad (2.1.16)$$

All these integrals are over the total surface of the ensemble scatterers. For the region  $r < r < r_1$  it can also be written similar expansions, but since at both sides secondary sources exist the proper form is

$$\bar{H}_I = \sum_{p,m} \left\{ (\alpha EI_{p,m}^1 \bar{M}_{p,m}^1 - \frac{i}{k} \alpha MI_{p,m}^1 \bar{N}_{p,m}^1) + (\alpha EI_{p,m}^2 \bar{M}_{p,m}^2 - \frac{i}{k} \alpha MI_{p,m}^2 \bar{N}_{p,m}^2) \right\} \quad (2.1.17)$$

But unfortunately, these multipole coefficients cannot be given by simple integrals as before. Therefore they should be obtained by boundary relationships. Taking the incident field  $\bar{H}_{in}$  can be written as an expansion of the form

$$\bar{H}_{in} = \sum_{p,m} (\alpha Ei_{p,m} \bar{M}_{p,m}^1 - \frac{i}{k} \alpha Mi_{p,m} \bar{N}_{p,m}^1) \quad (2.1.18)$$

As it was mentioned before, it is assumed that no interactions occur with the scattered fields. Then

$$\bar{n} \times (\bar{H}_{in} + \bar{H}_{sc}) = \bar{H}_I \quad (2.1.19)$$

Assuming that the series converge, a finite number of terms are enough to meet a certain precision. Smoothness assures lack of singularity for the domain  $\bar{r}'$  and (2.1.19) with (2.1.13-16) can be written as an array relation

$\alpha EI^1_{P_m}$	$\alpha MS_{P_m}$
$\alpha MI^1_{P_m}$	$\alpha ES_{P_m}$
$\alpha EI^2_{P_m}$	$\alpha MO_{P_m}$
$\alpha MI^2_{P_m}$	$\alpha EO_{P_m}$

The elements of the square matrices can be written as

$$IE_{\text{pmp'm'}}^{\text{ij}} = k^2 \oint_s (\bar{M}_{\text{p,m}}^{i*} \times \bar{M}_{\text{p',m'}}^j) \cdot \bar{n}' ds' \quad (2.1.21)$$

$$IM_{pmp'm'}^{ij} = -ik \phi_s (\bar{M}_{p,m}^{i*} \times \bar{N}_{p',m'}^j) \cdot \bar{n}' ds' \quad (2.1.22)$$

$$CE_{pmp'm'}^{ij} = -ik\phi_s (\bar{N}_{p,m}^{i*} \times \bar{M}_{p',m'}^j) \cdot \bar{n}' ds' \quad (2.1.23)$$

$$CM_{pmp'm'}^{ij} = - \oint_s (\bar{N}_{p,m}^{i*} \times \bar{N}_{p',m'}^j) \cdot \bar{n}' ds' \quad (2.1.24)$$

These entities have interesting properties and the essential ones can be derived from Green's second vector identity[8]and vector wave functions,these are

$$IE_{p,m,p',m'}^{i,j} = CM_{p,m,p',m'}^{i,j} \quad (2.1.25)$$

$$IM_{p,m,p',m'}^{i,j} = - CE_{p,m,e'm}^{j,i} \quad (2.1.26)$$

$$IE_{p,m,p',m'}^{J,i} = - (IE_{p',m',p,m}^{i,J})^* \quad (2.1.27)$$

$$IM_{p,m,p',m'}^{J,i} = - (IM_{p',m',p,m}^{i,J})^* \quad (2.1.28)$$

where  $i=1,2$  and  $j=1,2$  only.

These properties are very important for testing the numerical results and increasing the computational efficiency.

Continuity of the total electromagnetic field accross the spherical boundaries, orthogonality property of vector wave functions on a spherical surface give

$$\begin{bmatrix} \alpha MS_{p,m} \\ \alpha ES_{p,m} \\ \alpha MO_{p,m} \\ \alpha EO_{p,m} \end{bmatrix} = \begin{bmatrix} 0 & \delta_{pp'} \delta_{mm'} A_1^P & 0 & \delta_{pp'} \delta_{mm'} A_3^P \\ \delta_{pp'} \delta_{mm'} A_2^P & 0 & \delta_{pp'} \delta_{mm'} A_3^P & 0 \\ 0 & \delta_{pp'} \delta_{mm'} & 0 & \frac{\delta_{pp'} \delta_{mm'}}{A_1} \\ \delta_{pp'} \delta_{mm'} & 0 & \frac{\delta_{pp'} \delta_{mm'}}{A_1} & 0 \end{bmatrix} \begin{bmatrix} \alpha EI_{p,m}' \\ \alpha MI_{p,m}' \\ \alpha ET_{p,m}' \\ \alpha MI_{p,m}'' \end{bmatrix} - \begin{bmatrix} A_2^P \alpha NP_{p,m} \\ A_2^P \alpha Ei_{p,m} \\ \alpha Mi_{p,m} \\ \alpha El_{p,m} \end{bmatrix} \quad (2.1.29)$$

where

$$A_1^P = J_p(kr_0) / H_p^2(kr_0) \quad (2.1.30)$$

$$A_2^P = J_p(kr_1) / H_p^1(kr_1) \quad (2.1.31)$$

$$A_3^P = H_p^2(kr_1) / H_p^1(kr_1) \quad (2.1.32)$$

Equations (2.1.20) and (2.1.19) reduce the problem to a simple system of equations of the form

$$Y = S X = B X + U \quad (2.1.33)$$

where  $S$ ,  $B$  and  $U$  are named as scattering, boundary and excitation matrices, respectively. For example the  $U$  matrix entities for a circularly polarized plane wave progressing in the positive  $z$  direction is given by [8]:

$$\alpha Ei_{p,m} = -m i^{p+1} (4\pi (2p+1))^{1/2} \quad (2.1.34)$$

$$\alpha Mi_{p,m} = i^{p+1} (4\pi (2p+1))^{1/2} \quad (2.1.35)$$

where  $m=+1$  corresponds left hand polarization,  $m=-1$  right hand polarization. Although  $m=\pm 1$  this does not imply  $m'=\pm 1$ . The values of  $m'$  depend on the shape of the ensemble, too. Thus the system is coupled and cannot be partitionized into small systems. This is possible for rotational ensembles and for this case  $m'$  takes the values of  $m$  only.

## 2.2. MODIFICATION OF THE THEORY FOR INCLUSION OF THE ORIGIN

If the origin is included by one of the bodies, the previous formulation fails since for the region  $r < r_0$  s type expansion is not valid. In the sphere of radius  $r_0$  and on it the total field must vanish, then the following equations can be written considering the orthogonality of the spherical vector wave functions on a spherical surface :

$$\alpha E_{in} + ik \oint_s \frac{4\pi}{c} \bar{R}(F') \cdot \bar{N}_n^{2*}(F') dS' = 0 \quad (2.2.1)$$

$$\alpha M_{in} - k^2 \oint_s \frac{4\pi}{c} \bar{R}(F') \cdot \bar{M}_n^{2*}(F') dS' = 0 \quad (2.2.2)$$

With previous definitions, the boundary matrix B becomes :

$$B = \begin{vmatrix} 0 & \delta_{pp'} \delta_{mm'} & 0 & \delta_{pp'} \delta_{mm'} \\ \delta_{pp'} \delta_{mm'} & A_2^P & 0 & A_3^P \\ 0 & 0 & \delta_{pp'} \delta_{mm'} & 0 \\ \delta_{pp'} \delta_{mm'} & A_2^P & A_3^P & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{vmatrix} \quad (2.2.3)$$

On the other hand a careful investigation of Green's identities with exclusion of singularities [8] yield the following symmetry relations :

$$IE_{nn'}^{ij} = CM_{nn'}^{ij} ; \quad n=p,m \quad (2.2.4)$$

$$IM_{nn'}^{11} = -CE_{nn'}^{11}, \quad (2.2.5)$$

$$IM_{nn'}^{22} = -CE_{nn'}^{22} - 2\delta_{nn'} \quad (2.2.6)$$

$$IM_{nn'}^{ij} = -CE_{nn'}^{ij} - \delta_{nn'} ; \quad i \neq j \quad (2.2.7)$$

$$[IE_{nn'}^{ji}] = -[(IE_{nn'}^{ij})^*]^T \quad (2.2.8)$$

$$[IM_{nn'}^{11}] = -[(IM_{nn'}^{11})^*]^T \quad (2.2.9)$$

$$[IM_{nn'}^{22}] = -[(IM_{nn'}^{22})^*]^T - 2\delta_{nn'} \quad (2.2.10)$$

$$[IM_{nn'}^{ji}] = -[(IM_{nn'}^{ij})^*]^T - \delta_{nn'} ; \quad i \neq j \quad (2.2.11)$$

Again these equations can be used as a check for the results or a short cut for the algorithms. The coupledness of the equations is still present and to obtain uncoupled equations the configuration must be rotationaly again.

### 3. NUMERICAL ANALYSIS OF THE PROBLEM

#### 3.1. INTRODUCTION

Numerical part of the problem is the evaluation of the necessary functions, taking surface integrals and the solution of the scattering equations. Though it seems very easy, from the computer programming point of view there exist handicaps which limit the choice of the method.

The most essential dilemma of the problem is the question of need for time-core efficiency and the precision required for the results. Short-cut algorithms are necessary. Selection of the method is very important. On the other hand, since the core size of the available computer is limited, storage strategy is vital. Though external storage devices(disk and tape) are available excessive usage of them increases the execution time. Since at least at the primary steps, double precision is necessary, situation turns out to be a cross word puzzle, since double precision increases core demand and execution time considerably.

For toroidal bodies, there is not much trouble. If the lower illumination with a circularly polarized plane wave is considered, everything becomes easy. However generality of the problem reduces.

### 3.2 GENERATION OF ASSOCIATE LEGENDRE FUNCTIONS

The first step in the evaluation of the scattering integrals is the precise generation of the necessary functions. Associate Legendre functions have very nice recursive properties, their derivatives are also associatable to the [18] functions in a simple fashion. Following property of them is quite important:

$$Y_P^{-m} P_P^{-m}(\cos \theta) = Y_P^m P_P^m(\cos \theta) \quad (3.2.1)$$

For lower illumination of the toroidal configurations with circularly polarized plane waves,  $m=1$  case will yield the results of  $m=-1$  case with slight modifications.

First two values have to be evaluated from explicit expressions :

$$P_1^1(\cos \theta) = -\sin \theta \quad (3.2.2)$$

$$\frac{d}{d\theta} P_1^1(\cos \theta) = -\cos \theta \quad (3.2.3)$$

$$P_2^1(\cos \theta) = -3 \sin \theta \cos \theta \quad (3.2.4)$$

$$\frac{d}{d\theta} P_2^1(\cos \theta) = -3 (\cos^2 \theta - \sin^2 \theta) \quad (3.2.5)$$

Other ones are generated by the recursive formulas :

$$P_k^1(\cos \theta) = \frac{(2k-1)}{(k-1)} P_{k-1}^1(\cos \theta) \cos \theta - \frac{k}{k-1} P_{k-2}^1(\cos \theta) \quad (3.2.6)$$

$$\frac{d}{d\theta} P_k^1(\cos \theta) = k \cot \theta P_k^1(\cos \theta) - \frac{k+1}{\sin \theta} P_{k-1}^1(\cos \theta) \quad (3.2.7)$$

For a case where  $m$  takes other values too, there exist other recursion relations which may be found in any mathematical handbook (i.e., 'Mathematical Handbook' [20],  
-13-

Speigel, McGraw Hill Book Co., 1968)

For  $\theta = 0$ ,  $P_k^1(\cos\theta) = 0$ ,  $k = 1, 2, 3, \dots, n$ . But in the analysis of scattering cross-sections  $P_k^1(\cos\theta)/\sin\theta$  has to be evaluated. Since the computer cannot take the limit for 0/0 case it is necessary to modify the previous equations as :

$$Q_k(\theta) = P_k^1(\cos\theta)/\sin\theta \quad (3.2.8)$$

$$Q_1(\theta) = -1 \quad (3.2.9)$$

$$Q_k(\theta) = \frac{(2k-1)}{k} Q_{k-1}(\theta) \cos\theta - \frac{k}{k-1} Q_{k-1}(\theta) \quad (3.2.10)$$

$$Q_2(\theta) = -3 \cos\theta \quad (3.2.11)$$

$$\frac{d}{d\theta} P_1^1(\cos\theta) = -\cos\theta \quad (3.2.12)$$

$$\frac{d}{d\theta} P_2^1(\cos\theta) = -3(\cos^2\theta - \sin^2\theta) \quad (3.2.13)$$

$$\frac{d}{d\theta} P_k^1(\cos\theta) = k \cos\theta Q_k(\theta) - (k+1) Q_{k-1}(\theta) \quad (3.2.14)$$

### 3.3 GENERATION OF SPHERICAL BESSEL, NEUMAN, HANKEL FUNCTIONS

Spherical Bessel, Neuman, Hankel function also offer some recursive relations. For Bessel and Neumann functions first two values are calculated from :

$$J_0(x) = \sin x / x \quad (3.3.1)$$

$$Y_0(x) = -\cos x / x \quad (3.3.2)$$

$$J_1(x) = (J_0(x)/x + Y_0(x)) \quad (3.3.3)$$

$$Y_1(x) = (Y_0(x)/x - J_0(x)) \quad (3.3.4)$$

Both functions and any linear combination of them-such as Hankel functions-have the following recursive properties :

$$f_m(x) = \frac{2m-1}{x} f_{m-1}(x) - f_{m-2}(x) \quad (3.3.5)$$

$$\frac{d}{dx} x f_m(x) = x f_{m-1}(x) - m f_m(x) \quad (3.3.6)$$

By means of these equations all of the radial functions can be evaluated easily. In Fortran , since computation by means of proper equivalencing real and complex quantities, allows direct references to real and imaginary parts, it may be useful to generate only one of the hankel functions. For this again first two values have to be evaluated by means of explicit equations:

$$H_0^{(1)}(x) = \frac{e^{ix}}{i x} \quad (3.3.7)$$

$$H_1^{(1)}(x) = \left(-i + \frac{1}{x}\right) H_0^{(1)}(x) \quad (3.3.8)$$

Other ones are generated again by the previous recursion equations. Complex conjugation(which is not a time taking job ) will give other hankel function. Bessel and Neumann functions are used referring only one part of the hankel functions. [18]

For small values of argument these methods may fail due to enlargening errors during recursions. Miller method(reverse recursion and scaling method) and series solution may be useful but time consuming for this case. Since the position of the bodies is not critical,one can avoid small arguments.

### 3.4 INDICAL FUNCTIONS

There is no other way, except direct evaluation, for the indicial functions. Since this is done only once, no trouble arises. They are used in the expressions together and therefore it is good to evaluate the following expressions, instead :

$$\xi_k = \gamma_k^{\frac{1}{2}} \Delta_k = \left\{ \frac{2k+1}{4\pi} \right\}^{\frac{1}{2}} \quad (3.4.1)$$

$$J_k = \frac{\gamma_k^{\frac{1}{2}}}{\Delta_k} = \frac{\xi_k}{k(k+1)} \quad (3.4.2)$$

If they are stored arraywise, it is better to evaluate them once, and assign them during compilation. Thus, since the object form of a working program is used and no compilation is required they won't effect duration.

### 3.5 NECESSARY MAPPINGS FOR INTEGRATION

Assume a collection of bodies is present at a certain region of the frame concerned. The surface integral over the total surface can be separated into single surface integrals as

$$\oint \bar{A} \cdot d\bar{S} = \sum_{S_1} \oint \bar{A} \cdot d\bar{S} + \sum_{S_2} \oint \bar{A} \cdot d\bar{S} + \dots + \sum_{S_n} \oint \bar{A} \cdot d\bar{S} \quad (3.5.1)$$

where n is the total number of bodies. Therefore the problem of finding the scattering matrices reduces to taking the integrals of

$$\oint_{S_L} \bar{A} \cdot d\bar{S} \quad (3.5.2)$$

type and summing them. Every surface can be described by two independent variables, parametrically. For example:

$$x = f(u, v) \quad (3.5.3)$$

$$y = g(u, v) \quad (3.5.4)$$

$$z = h(u, v) \quad (3.5.5)$$

This is a mapping into  $(u, v)$  system. Then with proper substitutions integrals may be converted into a form like :

$$\oint_{S_i} \bar{A} \cdot d\bar{S} = \iint_{A_{uv}^i} \bar{a}(u, v) \cdot d\bar{S}(u, v) \quad (3.5.6)$$

where  $A_{uv}^i$  is the image of the actual surface  $S_i$ . But if the transformation is not one to one, to describe  $A_{uv}^i$  for the computer becomes very difficult (especially for a general program) and therefore the best is to find a nice transformation where  $A_{uv}^i$  is just a rectangle. This can be established for a convex body if the body is described with respect to an origin inside the body.

For surface element  $d\bar{S}$  one can make the following separation :

$$d\bar{S} = \bar{n} dS \quad (3.5.7)$$

where  $\bar{n}$  is the unit normal (outward) and  $dS$  is the unit surface element. On the other hand if the surface is described as  $r=r(u, v)$ , where  $r$  shows the magnitude of the position vector, one can write:

$$dS = \left| \left( \frac{\partial \bar{r}}{\partial u} du \right) \times \left( \frac{\partial \bar{r}}{\partial v} dv \right) \right| = \left| \frac{\partial \bar{r}}{\partial u} \times \frac{\partial \bar{r}}{\partial v} \right| du dv \quad (3.5.8)$$

$$\text{but } \pm \frac{\partial \bar{r}}{\partial u} \times \frac{\partial \bar{r}}{\partial v} = \bar{n} \left| \begin{array}{c} \frac{\partial \bar{r}}{\partial u} \times \frac{\partial \bar{r}}{\partial v} \end{array} \right| \quad (3.5.9)$$

$$\bar{n} = \left[ \frac{\partial \bar{r}}{\partial u} \times \frac{\partial \bar{r}}{\partial v} \right] / \left| \frac{\partial \bar{r}}{\partial u} \times \frac{\partial \bar{r}}{\partial v} \right| \quad (3.5.10)$$

Since these derivatives are tangent to the surface, their cross product will be in the direction of the normal. Then

$$\pm d\bar{s} = \pm \bar{n} ds = \frac{\partial \bar{r}}{\partial u} \times \frac{\partial \bar{r}}{\partial v} du dv \quad (3.5.11)$$

is found and (3.5.2) becomes

$$\oint_{S_i} \bar{A} \cdot d\bar{s} = \iint_{A_{uv}} \bar{a}(u, v) \cdot \left\{ \frac{\partial \bar{r}}{\partial u} \times \frac{\partial \bar{r}}{\partial v} \right\} du dv \quad (3.5.12)$$

Here it is necessary to take the cross multiplication which will produce outgoing normal direction. [16]

### 3.6. GENERAL STRUCTURES

The best way to express general convex structures is to describe the bodies with respect to their body frame spherically. Thus the surface of a certain body will be :

$$\bar{r} = r(\theta, \varphi) \bar{u}_r \quad (3.6.1)$$

This ensures smoothness if  $r$  is analytical with respect to  $\theta, \varphi$ .

Then

$$\frac{\partial \bar{r}}{\partial \theta} = \frac{\partial r}{\partial \theta} \bar{u}_r + r \bar{u}_\theta \quad (3.6.2)$$

$$\text{since } \frac{\partial \bar{u}_r}{\partial \theta} = \bar{u}_\theta \quad (3.6.3)$$

$$\frac{\partial \bar{r}}{\partial \varphi} = r \sin \theta \bar{u}_\varphi + \frac{\partial r}{\partial \varphi} \bar{u}_r \quad (3.6.4)$$

since  $\frac{\partial \bar{F}}{\partial \varphi} = \sin \theta \frac{\partial \bar{u}_r}{\partial \varphi}$  (3.6.5)

and thus

$$d\bar{S} = \left( \frac{\partial \bar{F}}{\partial \theta} \times \frac{\partial \bar{F}}{\partial \varphi} \right) d\theta d\varphi = \left( r^2 \sin \theta \bar{u}_r + r \frac{\partial r}{\partial \theta} \sin \theta \bar{u}_\theta + r \frac{\partial r}{\partial \varphi} \bar{u}_\varphi \right) d\theta d\varphi \quad (3.6.7)$$

or  $\oint \bar{A} \cdot d\bar{S} = \int_0^{2\pi} \int_0^\pi \bar{a}(\theta, \varphi) \cdot \left\{ r^2 \sin \theta \bar{u}_r + r \frac{\partial r}{\partial \theta} \sin \theta \bar{u}_\theta - r \frac{\partial r}{\partial \varphi} \bar{u}_\varphi \right\} d\theta d\varphi \quad (3.6.8)$

is obtained.

But unfortunately  $\bar{a}(\theta, \varphi)$  can best be evaluated with respect to prime coordinate frame (original one). Therefore one has to describe vectors in another system. A vectorial coordinate change requires two things, change of the point descriptions, base vector transformations. Assume the coordinates of the body frame is given as  $(x_c, y_c, z_c)$ , then

$$x_p = r_b \sin \theta_b \cos \varphi_b + x_c \quad (3.6.9)$$

$$y_p = r_b \sin \theta_b \sin \varphi_b + y_c \quad (3.6.10)$$

$$z_p = r_b \cos \theta_b + z_c \quad (3.6.11)$$

where the subscripts b, p show body and prime frames respectively.

On the other hand in terms of prime spherical coordinates are

$$r_p = (x_p^2 + y_p^2 + z_p^2)^{1/2} \quad (3.6.12)$$

$$\theta_p = \arccos(z_p / r_p) \quad (3.6.13)$$

$$\varphi_p = \arctg(y_p / x_p) \quad (3.6.14)$$

Thus prime coordinates can easily be found. To find a direct relation between vectorial components of the frames is a bit

= 1 / -

difficult. But one can use a rectangular system as an intermediate step. For rectangular systems if axes are kept parallel to each other, choosing a particular frame is immaterial for vectorial transformations.

$$\begin{vmatrix} \bar{a}_x \\ \bar{a}_y \\ \bar{a}_z \end{vmatrix} = \begin{vmatrix} \sin\theta_b \cos\varphi_b & \cos\theta_b \cos\varphi_b & -\sin\varphi_b \\ \sin\theta_b \sin\varphi_b & \cos\theta_b \sin\varphi_b & \cos\varphi_b \\ \cos\theta_b & -\sin\theta_b & 0 \end{vmatrix} \begin{vmatrix} \bar{a}_r \\ \bar{a}_\theta \\ \bar{a}_\psi \end{vmatrix} \quad (3.6.15)$$

gives body-frame cartesian components. For later reference we will use

$$S(\alpha, \beta) = \begin{vmatrix} \sin\alpha \cos\beta & \cos\alpha \cos\beta & -\sin\beta \\ \sin\alpha \sin\beta & \cos\alpha \sin\beta & \cos\beta \\ \cos\alpha & -\sin\alpha & 0 \end{vmatrix} \quad (3.6.16)$$

see that

(3.6.17)

thus the transformation is orthogonal and inverse of  $S$  is equal to its transpose. Using this, a second transformation can be written immediately

$$\begin{vmatrix} \bar{a}_r \\ \bar{a}_\theta \\ \bar{a}_\psi \end{vmatrix} = S^T(\theta_p, \varphi_p) \begin{vmatrix} \bar{a}_x \\ \bar{a}_y \\ \bar{a}_z \end{vmatrix} \quad (3.6.18)$$

Therefore

### 3.7 TOROIDAL STRUCTURES AND DEGENERATE SPHERICAL COORDINATES

With the name 'toroidal body', a revolutionary volume is understood, which is made revolving a plane figure around the z axis, which is chosen coincident with the direction of incidence. Although a toroidal ensemble cannot be assumed to be a general multibody structure, it presents very nice properties, which reduce the calculations considerably. For such a structure, the equations given in chapter 2 become uncoupled for  $m, m'$ . Thus a huge system of equations can be partitioned into smaller and simpler systems. If the incident field is expandable for a finite variation of  $m$ 's, as in case of a c.p. plane wave in z direction, the number of the partitioned systems are also finite. Regardless these facts, the previous method described in 3.6 is inadequate for solving toroidal scattering problems (for example from a ring) since with that transformation mapping is not one-to-one, for toroids. Therefore more or less a method to handle toroidal configurations will be useful.

Toroidal ensembles can best be treated by degenerate spherical coordinates. For this system surface integrals become simple integrals. Assume the characteristic curve of the toroid (cross section at only one side of a plane including z axis) can be described with a polar system (if the curve is convex) by a one to one mapping from  $\beta$  to  $\rho$ . Where  $\beta$  is the angular variable,  $\rho$  denotes the radial variable. Then the surface integrals become integrals of the form

$$\int_{\beta_{\min}}^{\beta_{\max}} (F_r(\beta) dS_r(\beta) + F_\theta(\beta) dS_\theta(\beta)) d\beta = \int_{\beta_{\min}}^{\beta_{\max}} f(\beta) d\beta \quad (3.7.1)$$

where the limits are shown in Fig.(3.7.1)

The necessity for a one-to-one correspondence is not a handicap, since any non-convex shape may be sectioned and may be described with respect to several suitable origins and limits. Fundamental equations of the degenerate spherical coordinates are given in terms of cartesian variables as

$$x_1 = (\alpha + \rho \cos \beta) \cos \varphi \quad (3.7.2)$$

$$x_2 = (\alpha + \rho \cos \beta) \sin \varphi \quad (3.7.3)$$

$$x_3 = (\alpha + \rho \sin \beta) \quad (3.7.4)$$

However this is not enough. Some other properties have to be known. The most importants are

$$h_\rho = \left\{ \sum_{i=1}^3 \left( \frac{\partial x_i}{\partial \rho} \right)^2 \right\}^{1/2} = 1 \quad (3.7.5)$$

$$h_\beta = \left\{ \sum_{i=1}^3 \left( \frac{\partial x_i}{\partial \beta} \right)^2 \right\}^{1/2} = \rho \quad (3.7.6)$$

$$h_\varphi = \left\{ \sum_{i=1}^3 \left( \frac{\partial x_i}{\partial \varphi} \right)^2 \right\}^{1/2} = \alpha + \rho \cos \beta \quad (3.7.7)$$

Remembering [15]

$$d\bar{r} = h_\rho d\rho \bar{u}_\rho + h_\beta d\beta \bar{u}_\beta + h_\varphi d\varphi \bar{u}_\varphi \quad (3.7.8)$$

$$\frac{\partial \bar{r}}{\partial \beta} = \frac{\partial \rho}{\partial \beta} \bar{u}_\rho + \rho \bar{u}_\beta \quad (3.7.9)$$

$$\frac{\partial \bar{r}}{\partial \varphi} = \frac{\partial \rho}{\partial \varphi} \bar{u}_\rho + (\alpha + \rho \cos \beta) \bar{u}_\varphi \quad (3.7.10)$$

Now taking the cross product

$$\bar{n} dS = \frac{\partial \bar{F}}{\partial \beta} \times \frac{\partial \bar{F}}{\partial \varphi} d\beta d\varphi \quad (3.7.11)$$

$$\bar{n} dS = [(a + \rho \cos \beta) \rho \bar{u}_\rho - (a + \rho \cos \beta) \frac{\partial \rho}{\partial \beta} \bar{u}_\beta - \rho \frac{\partial \rho}{\partial \varphi} \bar{u}_\varphi] d\beta d\varphi \quad (3.7.12)$$

is obtained. However for toroidal bodies further reductions are possible because

$$\bar{u}_\rho \cdot \bar{F} = 0 \quad (3.7.13)$$

$$\frac{\partial \rho}{\partial \varphi} = 0 \quad (3.7.14)$$

in short a modified surface element is produced as

$$\bar{n} dS' = 2\pi (a + \rho \cos \beta) (\rho \bar{u}_\rho - \frac{\partial \rho}{\partial \beta} \bar{u}_\beta) d\beta \quad (3.7.15)$$

Since the vector field to be integrated is known in terms of its spherical components, transformation matrix has to be known for vectorial conversions. This can be derived from F.3.7.3. and the following result is achieved :

$$\begin{vmatrix} \bar{u}_r \\ \bar{u}_\theta \\ \bar{u}_\varphi \end{vmatrix} = \begin{vmatrix} \sin(\beta + \theta) & \cos(\beta + \theta) & 0 \\ \cos(\beta + \theta) & -\sin(\beta + \theta) & 0 \\ 0 & 0 & 1 \end{vmatrix} \begin{vmatrix} \bar{u}_\rho \\ \bar{u}_\beta \\ \bar{u}_\varphi \end{vmatrix} \quad (3.7.16)$$

Now the surface element vector can be transferred into a spherical system as

$$dS_r = 2\pi (a + \rho \cos \beta) (\rho \sin(\beta + \theta) - \frac{\partial \rho}{\partial \beta} \cos(\beta + \theta)) d\beta \quad (3.7.17)$$

$$dS_\theta = 2\pi (a + \rho \cos \beta) (\rho \cos(\beta + \theta) + \frac{\partial \rho}{\partial \beta} \sin(\beta + \theta)) d\beta \quad (3.7.18)$$

in these equations following identity is used :

$$\Theta = \arctg \frac{a + \rho \cos \beta}{b + \rho \sin \beta} \quad (3.7.19)$$

Similarly

$$r = \{(a + \rho \cos \beta)^2 + (b + \rho \sin \beta)^2\}^{1/2} \quad (3.7.20)$$

By means of these transformations, integrals at the end become simple integrals. They can be evaluated by Gauss-Legendre type of quadrature. Related formulas for this time are simply

$$\int_{\beta_{\min}}^{\beta_{\max}} f(\beta) d\beta = \sum_i f(\beta_i) w'_i \quad (3.7.21)$$

where

$$\beta_i = \frac{\beta_{\max} - \beta_{\min}}{2} x_{qi} + \frac{\beta_{\max} + \beta_{\min}}{2} \quad (3.7.22)$$

$$w'_i = \frac{\beta_{\max} - \beta_{\min}}{2} w_{qi} \quad (3.7.23)$$

in these equations  $x$  and  $w$  stands for quadrature arguments and weights respectively.

### 3.8. SOLUTION OF THE SYSTEM OF EQUATIONS

After evaluation of the scattering matrice and the boundary matrice preceisely, one has to find the solution to the system given in chapter 2. It is better to find  $X$  from the equation

$$(S-B) X = U \quad (3.8.1)$$

than obtain  $Y$  from the boundary relations. Since the scattering integrals are complex a careful investigation is necessary. Solution of this matrice system is vital because this is the part which enforce the computer core capacity. Therefore a preceise, fast, less-core requiring algorithm has to be used. If the system is not an uncoupled one for  $m, m'$ , matrice sizes increase tre-

mendously. This is still a problem to be solved. If the equations are uncoupled, solutions for a sequence of equations of simpler or simpler form is necessary and can be solved with less trouble. If we can limit expansion of the incident wave over  $m$  to a finite number of values, problem reduces more. This is the case for a lower illumination of a toroidal system by a c.p. plane wave. Since we always want to increase the number of terms in the expansions to provide a better approximation, whatever the equation system is, it is a large one.

Several methods are employed during investigations. Some of these methods and their drawbacks are outlined below :

a) Separation of Real and Imaginary Parts:

Given a system of complex linear equations of the form

$$A X = B \quad (3.8.2)$$

It can be separated into real and imaginary parts in the form

$$\begin{vmatrix} A_r & -A_i \\ A_i & A_r \end{vmatrix} \begin{vmatrix} X_r \\ X_i \end{vmatrix} = \begin{vmatrix} B_r \\ B_i \end{vmatrix} \quad (3.8.3)$$

Thus a real system of equations is obtained. Although this seems as if an easy method, it is not efficient because it enlarges the system unnecessarily. Because the system size is now  $2n$  if there were  $n$  equations at the beginning. Core capacity requirement makes the usage of single precision necessary. Since the

largeness of the system brings enough error, single precision replacement makes the situation worse. For this reason, this is not a good method for solving such a system.

b) Crout Iterative Refinement Technique:

This is a rather simple technique. If A is a nonsingular complex matrix, it can be factorized in the form [19]

$$A = LU \quad (3.8.4)$$

where L is a lower triangle matrix, U is upper triangular. The elements of the matrices L and U of the crout factorization may be determined in the order r-th column of L, r-th row of U where r varies from 1 to n, by means of the equations :

$$\sum_{k=1}^{r-1} l_{ik} u_{kr} + l_{ir} = \alpha_{ir} \quad (i = r, \dots, n) \quad (3.8.5)$$

$$\sum_{k=1}^{r-1} l_{rk} u_{ki} + l_{rr} u_{ri} = \alpha_{ri} \quad (i = r+1, \dots, n) \quad (3.8.6)$$

After partitioning, to obtain the solution two steps are necessary. First the system

$$L y = b \quad (3.8.7)$$

is solved, then the subsidiary system

$$U x = y \quad (3.8.8)$$

is solved. Although this seems to be a rather simple algorithm, stability has to be assured by means of some pivoting. Pivoting

performed can best be described in terms of the usual algorithmic interpretation including the storage arrangement. There are n major steps and at the beginning of the rth step, the stored array is illustrated when n=4, r=3

$$\begin{bmatrix} l_{11} & u_{12} & u_{13} & u_{14} \\ l_{21} & l_{22} & u_{23} & u_{24} \\ l_{31} & l_{32} & a_{33} & a_{34} \\ l_{41} & l_{42} & a_{43} & a_{44} \\ \text{int}_1 & \text{int}_2 & \text{int}_3 & \text{int}_4 \end{bmatrix} \quad (3.8.9)$$

The rth step is as follows

i) Compute  $l_{ir} = a_{ir} - \sum_{k=1}^{r-1} l_{ik} u_{kr}$  and over write on  $a_{ir}$  ( $i=r, r+1, \dots, n$ )

ii) If  $\text{int}_r$  is the smallest integer for which  $|l_{\text{int}_r, r}| = \max_{i>r} |l_{ir}|$ , then store the integer  $\text{int}_r$  and interchange the whole rows of r and  $\text{int}_r$  in the current array.

iii) Compute  $u_{ri} = (a_{ri} - \sum_{k=1}^{r-1} l_{rk} u_{ki}) / l_{rr}$  and over write on  $a_{ri}$  ( $i=r+1, \dots, n$ )

After such n steps A is replaced by L and U and the n integers  $\text{int}_i$  ( $i=1, \dots, n$ ) give the details of the interchanges. If  $P_r$  denotes the permutation matrix, premultiplication by which gives interchange of rows r and  $\text{int}_r$ , then

$$P_n P_{n-1} \dots P_1 A = \tilde{A} = LU \quad (3.8.10)$$

in other words we have the orthodox Crout decomposition of A defined by Eq(3.8.10). The triangles L and U and the array of

integers int, enable us to solve the equations  $A x = b$  corresponding any right side  $b$ . For we have

$$P_n P_{n-1} \dots P_1 A x = b \quad (3.8.11)$$

$$\text{or } L U x = P_n \dots P_1 b \quad (3.8.12)$$

We have only to subject  $b$  to the appropriate interchanges to give say  $\tilde{b}$  and solve

$$Ly = \tilde{b} \quad (3.8.13)$$

$$Ux = y \quad (3.8.14)$$

In general the computer solution of  $A x = b$  can be refined, using the crout factorization, by means of the iterative procedure

$$x^{(0)} = 0 \quad (3.8.15)$$

$$r^{(s)} = b - Ax^{(s)} \quad (3.8.16)$$

$$L U d^{(s)} = P_n \dots P_1 r^{(s)} \quad (3.8.17)$$

$$x^{(s+1)} = x^{(s)} + d^{(s)} \quad (3.8.18)$$

Crout method is a precise method. But both factorizing and iterating schemes reserve large amount of core for their instructions. Time requirement is also important. Since undestroyed form of  $A$  has to be preserved, external devices are necessary to solve storage requirements. However Crout scheme requires less operations than Gauss elimination (without iterations.).

c) Gauss Elimination Method :

This is a well known method and needs no description. It applies equally well for Complex Systems too. However if some pivoting is not done, results are strange due to numerical instabilities.

In scattering algorithms two methods of pivoting is used, namely column-row pivoting and only row pivoting. Row pivoting gives satisfactory results if the system is not too large and some strong ill conditionedness does not occur. Other method is time consuming, but more precise. In pivoting square of the absolute value is used as a measure for significance.

Gauss elimination scheme allows iterative refinement too. Assume  $\epsilon$  denotes some deviation (in vector form) from the actual result due to operational truncations and numerical instabilities. If the result found is substituted in the equation, it will yield a difference  $\Delta b$  from the actual constant vector. But the following two equations show that  $\epsilon$  and  $\Delta b$  are related to each other with the same coefficient matrix. Therefore an iterative refinement technique is applicable until a certain tolerance is achieved.

$$A(x + \epsilon) = b + \Delta b \quad (3.8.19)$$

$$A\epsilon = \Delta b \quad (3.8.20)$$

If iterative refinement is used, Crout scheme becomes more economical from every respect. [9],[10],[12],[13],[14]

d) Gauss Seidel Iteration :

If the scattering equations are re-arranged such that x vector is now in the form :

$$X' = \begin{vmatrix} \alpha M I^1 \\ \cdots \\ \alpha E I^1 \\ \cdots \\ \alpha M I^2 \\ \cdots \\ \alpha E I^2 \end{vmatrix} \quad (3.8.21)$$

the iterative technique becomes applicable without any modification. This is due to great possibility of having larger terms on the diagonal. The simplest of the methods which can be applied is the Gauss-Seidel iterations and this can be applied with suitable error comparison-i.e. using the square of the absolute value-to complex systems also. But convergence cannot be guaranteed for every case. Otherwise it is very suitable for large system solving; it requires less time and has a good precision. [17]

e) Larger Systems :

It is necessary to develop non-iterative methods to solve large systems with less core requirement. The most general problem still requires this. Although not tried, one possibility is to partitionize the coefficient matrices and invert.

Steps are as follows : Assume A is the given matrice and it is partitioned  $A_{11}$  being a square matrice.

$$A = \begin{vmatrix} A_{11} & A_{12} \\ A_{21} & A_{22} \end{vmatrix} \quad (3.8.22)$$

The inverse of A can be written as :

$$A^{-1} = \begin{vmatrix} A_{11}^{-1} + X\theta^{-1}Y & -X\theta^{-1} \\ -\theta^{-1}Y & \theta^{-1} \end{vmatrix} \quad (3.8.23)$$

Where

$$X = A_{11}^{-1} A_{12} \quad (3.8.24)$$

$$Y = A_{21} A_{11}^{-1} \quad (3.8.25)$$

$$\theta = A_{22} - Y A_{12} \quad (3.8.26)$$

$$A_{22} - A_{21}X = A_{22} - A_{21}A_{11}^{-1}A_{12} \quad (3.8.27)$$

This method seems to be applicable only with some external storage devices for the computer available at METU Computation Center.

Consider Hankel functions for asymptotical values :

$$h_p^{(1)}(r) = (-i)^{p+1} \frac{e^{ir}}{r} \quad (3.9.7)$$

neglecting the terms varying with  $1/R^2$  compared with the terms varying with  $1/R$ , using the relation for toroidal lower illumination.

$$(E^{sc})_\theta = 2 \frac{e^{ir}}{r} \sum_p -\frac{\gamma_p^1}{\Delta_p} (-i)^{p+1} (\alpha E s_p^+ \frac{dP_p^1}{d\theta} + \alpha M s_p^+ \frac{P_p^1}{\sin\theta}) \quad (3.9.8)$$

$$(E^{sc})_\varphi = 2 \frac{e^{ir}}{r} \sum_p \frac{\gamma_p^1}{\Delta_p} (-i)^{p+1} (\alpha E s_p^+ \frac{P_p^1}{\sin\theta} + \alpha M s_p^+ \frac{dP_p^1}{d\theta}) \sin\varphi \quad (3.9.9)$$

and the definition for polar re-radiation diagram

$$\sigma(\theta, \varphi) = \frac{4\pi r^2}{\lambda^2} \frac{\bar{U}(\theta, \varphi) \cdot \bar{S}_{\perp}^{sc*}(\theta, \varphi)}{\bar{U}_z \cdot \bar{S}_{\parallel}^{av}} \quad (3.9.10)$$

a simple expression of the form

$$\sigma\{\theta, \varphi\} = \frac{4\pi}{\lambda^2} \left[ \cos^2 \varphi \left| \sum_p \bar{F}_p(\theta) \right|^2 + \sin^2 \varphi \left| \sum_p \bar{G}_p(\theta) \right|^2 \right] \quad (3.9.11)$$

is obtained. Where

$$\bar{F}_p(\theta) = \frac{-\gamma_p(-i)^{p+1}}{\Delta_p} \left( \alpha E s_p^+ \frac{dP_p^1}{d\theta} + \alpha M s_p^+ \frac{P_p^1}{\sin\theta} \right) \quad (3.9.12)$$

$$\bar{G}_p(\theta) = \frac{\gamma_p(-i)^{p+1}}{\Delta_p} \left( \alpha E s_p^+ \frac{P_p^1}{\sin\theta} + \alpha M s_p^+ \frac{dP_p^1}{d\theta} \right) \quad (3.9.13)$$

Specially for  $\varphi=0$  and  $\varphi=\frac{\pi}{2}$  we can write

$$\frac{\sigma(\theta, 0)}{\lambda^2} = \frac{1}{4\pi^2} \left| \sum_p \frac{\sqrt{2p+1}}{p^2+p} (-i)^{p+1} \left( \alpha E s_p^+ \frac{dP_p^1}{d\theta} + \alpha M s_p^+ \frac{P_p^1}{\sin\theta} \right) \right|^2 \quad (3.9.14)$$

$$\frac{\sigma(\theta, \frac{\pi}{2})}{\lambda^2} = \frac{1}{4\pi^2} \left| \sum_p \frac{\sqrt{2p+1}}{p^2+p} (-i)^{p+1} \left( \alpha M s_p^+ \frac{dP_p^1}{d\theta} + \alpha E s_p^+ \frac{P_p^1}{\sin\theta} \right) \right|^2 \quad (3.9.15)$$

these expressions for  $\theta=\pi$  give the back scattering cross-section.

### 3.10. PROGRAMMING

Several computer programs are prepared based on the methods described. For runs of these programs IBM-360/40 system of METU is used. This computer, although is a relatively fast computer , it has not enough core capacity to handle a general scattering problem without using some tricks and external devices; but this makes a considerable amount of increase in duration. But for lower illumination of symmetric structures a compleate subprogram is available to evaluate upto 8 multipole coefficients without any extra device or trick. After 8 coefficients multiphase programming and usage of disk become necessary.

The main subroutine which solves toroidal problem does not depend on the selection of the origin, with proper branchings origin inclusive and origin exclusive problems can be treated. Program works for single and multibody, different quadratures, without re-running infinitely many configurations. It can be used for a frequency sweep over some configurations. If required with a communication code S matrice, coefficients, polar diagram data, polar diagram plots are given one by one or all together.

Equation solving and Bessel,Legendre generating sections are sepearte moduls which can be replaced with others. For example instead of forward recursion, Miller method or series method can be used, Gauss Seidel may be used instead of Crout,etc.

Programming language is in general Fortran, IV 360 DOS level which is sufficient for a problem of scientific nature. There are many versions of the programs which are used for different purposes. They are run in their object deck forms to save compilation times (since the minimum compilation time recorded is 5 min.)

Other languages are also investigated at the beginning. Algol which provides a recursive flexibility cannot be used since the available system is not equipped with Algol compiler. Pl-I seems compatible with Fortran but available compiler is for a subset of it, therefore no attempt is made to write in Pl-I. Assembler language is not proper because of clerical tasks which are necessary, but when Fortran fails it can be mixed with Fortran with proper links. This is done at the primary steps where duration of the operations had to be measured. This can be done by the Assembler Language but not by Fortran of 360/40. Getime macro can be used for this [1] purpose, if this is done with Binary it can be transferred into Fortran being an integer variable. Thus the time is learned by Fortran. Several other tricks can be done, for example standard Fortran function REAL(.) can be established for complex double precision variables also (which is not available) with minimum operations.

With this form of the programs no attempt should be made to use with an arbitrary system, because some unconventional features are also used either for saving time or core. Suitable changes can be made easily however.

#### 4. EXPERIMENTAL INVESTIGATIONS

Some back-scattering measurements are made from a few configurations for which numerical values were present. Due to set up problems no experimental re-radiation diagram is obtained. Toroidal configurations were impossible to manufacture with the tools present.

The procedure to measure the back-scattering cross-section from a configuration for a certain frequency takes about four hours, since repetitions are necessary. Frequencies of the incident waves are chosen to be in the range of 8-11 GHZ. This is to minimize the Rayleigh distance of the radiating antennas. Otherwise in-room measurements are impossible and the ensemble has to be located very far from the antenna. The bodies were manufactured from brass. Tolerance of manufacturing is less than 0.05 mm. Only certain shapes can be made easily, such as spheres, prolate spheroids, eight like figures, etc. But their ensembles may be chosen in various forms. Some of the configurations reduce the total back-scattering cross sections below the proper measurement limit. This is due to tendency to cancel each other of the re-radiated fields of different objects. For example if a body is put in the shadow region of another one, sometimes the backscattering cross-section of the ensemble is less than the back-scattering cross sections of the isolated bodies.

Spherical TEM waves which are produced by a horn antenna may be assumed to be plane waves after Rayleigh distance and this may be used as a plane wave source. But if the region of interest is too long, decay in the magnitude disturbs the accuracy of the results. Phase variations may be reduced placing the bodies far from the horn antenna. Another possibility is to place the ensemble close to the aperture near the symmetry axis of the antenna. If the uniform illumination of the aperture can be obtained, plane wave assumption for small objects is valid. Both of these methods are used to check the accuracy of the results.

It is necessary to isolate the set-up with dissipative sheets to prevent unwanted scatterings. Thus an artificial free space is established. First the antenna is matched to empty range. Then a region where fields are less changing with distance is found. This can be found if the antenna is slightly mismatched. The reflection level from the ensemble is recorded. Then the same measurement is repeated with spheres of several diameters. The difference between the sphere and the ensemble scatterings are measured by means of a calibrated attenuator, obtaining the same amount of deflection of VSWR in every measurement.

Backscattering cross-sections of the spheres are calculated for every frequency and they are compiled as tables. From these tables actual back-scattering cross-section of the ensemble

can be calculated. Some data which is obtained in this fashion is shown in the tables I-III.

Evidently for elongated configurations, the results will be deviated from the calculated ones, because simulation of a plane wave will fail for such ensembles. This is observed for the objects separated too much from each other. Other important factor is the position of the standard spheres. It is difficult to replace the ensemble by a sphere, especially keeping the coincidence of the symmetry axis with the axis of the antenna. Therefore for different spheres, different results are obtained.

The separation of the bodies has to be established properly. Otherwise at high frequencies some troubles may occur due to increase in the error of the optical length of the separation. This is provided placing a dielectric separator of fixed thickness in between the bodies. This dielectric has the properties very close to the properties of the air.

The results which are few in quantity are satisfactory. Some precise measuring techniques seem to be necessary for further research in this field. Some deviations are observed for elongated bodies as expected. A more uniform source is necessary to investigate the full range of the numerical results.

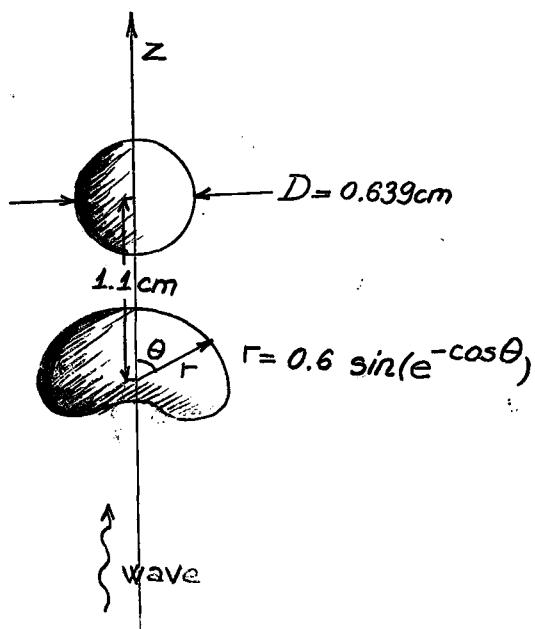


Fig 4.1. - Configuration (A)

Table I - Configuration (A), f= 10.0 GHZ, R=18 cm.

Sphere *	Diam. (cm)	Optical Radius	Sphere $\sigma_{db}$	Rel.to Sphere (db)	$\sigma_{db}$	$\sigma_b/\pi$
1	.639	.669	-3.1	11.5	8.4	2.2
2	.793	.830	1.46	7.75	8.2	2.1
3	.940	.982	3.99	4.0	8.	2.0
4	1.109	1.160	5.06	2.85	7.9	1.95
5	1.200	1.256	4.94	2.9	7.84	1.92
6	1.268	1.325	4.59	3.25	7.8	1.9

CALCULATED :  $\frac{\sigma_b}{\pi} = 1.94$  ,  $\sigma_b = 7.9$  db.

Table II - Configuration (A), f=9.6 GHZ, R = 18 cm.

Sphere *	Diam. (cm.)	Optical Radius	Sphere $\sigma_{db}$	Rel.to Sphere (db)	$\sigma_{db}$	$\sigma_b/\pi$
1	.639	.642	-3.709	11.9	8.2	2.2
2	.793	.796	1.025	6.5	7.525	1.8
3	.940	.945	3.832	3.5	7.532	1.8
4	1.109	1.112	5.307	2.0	7.307	1.7
5	1.200	1.205	5.414	1.9	7.314	1.7
6	1.268	1.272	5.244	2.0	7.244	1.68

CALCULATED :  $\sigma_b/\pi = 1.73$ ,  $\sigma_{db} = 7.4$  db

Table III - Configuration (A), f=9.2 GHZ, R=18 cm.

Sphere *	Diam. ( cm.)	Optical Radius	Sphere $\sigma_{db}$	Rel.to Sphere	$\sigma_{db}$	$\sigma_b/\pi$
1	.639	.615	-4.36	9.95	5.59	1.15
2	.793	.765	0.569	5.20	5.77	1.17
3	.940	.906	3.633	2.3	5.933	1.25
4	1.109	1.067	5.440	0.6	6.04	1.27
5	1.200	1.158	5.752	0.3	6.052	1.28
6	1.268	1.22	5.747	0.38	6.12	1.3

CALCULATED :  $\sigma_b/\pi = 1.505$ ,  $\sigma_{db} = 6.6$  db

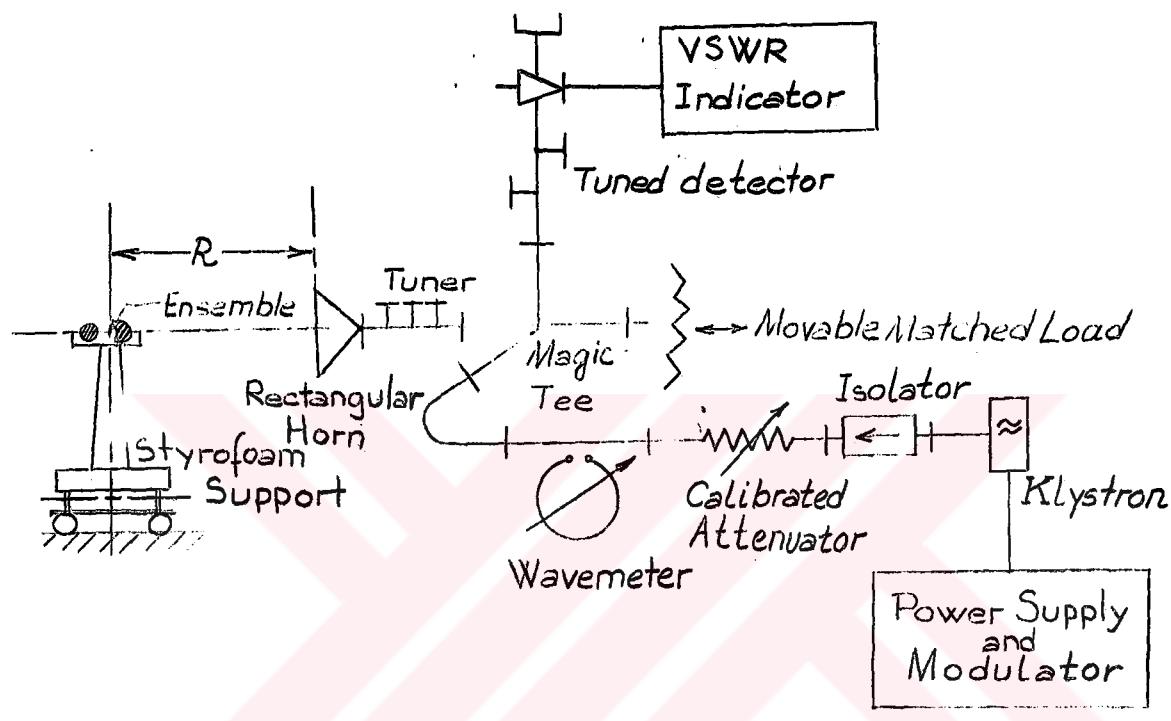


Fig. 4.2. Experimental Set-Up for  $G_f$  Measurements

## 5. GENERAL CONCLUSIONS

Numerical analysis of the problem of scattering of electromagnetic waves from arbitrary perfectly conducting ensembles show the fact the method is limited for such ensembles due to computational difficulties. A large storage is necessary to utilize the algorithms proposed for this problem. Usage of external storage devices may solve this problem however. But this is disadvantageous since the duration increase with such devices, especially if frequent references are made. This means the cost of the research increase. However even with the systems of larger core capacity to increase the number of multipole coefficients may be impossible after a certain limit.

Toroidal ensembles may be evaluated with less trouble if the direction of the incident field is coincident with the axis of revolution. Programs are made for this case requiring approximately the complete storage of a 64 K.byte computer for evaluation of 8 multipole coefficients. Execution of the program for a complete solution leading to re-radiation diagrams and plots of them takes 3 minutes for a certain configuration, with the system IBM 360 40-T. However for an arbitrary excitation of such an ensemble disk usage and multiphase programming may be necessary. For this case there is a chance of using the external

devices more efficiently due to the uncoupled property of the scattering equations. Thus the number of references to an external file reduces. Incidence with a different angle can be investigated expanding the wave with respect to the main coordinate system. This also requires external storage but with uncoupled equations.

Experiments show the difficulty to provide idealized measurement conditions. Since the dimensions increase with the second body due to the separation between the bodies, one has to find a more uniform wave source first, or produce smaller body dimensions. Then there is a limit for observable backscattering cross-sections with the used equipment and smaller bodies produce smaller back-scattering cross-sections. On the other hand to locate the bodies with the required precision is very difficult.

Results are satisfactory and show the validity of the method followed. The modification of the theory seems to be valid also. The most essential property of the formulation is the inclusion of the multiple scatterings without extra effort, since the problem is considered as a radiation due to induced surface sources. If some part of the ensemble touches one of the boundary separation spheres, along a continuous surface segment, the formulation fails since the expansions are no more valid with their previous forms. However, this is not a drawback, it is possible to choose a second origin which allows a valid expansion.

## APPENDIX 1. IE AND IM TYPE INTEGRANDS IN SPHERICAL COORDINATES

Throughout this section, distances will be normalized by  $k$ , thus every distance is optical. This helps simplifications and generalizations. Every free  $k$  in the expressions is associatable to a distance unit of the surface elements or the expressions.

### a) IE type

If the preceding argument is applied the integrand becomes

$$\overline{IE}_{pqmm'}^{\alpha\beta}(\bar{r}) = \overline{M}_{pm}^{\alpha*}(\bar{r}) \times \overline{M}_{qm}^{\beta}(\bar{r}) \quad (\text{A.1.1})$$

where  $\overline{M}_{pm}^{\alpha}$  can be written in normalized form as

$$\overline{M}_{pm}^{\alpha}(\bar{r}) = Z_p^{\alpha}(r) \overline{X}_{pm}^{\alpha}(\theta, \varphi) \quad (\text{A.1.2})$$

where  $Z$  and  $X$  are defined in chapter 2. Evidently

$$\{\overline{M}_{pm}^{\alpha}\}_r = 0 \quad (\text{A.1.3})$$

$$\{\overline{M}_{pm}^{\alpha}\}_{\theta} = Z_p^{\alpha}(r) \{\overline{X}_{pm}^{\alpha}(\theta, \varphi)\}_{\theta} \quad (\text{A.1.4})$$

$$\{\overline{M}_{pm}^{\alpha}\}_{\varphi} = Z_p^{\alpha}(r) \{\overline{X}_{pm}^{\alpha}(\theta, \varphi)\}_{\varphi} \quad (\text{A.1.5})$$

$$\{\overline{X}_{pm}^{\alpha}\}_r = 0 \quad (\text{A.1.6})$$

$$\{\overline{X}_{pm}^{\alpha}\}_{\theta} = -m \frac{\gamma_{pm}}{\Delta_{pm}} \frac{P_p^m(\cos\theta)}{\sin\theta} e^{im\varphi} \quad (\text{A.1.7})$$

$$\{\overline{X}_{pm}^{\alpha}\}_{\varphi} = -i \frac{\gamma_{pm}}{\Delta_{pm}} e^{im\varphi} \frac{d}{d\theta} P_p^m(\cos\theta) \quad (\text{A.1.8})$$

where  $P_p^m$  is the well known Associate Legendre Function, and

$$\gamma_p^m = \sqrt{\frac{(2p+1)}{4\pi} \frac{(p-m)!}{(p+m)!}} \quad (A.1.9)$$

$$\Delta_p = ((p+1)p)^{1/2} \quad (A.1.10)$$

are indicial functions. If the cross product is taken term by term

$$[\bar{IE}_{pqmm'}^{\alpha\beta}(\bar{r})]_r = m \frac{\gamma_p^m \gamma_q^{m'}}{\Delta_p \Delta_q \sin\theta} \left\{ P_p^m(\cos\theta) \frac{d}{d\theta} P_q^{m'}(\cos\theta) + P_q^{m'}(\cos\theta) \frac{d}{d\theta} P_p^m(\cos\theta) \right\} \{ i Z_p^{\alpha*}(\bar{r}) Z_q^{\beta}(\bar{r}) \} e^{i(m-m')\varphi} \quad (A.1.11)$$

$$[\bar{IE}_{pqmm'}^{\alpha\beta}(\bar{r})]_{\theta} = 0 \quad (A.1.12)$$

$$[\bar{IE}_{pqmm'}^{\alpha\beta}(\bar{r})]_{\varphi} = 0 \quad (A.1.13)$$

b) IM type

Again using the normalization and using

$$\bar{N}_{pm}^{\alpha}(\bar{r}) = \bar{\nabla} \cdot \bar{M}_{pm}(\bar{r}) \quad (A.1.14)$$

$$[\bar{N}_{pm}^{\alpha}]_r = i \Delta_p \frac{Z_p^{\alpha}(r)}{r} \gamma_p^m P_p^m(\cos\theta) e^{im\varphi} \quad (A.1.15)$$

$$[\bar{N}_{pm}^{\alpha}]_{\theta} = -\frac{1}{r} \{ \bar{X}_{pm}(\theta, \varphi) \}_{\theta} \frac{d}{dr} [r Z_p^{\alpha}(r)] \quad (A.1.16)$$

$$[\bar{N}_{pm}^{\alpha}]_{\varphi} = \frac{1}{r} \{ \bar{X}_{pm\varphi}(\theta, \varphi) \}_{\varphi} \frac{d}{dr} [r Z_p^{\alpha}(r)] \quad (A.1.17)$$

Then  $[\bar{IM}_{pqmm'}^{\alpha\beta}(\bar{r})]_r = -\frac{\gamma_p^m \gamma_q^{m'}}{r \Delta_p \Delta_q} \left\{ \frac{P_p^m(\cos\theta) P_q^{m'}(\cos\theta)}{\sin^2\theta} + \frac{\frac{dP_p^m(\cos\theta)}{d\theta} \frac{dP_q^{m'}(\cos\theta)}{d\theta}}{\Delta_p \Delta_q} \right\} \{ i Z_p^{\alpha*}(\bar{r}) \frac{d}{dr} Z_q^{\beta}(\bar{r}) \} e^{i(m-m')\varphi} \quad (A.1.18)$

$$[\bar{IM}_{pqmm'}^{\alpha\beta}(\bar{r})]_{\theta} = \frac{\Delta_q \gamma_p^m \gamma_q^{m'}}{\Delta_p r} P_q^m(\cos\theta) \frac{d(P_p^m(\cos\theta))}{d\theta} \{ i Z_p^{\alpha*}(\bar{r}) Z_q^{\beta}(\bar{r}) \} e^{i(m-m')\varphi} \quad (A.1.19)$$

$$[\bar{IM}_{pqmm'}^{\alpha\beta}(\bar{r})]_{\varphi} = m \frac{\gamma_p^m \gamma_q^{m'}}{\Delta_p r \sin\theta} Z_p^{\alpha*}(\bar{r}) Z_q^{\beta}(\bar{r}) P_p^m(\cos\theta) P_q^{m'}(\cos\theta) e^{i(m-m')\varphi} \quad (A.1.20)$$

## APPENDIX II. USAGE OF THE PROGRAM GIVEN IN THE APPENDIX

The given program package is one of the versions(CUBEX) which is used to investigate some standard configurations, for a range of frequencies. Data cards are arranged as follows :

The first card contains a description of the problem with a field of (1011,10X,2D30.15) where the parameters are:

1st integer	Number of surfaces.
2nd integer	0 if the origin is included.
3rd integer	1 if S matrix is required.
4th integer	1 if the program will continue on the pr.
5th integer	1 if the multipole coefficients are req.
6th integer	1 if the program will continue on the pr.
7th integer	1 if $\sigma(\theta, \psi)$ is required.
8th integer	1 if the plot of $\sigma(\theta, 0)$ is required.
9th integer	1 if the plot of $\sigma(0, \psi)$ is required.
10th integer	Number of frequency steps.

Other parameters are the starting frequency and the increment in terms of 4HZ. After the first card blocks for different bodies follow.

Cards for ith body is as follows :

- 1st card(32A1) Literal description of the body, 1/2
- 2nd card(2D3518) Body origin coordinates(cm)( $p=(x^2+y^2), z$ )
- 3rd card(2D3518)  $\beta_{\max}$ ,  $\beta_{\min}$
- 4th card(D35,18,44X,I1) Maximum(cm) size of the body and type of the crosssectional shape(number code as in the subroutine ARGON)

Then the quadrature information follows (if this is the first ensemble to be solved) :

1st Card (I4) Number of quadrature points.

2nd and the others (2D35.18) argument and weights for positive values of the argument only.

If a sequence of problems have to be solved, same type of cards follow, but no quadrature cards are necessary after other ensemble descriptions.

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```

IMPLICIT REAL*8(A-H,O-Z)
REAL*8 QA(9),QB(9),QSW(4)
CDMMDN MQX(4),RSW(4)
COMMNDN FLICK/AA(9),BB(9),TMX(9),TNN(9),ZBCX(4,9),RMAX,RMIN,NSF
EXTERNAL ARGON
1 READ(1,22,END=1453)NSF,ICODE,11,12,13,14,15,16,17,INKA,FSTR,FINC,
2 ((ZBCX(I,J),I=1,4),QA(J),QB(J),TMX(J),TNN(J),MQX(J),
2 J=1,NSF)
22 FORMAT(10I1,10X,2G30.15/(4A8/2G35.18/2G35.18/G35.18,44X,11))
FRAC=FSYR-FINC
DO 30 I=1,INKA
RMAX=0.D00
RMIN=0.D00
IF(ICODE.NE.0)RMIN=25.D00
FRAC=FRAC+FINC
FKAC=10.2D00*3.141592653589793/3.D00)*FRAC
DO 34 J=1,NSF
AA(J)=QA(J)*FKAC
BB(J)=QB(J)*FKAC
34 RSW(J)=QSW(J)*FKAC
CALL VENUS(ARGON,11,12,13,14,15,16,17)
WRITE(3,6)FRAC
6 FORMAT(//T8,3X,'FREQUENCY = ',F12.7,' GHZ')
36 CONTINUE
GO TO 16
453 STOP
END

```

```

SUBROUTINE ARGON(R,RT,COZE,SIZE,ZETA,L$F)
IMPLICIT REAL*8(A-H,D-Z)
COMMON MQX(4),R,SM(4)
BEK=R*SM(L$F)
MET=MQX(L$F)
GO TO (1,2,3,4,5,6,7,8),MET
C----- SPHERE
C----- 1 R=BEK
      RT=0.D00
      RETURN
C----- PROLATE SPHEROID
C----- 2 R=BEK/(2.D00-SIZE)
      RT=R*COZE/(2.D00-SIZE)
      RETURN
C----- VERTICAL EIGHT
C----- 3 R=BEK*(2.D00-COZE**2+SIZE**2)/3.D00
      RT=4.D00*BEK*SIZE*COZE/3.D00
      RETURN
C----- LONG BEAN
C----- 4 CST=DEXP(-SIZE)
      R=BEK*D$IN(CST)
      RT=-BEK*D$OS(CST)*CST*COZE
      RETURN
C----- HORIZONTAL EIGHT
C----- 5 R=BEK*(2.D00+COZE*COZE-SIZE*SIZE)/3.D00
      RT=-4.D00*BEK*SIZE*COZE/3.D00
      RETURN

```

## LIST OF ARGON

&lt; CONTINUED &gt;

PAGE 3

```
      RETURN          ARGO 37
C--- SHORT BEAN    ARGO 38
C--- C---          ARGO 39
C---   6 R=BEK*DSIN(2*D00-SIZE)/3.D00  ARGO 40
C---   RT=-BEK*DCOS(2*D00-SIZE)*COZE/3.D00  ARGO 41
C---   RETURN          ARGO 42
C--- C--- STAR OF FOUR LEAVES  ARGO 43
C--- C---          ARGO 44
C---   7 R=BEK*(2*D00+DCOS((4*D00*ZETA))/3.D00  ARGO 45
C---   RT=-4*D00*BEK*DSIN((4*D00*ZETA))/3.D00  ARGO 46
C---   RETURN          ARGO 47
C--- C---          ARGO 48
C---   8 R=BEK*(2*D00-SIZE)/3.D00  ARGO 49
C---   RT=-BEK*COZE/3.D00  ARGO 50
C---   RETURN          ARGO 51
C--- C---          ARGO 52
C---   END           ARGO 53
C---          ARGO 54
C---          ARGO 55
C---          ARGO 56
```

```

SUBROUTINE VENUS(ILKNUR, Q1, Q2, Q3, Q4, Q5, Q6, Q7)
IMPLICIT COMPLEX*16(A-H),REAL*8(O-Z)
REAL*8 XQ(200), YQ(200), KSI(5), ETA(5), AA(9), BB(9), AABS, BABS,
      COTA, COZE
      P(5), PT(5), PK(5),
      SIFIR(37), DOKSAN(37), SS(2,5,2,2,5,2,2)
      REAL TYPX(2), ADI(2), ADM(3), X0(74), X1(74), Z0(74), Z3(74), AB1(3),
      AB2(3), DRD(3), ARAS1(25), ARAS2(25), STAR(74)
      INTEGER I, TET(37), IPROB/0/, Q1, Q2, Q3, Q4, Q5, Q6, Q7
      COMPLEX*16 ADA(400), S(5,2,2,5,2,2), H1(5), H2(5), B(5,2,2), U(5,2,2),
      US(5,2), SEV, SEB, Z1(5,2), Z2(5,2), Y1(5,2), Y2(5,2)
      DIMENSION ZBCX(4,9)
      DIMENSION TMX(9), TNN(9)
      COMMAND/FLICK/AA,BB,TMX,TNN,ZBCX,RMAX,RMIN,NSF
      EQUIVALENCE (SS(1),S(1)),ADA(1)
      IF(IPROB)104,781,104
      TAB=3.141592653589793
      TAC=TAB/180.D00
      TAB=TAB*2.D00
C-----READ QUADRATURE POINTS
C-----READ(1,101)NQD,(XQ(I),YQ(I),I=1,NQD,2)
      101 FORMAT(14/(2G3.30))
      DO 102 I=2,NQD,2
      XQ(I)=-XQ(I-1)
      102 YQ(I)=YQ(I-1)
C-----PREPARE INDICAL FUNCTIONS
      STRG=4.*D00*3.141592653589793
      DO 103 I=1,5
      KSI(I)=DSQRT((2*I+1)/STRG)
      103 ETA(I)=KSI(I)/DFLOAT(I*I+1)
C-----START OF THE PROBLEM

```

```

C----- 10% CONTINUE
C----- IPROB=IPROB+1
C----- CLEAR UP THE INTEGRATION AREA
C----- DO 106 I=1,400
C----- 106 ADA(I)=0.D00,0.D00
C----- START OF INTEGRATION
C----- DO 110 LSF=1,NSF
      TAF=TAC*(TMX(LSF)-TMN(LSF))/2.D00
      TAK=TAC*(TMX(LSF)+TMN(LSF))/2.D00
      TAM=TAF*TAB
      ZAVL=AA(LSF)
      ZBVL=BB(LSF)
      DO 110 LQD=1,NQD
      ZETA=TAF*XQ(LQD)+TAK
      COZE=DCOS(ZETA)
      SIZE=DSIN(ZETA)
      WEB=TAM*YQ(LQD)
      CALL ILKNUR(AABS,BABS,COZE,SIZE,ZETA,LSF)
      ZZ=AABS*SIZE
      PP=AABS*COZE
      ZZ=ZBVL+ZZ
      PP=PP+ZAVL
      SEL=AABS*PP*WEB
      SEM=BABS*PP*WEB
      RR=DSQRT((ZZ**2+PP**2))
      IF(RR.GT.RMAX)RMAX=RR
      IF(RR.LT.RMIN)RMIN=RR
      COTA=ZZ/RR
      SOTA=PP/RR
      ZCS=COZE*COTA-SIZE*SOTA
      ZSS=COZE*SOTA+SIZE*COTA
      VENU 37
      VENU 38
      VENU 39
      VENU 40
      VENU 41
      VENU 42
      VENU 43
      VENU 44
      VENU 45
      VENU 46
      VENU 47
      VENU 48
      VENU 49
      VENU 50
      VENU 51
      VENU 52
      VENU 53
      VENU 54
      VENU 55
      VENU 56
      VENU 57
      VENU 58
      VENU 59
      VENU 60
      VENU 61
      VENU 62
      VENU 63
      VENU 64
      VENU 65
      VENU 66
      VENU 67
      VENU 68
      VENU 69
      VENU 70
      VENU 71
      VENU 72

```



## LIST OF VENUS

&lt; CONTINUED &gt;

PAGE 7

```

RENGIN=-DFLOAT( I/2+J/2 )
DO 688 K=1,5
DO 689 I=1,5
  B(I,2,1)=(0.000,0.000)
  B(I,2,2)=(0.000,0.000)
  GO TO 309
308 CALL FIDAN(H1,H2,5,RMIN)
  DO 112 I=1,5
    STRG=H2(I)
    B(I,2,1)=(1.000,0.000)
112   B(I,2,2)=H2(I)/STRG
309 CALL FIDAN(H1,H2,5,RMAX)
  DO 113 I=1,5
    STRG=H1(I)
    B(I,1,1)=STRG/H2(I)
113   B(I,1,2)=H2(I)/H1(I)
C----- GENERATE U MATRICE
C----- SEV=(0.000,1.000)
      DO 114 I=1,5
        SEB=SEV
        SEV=SEV*(0.000,1.000)
        VANTUZ=DSQRT((3.141592653589793*(8*I+4)))
        U(I,1,2)=SEB*VANTUZ
        U(I,2,2)=SEB*VANTUZ
        U(I,1,1)=B(I,1,1)*U(I,1,2)
        U(I,2,1)=B(I,1,1)*U(I,2,2)
114   IPAG=0
C----- OUTPUT SECTION
C----- IF(Q1=1)782,783,782
783 DO 115 KW=1,2
      DO 115 IA=1,2

```

```

VENU 109
VENU 110
VENU 111
VENU 112
VENU 113
VENU 114
VENU 115
VENU 116
VENU 117
VENU 118
VENU 119
VENU 120
VENU 121
VENU 122
VENU 123
VENU 124
VENU 125
VENU 126
VENU 127
VENU 128
VENU 129
VENU 130
VENU 131
VENU 132
VENU 133
VENU 134
VENU 135
VENU 136
VENU 137
VENU 138
VENU 139
VENU 140
VENU 141
VENU 142
VENU 143
VENU 144

```



```

DO 119 I=1,2
119 WRITE(3,120)(ADM(I),K=1,2),((U(L,M,I),M=1,2),L=1,5)
      WRITE(3,120)(ADM(3),K=1,2),((U(L,M),M=1,2),L=1,5)
120 FORMAT("//T24,*ALPHA E*,A4,T98,*ALPHA M*,A4,/
      1 (T2,2D25.13,25X,2D25.13))

```

## C BISTATIC CROSS SECTIONS

```

C
766 IF(Q4-1)125,788,125
768 STRG=1.D00/6.283185307179586
      RAA=3.141592653589793/36.D00
      CZERO=(0.D00,-1.D00)
      DO 641 I=1,5
      CNINE=CZERO
      CZERO=CZERO*(0.D00,-1.D00)
      SEV=DCMPLX(STRG *DSQRT(DFLOAT(2*I+1))/DFLOAT(I*(I+1)),0.D00)
      US(I,1)=US(I,1)*SEV*CNINE
      US(I,2)=US(I,2)*SEV*CZERO
      DO 122 I=1,37
      TETA=RAA*DFLOAT(I-1)
      COTA=DCOS(TETA)
      SOTA=DSIN(TETA)
      ITET(I)=3*(I-1)
      CALL BADE(IP,PT,PK,COTA,SOTA,5)
      CNINE=(0.D00,0.D00)
      CZERO=CNINE
      DO 121 K=1,5
      APT0=DCMPLX(PT(K),0.D00)
      APK0=DCMPLX(PK(K),0.D00)
      CZERO=US(K,2)*APK0+US(K,1)*APTO+CZERO
      CNINE=US(K,2)*APTO+US(K,1)*APK0+CNINE
      121 SIFIR(I)=CZERO*DCONJG(CZERO)
      DOKSAN(I)=CNINE*DCONJG(CNINE)
      X0(I)=SIFIR(I)*SOTA
      Z0(I)=SIFIR(I)*COTA
      X1(I)=DOKSAN(I)*SOTA

```

```

73(I)=DOKSAN(I)*COTA
LSF=I+37
X0(LSF)=-X0(I)
X2(LSF)=-X2(I)
Z0(LSF)=Z0(I)
522 Z3(LSF)=Z3(I)
C-- OUTPUT BISTATIC-CROSS SECTIONS
C-- IF(Q5=1)789,790,789
790 IPAG=IPAG+1
      WRITE(3,116)IPRUB,IPAG,RMAX,RMIN,NSF,(I,AA(I),I,BB(I),
1      I,TMX(I),I,TMN(I),I,(ZBCX(J,I),J=1,%),
2      I=1,NSF)
      WRITE(3,123)(ITET(I),SIFIR(I),DOKSAN(I),I=1,37),
3      FORMAT("//18,'THETA<DEG>',10X,'FDR<PHI=0>',10X,'FOR <PHI=90> • /'
4      178,3X,13,7X,D20.10)
789 IF(Q6=1)791,792,791
792 CALL ZEYNEP(X0,Z0,STAR,74,ORD,AB1,ARAS1,L)
791 IF(O7=1)25,794,125
794 CALL ZEYNEP(X1,Z3,STAR,74,ORD,AB2,ARAS2,L)
125 RETURN
DATA TYPX/8HREALIMAG/,ADI/8HIEIEIMIM/,ADM/12HI(1)I(2)$
1      AB1/12H XUNIT/,AB2/12H YUNIT/,ORD/2H
2      ARAS1/100HBSTATIC CROSS-SECTION FOR PHI=0 DEGREE
3      L, VENU 242
4      VENU 243
5      VENU 244
6      VENU 245
END
5 STAR/74*4H****/

```

```

SUBROUTINE ZEYNEP(X,Y,Z,N,ORD,ASB,ARAS,ICHART)
  DIMENSION X(N),Y(N),Z(N),ORD(3),ASB(3),ARAS(25),VAR(5)
  1001 FORMAT(1H2,T3,2A4,1H(,A4,1H))
  1002 FORMAT(1T2,E13.6,1H-)
  1003 FORMAT(1H+,T15,102A1)
  1004 FORMAT(T16,10I(1H),T16,5(1H),19X)
  1005 FORMAT(T13,S(EL3.6,7X),2A4,1H(,A4,1H))
  1006 FORMAT(//T10,6HCHART ,13,1H-,25A4)

C----- SORT ORDINATE VARIABLES IN DESCENDING ORDER
C----- 10 DO 11 I=1,N
      JASMIN=I
      COBRA=Y(I)
      DO 66 J=1,N
        IF(Y(J)-COBRA)66,66,64
  64  COBRA=Y(J)
      JASMIN=J
  66  CONTINUE
      IF(JASMIN-I)70,11,70
  70  Y(JASMIN)=Y(I)
      Y(I)=COBRA
      COBRA=X(JASMIN)
      X(JASMIN)=X(I)
      X(I)=COBRA
      COBRA=Z(JASMIN)
      Z(JASMIN)=Z(I)
      Z(I)=COBRA
  11  CONTINUE
C----- FIND BASE AND AMPLITUDE WIDTHS
C-----  YWIDTH=ABS(Y(1)-Y(N))
      XMIN=X(1)
      XMAX=X(1)
      DO 20 I=1,N

```

```

IF(X(I).GT.XMAX)XMAX=X(I)
IF(X(I).LT.XMIN)XMIN=X(I)
20 CONTINUE
XWIDTH=ABS(XMAX-XMIN)

C-----PREPARATION BEFORE ACTUAL PRINTING
C-----WRITE(3,1001)(ORD(I),I=1,3)
C-----START ACTUAL GRAPHING
C-----XPIITCH=XWIDTH/100.0
C-----YPIITCH=YWIDTH/50.0
C-----IF(XWIDTH.EQ.0.0)XPIITCH=0.01
C-----IF(YWIDTH.EQ.0.0)YPIITCH=0.02
JASMIN=1
DO 30 I=1,51
C-----SINGLE ROW PRINTING
C-----COBRA=Y(I)-(I-1)*YPIITCH
C-----WRITE(3,1002)COBRA
31  IF(ABS(Y(JASMIN)-COBRA).GT.(YPIITCH/2.0))GO TO 30
JESPA=(X(JASMIN)-XMIN)/XPIITCH+1
WRITE(3,1003)(BLANK,LALE=1,JESPA),Z(JASMIN)
IF(JASMIN.EQ.N)GO TO 30
JASMIN=JASMIN+1
GO TO 31
30 CONTINUE
C-----UNDERSCALING
C-----WRITE(3,1004)
DO 41 J=1,5
41  VAR(J)=XMIN+XPIITCH*(J-1)*20
WRITE(3,1005)(VAR(J),J=1,5),(ASB(L),L=1,3)

```

```
WRITE(3,1006)ICHART,TARAS(J),J=1,25)
RETURN
DATA BLANK/LH /
END
```

ZEYN 73  
ZEYN 74  
ZEYN 75  
ZEYN 76

```

SUBROUTINE CAPRI(A,B,N)
COMPLEX*16 A(N,N),B(N),E
REAL*8 DTMP,DMAX
NN=N-1
DO 106 K=1,NN
  KK=K+1
  LMAX=0
  DMAX=0.D00
  DO 102 J=K,N
    DTMP=A(J,K)*DCONJG(A(J,K))
    IF(DTMP-DMAX)102,101,101
    DMAX=DTMP
    LMAX=J
  102 CONTINUE
  103 E=B(LMAX)
  B(LMAX)=B(K)
  B(K)=E
  DO 104 L=K,N
    E=A(LMAX,L)
    A(LMAX,L)=A(K,L)
    104 A(K,L)=E
    DO 105 L=KK,N
      E=-A(L,K)/A(K,K)
      B(L)=B(L)+E*B(K)
    105
    DO 106 J=KK,N
      A(L,J)=A(L,J)+E*A(K,J)
    106 B(N)=B(N)/A(N,N)
    DO 108 L=1,NN
      K=N-L
      KK=K+1
      DO 107 J=KK,N
        B(K)=B(K)-A(K,J)*B(J)
      107
      108 B(K)=B(K)/A(K,K)
    RETURN
  END

```

```

CAPR 1 2
CAPR 2 3
CAPR 3 4
CAPR 4 5
CAPR 5 6
CAPR 6 7
CAPR 7 8
CAPR 8 9
CAPR 9 10
CAPR 10 11
CAPR 11 12
CAPR 12 13
CAPR 13 14
CAPR 14 15
CAPR 15 16
CAPR 16 17
CAPR 17 18
CAPR 18 19
CAPR 19 20
CAPR 20 21
CAPR 21 22
CAPR 22 23
CAPR 23 24
CAPR 24 25
CAPR 25 26
CAPR 26 27
CAPR 27 28
CAPR 28 29
CAPR 29 30
CAPR 30 31
CAPR 31 32
CAPR 32 33
CAPR 33 34
CAPR 34 35
CAPR 35 36

```

```
SUBROUTINE YAPRAK (H,G,N,X)
IMPLICIT COMPLEX*16(A-H),REAL*8 (0-Z)
COMPLEX*16 G(N,2),H(N,2)
SARAY(A)=A
DRILL(A)=DCMPLX(SARAY(A),0.D00)
C=DCMPLX(DSIN(X)/X,DCOS(X)/X)
H(1,2)=C*DCMPLX((1.D00/X,1.D00)
XX=X*X
H(2,2)=C*DCMPLX((3.D00-XX)/XX,3.D00/X)
G(2,2)=C*DCMPLX((XX-2.D00)/X,-2.D00)
G(2,2)=X*H(1,2)-2.D00*H(2,2)
DO 101 I=3,N
H(I,2)=(2*I-1)*H(I-1,2)/X-H(I-2,2)
103 G(I,2)=X*H(I-1,2)-1*H(I,2)
DO 102 I=1,N
H(I,1)=DRILL(H(I,2))
102 G(I,1)=DRILL(G(I,2))
RETURN
END
```

```
1      YAPR
2      YAPR
3      YAPR
4      YAPR
5      YAPR
6      YAPR
7      YAPR
8      YAPR
9      YAPR
10     YAPR
11     YAPR
12     YAPR
13     YAPR
14     YAPR
15     YAPR
16     YAPR
17     YAPR
18     YAPR
19     YAPR
```

```
SUBROUTINE BADE(P,PT,PK,COTA,SOTA,N)
IMPLICIT REAL*8 (A-H,O-Z)
REAL*8 P(N),PT(N),PK(N)
PK(1)=-1.000
P(1)=PK(1)*SOTA
PT(1)=-COTA
PK(2)=-3.000*COTA
P(2)=PK(2)*SOTA
PT(2)=-3.000*(COTA**2-SOTA**2)
DO 101 I=3,N
PK(I)=(2*I-1)*COTA*PK(I-2)/DFLOAT(I-1)
P(I)=PK(I)*SOTA
PT(I)=I*COTA*PK(I)-(I+1)*PK(I-2)
101 RETURN
END
```

```
SUBROUTINE FIDAN(H1,H2,N,X)
IMPLICIT COMPLEX*16(A-H),REAL*8(0-Z)
COMPLEX*16 H1(N),H2(N)
C=DCMPLX(DSIN(X)/X,DCOS(X)/X)
H2(1)=C*DCMPLX(1.000/X,1.000)
XX=X*X
H2(2)=C*DCMPLX((3.000-XX)/XX,3.000/X)
DO 101 I=3,N
  H2(I)=(2*I-1)*H2(I-1)/X-H2(I-2)
DO 102 I=1,N
  H1(I)=DCDNJG(H2(I))
102 RETURN
END
```

```
1 FIDA
2 FIDA
3 FIDA
4 FIDA
5 FIDA
6 FIDA
7 FIDA
8 FIDA
9 FIDA
10 FIDA
11 FIDA
12 FIDA
13 FIDA
```

K\*RMAX = 2.85  
 K\*RMIN = 0.0  
 #OF SF = 2

P1 = 0.0 Z1 = 0.0 BMX1 = 90.00 BMN1 = -90.00 R1 = LONG BEAN OF FOCAL MAXIMA D=6CM.  
 P2 = 0.0 Z2 = 2.21 BMX2 = 90.00 BMN2 = -90.00 R2 = SPHERE OF RADIUS 0.3195 CM

## REAL(IE11)

0.0	0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0	0.0

## IMAG(IE11)

-0.1514943766639D-01	0.1429783590106D-01	0.1271185216963D-01	0.4010127659350D-02	0.1089658140189D-02
0.1429783590106D-01	0.2147050045582D-01	0.1035135984882D-01	0.3734659796018D-02	0.9995905609435D-03
0.1271185216963D-01	0.1035135984882D-01	0.5376395127255D-02	0.1886586604034D-02	0.5066672363029D-03
0.4010127659350D-02	0.3734659796018D-02	0.1886586604034D-02	0.6637752376662D-03	0.1791115347504D-03
0.1089658140189D-02	0.9995905609435D-03	0.5066672363029D-03	0.1791115347504D-03	0.4867195991243D-04

## REAL(IE12)

0.8333626329013D-01	0.2402406149434D-01	0.3306097188328D 00	-0.5685398011349D 01	0.5475951667236D 02
0.2402406149434D-01	-0.2100470069449D-01	-0.5806285292997D-02	0.4222369755642D 00	-0.7958433748937D 01
-0.2773928115717D-02	-0.5806285283654D-02	-0.1915152486241D-01	-0.6398226205365D-02	0.4550824743069D 00
0.1423798852917D-03	-0.4567549649191D-02	-0.6398226157873D-02	-0.1220912329235D-01	-0.3603078173463D-02
-0.1461693695675D-04	-0.1205778331374D-02	-0.2331119220658D-02	-0.3603079267493D-02	-0.8209725159791D-02

## IMAG(IE12)

-0.1514943766639D-01	0.1429783590106D-01	0.1271185216963D-01	0.4010127659350D-02	0.1089658140189D-02
0.1429783590106D-01	0.2147050045582D-01	0.1035135984882D-01	0.3734659796018D-02	0.9995905609435D-03
0.1271185216963D-01	0.1035135984882D-01	0.5376395127255D-02	0.1886586604034D-02	0.5066672363029D-03
0.4010127659350D-02	0.3734659796018D-02	0.1886586604034D-02	0.6637752376662D-03	0.1791115347504D-03
0.1089658140189D-02	0.9995905609435D-03	0.5066672363029D-03	0.1791115347504D-03	0.4867195991243D-04

K\*RMAX = 2.85  
 K\*RMIN = 0.0  
 #OF SF = 2

P1 = 0.0 Z1 = 0.0 BMX1 = 90.00 BMN1 = -90.00 R1 = LONG BEAN OF FOCAL MAXIMA D=6CM.  
 P2 = 0.0 Z2 = 2.21 BMX2 = 90.00 BMN2 = -90.00 R2 = SPHERE OF RADIUS 0.3195 CM

## REAL(IE21)

-0.8333626329013D-01	-0.2402406149438D-01	0.2773928115717D-02	-0.1423798852917D-03	0.1461693695675D-04
-0.2402406149434D-01	0.2100470069449D-01	0.5806285283654D-02	0.4567549649191D-02	0.1205778331374D-02
-0.3306097188328D-00	0.5806285292997D-02	0.1915152486241D-01	0.6398226205365D-02	0.2391119220658D-02
0.5685398011349D-01	-0.4222369755642D 00	0.6398226205365D-02	0.1220912329235D-01	0.3603079267493D-02
-0.5475951667236D-02	0.7958433748937D 01	-0.4550824743069D 00	0.3603078173463D-02	0.8209725159791D-02

## IMAG(IE21)

-0.1514943766639D-01	0.1429783590106D-01	0.1271185216963D-01	0.4010127659350D-02	0.1089658140189D-02
0.1429783590106D-01	0.2147050045582D-01	0.1035135984882D-01	0.3734659796018D-02	0.9995905609435D-03
0.1271185216963D-01	0.1035135984882D-01	0.5376395127255D-02	0.1886586604034D-02	0.5066672363029D-03
0.4010127659350D-02	0.3734659796018D-02	0.1886586604034D-02	0.6637752376662D-03	0.1791115347504D-03
0.1089658140189D-02	0.9995905609435D-03	0.5066672363029D-03	0.1791115347504D-03	0.4867195991243D-04

## REAL(IE22)

0.360137233488D-17	-0.9652669876878D-12	0.3333836469685D 00	-0.5685540391235D 01	0.5475953128930D 02
0.9652669876878D-12	-0.5285485590867D-18	-0.9382926810272D-11	0.4268045252134D 00	-0.7957277970606D 01
-0.3333836469685D 00	0.9342939175951D-11	0.3264416866494D-17	-0.4749158178671D-10	0.4374135935275D 00
0.5685540391235D 01	-0.4268045252134D 00	0.4749158148354D-10	-0.9757819552370D-18	0.1094030710263D-08
-0.5475953128930D 02	0.7957227970606D 01	-0.4574135935275D 00	-0.1094030720671D-08	0.1036768327439D-17

## IMAG(IE22)

0.7033073968588D 00	0.9507193917475D 01	0.1320345297834D 02	0.3536691311666D 03	0.7386570153228D 03
0.5507193917475D 01	0.1789742237258D 02	0.27972387832D 03	0.1419310786827D 04	0.4157704479768D 05
0.1320345297834D 02	0.27972387832D 03	0.1613802951742D 04	0.3624219282889D 05	0.2938912048720D 06
0.3936691311666D 03	0.1419310786827D 04	0.3624219282889D 05	0.3160680670886D 06	0.8964882295077D 07
0.7386570153228D 03	0.4157704479768D 05	0.2938912048720D 06	0.8964882295077D 07	0.1084992196875D 09

K\*RMAX = 2.85  
K\*RMIN = 0.0  
#OF SF = 2

P1 = 0.0 Z1 = 0.0 BMX1 = 90.00 BMN1 = -90.00 R1 = LONG BEAN OF FOCAL MAXIMA 0.6CM.  
P2 = 0.0 Z2 = 2.21 BMX2 = 90.00 BMN2 = -90.00 R2 = SPHERE OF RADIUS 0.3195 CM

## REAL (IM11)

0.0	0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0	0.0

## IMAG (IM11)

-0.11530305230650 00	0.2221913169051D-01	0.4239083144190D-02	0.2532962082813D-03	0.8881900307252D-04
0.2221908186212D-01	-0.9370834645884D-02	-0.1857132447155D-02	-0.1335527926428D-02	-0.4985110690947D-03
0.4239145898031D-02	-0.1857258930323D-02	-0.3023883995218D-02	-0.1312382003318D-02	-0.4315394274939D-03
0.2531342619251D-03	-0.1335412502558D-02	-0.1312369280061D-02	-0.6362804265048D-03	-0.2069860196218D-03
0.8893292746105D-04	-0.4984906097682D-03	-0.431573367424D-03	-0.2069815160399D-03	-0.6908496277827D-04

## REAL (IM12)

-0.1860422167828D-00	-0.2802145716681D 00	0.2136978776086D 01	-0.1438011063478D 02	0.3116336549478D 03
-0.5480574729668D-01	-0.4051367673009D 00	-0.3091678714849D 00	0.1926084173348D 01	-0.3983720109490D 02
-0.1194819438107D-01	-0.1905728269115D-01	-0.4917973232279D 00	-0.222284083427D 00	0.5036634677874D 00
-0.3344728059779D-02	-0.6253047984524D-02	-0.1039988415317D-01	-0.5225042482064D 00	-0.1091677020923D 00
-0.1022786758938D-02	-0.1007319751926D-02	-0.2770476860297D-02	-0.7732560406109D-02	-0.5281715738513D 00

## IMAG (IM12)

-0.1153030523065D 00	0.2221913169051D-01	0.4239083144190D-02	0.2532962082813D-03	0.8881900307252D-04
0.2221908186212D-01	-0.9370834645884D-02	-0.1857132447155D-02	-0.1335527926428D-02	-0.4985110690947D-03
0.4239145898031D-02	-0.1857258930323D-02	-0.3023883995218D-02	-0.1312382003318D-02	-0.4315394274939D-03
0.2531342619251D-03	-0.1335412502558D-02	-0.1312369280061D-02	-0.6362804265048D-03	-0.2069860196218D-03
0.8893292746105D-04	-0.4984906097682D-03	-0.431573367424D-03	-0.2069815160399D-03	-0.6908496277827D-04

K\*RMAX = 2.85  
K\*RMIN = 0.0  
#OF SF = 2

P1 = 0.0 Z1 = 0.0 BMX1 = 90.00 BMN1 = -90.00 R1 = LONG BEAN OF FOCAL MAXIMA 0.6CM.  
P2 = 0.0 Z2 = 2.21 BMX2 = 90.00 BMN2 = -90.00 R2 = SPHERE OF RADIUS 0.3195 CM

## REAL (IM21)

-0.8139977832174D 00	0.5480579683031D-01	0.1194815481529D-01	0.3544761561030D-02	0.1022729869745D-02
0.2802143524676D 00	-0.5948632327041D 00	0.1905734356467D-01	0.6252885342125D-02	0.1007273100750D-02
-0.2136982424725D 01	0.3091672503726D 00	-0.5082026768062D 00	0.1039983903245D-01	0.2771094407364D-02
0.1438173688640D 02	-0.1926104870595D 01	0.2225267919921D 00	-0.4776957505829D 00	0.7731378051206D-02
-0.3117809034364D 03	0.3984977363765D 02	-0.5037669675560D 00	0.1091666058201D 00	-0.4718288373380D 00

## IMAG (IM21)

-0.1153030523065D 00	0.2221913169051D-01	0.4239083144190D-02	0.2532962082813D-03	0.8881900307252D-04
0.2221908186212D-01	-0.9370834645884D-02	-0.1857132447155D-02	-0.1335527926428D-02	-0.4985110690947D-03
0.4239145898031D-02	-0.1857258930323D-02	-0.3023883995218D-02	-0.1312382003318D-02	-0.4315394274939D-03
0.2531342619251D-03	-0.1335412502558D-02	-0.1312369280061D-02	-0.6362804265048D-03	-0.2069860196218D-03
0.8893292746105D-04	-0.4984906097682D-03	-0.431573367424D-03	-0.2069815160399D-03	-0.6908496277827D-04

## REAL (IM22)

-0.1000000000000D 01	-0.2254087748378D 00	0.2148926930910D 01	-0.1437676587314D 02	0.31163467776774D 03
0.2254086051709D 00	-0.1000000000000D 01	-0.2901105279272D 00	0.1934337098690D 01	-0.3983619378180D 02
-0.2148930619106D 01	0.2901099676814D 00	-0.1000000000003D 01	-0.2121269078389D 00	0.5064345421650D 00
0.1437839215835D 02	-0.1934357918580D 01	0.2121269078389D 00	-0.9999999987892D 00	-0.1014363239811D 00
-0.3117819262232D 03	0.39844766317900 02	-0.5065374444163D 00	0.10143404541400 00	-0.10000000411189D 01

## IMAG (IM22)

0.1860073464703D 01	0.1159524446553D 02	0.1248490933672D 03	0.1020639606649D 04	0.1896725495714D 05
0.1159532180408D 02	0.179424202233D 03	0.1677667466909D 04	0.2719685892675D 05	0.3363707551712D 06
0.1248448866019D 03	0.1677689186155D 04	0.3104144250627D 05	0.4301813838488D 06	0.9046023886456D 07
0.1020785542331D 04	0.2719643542169D 05	0.4301787121597D 06	0.9700679397287D 07	0.1791213562201D 09
0.1896476590716D 05	0.3363588219928D 06	0.9046584490880D 07	0.1791066609198D 09	0.4793349164486D 10

K\*RMAX = 2.85  
 K\*RMIN = 0.0  
 #OF SF = 2

P1 = 0.0 Z1 = 0.0 BMX1 = 90.00 BMN1 = -90.00 R1 = LONG BEAN OF FOCAL MAXIMA 0.6CM.  
 P2 = 0.0 Z2 = 2.21 BMX2 = 90.00 BMN2 = -90.00 R2 = SPHERE OF RADIUS 0.3195 CM.

## ALPHA EI(1)

0.3197070616155D 01 0.7392384296625D 01  
 -0.7809811162733D-01 0.6879934629353D 01  
 -0.992852824664D 01 -0.1636249095216D 00  
 -0.1023739141618D 01 -0.1055376138671D 02  
 0.8881010660451D 01 0.3850310344752D 00

## ALPHA MI(1)

0.1731658526074D 01 0.50941289551600 01  
 -0.8053437291969D 01 0.1869696759079D-01  
 -0.1077678726296D 00 -0.8799956841751D-01  
 0.1062530973881D 02 0.9203233943632D 00  
 0.3079156737492D 00 0.1423234804515D 02

## ALPHA EI(2)

0.1533944477862D 01 -0.3875255791874D 01  
 -0.1056056973400D 00 0.3458112875710D 00  
 -0.5941902723464D-02 0.1076829120492D-01  
 0.5222635079649D-03 -0.1642453755156D-02  
 -0.7184307222938D-05 0.3255204178401D-04

## ALPHA MI(2)

-0.7621971261342D 00 0.1776364468920D 00  
 -0.4629084937890D-01 0.1321003173645D 00  
 0.1046777349550D-01 -0.1189904684915D-01  
 -0.1748436701605D-03 0.1675412229315D-04  
 -0.6359091147730D-05 0.1094767343692D-04

## ALPHA ES

-0.1417053830480D 01 0.3711345886073D 01  
 0.1903476160144D 00 -0.6521133810749D 00  
 0.7981278143519D-02 -0.1463825634363D 00  
 -0.5911831227160D-02 -0.4313287271262D-01  
 -0.1937758826673D-02 -0.1331700803173D-01

## ALPHA MS

0.9348766484838D 00 -0.876737382927D 00  
 -0.1948216949821D 00 -0.1121587663840D 00  
 -0.1415191663032D 00 0.2254085626165D-01  
 -0.4027394978120D-01 0.1333839804747D-02  
 -0.1142695278500D-01 0.146496896829D-02

K\*RMAX = 2.85  
 K\*RMIN = 0.0  
 #OF SF = 2

P1 = 0.0 Z1 = 0.0 BMX1 = 90.00 BMN1 = -90.00 R1 = LONG BEAN OF FOCAL MAXIMA 0.6CM.  
 P2 = 0.0 Z2 = 2.21 BMX2 = 90.00 BMN2 = -90.00 R2 = SPHERE OF RADIUS 0.3195 CM.

## THETA&lt;DEG&gt;

## FOR&lt;PHI=0&gt;

## FOR&lt;PHI=90&gt;

0	0.3035762581D 00	0.3035762581D 00
5	0.3003537909D 00	0.3035772163D 00
10	0.2908064324D 00	0.3032241824D 00
15	0.2752988026D 00	0.3029388182D 00
20	0.2544498749D 00	0.3027905449D 00
25	0.2291380866D 00	0.3029532423D 00
30	0.2004887500D 00	0.3035568553D 00
35	0.1698321377D 00	0.3046435813D 00
40	0.1386266253D 00	0.3061380278D 00
45	0.1003505796D 00	0.3078388796D 00
50	0.8037660347D-01	0.3094363273D 00
55	0.3584997017D-01	0.3105550895D 00
60	0.3559336084D-01	0.3108179206D 00
65	0.2006082912D-01	0.3099200205D 00
70	0.9349239538D-02	0.3077018505D 00
75	0.3267818479D-02	0.3042074486D 00
80	0.1452195674D-02	0.2997176151D 00
85	0.3500734283D-02	0.2947516793D 00
90	0.9106150369D-02	0.2900366654D 00
95	0.1815577908D-01	0.2865472211D 00
100	0.3077901548D-01	0.2849228529D 00
105	0.4732968871D-01	0.2863708710D 00
110	0.6830323647D-01	0.2915646078D 00
115	0.9430173796D-01	0.3010475115D 00
120	0.1233716120D 00	0.31050565973D 00
125	0.1618461718D 00	0.33356266601D 00
130	0.2032257412D 00	0.3557704753D 00
135	0.2486208764D 00	0.3811693084D 00
140	0.29667063000D 00	0.4085329063D 00
145	0.3456315139D 00	0.4365965354D 00
150	0.3935183261D 00	0.4640288921D 00
155	0.4302694095D 00	0.4895463899D 00
160	0.4779081338D 00	0.5120020796D 00
165	0.5106787315D 00	0.5304379020D 00
170	0.5351451864D 00	0.5441169852D 00
175	0.5502528336D 00	0.5525248083D 00
180	0.5553603362D 00	0.5553603362D 00

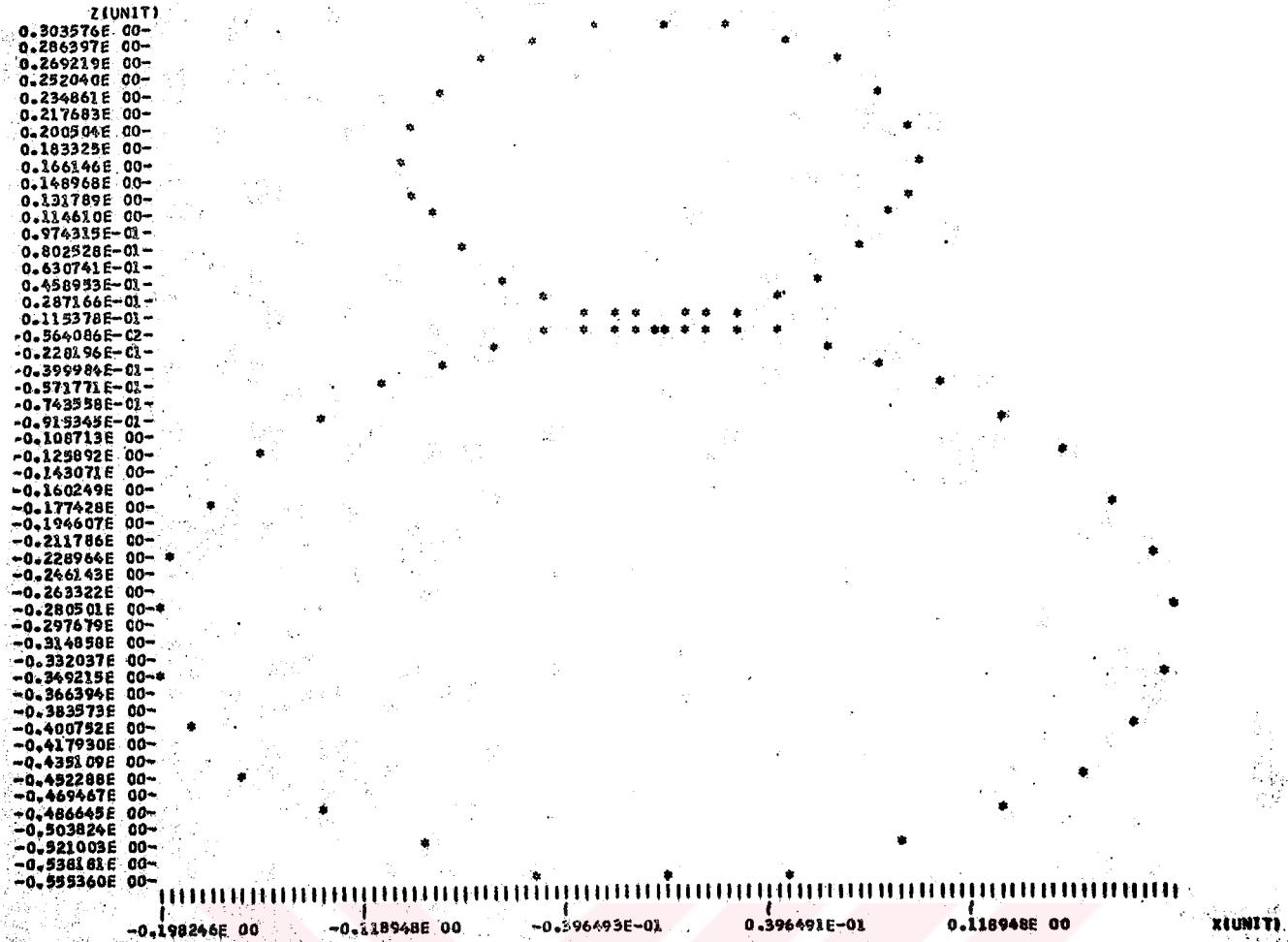


CHART 1-BISTATIC CROSS-SECTION FOR PHI=0 DEGREE

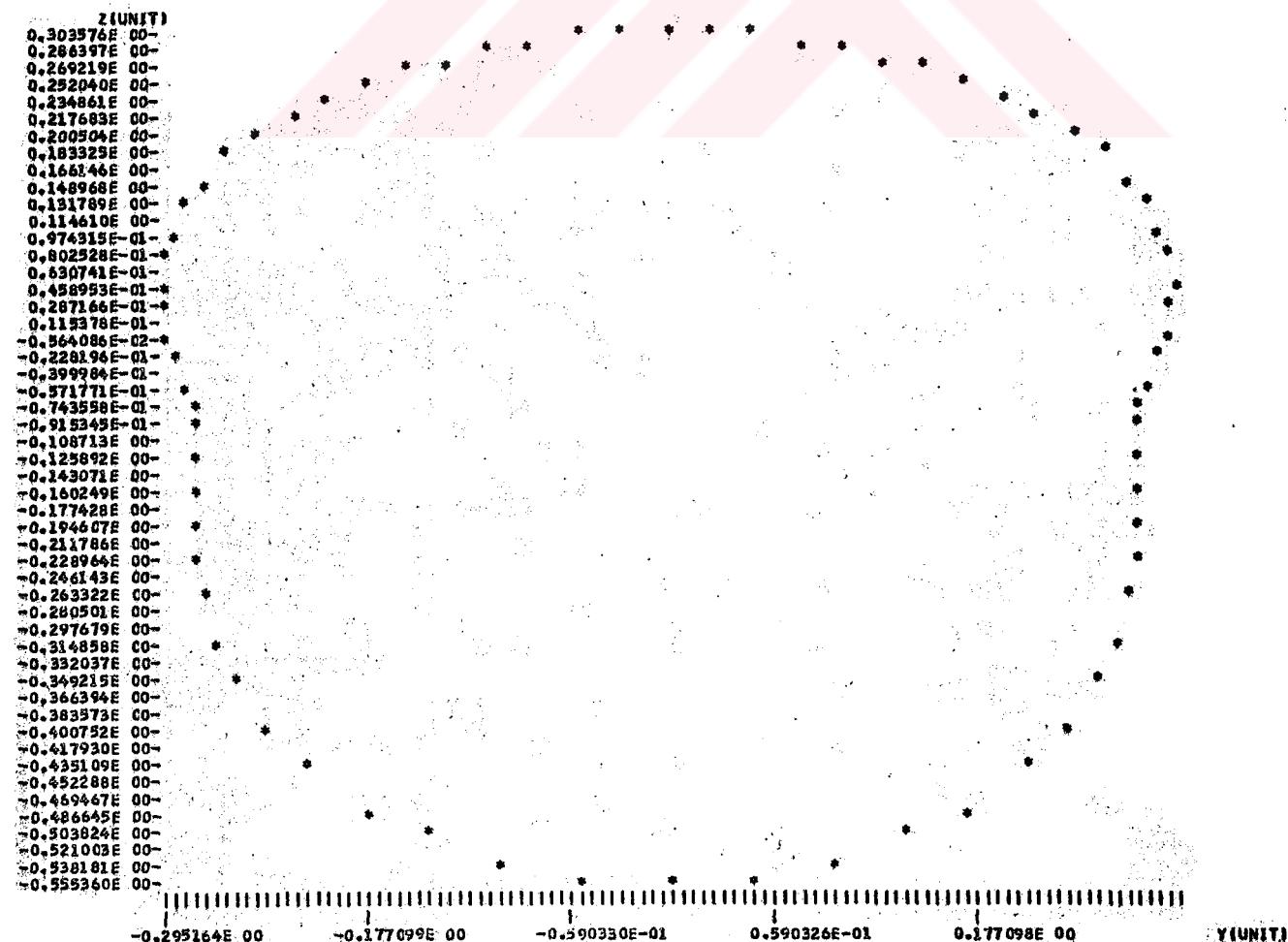


CHART 2-BISTATIC CROSS-SECTION FOR PHI=90 DEGREES