

SUBROUTINES FOR COMPUTING THE PARAMETERS OF THE
ELECTROMAGNETIC RADIATION SCATTERED BY A SPHERE

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TABLE OF CONTENTS

I.	Introduction	1
II.	Theory	6
	2.1 Stokes Parameters	6
	2.2 Expressions for Mie Scattering	11
III.	A Second Look at the Mie Expressions	17
	3.1 a_n and b_n	17
	3.2 A_n and W_n	20
	3.3 π_n and τ_n	25
IV.	Details about the Subroutines	29
	4.1 General	29
	4.2 Timing	32
	4.3 Storage Requirements	34
	4.4 Accuracy	34
	4.5 Sample Problem	35
V.	Conclusion	36
	Acknowledgments	37
	References	38
	Footnote	40
Appendix I:	Listing of Sample Problem	41
Appendix II:	Listing of DAMIE Subroutine	44
Appendix III:	Listing of DBMIE Subroutine	50
Appendix IV:	Sample Output using DAMIE; $m = 1.342 - 0.0i$, $r = 100\mu$, $\lambda = 0.4\mu$	56
Appendix V:	Sample Output using DBMIE; $m = 1.342 - 0.1i$, $r = 100\mu$, $\lambda = 0.4\mu$	61

I. INTRODUCTION

The passage of the electromagnetic radiation through a medium is generally accompanied by the removal of a fraction of the energy from the incident beam. This fraction may get partly absorbed within the medium, and may get partly scattered, i.e., reappear in same as well as other directions with or without change in frequency. Even when the scattering process does not involve any change in frequency (a case to which this report is restricted), the characteristics of the scattered radiation are determined by the wavelength (λ) of the incident radiation, refractive index ($m = n_1 - i n_2$) of the medium, and size as well as the shape of the discrete particles in the medium. Because of this, measurements and proper interpretation of the characteristics of the scattered radiation offer a very good opportunity for obtaining information about the state of the medium. Hence, numerical determination of the characteristics of the scattered radiation for a given model is of prime importance in several diversified fields such as planetary and atmospheric optics, astrophysics, laser, radar and searchlight applications, and chemistry. To this end, the first step is the evaluation of the characteristics of the radiation scattered by a single particle.

The numerical evaluation of the parameters of the radiation scattered by a particle very small compared to λ is very simple and straightforward (Rayleigh, 1871). The expressions for the radiation scattered by a sphere whose radius (r) is comparable to, or greater than, λ were first derived by Mie (1908) and as such,

this process of scattering is known as Mie scattering. The expressions for Mie scattering involve series whose terms contain Bessel functions of half-integral order (spherical Bessel function) with complex argument, and, first and second derivatives of the Legendre polynomials. Furthermore, the number of terms required for evaluating the series with reasonable accuracy is of the order of the size parameter $x = 2\pi r/\lambda$. Hence, the reliable Mie scattering computations for large spheres can be described as difficult, tedious and time consuming.

The derivations of the Mie expressions as well as a catalog of its available numerical solutions in some specific cases can be found in the treatise, "Light Scattering by Small Particles" by Van de Hulst (1957). Prior to this publication, numerical evaluation of the Mie expressions was confined to small values of x (< 10) for partly-absorbing ($n_2 > 0$) spheres. For non-absorbing spheres ($n_2 = 0$), the most outstanding work is that of Gumprecht and Slepcevich (1951, 1953), who carried their calculations to $x = 400$. As has been pointed out by Van-de Hulst, the work of these authors cannot be considered to be complete c. the laws of geometric and physical optics can be applied to the transfer problem in a sphere with some reasonable confidence only if x is of the order of 2000. Furthermore, the number of directions for which the radiation parameters are given by these authors is too small to give any insight into the fine structure of the radiation field. Besides, many fields of application demand extensive numerical data for values of refractive indices other than those used by Gumprecht and Slepcevich.

Since 1957, some notable progress has been made in extending the range of x for partly-absorbing spheres (e.g., Giese, 1961; Delmendjion, Clasen and Viezee, 1961; Irvine, 1963, 1965; Cunningham, Wells and Collins, 1966; Plass, 1966; and Kattawar and Plass, 1967). However, surprisingly enough, no results have been reported for large spheres ($x > 400$) which are partly-absorbing or non-absorbing (see Footnote on p. 40).

This absence of progress may be attributed partly to the limited interest of various workers. However, since Mie computations involve evaluation of series whose terms increase by several orders of magnitude with increase of n and change sign in rather uneven manner, it is essential to carry out basic arithmetic in double precision if x is large. The other factor is the large storage requirements. Computations of the radiation scattered in 37 different directions [0 (5) 180] by a sphere having size parameter $x = 1000$ with straightforward programming procedure can require use of more than one hundred thousand double precision words of storage, a requirement too large to be fulfilled by the main storage area of any modern computer. Hence, one is required to use magnetic tapes or disks.

The main difficulty in evaluation of the characteristics of the radiation scattered by a large, partly-absorbing sphere is the very rapid propagation of the errors when the function $A_n(mx)$ is computed using a recurrence relationship (sec. 3.2). A recurrence relationship provides a very powerful computing tool, especially in automatic work. However, since generation is carried out perforce

with rounded values, the errors may or may not grow relative to the size of the wanted function. If the errors do grow, the recurrence scheme is said to be unstable. When the upward recurrence scheme (i.e., starting with the value of $A_0(mx)$, one computes successively higher values of $A_n(mx)$ by making use of recurrence formulae) is unstable, in general the downward recurrence scheme is found to be very stable (Abramowitz and Stegun, 1964, p. XIII). It is not necessary to know the initial value for starting the downward recurrence scheme; one can make use of the ratio method first described by Miller (1952). Reference may also be made of an independent later work in this direction by Corbató and Uversky (1959), who have applied the downward recurrence procedure for generation of spherical Bessel functions on digital computers. Kattawar and Plass (1967) seem to be the first to use this downward recurrence procedure in Mie scattering calculations.

The purpose of this report is to describe and to present two FORTRAN subroutines with the help of which one can evaluate the characteristics of the electromagnetic radiation scattered by a sphere, in as many as 200 different directions. The subroutines are written using FORTRAN IV H language and IBM System/360 computers in mind.

In the first subroutine called DAMIE, all the functions are computed using the upward recurrence procedure and as such the storage requirement is very modest (15474 bytes for any value of the size parameter). However, because of

the use of upward recurrence, the scheme becomes unstable for $A_n(mx)$ when $n_2 > 0$, i.e., the spheres are partly-absorbing. From the discussion in sec. 3.2, it will be clear that further work is needed in understanding the problem of the error propagation in the case of function $A_n(mx)$. The only positive statement which can be made at this stage, is as follows. The final results obtained using DAMIE should not be accepted without checking them against those obtained using DBMIE if the parameter $n_2 x$ is about 50 or higher, or x is of the order of 100 or more.

In the second subroutine called DBMIE, the function $A_n(mx)$ is computed using downward recurrence procedure and hence, all its values have to be stored for further use. This does not lead to any significant increase in computer time but the storage requirement is very high. Since the recurrence is started with

$$A_N(mx) = 0.0 + 0.0i \quad (1)$$

where

$$N = 1.1(n_1^2 + n_2^2)^{1/2}x + 1, \quad (2)$$

a storage of 127340 bytes can accomodate x of about 20 if $n_1 = n_2 = 200$, a case for which some results have been presented by Van de Hulst (1957, p. 293). However, the results obtained using DBMIE are expected to be more reliable than those obtained with DAMIE under all foreseeable conditions.

II. THEORY

2.1 Stokes Parameters:

A beam of electromagnetic radiation is vectorial in nature as it requires more than one (actually four) parameters to fully describe its state. These parameters, as customarily understood, are as follows:

1. Specific Intensity (I): It is the amount of energy flowing normally through unit area in a cone of unit solid angle per unit time and per unit frequency interval. It is also referred to as Intensity, Radiance or Brightness.
2. Degree of Polarization (P): As a Nicol prism (analyser) is rotated around the axis of observation, the intensity of the source as seen through the prism, undergoes variations especially if the source consists of some scattered light, or alternately, if the beam happens to pass through some polarizing medium before reaching the Nicol prism. If this occurs, the intensity passes through a maximum value (I_{\max}) and a minimum value (I_{\min}) twice during a 360° rotation. The positions of I_{\max} and I_{\min} are at right angles to each other. The degree of polarization of the observed radiation is given by

$$P = \frac{I_{\max} - I_{\min}}{I_{\max} + I_{\min}} . \quad (3)$$

3. Deviation of Plane of Polarization (X): For several reasons, it may be convenient to refer to the directions of I_{\max} and that of I_{\min} with respect to two arbitrary directions (say e and r) which are mutually at right angles to each other. The $e-r$ plane is assumed to be at right angles to the direction of observation. If the direction of polarization, i.e., that of I_{\max} , makes an angle X , say with that of e , then the plane of polarization is said to deviate by an angle X .
4. Ellipticity of Polarization (Δ): The phenomenon of polarization is best understood by imagining two vibrations of complex amplitudes E_{\max} and E_{\min} along the direction of I_{\max} and I_{\min} respectively. Then, $I_{\max} = |E_{\max}|^2$ and $I_{\min} = |E_{\min}|^2$ if the constant of proportionality is assumed to be unity. If these vibrations maintain a constant phase difference (Δ) between them over millions of vibrations, the radiation is said to be elliptically polarized. This restriction (millions of vibrations) is necessary because one vibration takes of the order of 10^{-8} seconds and any available instrument averages over a period of the order of a second. The detection of this ellipticity requires the use of a wave plate (an optical component which introduces some arbitrary phase difference between the two vibrations) and also the Nicol prism.

This representation of the state of a beam of electromagnetic radiation by the customary parameters (I , P , X and Δ) though complete and easily understood, leads

to unwanted complications when one is interested, e.g., in obtaining the results of interaction of several beams on the same particle. This is because these customary parameters are of different dimensions and are non-additive. This difficulty can be bypassed by making use of the following alternate set of parameters first proposed by Stokes (1852):

In the Stokes representation, the intensity I is treated as a one-column matrix of four elements (or alternately, a vector with four elements). All the elements have the same dimensions and are also additive.

$$I = \begin{vmatrix} I \\ Q \\ U \\ V \end{vmatrix} \quad (4)$$

For the definition of these elements, let us assume two directions, e and r , at right angles to each other, such that the err plane is perpendicular to the direction of the propagation of the radiation under study. Let us further assume that the complex amplitudes of electrical vibrations are E_e^* and E_r^* , respectively, along the e - and r -directions. Their complex conjugates will be denoted by E_e and E_r . Again assuming the constant of proportionality to be unity, the four elements are defined as follows:

$$\left. \begin{aligned} I &= E_e E_e^* + E_r E_r^*, \\ Q &= E_e E_r^* - E_r E_e^*, \\ U &= E_e E_r + E_r E_e, \\ V &= I(E_e E_r^* - E_r E_e^*) \end{aligned} \right\} \quad (5)$$

and

This matrix, though symmetric, is less simple than the following somewhat less symmetric matrix which we shall also refer to as the Stokes matrix:

$$I = \begin{vmatrix} I_e \\ I_r \\ I_u \\ I_v \end{vmatrix} \quad (6)$$

where the elements of the two matrices are related as follows:

$$\left. \begin{aligned} I &= I_e + I_r, \\ Q &= I_e - I_r, \\ U &= I_u, \\ V &= I_v \end{aligned} \right\} \quad (7)$$

and

The following relationships can be shown to exist between the Stokes and the customary parameters of the beam of electromagnetic radiation (e.g., Chandrasekhar, 1950 ; Van de Hulst, 1957) :

$$P = \frac{\left[(I_e - I_r)^2 + I_u^2 + I_v^2 \right]^{1/2}}{I_e + I_r}, \quad (6)$$

and

$$\tan 2X = \frac{I_u}{I_e - I_r} \quad (7)$$

In order to relate β to the Stokes parameters, we have to define an angle θ given

by

$$\sin 2\beta = \frac{I_v}{\sqrt{(I_e - I_r)^2 + I_u^2 + I_v^2}}, \quad (8)$$

then

$$\tan \beta = \frac{\tan 2\beta}{\sin 2X} \quad (9)$$

If there is no elliptical polarization, i.e., $\beta = 0$, the parameter $I_v = 0$.

A partially polarized beam of radiation can then be represented by three parameters,

I_e , I_r , and I_u . If the direction of maximum should coincide with that of e or r , i.e., $X = 0^\circ$ or 90° , then $I_u = 0$ and

$$P = \frac{|I_e - I_r|}{|I_e + I_r|}. \quad (10)$$

A natural beam of radiation which is completely unpolarized has for its Stokes parameters: $1/2$, $1/2$, 0 , 0 .

2.2 Expressions for Mie Scattering

The expressions for the radiation scattered by a sphere of radius (r), and of material with a complex index of refraction (m), have been aptly derived by Van de Hulst (1957), and also by Born and Wolf (1964). We shall therefore enumerate the final results only.

Let I_i and I_s respectively represent the Stokes parameters of the radiation incident on and scattered by a sphere. Then,

$$I_s = F' \cdot I_i \quad (11)$$

where F' is a four-by-four matrix referred to as a "transformation matrix" by Van de Hulst (1957, p. 44). It has the following form

$$F' = \begin{vmatrix} M_2 & 0 & 0 & 0 \\ 0 & M_1 & 0 & 0 \\ 0 & 0 & S_{21} & -D_{21} \\ 0 & 0 & D_{21} & S_{21} \end{vmatrix}. \quad (12)$$

The matrix E' and hence the matrix \underline{S} are the functions of the following parameters:

$$x = 2\pi/\lambda \quad , \text{ where } \lambda \text{ is the wavelength of the incident radiation;}$$

$$\mu = n_1 / (n_1 + n_2) \quad , \text{ index of refraction of the material of the sphere with respect to its surroundings;}$$

and $\theta = \text{the angle between the direction of the incident and the scattered radiation.}$

- Van de Hulst has also noted that only three elements of E' are independent, the interrelationship being

$$S_{21}^2 + D_{21}^2 = M_2 M_1 \quad . \quad (15)$$

However, as S_{21} and D_{21} can be positive or negative, no useful purpose is served by using this relationship in actual computations.

In order to evaluate these elements, one first defines the complex amplitudes $S_1(x, m, \theta)$ and $S_2(x, m, \theta)$ for the scattered radiation.

$$S_1(x, m, \theta) = \sum_{n=1}^{\infty} \frac{(2n+1)}{n(n+1)} \left\{ a_n(x, m) \pi_n(\mu) + b_n(x, m) \tau_n(\mu) \right\} \quad (16)$$

and

$$S_2(x, m, \theta) = \sum_{n=1}^{\infty} \frac{(2n+1)}{n(n+1)} \left\{ b_n(x, m) \pi_n(\mu) + a_n(x, m) \tau_n(\mu) \right\} \quad (17)$$

where

$$\mu = \cos \theta \quad . \quad (18)$$

Then,

$$M_1 = S_1 S_1^* \quad ,$$

$$M_2 = S_2 S_2^* \quad ,$$

$$S_{21} = \pm (S_2 S_1^* + S_1 S_2^*) \quad ,$$

$$\text{and} \quad D_{21} = \frac{i}{2} (S_2 S_1^* - S_1 S_2^*) \quad .$$

The functions $a_n(x, m)$ and $b_n(x, m)$ are given by

$$a_n(x, m) = \frac{\tau_n(mx) \tau_n(x) - m \tau_n(mx) \tau_n'(x)}{\tau_n(mx) \tau_n(x) + m \tau_n(mx) \tau_n'(x)} \quad , \quad (20)$$

and

$$b_n(x, m) = \frac{m \tau_n'(mx) \tau_n(x) - \tau_n(mx) \tau_n'(x)}{m \tau_n(mx) \tau_n(x) + \tau_n(mx) \tau_n'(x)} \quad , \quad (21)$$

$\gamma_n'(z)$ and $\varepsilon_n'(x)$ are the respective derivatives of $\gamma_n(z)$ and $\varepsilon_n(x)$ with respect to z and x respectively.

$$\gamma_n'(z) = z j_n(z) \quad (22)$$

$$\varepsilon_n'(x) = x [j_n(x) - i \gamma_n(x)] \quad (23)$$

$$\gamma_n'(z) = z j_{n+1}(z) + n j_n(z) \quad (24)$$

$$\text{and } \varepsilon_n'(x) = x [j_{n+1}(x) - i \gamma_{n+1}(x)] - n [j_n(x) - i \gamma_n(x)] \quad (25)$$

j_n and γ_n are the spherical Bessel functions of the first and second kind, respectively. The following recurrence relationships and initial values for setting up the recurrence procedure can be found in any of the standard mathematical books (e.g., Antosiewicz, 1964).

$$f_{n+1}(z) = \frac{2n+1}{z} f_n(z) - f_{n-1}(z) \quad (26)$$

$$[f_n(z); j_n(z), \gamma_n(z)]$$

$$j_{-1}(z) = -\gamma_0(z) = \frac{\cos z}{z} \quad (27)$$

and

$$j_0(z) = \gamma_{-1}(z) = \frac{\sin z}{z} \quad (28)$$

The phase functions $\pi_n(\mu)$ and $\tau_n(\mu)$ appearing in eqs. (16) and (17) can be expressed in terms of the Legendre polynomials as follows:

$$\pi_n(\mu) = \frac{d P_n(\mu)}{d \mu} \quad (29)$$

and

$$\tau_n'(\mu) = \mu \pi_n(\mu) - (1-\mu^2) \frac{d \pi_n(\mu)}{d \mu} \quad (30)$$

Besides the transformation matrix F^* , other terms of considerable importance are the dimensionless constants referred to as "efficiency factors" by Van de Hulst (1957, p. 14 and p. 127). Q_e , the efficiency factor for extinction, which is the ratio of the total amount of energy removed from the incident beam to the geometric cross-section (πr^2) of the particle, can be obtained from the values of a_n and b_n only.

$$Q_e(x, m) = \frac{2}{\pi} \sum_{n=1}^{\infty} (2n+1) [\operatorname{Re}(a_n) + \operatorname{Re}(b_n)] \quad (31)$$

The symbol Re stands for the real part of the quantity in parentheses. Q_s , the efficiency factor for scattering, is given by the following expression:

$$Q_s(x, m) = \frac{2}{\pi} \sum_{n=1}^{\infty} (2n+1) [\|a_n\|^2 + \|b_n\|^2] \quad (32)$$

If there is no absorption, i.e., $n_2 = 0$, $Q_e = Q_s$. Otherwise, Q_a ,

the efficiency factor for absorption, is given by

$$Q_a = Q_e - Q_s. \quad (33)$$

Another dimensionless quantity of considerable interest is the so-called "asymmetry factor" represented by $\overline{\cos \theta}$.

$$\overline{\cos \theta} = \frac{4}{\pi^2 Q} \sum_{n=1}^{\infty} \left\{ \frac{n(n+2)}{n+1} \operatorname{Re} \left[a_n^{*} a_{n+1} + b_n^{*} b_{n+1} \right] + \frac{2n+1}{n(n+1)} \operatorname{Re} \left[a_n b_n^{*} \right] \right\} \quad (34)$$

As noted by Irvine (1963), and again by Kottawur and Plass (1967) (but not by Giese (1961)), asterisks appearing in eq. (34) have been omitted in Van de Hulst (1957).

Q_p , the efficiency factor for radiation pressure, can then be obtained using the following:

$$Q_p = Q_e - \overline{\cos \theta} + Q_s \quad (35)$$

III. A SECOND LOOK AT THE MIE EXPRESSIONS

Since the evaluation of the MIE amplitudes S_1 and S_2 can lead to the computations of a thousand or more terms by recurrence, it is advisable to give some thought to the question of the propagation of errors in the case of each of the four functions.

3.1 a_n and b_n

The expressions for the functions a_n and b_n (eqs. (20) and (21)) for complex values of the parameter m contain the spherical Bessel functions $j_n(mx)$ which in turn require hyperbolic sine and hyperbolic cosine functions $\sinh(n_2 x)$ and $\cosh(n_2 x)$. As is well-known, the values of these hyperbolic functions increase exponentially with $n_2 x$. Since the computers deal only with finite numbers, the evaluation of the expressions in the present form can result in an overflow if $n_2 x \sim 170$ and IBM System/360 computers are used.

This overflow can be avoided by dividing both numerator and denominator of the expression for a_n as well as that of b_n by $j_n(mx)$ as all the terms contain $j_n(mx)$ or $j_{n+1}(mx)$. Infeld (1947) who seems to be the first to recognize this, introduced the so-called logarithmic derivative of $\Psi_n(mx)$ and $\Xi_n(x)$ functions given by eqs. (22) and (23), respectively. Some of the properties of these logarithmic derivative functions have been discussed by Aden (1951). The function

$$\frac{d[\log \Upsilon_n(mx)]}{d(mx)} = \frac{\frac{d}{dx} \Upsilon_n(mx)}{\Upsilon_n(mx)} \quad (36)$$

is denoted as $\sigma_n(mx)$ by Aden (1951), as $A_n(mx)$ by Deirmendjian and Clasen (1962), and as $D_n(mx)$ function by Kattawar and Plass (1967). Using the notations of Deirmendjian and Clasen (1962), we have,

$$a_n(x, m) = \frac{\left\{ \frac{A_n}{m} + \frac{n}{x} \right\} \operatorname{Re}[W_n(x)] - \operatorname{Re}[W_{n-1}(x)]}{\left\{ \frac{A_n}{m} + \frac{n}{x} \right\} W_n(x) - W_{n-1}(x)} \quad , \quad (37)$$

and

$$b_n(x, m) = \frac{\left\{ m A_n(mx) + \frac{n}{x} \right\} \operatorname{Re}[W_n(x)] - \operatorname{Re}[W_{n-1}(x)]}{\left\{ m A_n(mx) + \frac{n}{x} \right\} W_n(x) - W_{n-1}(x)} \quad , \quad (38)$$

The logarithmic derivative of $\Upsilon_n(mx)$, i.e., the function $A_n(mx)$, has the following form:

$$A_n'(mx) = -\frac{n}{mx} + \frac{j_{n-1}(mx)}{j_n(mx)} = -\frac{n}{mx} + \frac{1}{\frac{n}{mx} - A_{n-1}(mx)} \quad , \quad (39)$$

The following initial value for setting up the recurrence can be obtained after making use of eqs. (27) and (28):

TABLE 1

Notations as used by Deirmendjian and Clasen (1962), and by Kattawar and Plass (1967).

Deirmendjian and Clasen	Kattawar and Plass
$\operatorname{Re}[W_n(x)]$	$\Upsilon_n(x)$
$W_n(x)$	$\sigma_n(x)$
$A_n(mx)$	$D_n(mx)$
$\frac{\operatorname{Re}[W_{n-1}(x)]}{\operatorname{Re}[W_n(x)]} - \frac{n}{x}$	$D_n(x)$
$\frac{W_{n-1}(x)}{W_n(x)} - \frac{n}{x}$	$G_n(x)$

$$A_n(mx) = \frac{J_{n-1}(mx)}{J_0(mx)} = \cot(mx) + \dots \quad (40)$$

The function $W_n(x)$ appearing in eqs. (37) and (38) is the same as $E_n(x)$ defined by eq. (22). Hence, after making use of eqs. (26) to (28), we have

$$W_n(x) = \frac{2n+1}{x} W_{n-1}(x) - W_{n-2}(x) \quad , \quad (41)$$

with

$$W_{-1}(x) = \cos x - i \sin x \quad , \quad (42)$$

and

$$W_0(x) = \sin x + i \cos x \quad . \quad (43)$$

At the first sight, modified expressions of $a_n(x, m)$ and $b_n(x, m)$ as given by Kattawar and Ploss (1967) appear to be different from those given by Deirmendjian and Clasen (1962). This is not really the case and one can check their equivalence by making use of Table 1.

3.2 A_n and W_n :

The initial value for setting up the recurrence procedure for computations of $A_n(mx)$ is given by eq. (40). With $m = n_1 - i n_2$, it can be written in either of the following forms:

$$A_0(mx) = \frac{\sin(2n_1x) + i \sinh(2n_2x)}{\cosh(2n_2x) - \cos(2n_1x)} \quad , \quad (44)$$

or

$$A_0(mx) = \frac{\sin(n_1x) \cos(n_1x) + i \sinh(n_2x) \cosh(n_2x)}{\sin^2(n_1x) + \sinh^2(n_2x)} \quad . \quad (45)$$

If $n_2x = 0$ and n_1x is an integer multiple of π , either of these forms will give rise to a division by zero and hence, a possible termination of the program, or wrong results. This is not a hypothetical case, as several situations can be visualized where $2n_1x/\pi$ can be an integer. Because of the generation of the round-off errors in a particular way, the expression for $A_0(mx)$ given by eq. (45) was found to be more suitable for computational purposes. However, it will be necessary to apply due caution if $2n_1x/\pi$ is very close to an integer, since round-off errors are generated differently in different computers.

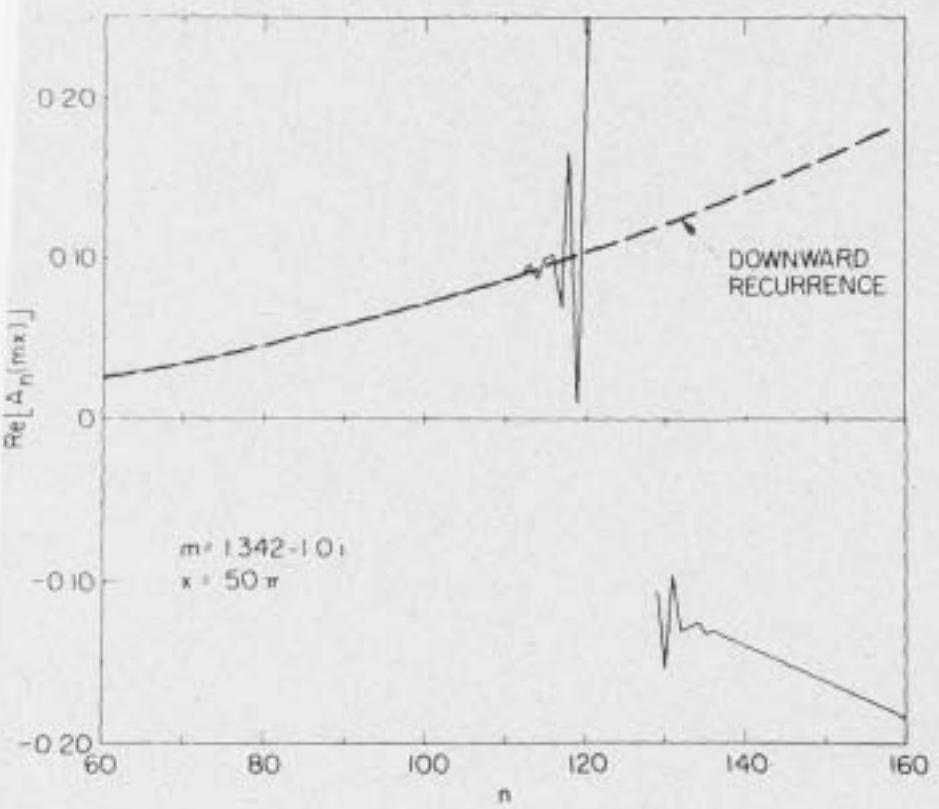
If $n_2 \neq 0$, a straightforward evaluation of $A_0(mx)$ will again run into trouble if $n_2x \geq 90$ or more. Since for these large values of n_2x , $\sinh(n_2x) \sim \cosh(n_2x) \sim \pm e^{n_2x}$, we may make use of the following approximation:

$$A_0(mx) = 0.0 + 1.0 i \quad (46)$$

A use of this approximation immediately leads to the question of propagation of errors, as the successive values of $A_n(mx)$ are computed by upward recurrence

FIGURE 1

Variation of real part of $A_n(mx)$ as a function of n , $m = 1,342 - 1.0 i$, $x = 50.0 \pi$. Broken curve represents values obtained using downward recurrence procedure. Thin solid curve represents values obtained using upward recurrence procedure.



[Eq. (39)]. As mentioned earlier (sec. 1), Kattawar and Plass (1967) have already pointed out that a numerical instability develops in this case. For setting up the downward recurrence procedure, they have suggested the use of $0.0 + 0.0 i$ as the initial value of $A_N(mx)$ for some value of $N > |mx|$. Since the computational error decreases very rapidly at each step of downward recurrence in this particular case, the calculations are insensitive to the assumed starting value and converge very rapidly to the correct value. The criterion for selecting downward recurrence at a value of N given by eq. (2) is rather strict.

The real part of the function $A_n(mx)$ for $m = 1,342 - 1.0 i$ and $x = 50.0 \pi$ as obtained after using the upward recurrence (thin solid curve) and downward recurrence (broken curve) procedures, is plotted in Figure 1 as a function of n . It can be seen that a numerical instability develops in the upward recurrence scheme when $n = 111$. Between 112 and 125, large oscillations develop. The amplitude of these oscillations decreases when n is about 130. For $n = 136$ and higher, a stabilization occurs, but the real part of $A_n(mx)$ achieves a value which is completely different from the correct value given by the broken curve.

The ultimate effect of this instability is to increase the values of the efficiency factor Q_3 [eq. (32)]. Eventually, Q_3 becomes greater than Q_e leading to negative values for Q_g [eq. (33)]. An example of this is provided in Table 2 where the values of Q_g as obtained using the upward recurrence procedur

TABLE 2

The values of the efficiency factor for absorption (Q_a) as a function of n_2 : $m = 1.342 - n_2 i$, $r = 10i$, $k = 0.4 \mu$, $x = 50.0 \pi$.

Column 2: Values obtained when function $A_n(mx)$ is computed using upward recurrence procedure (DAMIE).

Column 3: Values obtained when function $A_n(mx)$ is computed using downward recurrence procedure (DBMIE).

n_2	Q_a with DAMIE	Q_a with DBMIE
0.0	0.0000	0.0000
0.0001	0.0535	0.0535
0.001	0.4149	0.4149
0.01	0.9649	0.9649
0.1	0.9653	0.9653
0.2	0.9542	0.9542
0.3	0.9389	0.9390
0.4	0.4913	0.9211
0.5	-0.4925	0.9016
0.6	-0.5265	0.8808
0.7	-1.1592	0.8592
0.8	-2.3079	0.8369
0.9	-1.4379	0.8141
1.0	-1.3789	0.7910

(subroutine DAMIE), and as obtained using the downward recurrence procedure (subroutine DBMIE) for computations of $A_n(mx)$, are tabulated for several values of n_2 . The assumed values of n_1 and x are 1.342 and 50.0π , respectively. The values of Q_a as obtained from these subroutines agree in four significant figures for n_2 up to 0.2. For $n_2 = 0.3$, a difference of one unit is noticeable in the fourth significant place. Thereafter, the difference increases very rapidly, and for $n_2 = 0.5$, the subroutine DAMIE returns negative values for Q_a .

A definite guideline for prohibiting the use of DAMIE subroutine cannot be easily established as the amount of absolute error propagated depends both upon $n_2 x$ and x . It depends upon x because the number of terms required for the convergence of the series depends upon x . In the above example, where $x = 157.08$, the results of DAMIE and DBMIE show significant differences for $n_2 i > 47.1$. On the other hand, for another case for which m was assumed to be $1.70 - 1.84 i$, the results obtained with DAMIE and DBMIE were found to agree within six significant figures for x up to 40.0, i.e., $n_2 x = 73.6$.

The computations of the function $W_n(x)$ which always has a real argument, are straightforward. The problem of the propagation of error is not serious if all the basic arithmetic is done in double precision.

3.3 $\pi_n(u)$ and $\tau_n(u)$

The phase functions $\pi_n(u)$ and $\tau_n(u)$ given by eqs. (29) and (30) can be com-

after making use of the following recurrence relationship based on the properties of the Legendre polynomials and their derivatives:

$$\pi_n(\mu) = \frac{2n+1}{n+1} \mu \pi_{n-1}(\mu) + \frac{n}{n+1} \pi_{n-2}(\mu) , \quad (47)$$

$$\text{and } \tau_n(\mu) = \mu [\pi_n(\mu) - \pi_{n-2}(\mu)] - (2n+1)(1-\mu^2) \pi_{n-1}(\mu) + \tau_{n-2}(\mu) , \quad (48)$$

where

$$\left. \begin{aligned} \pi_0(\mu) &= 0 \\ \pi_1(\mu) &= 1 \\ \tau_0(\mu) &= 0 \\ \tau_1(\mu) &= \mu \end{aligned} \right\} \quad (49)$$

and

The straightforward upward recurrence poses no serious problem. The absolute error increases by a factor of about 10^6 as values $\pi_{3000}(\mu)$ and $\tau_{3000}(\mu)$ are obtained from those of $\pi_0(\mu)$, etc., given by eq. (49). Since this is a loss of six to seven significant figures, it is again necessary to carry out basic arithmetic in double precision.

The values of $\pi_n(1)$ and $\tau_n(1)$ obtained after making use of eqs. (47), (48), and (49), can be checked against those obtained from the following very simple expressions for the same (Van de Hulst, 1957):

$$\tau_n(1) = \tau_n(1) = \frac{1}{2} n(n+1) \quad (50)$$

For $\mu = 90^\circ$, i.e., $\mu = 0$, it can be shown that

$$\left. \begin{aligned} \pi_n(0) &= 0 \quad (\text{if } n \text{ is even}) \\ \text{and} \\ \pi_n(0) &= \frac{2(-1)^{(n-1)/2}}{\Gamma(\frac{n}{2}+1)} \quad (\text{if } n \text{ is odd}) \end{aligned} \right\} \quad (51)$$

The values of gamma functions for large values of n can be found in one of the tables in the series of Mathematical Tables published by the U. S. National Bureau of Standards (1951).

If the values of $\pi_n(\mu)$ and $\tau_n(\mu)$ for $\mu > 0$ have been already computed, it is not necessary to compute those of $\pi_n(-\mu)$ and $\tau_n(-\mu)$ as was done by Gumprecht and Slepcevich (1951). Instead, one can make use of the following relationships:

$$\pi_n(-\mu) = (-1)^{n-1} \pi_n(\mu) \quad (52)$$

and

$$\tau_n(-\mu) = (-1)^n \tau_n(\mu) . \quad (53)$$

These relationships can be derived easily by making use of the following well-known properties of the Legendre polynomials:

$$P_n(-\mu) = (-1)^n P_n(\mu) \quad ; \quad (\mu > 0) \quad (54)$$

and

$$P'_n(-\mu) = (-1)^{n+1} P'_n(\mu) \quad , \quad (\mu > 0) \quad (55)$$

where P'_n represents the derivative of P_n with respect to μ .

However, it appears that the above mentioned properties have gone unnoticed in this particular field. Van de Hulst (1957) has not given them explicitly, even though he does make use of this relationship for $\mu = -1$ at several places. He has not commented on the tables of Gumprecht and Slepcevich (1951) who have unnecessarily extended the length of the tables. The later researchers (e.g., Giese, 1961; Deirmendjian and Clasen, 1962; Giese, et al., 1962; Cunningham, et al., 1966; and Kattawar and Ploss, 1967) do not refer to this relationship either. The advantages of making use of these relationships should be more evident after studying the published data of various authors, and after recognizing the ease of programming for a supplementary angle (appendices II and III).

IV. DETAILS ABOUT THE SUBROUTINES

4.1 General

Two subroutines, called DAMIE and DBMIE, are written with the purpose of helping the user to obtain parameters of the radiation scattered by a sphere of a given size. The actual FORTRAN statements (FORTRAN IV language) are listed in appendices II and III respectively. A survey of the literature showed that the user is likely to be interested in computing the radiation field for the value of the size parameter $x (= 2\pi/\lambda) \sim 10^3$, and possibly for about the same number of directions of scattering. Such demands, along with the necessity of carrying out all computations in double precision, are bound to tax the storage capacities of any computer. In order to avoid any use of tapes or disks, it was decided to use the storage very sparingly. As mentioned in earlier sections, the main difference between the two subroutines is the manner in which one of the functions is evaluated. In DAMIE, the function $A_n(mx)$ is evaluated using an upward recurrence scheme, while in DBMIE a downward recurrence procedure has been used.

The CALL statement for either subroutine requires the following parameters in the order listed:

x : size parameter of the scattering sphere, $x > 0$.

n_1 : real part of the refractive index of the material of the sphere;
 $n_1 > 1.0$ because the present subroutines are not tested for values of $n_1 \leq 1.0$;

TABLE 3

Some symbols as used in the text and their equivalence in DAMIE and DBMIE subroutines.

TEXT	SUBROUTINES
x	X
n_1	RFR
n_2	RFI
π	THETD(J)
Q_e	QEXT
Q_s	QSCAT
$\frac{\cos \pi}{\cos \pi} + Q_s$	CTBRQS
$M_2^{(1)}$	ELTRMX (1,J,1)
$M_1^{(2)}$	ELTRMX (2,J,1)
$S_{21}^{(\pi)}$	ELTRMX (3,J,1)
$D_{21}^{(\pi)}$	ELTRMX (4,J,1)
$M_2^{(180 - \pi)}$	ELTRMX (1,J,2)
$M_1^{(180 - \pi)}$	ELTRMX (2,J,2)
$S_{21}^{(180 - \pi)}$	ELTRMX (3,J,2)
$D_{21}^{(180 - \pi)}$	ELTRMX (4,J,2)
m	RF
w_{n-2}, w_{n-1}, w_n	WM1, WFN(1), WFN(2)
A_{n-1}, A_n	ACAP(1), ACAP(2); ACAP(N) in DBMIE.
a_n, a_{n-1}	FNA, FNAP
b_n, b_{n-1}	FNB, FNBP
$\tau_{n-2}, \tau_{n-1}, \tau_n$	PI(1,J), PI(2,J), PI(3,J)
$\tau_{n-2}, \tau_{n-1}, \tau_n$	TAU(1,J), TAU(2,J), TAU(3,J)

n_2 : imaginary part of the refractive index of the material of the sphere,

$n_2 \geq 0.0$;

π : Angles between the directions of the scattered radiations and direction of the incident radiation for which computations are required; π should be entered in degrees and its value should not exceed 90.0° .

JX: Total number of π 's. Its value should not exceed 100 unless the dimensions in all related statements are appropriately changed. It must be greater than one.

It should be noted that the subroutines do not check for the conditions for x, n_1 and n_2 . If the conditions are not satisfied, the subroutines are likely to return meaningless results.

The following quantities are returned by the subroutines: Q_e : eq. (31); Q_s : eq. (32); $\frac{\cos \pi}{\cos \pi} + Q_s$: eq. (34); and the elements of the transformation matrix F (eq. 14) for all values of π as well as $180 - \pi$ (Table 3).

For the proper functioning of the subroutines, it is necessary to declare the following as REAL * 8 or as DOUBLE PRECISION in the main program:

X, RFR, RFI, QEXT, QSCAT, CTBRQS, THETD(100), and ELTRMX (4,100,2).

After studying the subroutines in conjunction with Table 3, the following should be evident:

1. For a given n , the values of the functions w_n, π_n, a_n and b_n are stored only if they are required at a later stage.
2. The above statement also applies to the function A_n which is computed in DAMIE by upward recurrence.

3. The real and imaginary parts of a complex quantity are obtained by means of EQUIVALENCE.
4. Storage for the elements of the transformation matrix (which are the quantities computed at the very end) is first utilized to store the values of the complex amplitudes S_1 and S_2 .
5. All the series are terminated when $|a_n|^2 + |b_n|^2 < 10^{-14}$.

4.2 Timing:

Average time t , in seconds, which the subroutine DAMIE takes to return the values of the four elements of the transformation matrix E for 182 values of θ , as well as the values of Q_e , Q_i and $\overline{\cos \theta} \cdot Q_i$, is given in Table 4 for several values of x . The computing facility used was an IBM System/360 Model 50 with FORTRAN IV H compiler at OPT = 0.

From the data presented in Table 4, it can be seen that the subroutine becomes more efficient with the increase of x in the range 1.0 to 100.0, since a ten-fold increase in x does not result in a ten-fold increase in t . It may be noted that the most time is consumed in computations of a_n and b_n , and hence, the subroutine will become less and less efficient as the computations are called for fewer and fewer values of θ .

Because of the use of downward recurrence procedure, the average time t for the subroutine DBMIE depends upon the values of n_1 , n_2 , and x . However, in several cases for which the timing tests were run, DBMIE was found to take on the average of 10 to 20% more time than DAMIE.

TABLE 4

Average time (t in seconds) which the subroutine DAMIE takes to return values of four elements of the transformation matrix for 182 values of θ , as well as values of Q_e , Q_i and $\overline{\cos \theta} \cdot Q_i$ for a sphere having size parameter x . Computing facility used: IBM System/360 Model 50, FORTRAN IV H compiler OPT = 0.

x	t in seconds
1.0	2.0
10.0	6.5
100.0	39.5
1000.0	350.

The output given in the appendices IV and V was obtained using FORTRAN IV G compiler, OPT = 0. This compiler seems to be more efficient for these subroutines as the computations for $x = 1571$ take about 300 seconds only.

4.3 Storage Requirements:

DAMIE : 15,474 bytes

DBMIE : 127,340 bytes

4.4 Accuracy:

Because of the nature of the problem, it is not possible to make a positive statement about the reliability of a given value. An overall figure of reliability of the first six to seven significant figures seems to be quite reasonable.

It may be added that these subroutines were used to compute the characteristics of the radiation field for several cases for which the published as well as unpublished numerical values were available, (e.g., Gumprecht and Slepcevich, 1953; Van de Hulst, 1957; Giese, 1961; Diermendjian and Clasen, 1962; Giese, et al., 1962; Irvine, 1963; Cunningham, et al., 1966; Pless, 1966; and Kattawar and Pless, 1967). As mentioned earlier, most of these results were obtained using single precision arithmetic and x up to 400 only. The results obtained using these subroutines and those obtained by the earlier authors were found to agree within the limits claimed by the various authors.

For a large sphere made of material with a refractive index of $1.342 + 0.0i$, Q_s , the efficiency factor for scattering, can be calculated from the following "best" formula suggested by Van de Hulst (1957, pp. 264, 265).

$$Q_s = 2.0 - 7.680 \sin(0.684x) - 1.84x^{-(2/3)} + \text{"ripple"}$$

(56)

For $x = 500.0 \pi$ (a case for which the results are presented in Appendix IV),

the second term is zero, and hence Q_s can be computed using the remaining three terms only. The value of Q_s so obtained is $2.0136 + \text{"ripple"}$, which compares very well with the value of 2.0129 given in the Appendix.

4.5 Sample Problem:

The FORTRAN statements for a sample problem with which one calls the DAMIE subroutine are listed in Appendix I. The same listing can be used for calling DBMIE after changing the CALL DAMIE ----- to CALL DBMIE ----- in the statement numbered DSMIE 036.

The sample problem program reads in the values of n_1 , n_2 , x , JX (number of values for which computations are required), AJX (interval $\Delta\theta$) and MMM. All the values except that of integer JX should be punched on the card in double precision format. If the last quantity (MMM) is equal to zero, the value of x remains unchanged. If $MMM \neq 0$, x is changed to $-2\pi x/(MMM \times 10^{-3})$. The TIME subroutine called at stages DSMIE 035 and DSMIE 039 helps to read the clock in milliseconds. If this subroutine is not available, these two statements, as well as the statements numbered DSMIE 019, DSMIE 040, and DSMIE 065 can be omitted.

Besides other related information and the quantities Q_e , Q_s , Q_o and $\cos(\theta)$, the sample program also prints out the values of all the four elements of the matrix F as well as the intensity and the degree of polarization of the radiation scattered by a sphere illuminated with natural light.

V. CONCLUSION

The problem of numerical evaluation of the characteristics of the electromagnetic radiation scattered by a sphere (Mie scattering) has been analyzed in great detail in previous sections. It has been shown that the question of numerical instability (Kattawar and Plass, 1967) when computing one of the several functions by upward recurrence is an important one indeed if the sphere is partly-absorbing, and is of moderate size (size parameter $x = 100$ or more). Following this, two subroutines are presented.

The subroutine DBMIE is expected to give reliable results under all foreseeable conditions. On the other hand, the subroutine DAMIE is expected to give reliable results only when the moderate and large spheres are made of non-absorbing material. For a partly-absorbing sphere, the results obtained using DAMIE can be expected to be reliable only if the absorption and the size parameters are of moderate magnitude ($n_2^a = 80$, $x = 100$). The present series of tests are insufficient to lay down a more specific criterion for limiting the use of DAMIE. Apart from providing a concise subroutine which can be used in several cases with great ease, the other reason for providing DAMIE is to permit the user to check the reliability of some of the vast amount of published and unpublished numerical data on this subject.

These subroutines are fairly efficient from the point of view of the storage requirements and timing considerations. However, since they are basically general purpose subroutines, more efficient versions can be obtained after appropriate changes. One may be able to relax some of the stringent criteria in some limited applications, and after checks in extreme cases, one may be able to carry out all the basic arithmetic in single precision.

In the end, it should be added that though these subroutines are checked in several alternate ways, they should be still be treated as experimental ones. Due precaution must be exercised before accepting some unusual results in new cases.

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Very recently, Fahlen and Bryant (1968) have reported some results for the radio-frequency back-scattered by a large sphere ($\chi = 3000.0$).

Listing of Sample Problem

Fahlen, T.S. and H. C. Bryant, Journ. Opt. Soc. Amer., Vol. 58, 1968,
pp. 304-310.

APPENDIX II

Listing of DAMIE Subroutine

```

TITLE: REAL PART OF FNPBP, TBL111: IMAGINARY PART OF FNPBP,
FNPBP C: FNPBP AND THE PRECEDING VALUES OF FNA C: FNB RESPECTIVELY;
EQUIVALENT: TBL111, TBL111, TBL111, TBL111, TBL111
EQUIVALENT: FNAP, TBL111, TBL111, TBL111
ID: C 28 100, 130 7 GO TO 20
WHILE1661:
WRITE1661
CALL1661
END IF = UCLMPLEX(REAL,=RFL1)
REAL = 1.000/END
REAL = 1.000/R
REAL = REAL * R
GO TO 4 = 1.00
D 1 TBL111,23 = 0.000 + THETBL111 = DABSETHLDE23
D 1 TBL111,23 = 0.000 + GO TO 23
STHETBL111 = 1.000
S12HTBL111 = 0.000
GO TO 33
C 23 TBL111,23 = 0.000 + GO TO 25
TBL111,23 = 3.1415926535897932 * THETBL111+0.000
STHETBL111 = 0.000111111
S12HTBL111 = 1.000 - S12HTBL111**2
GO TO 32
C 25 TBL111,23 = 0.000 + GO TO 26
S12HTBL111 = 0.000
STHETBL111 = 1.000
GO TO 32
26 WHILE1661: THETBL111
WRITE1661
CALL1661
END IF = UCLMPLEX(REAL,=RFL1)
REAL = 1.000
END IF = UCLMPLEX(TBL111,-RFL1)
TBL111,23 = 0.000
PEI2,23 = 1.000
TAU2,23 = 0.000
TAU2,23 = S12HTBL111
33 LUMTBL111
TBL111 = UCLMPLEX
TBL111 = UCLMPLEX(TBL111,-RFL1)
S12HTBL111 = UCLMPLEX(TBL111,TBL111)
S12HTBL111 = REAL * S12HTBL111 - S12HTBL111
TBL111 = REAL * R
D 1 TBL111,23 = 0.000 + GO TO 40
TBL111 = 3.1400 * UCLMPLEX(TBL111)
S12HTBL111 = 0.000111111
TBL111 = TBL111 + TBL111
TBL111 = TBL111 - TBL111

```

DANIE049
 DANIE050 0057 T111 = T121***2
 DANIE051 0058 T121 = T111 * T121
 DANIE052 0059 T111 = T121
 DANIE053 0060 T111 = HFR * R
 DANIE054 0061 T111 = JAIN1***1
 DANIE055 0062 F111 = HOU1***1
 DANIE056 0063 T111 = T111 + T111***2
 DANIE057 0064 T111 = T111 + T111
 DANIE058 0065 ACAP111 = DUMPLITE111, T121/T111
 DANIE059 0066 GU T1 53
 DANIE060 0067 40 ACAP111 = DUMPLAT 0.000,1.000 1
 DANIE061 0068 WRITE111, T111
 DANIE062 0069 WRITE11101
 DANIE063 0070 30 ACAP111 = - HFR + TELGU1***1 - ALAP1111
 DANIE064 0071 T111 = ACAP111 + HFR + RR
 DANIE065 0072 T111 = ALAP111 + RR + RR
 DANIE066 0073 --> FRA = T111***1 - TAC111***1, CLEMENT111 - MEL111
 DANIE067 0074 FRA = T111***1 - TAC111***1 + MIN111 - MEL111
 DANIE068 0075 FRA = T111
 DANIE069 0076 FRA = T111
 DANIE070 0077 T111 = 1.5000
 DANIE071 L FROM REAL TO THE STATEMENT NUMBER NO1, ELEMENT111,2,31 HAS THE
 DANIE072 L FOLLOWING MEANING:
 DANIE073 L ELEMENT111,2,31 REAL PART OF THE FIRST COMPLEX AMPLITUDE.
 DANIE074 L ELEMENT111,2,31 IMAGINARY PART OF THE FIRST COMPLEX AMPLITUDE.
 DANIE075 L ELEMENT111,2,32 REAL PART OF THE SECOND COMPLEX AMPLITUDE.
 DANIE076 L ELEMENT111,2,32 IMAGINARY PART OF THE SECOND COMPLEX AMPLITUDE.
 DANIE077 L K = 1,2 FOR THE T111 AND K = 2,1 FOR T121,0 - INC1011
 DANIE078 L 08 INITIATION OF THE COMPLEX AMPLITUDE1 VAN DE HOEST,F-125.
 DANIE079 0110 T111 = T111 * T111
 DANIE080 0111 T111 = T111 * T111
 DANIE081 0112 T111 = T111 * T111
 DANIE082 0113 T111 = T111 * T111
 DANIE083 0114 DI 0 4 = 1.128
 DANIE084 0115 ELTHR111,2,11 = T1111 * P112,21 + T1111 * TAU12,21
 DANIE085 0116 ELTHR111,2,11 = T1121 * P112,21 + T1111 * TAU12,21
 DANIE086 0117 ELTHR111,2,11 = T1111 * P112,21 + T1111 * TAU12,21
 DANIE087 0118 ELTHR111,2,11 = T1111 * P112,21 + T1121 * TAU12,21
 DANIE088 0119 ELTHR111,2,21 = T1121 * P112,21 - T1111 * TAU12,21
 DANIE089 0120 ELTHR111,2,21 = T1121 * P112,21 - T1121 * TAU12,21
 DANIE090 0121 ELTHR111,2,21 = T1121 * P112,21 - T1111 * TAU12,21
 DANIE091 0091 50 CONTINUE
 DANIE092 0072 QSL1 = 2.000 + 1. T1111 + T11111
 DANIE093 0073 QSL1 = T1111***2 + T1121***2 + T1111***2 + T1121***2,0,7500
 DANIE094 0074 CTM425 = 0.300
 DANIE095 0075 R = 1
 DANIE096 0076 05 T111 = 286 - 1

```

T131 = 4 - 3
T132 = 2 * 3 + 1
m1 = T132/3 = 1.333
P133=J1 = ((11114*E12/2)*LSTH(J1)) *NP131,J1/E12
T134=J1 = -(11112*P131,J1-P131,J1-T131*5/12*H12*P12*J1)
L1=T134,J1
CONTINUE
WNL = NP131,J1
WNL1 = WNL*Z1
WNL2 = 11118*WNL*NP131,J1 = WNL
ACAP123 = ALAP123
ACAP123 = -1.4 * WNL + 11.000/INWNL*ACAP123
FC1 = ACAP123*WNL + INWL
FC2 = ALAP123*WNL + INWL
PN1 = T131*T131-T133*T133*NP132 = -NP131,J1
PN2 = T131*T131-T133*T132*NP132 = -NP131,J1
PN3 = 0
PN4 = T131*T131*NP132
PN5 = T131*T131*NP132 + 1.000*NP132
LT132 = WNL*NP131,J1 + T132*NP132*Z1 + T133*NP131,J1 + T134*NP132*Z1 + T135*NP131,J1 + T136*NP132*Z1 + T137*NP131,J1 + T138*NP132*Z1
LT133 = -WNL + T131*T132*NP132
LT134 = T131*T132 + T132*T132*Z1 + T133*T132
NP131,J1 = WNL + PN1 * P131
T131 = T131*T131
P = T132*Z1
H1 = H1*Z1
H2 = H2*Z1
H3 = H3*Z1
LT135 = LT135*NP131,J1 + T135*T135*NP132 + T136*T135*Z1 + T137*T135*NP132 + T138*T135*Z1
LT136 = LT136*NP131,J1 + T136*T136*NP132 + T137*T136*Z1 + T138*T136*NP132
LT137 = LT137*NP131,J1 + T137*T137*NP132 + T138*T137*Z1 + T139*T137*NP132
LT138 = LT138*NP131,J1 + T138*T138*NP132 + T139*T138*Z1 + T140*T138*NP132
LT139 = LT139*NP131,J1 + T139*T139*NP132 + T140*T139*Z1 + T141*T139*NP132
LT140 = LT140*NP131,J1 + T140*T140*NP132 + T141*T140*Z1 + T142*T140*NP132
NP132 = 0
LT141 = LT141*NP131,J1 + T141*T141*NP132 + T142*T141*Z1 + T143*T141*NP132 + T144*T141*Z1
LT142 = LT142*NP131,J1 + T142*T142*NP132 + T143*T142*Z1 + T144*T142*NP132
LT143 = LT143*NP131,J1 + T143*T143*NP132 + T144*T143*Z1 + T145*T143*NP132
LT144 = LT144*NP131,J1 + T144*T144*NP132 + T145*T144*Z1 + T146*T144*NP132
NP133 = 0
LT145 = LT145*NP131,J1 + T145*T145*NP132 + T146*T145*Z1 + T147*T145*NP132 + T148*T145*Z1
LT146 = LT146*NP131,J1 + T146*T146*NP132 + T147*T146*Z1 + T148*T146*NP132
LT147 = LT147*NP131,J1 + T147*T147*NP132 + T148*T147*Z1 + T149*T147*NP132
LT148 = LT148*NP131,J1 + T148*T148*NP132 + T149*T148*Z1 + T150*T148*NP132
LT149 = LT149*NP131,J1 + T149*T149*NP132 + T150*T149*Z1 + T151*T149*NP132
LT150 = LT150*NP131,J1 + T150*T150*NP132 + T151*T150*Z1 + T152*T150*NP132
NP134 = 0
LT151 = LT151*NP131,J1 + T151*T151*NP132 + T152*T151*Z1 + T153*T151*NP132 + T154*T151*Z1
LT152 = LT152*NP131,J1 + T152*T152*NP132 + T153*T152*Z1 + T154*T152*NP132
LT153 = LT153*NP131,J1 + T153*T153*NP132 + T154*T153*Z1 + T155*T153*NP132
LT154 = LT154*NP131,J1 + T154*T154*NP132 + T155*T154*Z1 + T156*T154*NP132
NP135 = 0
LT155 = LT155*NP131,J1 + T155*T155*NP132 + T156*T155*Z1 + T157*T155*NP132 + T158*T155*Z1
LT156 = LT156*NP131,J1 + T156*T156*NP132 + T157*T156*Z1 + T158*T156*NP132
LT157 = LT157*NP131,J1 + T157*T157*NP132 + T158*T157*Z1 + T159*T157*NP132
LT158 = LT158*NP131,J1 + T158*T158*NP132 + T159*T158*Z1 + T160*T158*NP132
LT159 = LT159*NP131,J1 + T159*T159*NP132 + T160*T159*Z1 + T161*T159*NP132
LT160 = LT160*NP131,J1 + T160*T160*NP132 + T161*T160*Z1 + T162*T160*NP132
NP136 = 0
LT161 = LT161*NP131,J1 + T161*T161*NP132 + T162*T161*Z1 + T163*T161*NP132 + T164*T161*Z1
LT162 = LT162*NP131,J1 + T162*T162*NP132 + T163*T162*Z1 + T164*T162*NP132
LT163 = LT163*NP131,J1 + T163*T163*NP132 + T164*T163*Z1 + T165*T163*NP132
LT164 = LT164*NP131,J1 + T164*T164*NP132 + T165*T164*Z1 + T166*T164*NP132
NP137 = 0
LT165 = LT165*NP131,J1 + T165*T165*NP132 + T166*T165*Z1 + T167*T165*NP132 + T168*T165*Z1
LT166 = LT166*NP131,J1 + T166*T166*NP132 + T167*T166*Z1 + T168*T166*NP132
LT167 = LT167*NP131,J1 + T167*T167*NP132 + T168*T167*Z1 + T169*T167*NP132
LT168 = LT168*NP131,J1 + T168*T168*NP132 + T169*T168*Z1 + T170*T168*NP132
LT169 = LT169*NP131,J1 + T169*T169*NP132 + T170*T169*Z1 + T171*T169*NP132
LT170 = LT170*NP131,J1 + T170*T170*NP132 + T171*T170*Z1 + T172*T170*NP132
NP138 = 0
LT171 = LT171*NP131,J1 + T171*T171*NP132 + T172*T171*Z1 + T173*T171*NP132 + T174*T171*Z1
LT172 = LT172*NP131,J1 + T172*T172*NP132 + T173*T172*Z1 + T174*T172*NP132
LT173 = LT173*NP131,J1 + T173*T173*NP132 + T174*T173*Z1 + T175*T173*NP132
LT174 = LT174*NP131,J1 + T174*T174*NP132 + T175*T174*Z1 + T176*T174*NP132
NP139 = 0
LT175 = LT175*NP131,J1 + T175*T175*NP132 + T176*T175*Z1 + T177*T175*NP132 + T178*T175*Z1
LT176 = LT176*NP131,J1 + T176*T176*NP132 + T177*T176*Z1 + T178*T176*NP132
LT177 = LT177*NP131,J1 + T177*T177*NP132 + T178*T177*Z1 + T179*T177*NP132
LT178 = LT178*NP131,J1 + T178*T178*NP132 + T179*T178*Z1 + T180*T178*NP132
LT179 = LT179*NP131,J1 + T179*T179*NP132 + T180*T179*Z1 + T181*T179*NP132
LT180 = LT180*NP131,J1 + T180*T180*NP132 + T181*T180*Z1 + T182*T180*NP132
NP140 = 0
LT181 = LT181*NP131,J1 + T181*T181*NP132 + T182*T181*Z1 + T183*T181*NP132 + T184*T181*Z1
LT182 = LT182*NP131,J1 + T182*T182*NP132 + T183*T182*Z1 + T184*T182*NP132
LT183 = LT183*NP131,J1 + T183*T183*NP132 + T184*T183*Z1 + T185*T183*NP132
LT184 = LT184*NP131,J1 + T184*T184*NP132 + T185*T184*Z1 + T186*T184*NP132
NP141 = 0
LT185 = LT185*NP131,J1 + T185*T185*NP132 + T186*T185*Z1 + T187*T185*NP132 + T188*T185*Z1
LT186 = LT186*NP131,J1 + T186*T186*NP132 + T187*T186*Z1 + T188*T186*NP132
LT187 = LT187*NP131,J1 + T187*T187*NP132 + T188*T187*Z1 + T189*T187*NP132
LT188 = LT188*NP131,J1 + T188*T188*NP132 + T189*T188*Z1 + T190*T188*NP132
LT189 = LT189*NP131,J1 + T189*T189*NP132 + T190*T189*Z1 + T191*T189*NP132
LT190 = LT190*NP131,J1 + T190*T190*NP132 + T191*T190*Z1 + T192*T190*NP132
NP142 = 0
LT191 = LT191*NP131,J1 + T191*T191*NP132 + T192*T191*Z1 + T193*T191*NP132 + T194*T191*Z1
LT192 = LT192*NP131,J1 + T192*T192*NP132 + T193*T192*Z1 + T194*T192*NP132
LT193 = LT193*NP131,J1 + T193*T193*NP132 + T194*T193*Z1 + T195*T193*NP132
LT194 = LT194*NP131,J1 + T194*T194*NP132 + T195*T194*Z1 + T196*T194*NP132
NP143 = 0
LT195 = LT195*NP131,J1 + T195*T195*NP132 + T196*T195*Z1 + T197*T195*NP132 + T198*T195*Z1
LT196 = LT196*NP131,J1 + T196*T196*NP132 + T197*T196*Z1 + T198*T196*NP132
LT197 = LT197*NP131,J1 + T197*T197*NP132 + T198*T197*Z1 + T199*T197*NP132
LT198 = LT198*NP131,J1 + T198*T198*NP132 + T199*T198*Z1 + T200*T198*NP132
LT199 = LT199*NP131,J1 + T199*T199*NP132 + T200*T199*Z1 + T201*T199*NP132
LT200 = LT200*NP131,J1 + T200*T200*NP132 + T201*T200*Z1 + T202*T200*NP132
NP144 = 0
LT201 = LT201*NP131,J1 + T201*T201*NP132 + T202*T201*Z1 + T203*T201*NP132 + T204*T201*Z1
LT202 = LT202*NP131,J1 + T202*T202*NP132 + T203*T202*Z1 + T204*T202*NP132
LT203 = LT203*NP131,J1 + T203*T203*NP132 + T204*T203*Z1 + T205*T203*NP132
LT204 = LT204*NP131,J1 + T204*T204*NP132 + T205*T204*Z1 + T206*T204*NP132
NP145 = 0
LT205 = LT205*NP131,J1 + T205*T205*NP132 + T206*T205*Z1 + T207*T205*NP132 + T208*T205*Z1
LT206 = LT206*NP131,J1 + T206*T206*NP132 + T207*T206*Z1 + T208*T206*NP132
LT207 = LT207*NP131,J1 + T207*T207*NP132 + T208*T207*Z1 + T209*T207*NP132
LT208 = LT208*NP131,J1 + T208*T208*NP132 + T209*T208*Z1 + T210*T208*NP132
LT209 = LT209*NP131,J1 + T209*T209*NP132 + T210*T209*Z1 + T211*T209*NP132
LT210 = LT210*NP131,J1 + T210*T210*NP132 + T211*T210*Z1 + T212*T210*NP132
NP146 = 0
LT211 = LT211*NP131,J1 + T211*T211*NP132 + T212*T211*Z1 + T213*T211*NP132 + T214*T211*Z1
LT212 = LT212*NP131,J1 + T212*T212*NP132 + T213*T212*Z1 + T214*T212*NP132
LT213 = LT213*NP131,J1 + T213*T213*NP132 + T214*T213*Z1 + T215*T213*NP132
LT214 = LT214*NP131,J1 + T214*T214*NP132 + T215*T214*Z1 + T216*T214*NP132
NP147 = 0
LT215 = LT215*NP131,J1 + T215*T215*NP132 + T216*T215*Z1 + T217*T215*NP132 + T218*T215*Z1
LT216 = LT216*NP131,J1 + T216*T216*NP132 + T217*T216*Z1 + T218*T216*NP132
LT217 = LT217*NP131,J1 + T217*T217*NP132 + T218*T217*Z1 + T219*T217*NP132
LT218 = LT218*NP131,J1 + T218*T218*NP132 + T219*T218*Z1 + T220*T218*NP132
LT219 = LT219*NP131,J1 + T219*T219*NP132 + T220*T219*Z1 + T221*T219*NP132
LT220 = LT220*NP131,J1 + T220*T220*NP132 + T221*T220*Z1 + T222*T220*NP132
NP148 = 0
LT221 = LT221*NP131,J1 + T221*T221*NP132 + T222*T221*Z1 + T223*T221*NP132 + T224*T221*Z1
LT222 = LT222*NP131,J1 + T222*T222*NP132 + T223*T222*Z1 + T224*T222*NP132
LT223 = LT223*NP131,J1 + T223*T223*NP132 + T224*T223*Z1 + T225*T223*NP132
LT224 = LT224*NP131,J1 + T224*T224*NP132 + T225*T224*Z1 + T226*T224*NP132
NP149 = 0
LT225 = LT225*NP131,J1 + T225*T225*NP132 + T226*T225*Z1 + T227*T225*NP132 + T228*T225*Z1
LT226 = LT226*NP131,J1 + T226*T226*NP132 + T227*T226*Z1 + T228*T226*NP132
LT227 = LT227*NP131,J1 + T227*T227*NP132 + T228*T227*Z1 + T229*T227*NP132
LT228 = LT228*NP131,J1 + T228*T228*NP132 + T229*T228*Z1 + T230*T228*NP132
LT229 = LT229*NP131,J1 + T229*T229*NP132 + T230*T229*Z1 + T231*T229*NP132
LT230 = LT230*NP131,J1 + T230*T230*NP132 + T231*T230*Z1 + T232*T230*NP132
NP150 = 0
LT231 = LT231*NP131,J1 + T231*T231*NP132 + T232*T231*Z1 + T233*T231*NP132 + T234*T231*Z1
LT232 = LT232*NP131,J1 + T232*T232*NP132 + T233*T232*Z1 + T234*T232*NP132
LT233 = LT233*NP131,J1 + T233*T233*NP132 + T234*T233*Z1 + T235*T233*NP132
LT234 = LT234*NP131,J1 + T234*T234*NP132 + T235*T234*Z1 + T236*T234*NP132
NP151 = 0
LT235 = LT235*NP131,J1 + T235*T235*NP132 + T236*T235*Z1 + T237*T235*NP132 + T238*T235*Z1
LT236 = LT236*NP131,J1 + T236*T236*NP132 + T237*T236*Z1 + T238*T236*NP132
LT237 = LT237*NP131,J1 + T237*T237*NP132 + T238*T237*Z1 + T239*T237*NP132
LT238 = LT238*NP131,J1 + T238*T238*NP132 + T239*T238*Z1 + T240*T238*NP132
LT239 = LT239*NP131,J1 + T239*T239*NP132 + T240*T239*Z1 + T241*T239*NP132
LT240 = LT240*NP131,J1 + T240*T240*NP132 + T241*T240*Z1 + T242*T240*NP132
NP152 = 0
LT241 = LT241*NP131,J1 + T241*T241*NP132 + T242*T241*Z1 + T243*T241*NP132 + T244*T241*Z1
LT242 = LT242*NP131,J1 + T242*T242*NP132 + T243*T242*Z1 + T244*T242*NP132
LT243 = LT243*NP131,J1 + T243*T243*NP132 + T244*T243*Z1 + T245*T243*NP132
LT244 = LT244*NP131,J1 + T244*T244*NP132 + T245*T244*Z1 + T246*T244*NP132
NP153 = 0
LT245 = LT245*NP131,J1 + T245*T245*NP132 + T246*T245*Z1 + T247*T245*NP132 + T248*T245*Z1
LT246 = LT246*NP131,J1 + T246*T246*NP132 + T247*T246*Z1 + T248*T246*NP132
LT247 = LT247*NP131,J1 + T247*T247*NP132 + T248*T247*Z1 + T249*T247*NP132
LT248 = LT248*NP131,J1 + T248*T248*NP132 + T249*T248*Z1 + T250*T248*NP132
LT249 = LT249*NP131,J1 + T249*T249*NP132 + T250*T249*Z1 + T251*T249*NP132
LT250 = LT250*NP131,J1 + T250*T250*NP132 + T251*T250*Z1 + T252*T250*NP132
NP154 = 0
LT251 = LT251*NP131,J1 + T251*T251*NP132 + T252*T251*Z1 + T253*T251*NP132 + T254*T251*Z1
LT252 = LT252*NP131,J1 + T252*T252*NP132 + T253*T252*Z1 + T254*T252*NP132
LT253 = LT253*NP131,J1 + T253*T253*NP132 + T254*T253*Z1 + T255*T253*NP132
LT254 = LT254*NP131,J1 + T254*T254*NP132 + T255*T254*Z1 + T256*T254*NP132
NP155 = 0
LT255 = LT255*NP131,J1 + T255*T255*NP132 + T256*T255*Z1 + T257*T255*NP132 + T258*T255*Z1
LT256 = LT256*NP131,J1 + T256*T256*NP132 + T257*T256*Z1 + T258*T256*NP132
LT257 = LT257*NP131,J1 + T257*T257*NP132 + T258*T257*Z1 + T259*T257*NP132
LT258 = LT258*NP131,J1 + T258*T258*NP132 + T259*T258*Z1 + T260*T258*NP132
LT259 = LT259*NP131,J1 + T259*T259*NP132 + T260*T259*Z1 + T261*T259*NP132
LT260 = LT260*NP131,J1 + T260*T260*NP132 + T261*T260*Z1 + T262*T260*NP132
NP156 = 0
LT261 = LT261*NP131,J1 + T261*T261*NP132 + T262*T261*Z1 + T263*T261*NP132 + T264*T261*Z1
LT262 = LT262*NP131,J1 + T262*T262*NP132 + T263*T262*Z1 + T264*T262*NP132
LT263 = LT263*NP131,J1 + T263*T263*NP132 + T264*T263*Z1 + T265*T263*NP132
LT264 = LT264*NP131,J1 + T264*T264*NP132 + T265*T264*Z1 + T266*T264*NP132
NP157 = 0
LT265 = LT265*NP131,J1 + T265*T265*NP132 + T266*T265*Z1 + T267*T265*NP132 + T268*T265*Z1
LT266 = LT266*NP131,J1 + T266*T266*NP132 + T267*T266*Z1 + T268*T266*NP132
LT267 = LT267*NP131,J1 + T267*T267*NP132 + T268*T267*Z1 + T269*T267*NP132
LT268 = LT268*NP131,J1 + T268*T268*NP132 + T269*T268*Z1 + T270*T268*NP132
LT269 = LT269*NP131,J1 + T269*T269*NP132 + T270*T269*Z1 + T271*T269*NP132
LT270 = LT270*NP131,J1 + T270*T270*NP132 + T271*T270*Z1 + T272*T270*NP132
NP158 = 0
LT271 = LT271*NP131,J1 + T271*T271*NP132 + T272*T271*Z1 + T273*T271*NP132 + T274*T271*Z1
LT272 = LT272*NP131,J1 + T272*T272*NP132 + T273*T272*Z1 + T274*T272*NP132
LT273 = LT273*NP131,J1 + T273*T273*NP132 + T274*T273*Z1 + T275*T273*NP132
LT274 = LT274*NP131,J1 + T274*T274*NP132 + T275*T274*Z1 + T276*T274*NP132
NP159 = 0
LT275 = LT275*NP131,J1 + T275*T275*NP132 + T276*T275*Z1 + T277*T275*NP132 + T278*T275*Z1
LT276 = LT276*NP131,J1 + T276*T276*NP132 + T277*T276*Z1 + T278*T276*NP132
LT277 = LT277*NP131,J1 + T277*T277*NP132 + T278*T277*Z1 + T279*T277*NP132
LT278 = LT278*NP131,J1 + T278*T278*NP132 + T279*T278*Z1 + T280*T278*NP132
LT279 = LT279*NP131,J1 + T279*T279*NP132 + T280*T279*Z1 + T281*T279*NP132
LT280 = LT280*NP131,J1 + T280*T280*NP132 + T281*T280*Z1 + T282*T280*NP132
NP160 = 0
LT281 = LT281*NP131,J1 + T281*T281*NP132 + T282*T281*Z1 + T283*T281*NP132 + T284*T281*Z1
LT282 = LT282*NP131,J1 + T282*T282*NP132 + T283*T282*Z1 + T284*T282*NP132
LT283 = LT283*NP131,J1 + T283*T283*NP132 + T284*T283*Z1 + T285*T283*NP132
LT284 = LT284*NP131,J1 + T284*T284*NP132 + T285*T284*Z1 + T286*T284*NP132
NP161 = 0
LT285 = LT285*NP131,J1 + T285*T285*NP132 + T286*T285*Z1 + T287*T285*NP132 + T288*T285*Z1
LT286 = LT286*NP131,J1 + T286*T286*NP132 + T287*T286*Z1 + T288*T286*NP132
LT287 = LT287*NP131,J1 + T287*T287*NP132 + T288*T287*Z1 + T289*T287*NP132
LT288 = LT288*NP131,J1 + T288*T288*NP132 + T289*T288*Z1 + T290*T288*NP132
LT289 = LT289*NP131,J1 + T289*T289*NP132 + T290*T289*Z1 + T291*T289*NP132
LT290 = LT290*NP131,J1 + T290*T290*NP132 + T291*T290*Z1 + T292*T290*NP132
NP162 = 0
LT291 = LT291*NP131,J1 + T291*T291*NP132 + T292*T291*Z1 + T293*T291*NP132 + T294*T291*Z1
LT292 = LT292*NP131,J1 + T292*T292*NP132 + T293*T292*Z1 + T294*T292*NP132
LT293 = LT293*NP131,J1 + T293*T293*NP132 + T294*T293*Z1 + T295*T293*NP132
LT294 = LT294*NP131,J1 + T294*T294*NP132 + T295*T294*Z1 + T296*T294*NP132
NP163 = 0
LT295 = LT295*NP131,J1 + T295*T295*NP132 + T296*T295*Z1 + T297*T295*NP132 + T298*T295*Z1
LT296 = LT296*NP131,J1 + T296*T296*NP132 + T297*T296*Z1 + T298*T296*NP132
LT297 = LT297*NP131,J1 + T297*T297*NP132 + T298*T297*Z1 + T299*T297*NP132
LT298 = LT298*NP131,J1 + T298*T298*NP132 + T299*T298*Z1 + T300*T298*NP132
LT299 = LT299*NP131,J1 + T299*T299*NP132 + T300*T299*Z1 + T301*T299*NP132
LT300 = LT300*NP131,J1 + T300*T300*NP132 + T301*T300*Z1 + T302*T300*NP132
NP164 = 0
LT301 = LT301*NP131,J1 + T301*T301*NP132 + T302*T301*Z1 + T303*T301*NP132 + T304*T301*Z1
LT302 = LT302*NP131,J1 + T302*T302*NP132 + T303*T302*Z1 + T304*T302*NP132
LT303 = LT303*NP131,J1 + T303*T303*NP132 + T304*T303*Z1 + T305*T303*NP132
LT304 = LT304*NP131,J1 + T304*T304*NP132 + T305*T304*Z1 + T306*T304*NP132
NP165 = 0
LT305 = LT305*NP131,J1 + T305*T305*NP132 + T306*T305*Z1 + T307*T305*NP132 + T308*T305*Z1
LT306 = LT306*NP131,J1 + T306*T306*NP132 + T307*T306*Z1 + T308*T306*NP132
LT307 = LT307*NP131,J1 + T307*T307*NP132 + T308*T307*Z1 + T309*T307*NP132
LT308 = LT308*NP131,J1 + T308*T308*NP132 + T309*T308*Z1 + T310*T308*NP132
LT309 = LT309*NP131,J1 + T309*T309*NP132 + T310*T309*Z1 + T311*T309*NP132
LT310 = LT310*NP131,J1 + T310*T310*NP132 + T311*T310*Z1 + T312*T310*NP132
NP166 = 0
LT311 = LT311*NP131,J1 + T311*T311*NP132 + T312*T311*Z1 + T313*T311*NP132 + T314*T311*Z1
LT312 = LT312*NP131,J1 + T312*T312*NP132 + T313*T312*Z1 + T314*T312*NP132
LT313 = LT313*NP131,J1 + T313*T313*NP132 + T314*T313*Z1 + T315*T313*NP132
LT314 = LT314*NP131,J1 + T314*T314*NP132 + T315*T314*Z1 + T316*T314*NP132
NP167 = 0
LT315 = LT315*NP131,J1 + T315*T315*NP132 + T316*T315*Z1 + T317*T315*NP132 + T318*T315*Z1
LT316 = LT316*NP131,J1 + T316*T316*NP132 + T317*T316*Z1 + T318*T316*NP132
LT317 = LT317*NP131,J1 + T317*T317*NP132 + T318*T317*Z1 + T319*T317*NP132
LT318 = LT318*NP131,J1 + T318*T318*NP132 + T319*T318*Z1 + T320*T318*NP132
LT319 = LT319*NP131,J1 + T319*T319*NP132 + T320*T319*Z1 + T321*T319*NP132
LT320 = LT320*NP131,J1 + T320*T320*NP132 + T321*T320*Z1 + T322*T320*NP132
NP168 = 0
LT321 = LT321*NP131,J1 + T321*T321*NP132 + T322*T321*Z1 + T323*T321*NP132 + T324*T321*Z1
LT322 = LT322*NP131,J1 + T322*T322*NP132 + T323*T322*Z1 + T324*T322*NP132
LT323 = LT323*NP131,J1 + T323*T323*NP132 + T324*T323*Z1 + T325*T323*NP132
LT324 = LT324*NP131,J1 + T324*T324*NP132 + T325*T324*Z1 + T326*T324*NP132
NP169 = 0
LT325 = LT325*NP131,J1 + T325*T325*NP132 + T326*T325*Z1 + T327*T325*NP132 + T328*T325*Z1
LT326 = LT326*NP131,J1 + T326*T326*NP132 + T327*T326*Z1 + T328*T326*NP132
LT327 = LT327*NP131,J1 + T327*T327*NP132 + T32
```

APPENDIX II

Listing of DBMIE Subroutine

L. TAUELL,J,J = THE PRECEDING VALUES OF RNA & FNU RESPECTIVELY.
 LWHITE1,J,J = TAU111,TAU112,TAU113,FNU1,TU111
 LWHITE100,J,J = TAU111,TAU112,TAU113,FNU1,TU111
 IF E 24 .LE. 100 F GO TO 20
 WHITE1,J,J
 WHITE100,J,J
 CALL 1,J,J
 20 IR = 100*TAUELL,J,J+FNU1
 IRB = 1.00000E
 RA = 1.00074
 RHEX = RHE + RA
 TLL1 = 1.444234E+04**21
 TLL1 = 0.000111111
 RMAX = 1.1000 * TLL1
 IF E RMAX .LE. 0.999 F GO TO 21
 WHITE1,J,J
 CALL 1,J,J
 CL RHE2 < TLL1
 DO 1 RHE1 =>1, 150 F GO TO 22
 RHE1 = 150
 RHE2 = 115
 CL ACAPINH1 + 1, 1 => 1.0000, 0.000 :
 DO 21 N = 1,NMAX
 NN = RHE1 - N + 1
 ACAPINH1 = 1.64*11 * RHEX + 1.0000*ENH1*ENHAX + ACAPINH1+1
 2.3 CONTINUE
 DO 32 J = 1,JX
 IF E THETD1,J,J <= 0.000 F THEFD1,J,J = QABS(THETD1,J,J)
 DO 33 IFETD1,J,J <= 0.000 F GO TO 24
 CTHF1,J,J = 1.000
 STHF1,J,J = 0.000
 GO TO 32
 25 IF E THETD1,J,J <= 90.000 F GO TO 25
 TLL1 = 1.31415926535897932 * THETD1,J,J/180.000
 LSTHT1,J,J = 0.000111111
 STHT1,J,J = 1.000 - LSTHT1,J,J**2
 GO TO 30
 26 IF E THETD1,J,J <= 90.000 F GO TO 26
 LSTHT1,J,J = 0.000
 STHT1,J,J = 1.000
 GO TO 32
 28 WHITE100,J,J THETD1,J,J
 WHITE100,J,J
 CALL 1,J,J
 30 CONTINUE
 DO 35 J = 1,JX
 P111,J,J = 0.000
 P112,J,J = 1.000
 DBMIE049
 DBMIE050
 DBMIE051
 DBMIE052
 DBMIE053
 DBMIE054
 DBMIE055
 DBMIE056
 DBMIE057
 DBMIE058
 DBMIE059
 DBMIE060
 DBMIE061
 DBMIE062
 DBMIE063
 DBMIE064
 DBMIE065
 DBMIE066
 DBMIE067
 DBMIE068
 DBMIE069
 DBMIE070
 DBMIE071
 DBMIE072
 DBMIE073
 DBMIE074
 DBMIE075
 DBMIE076
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 DBMIE078
 DBMIE079
 DBMIE080
 DBMIE081
 DBMIE082
 DBMIE083
 DBMIE084
 DBMIE085
 DBMIE086
 DBMIE087
 DBMIE088
 DBMIE089
 DBMIE090
 DBMIE091
 DBMIE092
 DBMIE093
 DBMIE094
 DBMIE095
 DBMIE096

0058 TAU112,J,J = 1.000
 TAU112,J,J = 0.000111111
 35 CONTINUE
 TLL1 = 0.000111111
 TLL2 = 0.000111111
 RHE1 = 0.000111111
 RHE2 = 0.000111111
 TLL1 = ACAPINH1 * RHE + RA
 PNA = TLL1*TLL11 - TAU111*TLL1*PNA1 - RHE111
 FNU = TLL2*TLL21 - TAU112*TLL2 * RHE112 - RHE111
 PNA = FNU
 FNUP = FNU
 TLL1 = 1.0000
 L. FROM HERE TO THE STATEMENT NUMBER 40, ELTHM1,J,J,F HAS THE
 FOLLOWING MEANING:
 ELTHM1,J,J,F1 REAL PART OF THE FIRST COMPLEX AMPLITUDE,
 ELTHM1,J,J,F2 IMAGINARY PART OF THE FIRST COMPLEX AMPLITUDE,
 ELTHM1,J,J,F3 REAL PART OF THE SECOND COMPLEX AMPLITUDE,
 ELTHM1,J,J,F4 IMAGINARY PART OF THE SECOND COMPLEX AMPLITUDE,
 N = 1 FOR THE TLL1,J,J AND N = 2 FOR 180.0 - THETD1,J,J
 DEFINITION OF THE COMPLEX AMPLITUDES: VAN DE HULST,P.125.
 TLL1 = TLL1 + TLL1
 TLL2 = TLL1 * TLL2
 TLL3 = TLL1 * TLL1
 TLL21 = TLL1 * TLL21
 00 60 J = 1,JX
 ELTHM1,J,J,F1 = TLL1 * P111,J,J + TLL11 * TAU112,J,J
 ELTHM1,J,J,F2 = TLL2 * P112,J,J + TLL21 * TAU122,J,J
 ELTHM1,J,J,F3 = TLL1 * P111,J,J + TLL11 * TAU112,J,J
 ELTHM1,J,J,F4 = TLL2 * P112,J,J - TLL21 * TAU122,J,J
 ELTHM1,J,J,F1 = TLL1 * P112,J,J - TLL11 * TAU112,J,J
 ELTHM1,J,J,F2 = TLL2 * P111,J,J + TLL21 * TAU122,J,J
 ELTHM1,J,J,F3 = TLL1 * P111,J,J - TLL11 * TAU112,J,J
 ELTHM1,J,J,F4 = TLL2 * P112,J,J - TLL21 * TAU122,J,J
 00 CONTINUE
 QCAT = 2.000 + 1 TLL11 + TLL111
 QCAT = TLL11**2 + TLL21**2 + TLL11**2*10.7500
 CFBQS = 0.000
 N = 2
 45 TLL1 = 2*N - 1
 TLL2 = N - 1
 TLL3 = 2 * N + 1
 DO 70 J = 1,JX
 P111,J,J = TLL1**P111,J,J + CTHF1,J,J*TLL11
 P112,J,J = TLL2**P112,J,J + STHF1,J,J*TLL21
 TAU112,J,J = CTHF1,J,J*TLL11 + P113,J,J - P111,J,J*TLL11
 TAU112,J,J = TLL11**P111,J,J + TLL21**P112,J,J + TLL11**P112,J,J

APPENDIX IV

Sample Output Using DAMIE Subroutine
 $m = 1.342 - 0.0i$, $r = 100.0 \mu$, $\lambda = 0.4 \mu$

ELEMENTS OF THE TRANSFORMATION MATRIX FOR A SPHERE WITH SIZE PARAMETER = 100.0

REFRACTIVE INDEX, KERL = 0.135200 01 TRANSMISSION = 0.0

WAVELET	M SUB 2	N SUB 1	S SUB 21	O SUB 21	INTENSITY	POLARIZATION
0.00000	0.1352000 13	0.1352000 13	0.1352000 13	-0.0i	0.1352000 13	0.0
1.00000	0.1625000 09	0.1625000 09	0.1625000 09	0.1047210 03	0.1625000 09	0.00100
2.00000	0.18673500 08	0.18673500 08	0.18673500 08	-0.2245870 07	0.18673500 08	-0.01470
3.00000	0.1982610 07	0.1982610 07	0.1982610 07	-0.5785350 06	0.1982610 07	-0.01570
4.00000	0.2579132 07	0.1979132 07	0.2233730 07	-0.3031660 06	0.2233730 07	-0.15510
5.00000	0.3599010 07	0.1775250 07	0.3079420 07	-0.1403790 06	0.3079420 07	0.04320
6.00000	0.41173190 06	0.1155630 06	0.1164140 06	-0.4184270 06	0.1164140 06	-0.00707
7.00000	0.12327500 06	0.1134660 06	0.1207690 06	0.4184180 06	0.1207690 06	0.07950
8.00000	0.48618610 05	0.42648610 05	0.4270400 05	-0.3130760 06	0.42648610 05	-0.21390
9.00000	0.11171250 05	0.1134430 05	0.1301210 05	-0.2930500 06	0.1134430 05	0.03420
10.00000	0.10763500 04	0.1054580 04	0.113120 04	-0.1159810 05	0.1054580 04	-0.01840
11.00000	0.17031920 04	0.165980 04	0.1805810 04	-0.5665160 05	0.1805810 04	-0.18350
12.00000	0.25599710 03	0.1891240 03	0.2451030 03	-0.1294280 04	0.2451030 03	-0.03550
13.00000	0.18456800 03	0.1752010 03	0.1847570 03	-0.1415140 04	0.1847570 03	0.00550
14.00000	0.07298620 02	0.0749820 02	0.0748240 02	-0.2626280 03	0.0748240 02	-0.00460
15.00000	0.06103790 02	0.0615410 02	0.0608940 02	-0.4764640 03	0.0608940 02	-0.02120
16.00000	0.05706480 02	0.0582360 02	0.05640150 02	-0.1517790 03	0.0582360 02	-0.02220
17.00000	0.06159360 02	0.0614960 02	0.06224860 02	-0.2830700 03	0.06224860 02	0.01240
18.00000	0.07118520 02	0.0746490 02	0.07310610 02	-0.5178180 03	0.07310610 02	0.02490
19.00000	0.16192700 02	0.1611750 02	0.17573890 02	-0.2626280 03	0.17573890 02	-0.09660
20.00000	0.04065620 02	0.0414890 02	0.03743050 02	-0.3140780 03	0.0414890 02	-0.12950
21.00000	0.14464520 02	0.1454240 02	0.14266900 02	-0.3020860 03	0.1454240 02	-0.08140
22.00000	0.14609950 02	0.1448090 02	0.1473120 02	-0.3834220 03	0.1448090 02	-0.02230
23.00000	0.13730190 02	0.12972100 02	0.13252350 02	-0.1686880 03	0.13331160 02	-0.11330
24.00000	0.12690700 02	0.12619700 02	0.12842600 02	-0.2833590 03	0.12842600 02	-0.13750
25.00000	0.13774150 02	0.14078100 02	0.1386010 02	-0.2153780 03	0.14078100 02	-0.09770
26.00000	0.14926430 02	0.15674520 02	0.15286370 02	-0.9579510 03	0.15674520 02	0.07050
27.00000	0.13088180 02	0.12942720 02	0.2990640 02	-0.2992590 03	0.12942720 02	-0.04270
28.00000	0.13152690 02	0.12775550 02	0.1293150 02	-0.3954590 03	0.12775550 02	-0.06160
29.00000	0.13818490 02	0.13563760 02	0.13655680 02	-0.48191010 03	0.13563760 02	-0.08960
30.00000	0.12266800 02	0.11993460 02	0.1089040 02	-0.2090680 03	0.11993460 02	-0.17440
31.00000	0.13054480 02	0.12819690 02	0.2937470 02	-0.2431940 03	0.12819690 02	-0.08170
32.00000	0.13383050 02	0.13218150 02	0.1312670 02	-0.1705860 03	0.13218150 02	-0.11220
33.00000	0.11818430 02	0.1044860 02	0.1125900 02	-0.1170530 03	0.1044860 02	-0.31110
34.00000	0.12525590 02	0.12399610 02	0.12430060 02	-0.3883370 03	0.12399610 02	-0.02950
35.00000	0.12588040 02	0.12694690 02	0.12623750 02	-0.2998900 03	0.12694690 02	0.02020
36.00000	0.11478110 02	0.11153100 02	0.10394420 02	-0.3047470 03	0.11153100 02	-0.18920
37.00000	0.11152560 02	0.10760100 02	0.1030860 02	-0.2330750 03	0.10760100 02	-0.15150
38.00000	0.11996610 02	0.11477700 02	0.1141580 02	-0.4903370 03	0.11477700 02	-0.11090
39.00000	0.11865310 02	0.11921130 02	0.1167820 02	-0.3255450 03	0.11921130 02	0.11630
40.00000	0.11829710 02	0.11024640 02	0.1106470 02	-0.1578340 03	0.11024640 02	0.14440
41.00000	0.11895910 02	0.11081740 02	0.1187750 02	-0.5783520 03	0.1187750 02	-0.03030
42.00000	0.11628300 02	0.11989110 02	0.11596670 02	-0.1958450 03	0.11989110 02	-0.01110
43.00000	0.11480370 02	0.1111540 02	0.1136500 02	-0.2936470 03	0.1111540 02	-0.03990
44.00000	0.11152740 02	0.1135750 02	0.11079760 02	-0.2316890 03	0.1135750 02	-0.14290
45.00000	0.11246710 02	0.11043170 02	0.1114290 02	-0.2403520 03	0.11043170 02	-0.08780
46.00000	0.11044320 02	0.11930730 02	0.11651950 02	-0.2822380 03	0.11930730 02	-0.13667
47.00000	0.10758720 02	0.10359150 02	0.1298260 02	-0.3559960 03	0.10359150 02	-0.15861
48.00000	0.10262640 02	0.8746330 02	0.9271330 02	-0.1943670 03	0.9271330 02	-0.07460

P1+0.00	0.16000000	0.0	0.11111100	0.0	-0.16555800	0.05	0.21282300	0.0	0.58999900	0.0	0.1
P2+0.00	0.12000000	0.0	0.09111100	0.0	-0.20312700	0.05	-0.47823900	0.0	0.38000000	0.0	0.1
P3+0.00	0.09000000	0.0	0.06025400	0.0	-0.06974800	0.05	-0.47810000	0.05	0.91429900	0.05	0.1
P4+0.00	0.11500000	0.0	0.12268100	0.0	-0.13908800	0.06	0.27652800	0.0	0.13932700	0.0	0.1
P5+0.00	0.14500000	0.0	0.18783200	0.0	-0.21237200	0.07	-0.31552800	0.05	0.95413900	0.0	0.1
P6+0.00	0.20500000	0.0	0.20163000	0.0	-0.15298300	0.06	-0.12970000	0.0	0.16020700	0.0	0.1
P7+0.00	0.22000000	0.0	0.19423600	0.0	-0.72267600	0.04	-0.17532300	0.0	0.19760400	0.05	0.1
P8+0.00	0.18000000	0.0	0.09696700	0.0	-0.53916600	0.05	0.79499900	0.03	0.54804700	0.05	0.1
P9+0.00	0.10000000	0.0	0.12883600	0.0	-0.10405000	0.06	-0.11143900	0.04	0.10646300	0.0	0.1
P10+0.00	0.10000000	0.0	0.10920300	0.07	-0.10920300	0.07	-0.10920300	0.07	0.10920300	0.0	0.1

EFFICIENCY FACTOR FOR SCATTERING = 0.2012940 01

EFFICIENCY FACTOR FOR SCATTERING = 0.2012940 01

EFFICIENCY FACTOR FOR ABSORPTION = 0.0

REFRACTIVE FACTOR = 1.68033400 00

TOTAL TIME FOR THIS CASE IN SECONDS = 249.484

APPENDIX V

Sample Output Using D8MIE Subroutine

$m = 1.342 - 0.1i$, $r = 100.0 \mu$, $\lambda = 0.4 \mu$

REF ID: A

J. V. DAVE, J. CANOSA

1974
Subroutine for Computing the Parameters
of the Electromagnetic Radiation
emitted by a Sphere

IBM RESEARCH LABORATORY

IBM ALTO SCIENTIFIC CENTER

REPORT NUMBER

320-3332

NUMBER OF PAGES

42

NO. OF REFERENCES

37

DATE COMPUTED

May 1974

SUBJECT

DATE OF INITIAL PRINTING

March 1968

DATE OF LAST PRINTING

April 1974

Two experimental IBM subroutines with which one can compute the monochromatic radiation emitted by a sphere, are described in great detail. The index of refraction, n_1 , of the material of the sphere is assumed to have the form $n_1 = n_0 + \alpha_1 r$ with $\alpha_1 < 0$. The formulae used for these subroutines were first derived by G. Mie (1908) and as such this scattering process is referred to as Mie scattering in the scientific literature.

A listing of these subroutines along with sample program outputs are provided in the Appendixes.

KEY TERMS FOR THE IBM SUBJECT INDEX

Scattering
Radiation
Optics
Atmosphere

IBM - Earth and Atmospheric

20 - Physics

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The availability of reprints is correct as of the printing date of this report.

* Reprints are not available from the Scientific Center.