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Application of Cauchy's Residue Theorem in Evaluating the Poles and Zeros of Complex Meromorphic Functions and Apposite Computer Programs

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Abstract

This note addresses itself to the problem of locating the zeros and poles of a complex meromorphic function M(s) in a specified rectangular or square region of the complex s-plane. It is assumed that M(s) has to be computed numerically as for example, a dispersion relation in plasma physics or ii) the system determinant of the matricized integral equation while employing the singularity expansion method (SEM)[1] to solve electromagnetic scattering problems. The procedure developed here eliminates the usual 2-dimensional search and replaces it with a direct constructive method for determining the poles of M(s) based on an application of Cauchy's residue theorem. The zeros of M(s) are easily found by applying the procedure to the reciprocal function 1/M(s). Two examples, i.e., 1) ratios of polynomials and 2) input impedance of a biconical antenna, are numerically illustrated.

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List of Principal Symbols

1. s = Complex variable =
$$s_R + s_I = \Omega + j\omega$$

$$M(s) = A_1(s)/A_2(s)$$
4. $N_a(f,C)$ = Argument number of the function $f(s)$ in a prescribed counterclockwise or

5.
$$N_O(f,C)$$
 = Number of zeros of $f(s)$ in C

6.
$$N_{p}(f,C)$$
 = Number of poles of $f(s)$ in C

8.
$$(z_1, z_2, \dots)$$
Locations of zeros and poles of a meromorphic function (p_1, p_2, \dots)

9.
$$M^{\text{nor}}(s)$$
 = Normalized version of $M(s)$

^{*} In addition to being analytic on the contour C which encloses the domain D, A(s) and M(s) are required not to vanish on C.

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I. Introduction

The problem of locating the poles and zeros of complex functions in a finite domain of the complex plane, occurs in many scientific disciplines e.g., dispersion relations in plasma physics, the singularity expansion method [1] in electromagnetic scattering or antenna problems. In an earlier note [2] Singaraju, et al. described a technique of locating the zeros of analytic functions in a given region of the complex plane. This note also included relevant computer programs and illustrative examples by way of i) polynomials ii) product of polynomials and exponentials and iii) determination of natural frequencies of a thin straight wire.

A sequential extension of the above mentioned work [2] is, of course, a method of locating poles and zeros of meromorphic functions when they coexist in a given contour. It is interesting to note that the word "meromorphic" is derived [3] from the Greek $\mu\epsilon\rho\sigma s = fraction$ and $\mu\sigma\rho\phi\eta = form$, and means "like a fraction." In keeping with the origin of the word "meromorphic," the complex function M(s) considered in this note will be a ratio of two entire functions of the complex variable s. The problem at hand can now be defined in terms of given and required quantities, as follows:

Given:

- i) A numerical way of evaluating a meromorphic complex function M(s) of a complex variable s,
- ii) A rectangular or a square region in the finite complex s-plane.

Find:

i) All the zeros and poles of M(s) in the given region.

Remark:

i) Typically, evaluation of M(s) is expensive in terms of computer time and hence it is desirable to optimize the number of M(s) computations.

II. A Review of SGB Technique for Finding the Zeros of Analytic Functions

In a recent note [2], Singaraju, Giri and Baum described a technique of locating the zeros of an analytic function A(s) in a finite domain D of the complex s-plane. This work, referred to as the SGB Technique also includes a family of computer programs titled SEARCH. This technique is based on the "principle of the argument" and a generalization thereof. The principle of argument for an analytic function is given by [4]

$$\frac{1}{2\pi i} \oint \frac{A'(s)}{A(s)} ds = N_o(A,C) \qquad (2.1)$$

where

A(s) = Analytic function* of s in a domain D enclosed by a simple contour C,

 $N_{O}(A,C)$ = Number of zeros of the analytic function A'(s) inside the contour C.

Equation (2.1) is a special case of

$$\frac{1}{2\pi i} \oint_{C} A_{m}(s) \frac{A'(s)}{A(s)} ds = \sum_{\alpha=1}^{N_{O}} A_{m}(s_{\alpha})$$
 (2.2)

obtained by setting $A_m(s) = 1$. In equation (2.2) $A_m(s)$ is an analytic (at least in and on C) multiplier function and s_{α} are the zeros of A(s) in C. If we choose $A_m(s) = s^n$ and consider n to take integer values ranging from 0 to N_C , we have

$$c_n = \frac{1}{2\pi i} \int_{C} s^n \frac{A'(s)}{A(s)} ds;$$
 for $n = 0, 1, 2, ... N_o$
(2.3)

which leads to

* A(s) is required not to vanish on the contour C.

$$C_0 = 1 + 1 + 1 + 1 + 1 + \dots + 1 = N_0$$
 (2.4.0)

$$c_1 = s_1 + s_2 + s_3 + s_4 \dots + s_N$$
 (2.4.1)

$$c_2 = s_1^2 + s_2^2 + s_3^2 + s_4^2 \dots + s_N^2$$
 (2.4.2)

$$C_{N_{0}} = s_{1}^{N_{0}} + s_{2}^{N_{0}} + s_{3}^{N_{0}} + s_{4}^{N_{0}} + s_{N_{0}}^{N_{0}}$$
 (2.4.N₀)

After determining the moments (C_n) SEARCH proceeds to locate the zeros in the given contour C (if any) by solving the above system of equations. By way of an interesting example, SEARCH has been used in easing the chase for those "elusive and ubiquitous" [5] SEM poles of a thin wire.

Above is a rather brief description of the underlying basis of the SGB technique and the interested reader is referred to Mathematics Note 42 [2] for all of the details regarding the working, limitations and use of the relevant computer programs. A logical extension of this work is of course a method of locating the zeros and poles of a meromorphic function in a finite region of the complex plane. In the following section, a method is developed which determines only the poles of a meromorphic function in a region regardless of whether or not there are zeros in that region. By applying the pole finding method of section III to the given function as well as its reciprocal, the zeros and poles of the given function in the given region are successfully located.

III. Poles and Zeros of Meromorphic Functions

A. Pole finder

In this section we shall develop a procedure to determine the number of poles [N_p(M,C)] and their locations of a meromorphic function M(s) in a given contour C. This procedure is independent of the presence or absence of zeros in C and also the actual shape of the contour C itself. However, for purposes of illustration and numerical ease, we shall consider the contour C to be a square as in figure 3.1 which in some special cases may be rectangular. We will also stipulate that the side of the square is equal to or not very different from unit length in the normalized s plane of figure 3.1.

The meromorphic function is representable by a ratio of two entire functions $E_1(s)$ and $E_2(s)$ as

$$M(s) = \frac{E_{1}(s)}{E_{2}(s)}$$

$$= \frac{A_{1}(s) A_{1}(s)}{A_{2}(s) A_{2}(s)}$$
(3.1)

E₁(s) and E₂(s) are in turn written as a product of an interior and an exterior analytic function. The subscripts "interior" and "exterior" are with reference to the contour C of figure 3.1. The "exterior" functions are required to be analytic in and on C, and not vanish on C. Our procedure of finding poles inside contour C allows for other types of singularities like essential or branch point to occur outside and sufficiently away from the contour C. The poles of M(s) within C are of course the zeros of the following equation

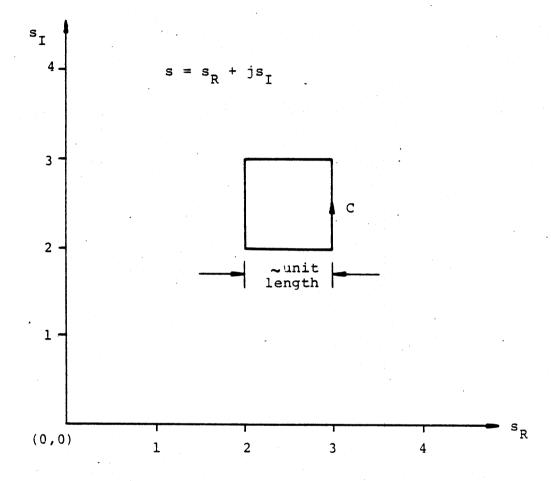


Figure 3.1 Normalized s-plane showing a simple square contour C.

$$A_{2int}(s) = 0$$
 (3.2)

Any analytic function can be represented by a suitable polynomial so that,

$$A_{lint}(s) = \prod_{i=1}^{N_{o}} (s - z_{i})$$
 (3.3)

$$A_{lext}(s) = E_{l}(s)/A_{lint}(s)$$
 (3.4)

$$A_{2_{int}}(s) = \prod_{k=1}^{N_p} (s - p_k)$$
 (3.5)

$$A_{2}_{ext}(s) = E_{2}(s)/A_{2}_{int}(s)$$
 (3.6)

leading to

$$M(s) = \frac{\begin{bmatrix} N_o \\ T \\ i=1 \end{bmatrix} (s-z_i) \begin{bmatrix} A_{1_{ext}}(s) \\ A_{2_{ext}}(s) \end{bmatrix}}{\begin{bmatrix} N_p \\ T \\ k=1 \end{bmatrix} (s-p_k) \begin{bmatrix} A_{2_{ext}}(s) \\ A_{2_{ext}}(s) \end{bmatrix}}$$
(3.7)

where

where

 z_i 's are zeros of M(s) within C,

 p_k 's are poles of M(s) within C,

 N_{O} = number of zeros of M(s) within c,

and

 N_{p} = number of poles of M(s) within C.

Our pole finding scheme determines N_p and subsequently p_k for k = 1, 2, N_p .

i) Finding the number of poles N_{p}

Besides the function values, what distinguishes a pole from a zero is the concept of residue and Cauchy's residue theorem. If the poles $(p_k$'s) are simple then the corresponding residues are given by

the corresponding residues are given by

$$R_{m} = \underset{s \to p_{m}}{\text{Lim}} \left[(s - p_{m}) M(s) \right]$$

$$= \underset{s \to p_{m}}{\text{Lim}} \left[(s - p_{m}) \underbrace{\prod_{i=1}^{N_{O}} (s - z_{i})}_{N_{O}} A_{1} \underbrace{(s)}_{ext} \right]$$

$$= \underset{s \to p_{m}}{\text{Lim}} \left[(s - p_{m}) \underbrace{\prod_{i=1}^{N_{D}} (s - p_{k})}_{K=1} A_{2} \underbrace{(s)}_{ext} \right]$$

$$= \begin{bmatrix} \prod_{i=1}^{N_{O}} (p_{m} - z_{i}) & A_{1_{ext}} (p_{m}) \\ \hline N_{p} & \\ \prod_{k=1}^{N_{p}} (p_{m} - p_{k}) & A_{2_{ext}} (p_{m}) \\ k \neq m \end{bmatrix} ; \text{ for } m = 1, 2...N_{p}$$
(3.8)

However, for the present purpose, R_{m} 's are of no interest and will eventually be eliminated.

We will now define residue moments by

$$D_{n} \stackrel{\triangle}{=} \frac{1}{2\pi j} \oint_{C} s^{n} M(s) ds$$
; for $n = 0, 1, 2, ... 2N_{p}$ (3.9)

In terms of the residues of equation (3.8), the residue moments are given by

$$D_0 = R_1 + R_2 + \dots + R_{N_p} = \sum_{q=1}^{N_p} R_q$$
 (3.10.0)

$$p_1 = p_1 R_1 + p_2 R_2 + \dots + p_{N_p} R_{N_p} = \sum_{q=1}^{N_p} p_q R_q$$
 (3.10.1)

$$D_{N_{p}} = p_{1}^{N_{p}} R_{1} + p_{2}^{N_{p}} R_{2} + \cdots + p_{N_{p}}^{N_{p}} R_{N_{p}} = \sum_{q=1}^{N_{p}} p_{q}^{N_{p}} R_{q}$$
 (3.10.N_p)

$$D_{2N_p} = p_1^{2N_p} R_1 + p_2^{2N_p} R_2 + \dots + p_{N_p}^{2N_p} R_{N_p} = \sum_{q=1}^{N_p} p_q^{2N_p} R_q \quad (3.10.2N_p)$$

The above system of equations can also be written compactly as

$$D_{n} = \sum_{q=1}^{N_{p}} p_{q}^{n} R_{q}; \text{ for } n = 0, 1, 2, ..., 2N_{p}$$
 (3.11)

Let us recall that we are trying to determine the order and the zeros of the polynomial A₂ (s) which give the unmber and locations of the poles of M(s) within C.

$$A_{2int}(s) = \prod_{k=1}^{N_p} (s - p_k) \equiv \sum_{k=0}^{N_p} a_k s^k$$
 (3.12)

where the coefficient a_N of the highest degree term may be set equal to 1 without any loss of generality. We shall now eliminate the residues $(R_q$'s) from the system of equations (3.10) by making use of equation (3.12). To achieve this, consider the first (N_p+1) number of equations in (3.10) starting with (3.10.0) and ending with (3.10.N_p). Multiplying these equations respectively by the coefficients a_0 to a_N , would yield

$$a_{0}^{D_{0}} + a_{1}^{D_{1}} + a_{2}^{D_{2}} + \dots + a_{N_{p}-1}^{D_{N_{p}-1}} + a_{N_{p}}^{D_{N_{p}}}$$

$$= R_{1} \left(a_{0} + a_{1}^{p_{1}} + a_{2}^{p_{1}^{2}} + \dots + a_{N_{p}-1}^{p_{p}-1} + a_{N_{p}}^{p_{p}-1} + a_{N_{p}}^{p_{p}} \right)$$

$$+ R_{2} \left(a_{0} + a_{1}^{p_{2}} + a_{2}^{p_{2}^{2}} + \dots + a_{N_{p}-1}^{p_{p}-1} + a_{N_{p}}^{p_{p}-1} + a_{N_{p}}^{p_{p}} \right)$$

$$\vdots$$

$$\vdots$$

$$+ R_{N_{p}} \left(a_{0} + a_{1}^{p_{N_{p}}} + a_{2}^{p_{N_{p}}} + \dots + a_{N_{p}-1}^{p_{N_{p}-1}} + a_{N_{p}}^{p_{p}-1} + a_{N_{p}}^{p_{N_{p}}} \right)$$

$$= \sum_{q=1}^{N_p} R_q \left[\sum_{k=0}^{N_p} a_k p_q^k \right]$$

$$= \sum_{q=1}^{N_p} \left[R_q A_{2_{int}} (p_q) \right]$$

$$= 0 (3.13)$$

because p_q for $q = 1, 2, ... N_p$ are the zeros of A_2 (s) = 0. Thus we have

$$a_0^{D_0+a_1^{D_1+a_2^{D_2}}} + \dots + a_{N_p^{D_N}} = 0$$
 (3.14)

Continuing this above procedure of successively multiplying a set of $(N_p + 1)$ equations from the system of equation (3.10) and using equation (3.12) will eliminate all of the residues $(R_q$; for $q = 1, 2, ...N_p)$ and lead to the following matrix equation

residues
$$(R_q; \text{ for } q = 1, 2, ... N_p)$$
 and lead to the following matrix equation

$$\begin{bmatrix}
D_0 & D_1 & D_2 & ... & ... D_{N_p} \\
D_1 & D_2 & D_3 & D_{N_p+1} \\
D_2 & D_3 & D_4 & D_{N_p+2} \\
... & ... & ... \\
... & ... & ... \\
D_{N_p} & D_{N_p+1} & D_{N_p+2} & ... & ... D_{2N_p}
\end{bmatrix} \begin{bmatrix}
a_0 \\
a_1 \\
a_2 \\
... \\
a_{N_p} & ... \\
0
\end{bmatrix} (3.15a)$$

or ·

$$\sum_{n=0}^{N_p} D_{m+n} a_n = 0; \text{ for } m = 0, 1, 2 \dots N_p$$
 (3.15b)

In this matrix equation, D's are the residue moments defined by equation (3.9) and a's are the coefficients of the polynomial A_2 (s), the zeros of which, are the poles of given meromorphic function M(s) within the contour C. From equation (3.15), we observe the following

a) If
$$N_p = 0$$
, then $D_i = 0$; for $i = 0, 1, 2, ...$ (3.16)

b) If
$$N_p = 1$$

$$D_i + a_0 D_{i-1} = 0 ; \text{ for } i = 1, 2, 3, ...$$
 (3.17)

c) If
$$N_p = 2$$

$$D_i + a_1 D_{i-1} + a_0 D_{i-2} = 0 ; \text{ for } i = 2, 3, 4..$$
(3.18)

d) If
$$N_p = 3$$

$$D_i + a_2 D_{i-1} + a_1 D_{i-2} + a_0 D_{i-3} = 0 ;$$
for $i = 3, 4, 5 ... (3.19)$

e) If
$$N_p = 4$$

$$D_i + a_3 D_{i-1} + a_2 D_{i-2} + a_1 D_{i-3} + a_0 D_{i-4} = 0 ;$$
For $i = 4, 5, 6, 7, ...(3.20)$

... etc.

Put differently, N_p will be = R - 1, where R = Rank of the infinite version of the D matrix of equation (3.15). However equations (3.16) thru (3.20 ...) are more useful in determining N_p , because as a by-product they yield the

coefficients a_k for $k = 0, 1, 2, ..., N_p$, as well. This will be illustrated as follows, for example, if $N_p = 3$, equation (3.19) for i = 3, 4 and 5 will give

$$D_{3} + a_{2}D_{2} + a_{1}D_{1} + a_{0}D_{0} = 0$$

$$D_{4} + a_{2}D_{3} + a_{1}D_{2} + a_{0}D_{1} = 0$$

$$D_{5} + a_{2}D_{4} + a_{1}D_{3} + a_{0}D_{2} = 0$$
(3.21)

which may be used in solving for a_0 , a_1 and a_2 . These a's may then be used in

$$D_{i} + a_{2}D_{i-1} + a_{1}D_{i-2} + a_{0}D_{i-3} = 0$$
;
for $i = 6, 7, 8 \dots$ (3.22)

to ensure that N is indeed 3. With the value of N and the coefficients of A (s) polynomial known, it is a simple matter to solve for the locations p_k of the poles of the given meromorphic function M(s) within the contour C being considered.

It is emphasized, at this stage that there are a few numerical pitfalls in implementing this scheme and section IV will address these problems specifically.

B. Zero finder

We still need to find the number $N_{\rm O}$ and locations $z_{\rm i}$, $i=1,\,2,\,\ldots\,N_{\rm O}$ of the zeros of the given meromorphic function M(s) in the given contour C. This is a rather trivial numerical exercise by virtue of the fact that the pole finder described above is independent of the presence or absence of zeros. In view of this, if we worked with the reciprocal function $M^{-1}(s)$, the zeros which now become poles inside contour C, are easily determined by using the pole finding scheme.

IV. Numerical Implementation and Results

In this section, we deal with the numerical implementation of the pole and zero finding schemes described in the preceding section.

Given the function and a rectangular region \mathcal{C} the normalized complex plane, we initially divide the region into a number of subcontours of approximately unit sized square regions (see Figure 4.1). Improved accuracy is obtained by centering each subcontour around the point 1 + j0 in the complex plane, via a simple change of variable. The need for this change of variable is explained in detail, later in this section, while describing the subroutine RESIDUE in which all of the residue moments of equation (3.9) are computed. Function values are computed at locations on the subcontour, determined by a 40-point Gaussian quadrature integration scheme and these values are stored in a complex array. In a sequential fashion, the stored function values are recalled and normalized for each of the subcontours of approximate size unity per side. The exponential normalization of the function, intended to improve the quality and accuracy of the location of the singularities (poles), is well described in the previous work under section III C of reference [2]. It is noted that the normalization function is an entire function with no zeros or poles in or on the subcontour, so that the singularities of the original function M(s) are undisturbed. reference to the typical subcontour $C_{m,n}$ shown in Figure 4.1, the entire function E(s) used in the process of normalization is given by

$$E(s) = u e^{VS}$$
 (4.1)

where u and v are real constants given by

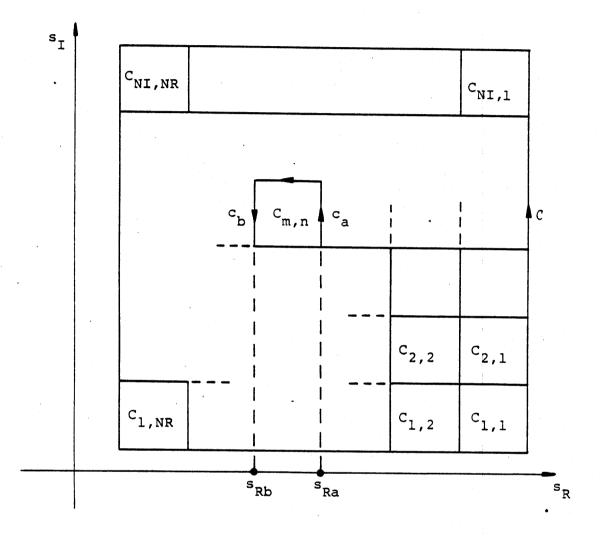


Figure 4.1 Division of the given rectangular domain into smaller rectangular (or square) subcontours

Note. $C_{m,n}$ is a typical subcontour and c_a and c_b are the parts of $C_{m,n}$ with constant real parts s_{Ra} and s_{Rb} respectively.

$$v = \frac{1}{(s_{Ra} - s_{Rb})} \ln \left(\frac{\text{average of } |M(s)| \text{ on } c_a}{\text{average of } |M(s)| \text{ on } c_b} \right)$$
 (4.2)

$$u = \exp(-vs_{Ra}) \times (average of |M(s)| on c_a)$$
or
$$(4.3)$$

= $\exp(-vs_{Rb}) \times (average of |M(s)| on c_b)$

Using the above entire function, the normalized function $M^{\text{nor}}(s)$ is obtained by using

$$M^{\text{nor}}(s) = M(s)/E(s)$$
 (4.4)

With this normalization scheme, the average magnitude of M(s) on c_a and c_b is unity and in general, does not depart significantly from unity on the rest of the subcontour. Also, effects of large phase variations in M(s) tend to become reduced when one works with $M^{nor}(s)$ instead. It is emphasized that, E(s) being an entire function does not disturb the singularities of the given function M(s).

A. Description of the Computer Programs

The following subroutines and function subprograms were written and they are useful in numerically evaluating the locations of poles and zeros of the given function M(s) which is meromorphic in a given region of the finite complex s-plane.

CONTOUR*, RESIDE, POLECHK, DETER*, ANGLER*,

POLY1, POLY2, POLY3, CFCTS

(* from Reference 2)

In what follows, we shall briefly describe each of these subroutines. Their listings are included in Appendix A.

a) Subroutine CONTOUR (CSL, CSF, NR, NI, KDM)

This program divides up the given scan area enclosed by the contour into a specified number of rectangular/square subcontours and calls subroutine RESIDUE once for each of the contours. The arguments appearing in this subroutine are as follows.

CSL : Coordinates of the upper left corner of C,

CSF : Coordinates of the lower right corner of C,

NR : Number of major divisions of the real axis within \mathcal{C} .

NI : Number of major divisions of the imaginary axis within C.

KDM: A multiplicative factor for Gaussian integration, i.e., the number of points in the integration is given by 40 × KDM per side of the subcontour. Although the largest allowable value is 4 because of the present dimensioning of the arrays in RESIDUE, KDM = 1 should be adequate in most cases.

This subroutine also summarizes the results by listing the location of all the poles and zeros found as well as the function values at the zeros and the reciprocal of the function value at the poles. These function values may be used by the user in judging the quality of pole-zero locations which are numerically determined.

b) Subroutine RESIDUE (CFCTS, CSM, CSMI, KDM)

This is the core subroutine in the entire package and essentially does the following:

 makes a change of variable so that the subcontour is now centered around the point (1 + j0) in the complex plane

- 2) computes the function value on this new subcontour at locations required by a 40-point Gaussian integration procedure, (the locations and the function values are stored in complex arrays CS and CF respectively),
- 3) normalizes the function values and computes the argument number and nine residue moments D_n for $n=0,1,\ldots,8$, of equation (3.9),
- 4) calls subroutine POLECHK which determines appotential value of the number of poles $N_{\rm p}$, the coefficients of the denominator polynomial and the pole locations as well,

and 5) goes through a similar procedure, working with the reciprocal function, to determine the location of zeros.

Various arguments of this subroutine are described below

CFCTS : User supplied function subprogram,

CSM : Coordinates of the upper left corner of the subcontour $C_{m,n}$,

CSMI : Coordinates of the lower right corner of the subcontour $C_{m,n}$,

KDM : Same as in subroutine CONTOUR.

In computing the residue moments, given by

$$D_{n} = \frac{1}{2\pi j} \int_{C_{m,n}}^{\infty} s^{n} M(s) ds$$
 (4.5)

a change of variable of the following form was found to improve the accuracy of the above integration,

 $z = s - s_C - 1$ (4.6)

where,

 s_c = coordinates of the center point of the subcontour $c_{m,n}$.

The reason for this change of variable is that when the subcontour $C_{m,n}$ is located away from the origin, the numerical value of the factor s^n in the integrand can become quite large compared with the average magnitude of unity for the normalized M(s) around the subcontour $C_{m,n}$. With the change of variable given by equation (4.6), the new subcontour in the z-plane is centered around the point 1+j0 so that the entire integrand will now have an average magnitude of unity resulting in improved accuracy for the numerical evaluation of the residue moments.

c) Subroutine POLECHK (DO, D, DAO, DA, NP)

This subroutine accepts the nine residue moments as input and determines a potential value for the number of poles $N_{\rm p}$ and the coefficients of the denominator polynomial as well. The various arguments of this subroutine are:

- DO : Zeroth residue moment D₀ which is simply the sum of the residues,
- D: A complex array which contains the residue mements D_n for n = 1, 2, ..., 8,
- DAO : Constant term in the denominator polynomial (a₀ of equation (3.12)),
- DA : A complex array which contains the remaining coefficients of the denominator polynomial,
- NP : Most likely value of the number of poles $N_{\mbox{\scriptsize p}}$ in the subcontour.

d) Subroutine DETER (CM, CB, CW, CV, MS, CD)

This subroutine finds the determinant of a given square matrix. A description of the various arguments of this subroutine is given below.

CM : Complex array continuing the matrix

elements,

CB, CW and CV : Complex working arrays, local to the

subroutine,

MS : Size of the input square matrix,

CD : Computed value of the determinant.

e) FUNCTION ANGLER (X, Y)

This function subprogram computes the phase $\, \varphi \,$ of a complex number in radians such that $\, 0 \, \leq \, \varphi \, \leq \, 2\pi \,$. The arguments are:

X : Real part of the complex number,

Y : Imaginary part of the complex number,

ANGLER: Phase ϕ of the complex number X + j Y

such that $0 \le \phi \le 2\pi$.

It was important to determine the phase in this range of $0 \le \varphi \le 2\pi$ for obtaining the argument number, rather than the commonly available range of $-\pi \le \varphi \le \pi$.

f) Subroutines POLY1 (C0, C1, CLIN)

POLY2 (CO, Cl, C2, CQUAD)

POLY3 (CO, C1, C2, C3, CUBE)

These three subroutines respectively solve for 1, 2 and 3 roots of the following linear, quadratic and cubic polynomials

$$C_0 + C_1 s = 0 (4.7a)$$

$$C_0 + C_1 s + C_2 s^2 = 0$$
 (4.7b)

$$C_0 + C_1 s + C_2 s^2 + C_3 s^3 = 0$$
 (4.7c)

when the appropriate coefficients are fed in. The arguments appearing in the three subroutines are:

CO, Cl, C2, C3 : Coefficients of the polynomial,

CLIN : Root of the linear equation (4.7a),

CQUAD : Two roots of the quadratic equation (4.7b),

CUBE : Three roots of the cubic equation (4.7c).

g) COMPLEX FUNCTION CFCTS (CS, CSHIFT)

This complex function subprogram numerically evaluates the meromorphic function M(s) for a prescribed s. The two arguments and the result of the function subprogram are

CS : Complex value of the variable s,

CSHIFT : Complex constant to facilitate a change of

variable (can be set equal to zero),

CFCTS : Function value.

This completes a brief description of the various subroutines and function subprograms in this package. It is recalled that, in the subroutine RESIDUE, we computed the nine residue moments D_n for $n=0,1,\ldots,8$, which introduces a limitation of no more than 3 poles in a subcontour which is cf approximate size unity square. Also when the number of zeros and poles $(N_0 + N_p)$ in any subcontour is more than 4, because of large changes in the function value in a small region, the residue moment calculations may not be sufficiently

accurate to determine the exact value of N_p . We have encountered no difficulty when $(N_0 + N_p) \le 4$ per subcontour and some modifications to improve the accuracy may be required for the rare occurrence in physical problems, when the poles and zeros are more closely bunched together. When $(N_0 + N_p) > 4$ per subcontour, another possibility is to divide the subcontour of unit size into smaller subcontours. This has not been automated in the present family of computer programs.

In the following two subsections, we consider two examples that validate the application of this pole finding scheme to determine the poles and zeros of given complex functions that are meromorphic in a specified region of the finite complex plane.

B. Ratio of Polynomials (Example 1)

Consider a meromorphic function $M_1(s)$ given by,

$$M_{1}(s) = \frac{\prod_{p=1}^{16} (s-z_{p})}{\prod_{q=1}^{10} (s-p_{q})}$$
(4.8)

which is readily seen to have 16 zeros and 10 poles in the complex s-plane. The locations of zeros indicated by z, and poles indicated by p are shown in Figure 4.2. The zeros are given by;

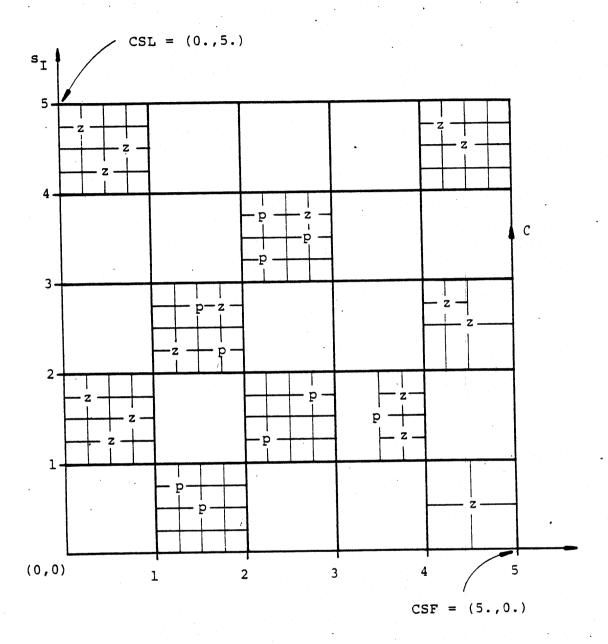


Figure 4.2. The pole-zero configuration of the given ratio of polynomials (example 1).

Note: 1) C is the given square scan area, uniquely specified by the points CSF and CSL

- 2) $z \rightarrow location of a zero$
- 3) $p \rightarrow location of a pole$

$$z_1 = 4.50 + j0.50, \ z_2 = 4.25 + j4.75, \ z_3 = 3.75 + j1.25, \ z_4 = 3.75 + j1.75$$

$$z_5 = 4.50 + j2.50, \ z_6 = 4.25 + j2.75, \ z_7 = 2.75 + j3.75, \ z_8 = 1.25 + j2.25$$

$$z_9 = 1.75 + j2.75, \ z_{10} = 4.50 + j4.50, \ z_{11} = 0.75 + j1.50, \ z_{12} = 0.50 + j1.25$$

$$z_{13} = 0.25 + j1.75, \ z_{14} = 0.50 + j4.25, \ z_{15} = 0.75 + j4.50, \ z_{16} = 0.25 + j4.75$$

$$(4.9)$$

The poles are given by:

$$p_1 = 3.50+j1.50$$
, $p_2 = 1.75+j2.25$, $p_3 = 1.50+j2.75$, $p_4 = 2.25+j3.25$
 $p_5 = 2.75+j3.50$, $p_6 = 2.25+j3.75$, $p_7 = 1.25+j0.75$, $p_8 = 2.75+j1.75$
 $p_9 = 1.50+j0.50$, $p_{10} = 2.25+j1.25$ (4.10)

In Figure 4.2, the scan area which is a square of side 5 units with its lower left corner as the origin of the complex s-plane is indicated by the counterclockwise contour \mathcal{C} . This contour is divided into 25 subcontours and the pole-zero locations are indicated in the various subcontours. The formal input variables CSL and CSF that uniquely specify the scan. area are also shown in the Figure 4.2. The results of using this function in order to recover the poles and zeros are presented in Table 1, in the same format as the subroutine CONTOUR would summarize. Comparing the results of Table 1 with those of equations (4.9) and (4.10), it is seen that the poles and zeros of this ratio of polynomial given by $M_1(s)$ are recovered with a high degree of accuracy. With the view of introducing large phase variations in $M_1(s)$, it was multiplied by an entire function of the form eAS. Define a new function N₁(s)

$$N_1(s) = e^{AS} M_1(s)$$
 (4.11)

where $M_1(s)$ is given by equation (4.8). The real

SHMMARY	OF	RESULTS
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				DOLLHILL	*	* *	* * * * * * * * *	. * * * * * * * * * * * * * * * * * * *
*	*	* * * * *	. * * * * * * *	* * * * * * * * * * * * * * * * * * *	* * ^ ^ "		Real part	Imag. part
			Real part	Imag. pare		=	51481437E-16	79124576E-16
	1	POLE AT	.35000000E+01	.15000000E+01	·	=	16026682E-15	52103539E-16
	2	POLE AT	.27500000E+01	.17500000E+01		=	27757148E-16	70485840E-17
	3	POLE AT	.22500000E+01	.12500000E+01	1/F (POLE)	=	.34166970E-10	.48627420E-10
	4	POLE AT	.22500000E+01	.37499999E+01	1/F (POLE)		.25440087E-10	.64503667E-11
	5	POLE AT	.22500000E+01	.32500000E+01	1/F (POLE)	=	.63817493E-10	92700647E-10
	6	POLE AT	.27499999E+01	.35000000E+01	1/F (POLE)		.85008163E-17	32501785E-17
	7	POLE AT	.15000000E+01	.50000000E+00	•	=	.15115588E-16	46924316E-17
	8	POLE AT	.12500000E+01	.75000000E+00	1/F (POLE)	=	18510407E-14	.46241384E-15
	9	POLE AT	.17500000E+01	.22500000E+01	1/F (POLE)		04130959F-14	.24127978E-14
	10	POLE AT	.15000000E+01	.27500000E+01	1/F (POLE)	* *	* * * * * * * *	* * * * * * * * *
,	* *	* * * *	* * * * * * * * * * * * * * * .4500000E+01	.50000000E+00	F(ZERO)	=	33974973E-10	.37801892E-10
		ZERO AT	.45000000E+01	.25000000E+01	F(ZERO)	=	.71209024E-11	91285320E-11
	2	ZERO AT		.27500000E+01	F(ZERO)	=	.14853889E-11	10834960E-10
2	3		.42500000E+01	.45000000E+01	F(ZERO)	=	.60244057E-11	.38265879E-11
8	4		.45000000E+01	.47500000E+01	F(ZERO)	=	.31923752E-11	.52167555E-11
	5		.42500000E+01	.17500000E+01	F(ZERO)	=	34168506E-11	.23797403E-09
	6		.37500000E+01	.12500000E+01	F (ZERO)	<u>.</u>	27022936E-09	.28549864E-09
	7	ZERO AT	.37500000E+01	.37500000E+01	F(ZERO)	= .	10640196E-09	31899668E-09
	8	ZERO AT	.27500000E+01	.27500000E+01	F(ZERO)	=	32643289E-08	54749864E-09
	. 9	ZERO AT	.17500000E+01	.22500000E+01	F(ZERO)	=	48766628E-09	.13125755E-09
	1	O ZERO AT	.12500000E+01	.17500000E+01	F(ZERO)	=	25715938E-06	.28969053E-06
	1	1 ZERO AT			F(ZERO)	=	55518217E-06	.30493810E-06
	1	2 ZERO AT	· · · · · · · · · · · · · · · · · · ·	.12500000E+01	F(ZERO)	=	48376999E-06	.23146903E-06
	1	3 ZERO AT		.15000000E+01	F (ZERO)	==	.30761459E-07	.49960265E-07
	1	4 ZERO AT		.47500000E+01	F (ZERO)	. =	.31526461E-08	.25607578E-07
	1	5 ZERO AT		.42500000E+01	F (ZERO)	-	.22238213E-07	.14267090E-07
	1	6 ZERO AT	.75000000E+00	.45000000E+01	F (ZERO)	* *	* * * * * * * *	* * * * * * * *
	*	* * * * *	**********		•			

Table 1. The poles and zeros of $M_1(s)$ found by the present method

constant A in the exponent was varied and the results (pole-zero locations) did not vary significantly in regions where the singularities were not dense. This can be attributed to the efficient exponential normalization procedure, described earlier. An improved scheme of determining $N_{\rm p}$ will be necessary for highly oscillatory functions.

C. Input Impedance of a Biconical Antenna (Example 2)

The input impedance of a biconical antenna may be written as [6,7],

$$Z_{in}(kl) \simeq Z_{c} \left[\frac{e^{jkl} + T(kl) e^{-jkl}}{e^{jkl} - T(kl) e^{-jkl}} \right]$$
(4.12)

where

 ℓ \equiv slant height of the cone

Z = characteristic impedance of a symmetrical bicone

$$= \frac{z_0}{\pi} \ln \left(\cot \frac{\theta}{2}\right) \simeq \frac{z_0}{\pi} \ln \left(\frac{2}{\theta}\right) \text{ for small angles}$$

Z = characteristic impedance of free space

 θ = half angle of the bicone in radians

 $k \equiv free space propagation constant$

T(kl) = effective terminal reflection coefficient

Rewriting the input impedance in the normalized Laplace transform variable plane

$$s = \frac{s\ell}{\pi c} \tag{4.13}$$

$$\tilde{z}_{in}(s) = \left(\frac{z_{in}}{z_{c}}\right) = \left\{\frac{e^{2\pi\delta} + \tilde{T}(s)}{e^{2\pi\delta} - \tilde{T}(s)}\right\}$$
(4.14)

where

$$k\ell = -j\pi\delta \tag{4.15a}$$

$$\tilde{\mathbf{T}}(\delta) = \left\{ \frac{1 - \tilde{\mathbf{y}}(\delta)}{1 + \tilde{\mathbf{y}}(\delta)} \right\} \tag{4.15b}$$

 $\tilde{y}(s)$ = normalized terminal admittance = $Z_C \tilde{Y}(s)$

$$\frac{1}{4\pi} \left(\frac{Z_{O}}{Z_{C}} \right) \left[2 \operatorname{Ein}(2\pi\delta) + e^{2\pi\delta} \left\{ \ln(2) + \operatorname{Ein}(2\pi\delta) - \operatorname{Ein}(4\pi\delta) \right\} + e^{-2\pi\delta} \left\{ -\ln(2) + \operatorname{Ein}(-2\pi\delta) \right\} \right]$$

$$(4.15c)$$

and Ein(z), following the notation of Ref. [8], is the exponential integral given by

Ein(z) =
$$\int_{0}^{z} \frac{1 - e^{-t}}{t} dt$$
 (4.16)

Equation (4.14) was used as the input mermorphic function and its zeros and poles were accurately determined and they were found to be in excellent agreement (5 places or better) with earlier results [7] which employed a rather tedious 2-dimensional search method. The results for the case of θ = .001° is shown plotted in figure 4.3.

In concluding this section, we note that the two examples considered here have shown that an application of Cauchy's residue theorem is very useful in determining the pole and zero locations of a complex function of a complex variable which is meromorphic in a given region.

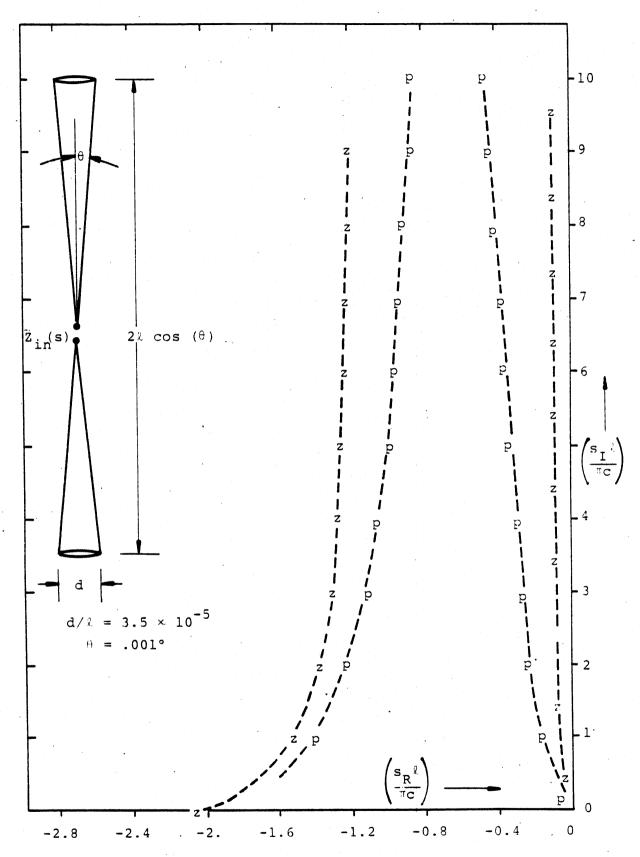


Figure 4.3 The pole-zero configuration of the input impedance $\tilde{z}_{in}(s)$ of a biconical antenna (Example 2)

V. Summary

In this note, we have developed a pole finding procedure based on an application of Cauchy's residue theorem. The validity of the procedure in determining the poles and zeros of a meromorphic function has been demonstrated with two numerically illustrated examples of i) a ratio of polynomials, and ii) the input impedance of a biconical antenna. It is to be emphasized that a major problem in this procedure lies in unambigiously in a given contour. determining the number of poles N_n This is further discussed in Appendix B. In view of this, it is expected that this procedure may be successfully applied in physical problems where the user has some a priori knowledge of the behaviour of the meromorphic function. In other instances, modifications in terms of improving the accuracy with which the residue moments are determined, may be required to circumvent certain problems e.g., highly oscillatory functions or functions with dense population of poles and zeros.

This work is a sequel to an earlier work [2] which concerned itself with the numerical evaluation of the zeros of an analytic function.

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APPENDIX A

Computer Program Listings

The following TEST Program illustrates the way in which this package entitled CONTOUR may be called as a subroutine.

PROGRAM TEST (INPUT, OUTPUT)
IMPLICIT COMPLEX (C)
CJMMCN NOZ, CZ(30), CFZ(30), NGP, CP(30), CFP(30), IWARN
CSL=(0.,5.)
CSF=(5.,0.)
NR=5
NI=5
KDM=1
CALL CONTOUR(CSL, CSF, NR, NI, KDM)
END

```
SUBROUTINE CONTOUR(CSL,CSF,NR,NI,KOM)
   IMPLICIT COMPLEX (C)
   COMMON NUZ, CZ (30), CFZ (30), NOP, CP (30), CFP (30), IHARN
   EXTERNAL CFCTS
   NUZ = C
   NOP=0
   I WARN=0
   RCSF=REAL (CSF)
   SCIF=4IMAG(CSF)
   RCSL =REAL(CSL)
   SCIL=AIMAG(CSL)
   RINC = (RCSL-RCSF) / FLOAT (NR)
   SCIINC=(SCIL-SCIF)/FLOAT(NI)
   ND4= (40 *K DM) * 4
   PRINT 810, RCSF, SCIF, RCSL, SCIL, NO4, NR. NI
810 FORMAT (1H1,2X,56HCOORDINATES OF THE LOWER RIGHT CORNER OF SCAN AR
   1EA ARE (,1X,F12.8,1H,,F12.8,1X,1H),/,3X,56HCDORDINATES OF THE UPPE
   2R LEFT CORNER OF SCAN AREA ARE (.1X.F12.8,1H.,F12.8,1X,1H).//.3X,
   343HTOTAL NUMBER OF PUINTS USED PER CONTOUR ARE, 2X, 14, //, 3X, 43HNUMB
   4ER OF DIVISIONS ALONG THE REAL AXIS ARE, 2X, 12, 5X, 44HNUMBER OF DIVI
   5SIONS ALONG THE IMAG. AXIS ARE, 2X, 14, //, 3X, 100H***************
   7************
    00 72C J=1.NR
    1-1=11
    SRMI=RCSF+RINC*FLOAT(J)
    SRM=RCSF+RINC*FLOAT(JJ)
    00 710 I=1,NI
    11=1-1
    SIM=SCIF+SCIINC*FLOAT(I)
    SIMI=SCIF+SCIINC*FLOAT(II)
    CSM=CMPLX(SRMI,SIM)
    CSMI=CMPLX(SRM.SIMI)
    CALL RESIDUE(CFCTS, CSM, CSMI, KDM)
    CONTINUE
710
    CONTINUE
720
    PRINT 730. IWARN
    FORMAT (1H1,50X, #SUMMARY OF RESULTS#,/,3X, #NO. OF WARNING MESSAGES
730
    2 = *, [3]
     PRINT 747
     IF (NOP.EQ. 0) GO TO 745
     DO 740 M=1.NOP
     PRINT 760, M, CP(M), CFP (M)
    CONTINUE
740
     IF (NOP.LE.30) GO TO 745
     PRINT 742 , NOP
742 FORMAT (1HO, //, 3X, # FOUND #, 2X, 13, 2X, #POLES. INCREASE DIMENSION OF
    S CP AND CFP ARRAYS ACCORDINGLY#1
745 CONTINUE
     IF (NUZ.EQ.0) GD TG 780
     FORMAT (1HC;/;3X, *****************************
     PRINT 747
    747
     00 750 M=1,NOZ
     PRINT 770.M.CZ(M) (CFZ(M)
    CONT INUE
750
     IF (NOZ.LE.30) GO TO 755
```

PRINT 752, NOZ FORMAT (1HO,//,3X, #1 FOUND #, 2X, 13, 2X, #ZEROS. INCREASE DIMENSION OF \$ CZ AND CFZ ARRAYS ACCORDINGLY#) CONTINUE PRINT 747 FORMAT (1H0,3x,13,2x, #POLE AT#, 2E20.8,8x, #1/F(POLE) 760 FORMAT (1HO, 3X, 13, 2X, #ZERO AT#, 2E20.8, 8X, #F(ZERO) 770 RETURN 780 PRINT 750 FORMAT (1HO, 10X, #SORRY .. RESIDUE. . COULD NOT FIND ANY ZEROS OR 790 SPOLES IN THE GIVEN SCAN AREA +,//) RETURN END

```
SUBROUTINE RESIDUE(CFCTS.CSM.CSMI.KDM)
 IMPLICIT COMPLEX (C)
COMMON NOZ, CZ (30), CFZ (30), NGP, CF(30), CFP(30), IWARN
 DIMENSION CS(641), ARG(641), CF(641), CSTO(641)
 DIMENSION X(40).W(40).CD(8).CA(3).CQUAD(2).CUBE(3)
 DATA NP/40/, X/-.998237709710555,-.990726238699457,-.9772599499837
174, -.957916819213792, -.932812808278677, -.902098806968874, -.8659595
203212260,-.824612230833312,-.778305651426519,-.727318255189927,-.6
371956684614180,-.612553889667980,-.549467125095128,-.4830758016861
479,-.413779204371605,-.341994090825758,-.268152185007254,-.1926975
580701371,-.116084070675255,-.387724175060508E-1,.387724175060508E-
61,.116084070675255,.192697580701371,.268152185007254,.341994090825
7758, 413779204371605, 483075801686179, 549467125095128, 6125538896
867980, 671956684614180, 727318255189927, 778305651426519, 82461223
90833312,.865959503212260,.902098806968874,.932812808278677,.957916
$819213792,.977259949983774,.990726238699457,.998237709710559/
 DATA W/. 45212770985332E-2,.104982845311528E-1,.164210583819079E-1
1. .22 2458491941670E-1.. 279370069800234E-1.. 334601952825478E-1.. 3878
221679744720E-1,.438709081856733E-1,.486958076350722E-1,.5322784698
339368E-1,.574397690993916E-1,.613062424929289E-1,.648040134566010E
4-1,.679120458152339E-1,.706116473912868E-1,.728865823958041E-1,.74
57231690579683E-1,.761103619006262E-1,.770398181642480E-1,.77505947
69784248E-1,.775059479784248E-1,.770398181642480E-1,.76110361900626
72E-1, .747231690579683E-1, .728865823958041E-1, .706116473912868E-1, .
8679120458152339E-1, .648040134566010E-1, .613062424929289E-1, .574397
9690993916E-1,.532278469839368E-1,.486958076350722E-1,.438709081856
$733E-1,.387821679744720E-1,.334601952825478E-1,.279370069800234E-1
$,.222458491941670E-1,.164210583849079E-1,.104982845311528E-1,.4521
 $27709853319E-2/
  IF (KDM.EQ.O) KDM=1
 CZERG=(0.,0.)
 CAUX 1=CSM
 CAUX2=CSMI
 ND=KDM*NP
  NC1=NC+1
  ND2=2*ND
  ND 3= 3*ND
  4* QN =4 CN
  NM1=ND4+1
  NMM1=ND4-1
  NDP=NC2+1
  C1=(1.,0.)
  RCSF=REAL(CSMI)
  RCSL=REAL(CSM)
  SCIF = AIMAG(CSMI)
  SCIL=AIMAG(CSM)
  PI=3.14159265
  P13=1.5*PI
  IP I = 2. *P I
  HPI= .5*PI
  NR = 1
  N 1=1
  [CKL=1
   ICKU=2
  CONT INUE
   DO 655 ICK=ICKL, ICKU
```

RINT=(RCSL-RCSF)/FLCAT(NR)

```
SCINT=(SCIL-SCIF)/FLOAT(NI)
      00 650 JT=1.NR
      JJT=JT-1
      SRMI =RCSF+RINT*FLCAT(JT)
      SRM=RCSF+RINT*FLOAT(JJT)
      SAUX 1 = SRMI
      SAUX 2= SRM
      DO 640 IT=1,NI
      IIT=IT-1
      SIMI=SCIF+SCINT*FLOAT(IIT)
      SIM=SCIF+SCINT*FLOAT(IT)
      IF (ICK.GT.1) GG TO 20
 10
      CONTINUE
      PRINT 660, SRM, SIMI, SRMI, SIM
      PRINT 150
      GO TO 25
      PRINT 200
20
 25
      CONTINUE
C
C
      CHANGE OF VARIABLE
C
      SMR=(SRM+SRMI)/2.
      SMI=(SIM+SIM[)/2.
      CSC=CMPLX(SMR,SMI)
      CSC=CSC-(1.,0.)
      CSM=CMPLX(SRM,SIMI)-CSC
      CSMI = CYPLX(SPMI, SIM) - CSC
      SRM=REAL(CSMI)
      SRMI = REAL (CSM)
      SIMI = AI MAG (CSMI)
      SIM=AIMAG(CSM)
      DELX=(SRM-SRMI)/(2.*FLOAT(KDM))
      DELY=(SIM-SIMI)/(2.*FLOAT(KDM))
      DO 140 K=1,4
      KK=K-1
      KKK=KK+KDM+NP
      IF (K-2) 40,50,30
   30 IF (K-3) 50,60,70
   40 YU=SIM.
      YL=SIMI
      XC=SRM
      GO TO 110
   50 XU=SRMI
      XL=SRM
       YC=SIM
       GO TO 80
   60 YU=SIMI
       YL = SIM
       XC=SRMI
       GC TC 110
   70 XU=SRM
       XL=SRMI
       YC=SIMI
    80 DL=(XU-XL1/FLOAT(KDM)
       DO 100 L=1,KDM
       LL=L-1
       LLL=LL *NP+KKK
       XLOW=XL+FLOAT(LL)*DL
```

```
XUP=XLOW+OL
    DLX=(XUP-XLOW)/2.
    PLX=(XUP+XLOW)/2.
    DB 90 M=1,NP
    MM=M+LLL
    CS(MM)=CMPLX(DLX+X(M)+PLX,YC)
 90 CENTINUE
JUNITHDD COL
    GO TO 140
110 DL=(YU-YL)/FLGAT(KDM)
    DO 130 L=1,KDM
    LL=L-1
    LLL=LL*NP+KKK
    YLOW=YL+FLOAT(LL) *DL
    YUP=YLOW+DL
    DLY=(YUP-YLCW)/2.
    PLY= (YUP+YLOW) /2.
    DQ 120 M=1.NP
     MM=M+LLL
    CS(MM)=CMPLX(XC,DLY+X(M)+PLY)
120 CONTINUE
130 CONTINUE
140 CONTINUE
    FORMAT(1HO, 2X, #WORKING WITH THE FUNCTION F(S) TO EXTRACT POLES....
    2 . . . . * 1/1
    FORMAT (1HO+2X, #WORKING WITH THE RECIPROCAL OF F(S) TO EXTRACT ZER
200
    205 . . . . . . . . . . . . / )
     CONTINUE
220
     00 250 K=1,NO4
     IF (ICK.EQ.1) CF(K)=CFCTS(CS(K),CSC)
     IF(ICK.EQ.2) GC TC 225
     CSTO(K)=CF(K)
     GO TO 250
    CF(K)=1./CSTO(K)
225
     CONT INUE
250
     CS(NM1)=CS(1)
     A SUM = 0.
     DG 300 K=1,ND
     ASUM=ASUM+CABS (CF (K))
 300 CONTINUE
     A 1=ASUM/FLOAT (ND)
     A SUM=0 .
     DO 310 K=NDP+ND3
      ASUM=ASUM+CABS(CF(K))
 310 CONTINUE
      A 2=A SUM/FLOAT (NC)
      A3=ALUG(A1/A2)/(SRM-SRMI)
      A4=A1*EXP(-A3*SRM)
      CSUM=( 0. , 0.)
      DO 320 K=1,ND4
      CSUM=CSUM+CF(K)
      CF(K)=CF(K)/(A4+CEXP(A3+CS(K)))
  320 CONTINUE
      CAVE =CSUM/FLOAT(NC4)
      AVEE = CABS(CAVE)
      CONT INUE
 325
      00 330 L=1.ND4
      RF=REAL(CF(L))
```

```
Z=AIMAG(CF(L))
    ARG(L)=ANGLER(RF,Z)
330 CONTINUE
    10 = 0
    OVF=ARG(1)
    OV=ARG(1)
    DVL=ARG(ND4)
    DO 390 K=1, NMM1
    IF (OV.GT.PI3.AND.OV.LT.TPI) GG TG 340
    IF (UV.GT.O..AND.DV.LT.HPI) GO TO 350
    GG TO 380
340 IF (ARG(K+1).GT.O..AND.ARG(K+1).LT.HPI) GO TO 360
    GO TO 380
350 IF (ARG(K+1).GT.P13.AND.ARG(K+1).LT.TP1) GO TO 370
    GO TO 380
360 IU=IC+1
    GO TO 380
370 IO=IO-1
380 OV= ARG(K+1)
    ARG(K+1)=IO*T PI+ARG(K+1)
390 CONTINUE
    IF (OVL.GT.PI3.AND.OVL.LT.TPI) GO TO 400
    IF (OVL.GT.O..AND.OVL.LT.HPI) GO TO 410
    GO TO 440
400 IF (OVF.GT.O..AND.OVF.LT.HPI) GC TO 420
    GO TO 440
410 IF (OVF.GT.PI3.AND.OVF.LT.TPI) GO TO 430
    GO TO 440
420 [G=IO+1
    GD TO 440
430 [C=IO-1
440 CONTINUE
     ARG(NM1)=FLOAT(IU)*TPI+ARG(1)
     IF (ICK.EQ.1) IAN1= IO
     IF (ICK.LT.2) GB TC 444
     [AN2=10
     ICHECK=IAN1+IAN2
     IF (ICHECK.EQ.O) GO TC 444
     IWARN= IWARN+1
     PRINT 442, IWARN, IAN1, IAN2
442 FORMAT [1HC, //, 3X, #WARNING NUMBER = #, [3, //, 3X,
    2 ARGUMENT NUMBER OF F(S)
                                  = \pm .13./.3X.
    3 # ARGUMENT NUMBER OF 1/F(S) = #, 13,/)
444 CONTINUE
     FINDING THE RESIDUE MOMENTS DO THRU D8.
     DO 520 L=1.9
     CON1=(C.,C.)
     CON2= (0.,0.)
     CON3=(0.,C.)
     CON4=10.,0.1
     LM1=1.-1
     DG 500 K=1.4
     KK=(K-1)*KDM*N?
     IF (K-2) 450, 465, 445
     IF (K-3) 465,480,495
445
450
     DO 460 M=1,KDM
     MM = (M-1) * NP
     DO 455 N=1,NP
```

```
NN=KK+MM+N
     CON1=CON1+((CS(NN)**LN1))*W(N)*CF(NN)
455
     CONTINUE
     CONTINUE
460
     GO TO 498
465
     DO 475 M=1.KDM
     MM=(M-1)*NP
     00 470 N= 1, NP
     NN=KK+MM+N
     CON2=CON2+(CS(NN) ** LM1) *W(N) *CF(NN)
470
     CONTINUE
     CONTINUE
475
     GO TO 498
480
     DO 490 M=1.KDM
     MM= (M-1) *NP
     DO 485 N=1,NP
     NN=KK+MM+N
      CON3 = CON3 + ( (CS(NN) * * LM1)) * W(N) * CF(NN)
485
     CONTINUE
     CONT INUE
490
      GO TO 498
      DO 497 M=1.KDM
495
      MM = (M-1) + NP
      DO 456 N=1.NP
      NN=KK+MM+iV
      CCN4=CON4+((CS(NN)++LM1))+W(N)+CF(NN)
496
      CCNTINUE
      CONTINUE
497
498
      CONTINUE
 500
      CONTINUE
      CON1=CON1+DELY+(0.,1.)
      CCN2=-CON2*DELX
      CON3 = CON3 + OELY + (0., -1.)
      CUN4=CUN4*DELX
      IF (L.GT.1) GG TO 510
      CD0=(CON1+CGN2+CON3+CON4)/(TPI*(0 -+ 1 -))
      GO TO 520
      CD(LM1)=(CON1+CON2+CON3+CON4)/(TPI*(G.,1.))
 510
      CONTINUE
 520
      ALL THE 9 RESIDUE MOMENTS ARE NOW COMPUTED.
C
      CALL POLECHK (CDO, CD, CAO, CA, NQ)
      NP1=NC+1
      IF (ICK.EQ.L) NQ1=NQ
      IF (ICK.EQ.2) NQ2=NQ
      IF (NO-4) 521,570,570
 521 IF (ICK.LT.2) GD TO 525
      NCK=NQ2-NQ1-IAN1
      IF (NCK.EQ.0) GO TO 525
       IWARN= [WARN+1
      PRINT 522, IWARN, NQ2, NQ1, IAN1
 522 FORMAT (1HC,//,3x, *WARNING NUMBER = *,13,//,3x, *NUMBER OF ZEROS NO
      22 = #, 13, 5x, #NUMBER OF POLES NQ1 = #, 13, /, 3X,
      3 + (NQ2-NQ1) DOES NOT EQUAL THE ARGUMENT NUMBER IAN1 = +, 13,/)
 525
      CONTINUE
       GO TC 528
       IF (NG-4) 528,570,570
 526
 528
      CONTINUE
       GO TO (530,540,550,560),NP1
```

```
PRINT 590, (ICK, NQ, CAO, (CA(K), K=1,3))
530
     IF (ICK.EQ.2) PRINT 620
     GO TO 580
     PRINT 590, (ICK, NQ, CAO, (CA(K), K=1,3))
540
     CALL POLY 1(CAO,C1,CL[N)
     IF (ICK.EQ.2) CLIN=CLIN+CSC
     IF (ICK.EQ.2) GO TO 545
     PRINT 600 CLIN
     NOP=NOP+1
     CP(NOP)=CLIN
     CFP(NOP) = 1./CFCTS(CLIN, CZERO)
     GO TO 580
     PRINT 610, CLIN
545
     NOZ=NOZ+1
     CZ(NOZ)=CLIN
     CFZ(NOZ)=CFCTS(CL IN, CZERO)
      PRINT 620
      GD TO 580
      PRINT 590, (ICK, NQ, CAO, (CA(K), K=1,3))
550
      CALL POLY 2(CAO, CA(1),C1,CQUAD)
      IF (ICK.EQ.1) GU TO 551
      CQUAC(1)=CQUAD(1)+CSC
      CQUAD(2) = CQUAD(2) + CSC
      GO TO 555
      CONTINUE
551
      PRINT 600, (CQUAD(K), K=1,2)
      DO 552 L=1.2
      NOP=NOP+1
      CP(NOP) = CQUAD(L)
      CFP(NOP)=1./CFCTS(CQUAD(L),CZERO)
      CCNTINUE
      GO TO 580
      PRINT 610, (CQUAD(K), K=1,2)
555
      DO 557 L=1.2
      NOZ=NOZ+1
      CZ(NOZ)=CQUAD(L)
      CFZ(NOZ) = CFCTS(CQUAD(L), CZERO)
      CONTINUE
 557
      PRINT 620
      GO TO 580
      PRINT 590, (ICK, NQ, CAO, (CA(K), K=1,3))
 560
      CALL POLY 3(CAO, CA(1), CA(2), C1, CUBE)
      IF (ICK.EQ.1) GC TO 561
      CUBE(1)=CUBE(1)+CSC
      CUBE(2)=CUBE(2)+CSC
      CUBE (3) = CUBE (3)+CSC
      GO TO 565
 561 - CONTINUE
      PRINT 600, (CUBE(K), K=1,3)
      DG 562 L=1.3
      NOP=NOP+1
      CP(NCP) = CUBE(L)
      CFP(NOP)=1./CFCTS(CUBE(L),CZERO)
      CONTINUE
 562
      GO TO 580
      PRINT 610, (CUBE(K), K=1,3)
       00 567 L=1.3
       NCZ=NOZ+1
```

```
CZ(NGZ)=CUBE(L)
     CFZ(NOZ)=CFCTS(CUBE(L),CZERO)
567
     CONTINUE
     PRINT 620
     GO TC 580
     IF (ICK.EQ.2) GO TO 575
570
     PRINT 572
    FORMAT (1HC,2X, #CAUTION ..... #,/,3X, #THIS CONTOUR HAS MORE THAN 3
572
    $ POLES. # . / . 3X . * THE USER IS URGED TO REWORK THIS CONTOUR # )
     GO TO 630
     CONT INUE
575
     PRINT 577
576
     FORMAT (1HC,3x, #CAUTION ..... + //,3X, #THIS CONTOUR HAS MORE THAN 3
577
    $ ZEROS. #, /, 3X, #THE USER IS URGED TO REWORK THIS CONTOUR #)
     GO TO 630
     CONTINUE
580
     FORMAT (1H0,3X, #ICK = #, 11,3X, #NP = #,11,3X,8E14.5,/)
590
     FORMAT (1H0,3X, +POLE AT+,2X, 2E20.8,/)
600
     FORMAT (1H0,3X, #ZERO AT#, 2X, 2E2C.8, /)
610
      FORMAT (1HO+1X+/////)
620
     CONTINUE
€30
      SRM I = SAUX I
      SRM=SAUX2
      CONTINUE
640
      CONTINUE
650
      CONT INUE
655
      FORMAT [1H0.2X.44HCOGRDINATES OF THE LOWER RIGHT CORNER ARE
 660
     1F12.8.1H., F12.8.1X.1H1./, 3X.44HCOORDINATES OF THE UPPER LEFT CORNE
     2R ARE (,1X,F12.8,1H,,F12.8,1X,1H))
      CSM=CAUX1
      CSMI=CAUX2
      RETURN
      END -
```

```
SUBROUTINE POLECHK(DO,D,DAO, DA, AP)
      IMPLICIT COMPLEX (D)
      DIMENSION D(8),CA(3),AD(8),D3(3,3)
      DIMENSION D3A(3,3),DW(3),DV(3)
C
      THIS SUBROUTINE FINDS A POTENTIAL VALUE FOR NP = NUMBER
C
      OF POLES IN THE GIVEN CONTOUR C.
Ċ
      IT ALSO DETERMINES THE COEFFICIENTS DA(4) OF THE
C
      DENGMINATOR POLYNOMIAL.
C
      NP=0
      1.0..0.1
      DO 10 IC=1,3
      JA(IC)=(0.,0.)
 10
      CONTINUE
      EPS=1.E-02
       ADO= CABS (-CO)
       00 2C [=1,8
       AD(I)=CABS(D(I))
 20
      CONT INUE
       PRINT 25, ADO, (AC(I), I=1,8)
      FURMAT (1HC,2x, #MAGNITUDES OF THE RESIDUE MOMENTS DO THRU D8#,//,3
 25
      $X.9E14.5./)
       IF (ACO.LE.EPS) GB TO 40
       CHECKING TO SEE IF NP = 1
C
       DA0 = -D(1)/D0
       00 30 1=1,7
       IP 1= I+1
       DQTY=-D(IP1)/D(I)
       TR=ABS(REAL(DAO-DGTY))
       TI=ABS(AIMAG(DAO-DQTY))
       IF (TR.GT.EPS.AND.TI.GT.EPS) GO TO 40
       CONTINUE
  30
       NP=1
       RETURN
  40
       CONTINUE
       IF (ADD.LE.EPS.AND.AD(1).LE.EPS) GO TO 60
       CHECKING T SEE IF NP = 2 .....
C
       DET=(D0*D(2))-(D(1)**2)
       JAO=((D(1)*D(3))-(D(2)**2))/DET
       DA1=((D(1)*D(2))-(D0*D(3)))/CET
       D4(1)=0A1
       DO 50 J=2.6
       JP1=J+1
       JP 2= J+2
       DQTY=C(JP2)+(D(JP1)*DA1)+(D(J)*DA0)
       TR=ABS(REAL(DCTY))
       TI=ABS(AIMAG(DQTY))
        IF (TR.GT.EPS.AND.TI.GT.EPS) GO TO 60
       CONTINUE
  50
       NP=2
        RETURN
       CONTINUE
  60
        IF (ADO.LE.EPS.ANO.AD(1).LE.EPS.AND.AD(2).LE.EPS) GO TO 180
        CHECKING TO SEE IF NP = 3 .....
 C
        00 90 IROW=1,3
        DO 80 JCOL=1,3
```

```
IND= IROW- 2+JCOL
     IF (IROW.EQ.1.AND.JCOL.EQ.1) GO TO 70
     D3(IROW, JCOL) =D(IND)
     GO TO 80
     D3(1,1)=00
73
     CONTINUE
80
     CONT INUE
9.3
     CALL DETER(D3, D3A, DW, CV, 3, DENGM)
     00 100 IROW=1,3
     IND= IRUW+2
     D3(IROW.1)=-D(IND)
     CUNTINUE
100
     CALL DETERIDS, DSA, DW, DV, 3, DNO ) .
     DC 130 [ROW=1,3
     DU 120 JCUL=1.2
      IF (JCOL.EQ.2) GO TO 110
      IND1 = IRCW-2+JCCL
      IF (IND1.EQ.0) GO TO 105
      D3(IROW, JCOL)=D(INDL)
     GC TO 120
     D3(1,1)=DC
105
      GO TO 120
110
      IND2=IRGW+2
      D3(IROW, JCOL) =-O(IND2)
      CONTINUE
120
      CONTINUE
 130
      CALL DETER(D3,D3A,DW,DV,3,DN1)
      00 160 IRCW=1.3
      DO 150 JCOL=2,3
      IF (JCOL.EQ.3) GO TO 140
      IND1=IROW-2+JCOL
      D3(IROW, JCOL) =D(INDL)
      GO TO 150
      IND2=IROW+2
 140
      D3(IRCW, JCCL) =-C(IND2)
      CONT INUE
 150
      CONTINUE
 160
      CALL DETER(D3,D3A,DW,DV,3,DN2)
      DAO=DNO/DENOM
      DA(1)=DN1/DENCM
      DA(2)=DN2/DENGM
       DO 170 J=3.5
       JP1=J+1
       JP 2= J+2
       DQTY=D(JP3)+(DA(2)*D(JP2))+(DA(1)*D(JP1))+(DAO*D(J))
       JP3=J+3
       TR=ABS(REAL(DQTY))
       TI=ABS(AIMAG(DCTY))
       IF (TR.GT.EPS.AND.TI.GT.EPS) GO TO 180
       CONTINUE
  170
       NP= 3
       RETURN
       CONTINUE
  180
        [F (ADO.GT.EPS) GO TO 200
       DO 190 ID=1.3
        IF (AD(ID).GT.EPS) GO TO 200
       CONTINUE
  190
        GB TC 340
```

```
200
      CONTINUE
 340
      CONTINUE
      DO 350 ID=4.8
      IF (AD(ID).GT.EPS) GO TO 370
350
      CONT INUE
      ALL MCMENTS ARE LESS THAN EPSILON.
C
      SO NP=0 FOR THIS CONTOUR.
C
      NP=0
      DAQ=[0.,0.]
      00 360 I=1,3
      DA(I)=(0.,0.)
 360
      CONTINUE
      RETURN
      CONTINUE
 370
      NP=4
      RETURN
      END
```

```
SUBROUTINE DETERICM, CB, CW, CV, MS, CD)
      THIS SUBROUTINE COMPUTES THE VALUE OF THE DETERMINANT
      CD OF A SQUARE MATRIX CM OF SIZE MS
C
C
      INPUTS. 1) COMPLEX MATRIX ELEMENTS CM
Ċ
               21SIZE OF THE SQUARE MATRIX MS
C
      OUTPUT. 11 COMPLEX VALUE OF DETERMINANT CO
Č
C
       IMPLICIT COMPLEX (C)
      DIMENSION CM(MS. MS), CB(MS. MS), CW(MS), CV(MS)
C
       REGIN TRIANGULARIZATION
C
       MCVE MATRIX CM TO CB SO THAT INPUT MATRIX WILL NOT
       BE DESTROYED
       DO 15 [=1,MS
       00 10 J=1. MS
       CB([,J) =CM([,J)
 10
       CONTINUE
 15
       DG 60 J=1.MS
       0=0.
       CALCULATE NORM SQUARED OF COLUMN J
C
       DG 20 K=J,MS
       Q=Q+CABS(CB(K,J)) **2
 20
       IF (C.GT.G.) GB TO 21
       IF THE NORM IS ZERO, THE MATRIX IS SINGULAR
C
       ISW=3
       PRINT 9
       FORMAT (1HO. #THE MATRIX IS SINGULAR #.//)
  9
       RETURN
       CALCULATE THE CIAGONAL ELEMENT OF THE MATRIX T. (CH(J))
 C
       CALCULATE THE DIAGONAL ELEMENT OF MATRIX U. (CB(J,J)) CALCULATE THE ELEMENT OF VECTOR V
 C
 C
       BEGIN ITERATION
 C
       BSQ=CABS(CB(J,J))**2
  21
        IF (BSQ.EQ.O.) C#(J)=SQRT(Q)
        IF (BSQ.GT.O.) CW(J)=SQRT(Q/BSQ)+CB(J,J)
        CB(J+J) = CB(J+J) + CW(J)
        CV(J)=-CB(J,J) *CONJG(CW(J))
        [F (J.EQ.MS) GG TO 60
        18=J+1
        DG 50 I= IB, MS
        CS=(0.,0.)
        nn 30 K=J.MS
        CS=CS+CB(K, I) *CONJG(CB(K,J))
   30
        CS=CS/CV(J)
        DO 40 K=J.MS
        CB(K, I)=CB(K, I)+CS*CB(K,J)
   40
        CONTINUE
   50
        CONTINUE
         CD=(1..0.)
         00 61 I=1, MS
         CD=CD*Ch(I)
   61
         RETURN
         END
```

FUNCTION ANGLER (X,Y) PI=3.14159265 [F (X) 90,10,50 10 IF (Y) 30,20,40 20 ANGLER=0. RETURN 30 ANGLER = 1.5 *PI RETURN 40 ANGLER =PI*.5 RETURN 50 [F (Y) 80,60,70 60 ANGLER=0. RETURN 70 ANGLER=ATAN(Y/X) RETURN 80 ANGLER = -ATAN(-Y/X)+2. *PI RETURN 90 XN=-X IF (Y) 120, 10C, 110 100 ANGLER=PI RETURN 110 ANGLER=PI-ATAN(Y/XN) RETURN 120 ANGLER=PI+ATAN (-Y/XN) RETURN END

SUBROUTINE POLY 1(CO.CI.CLIN)
IMPLICIT COMPLEX (C)

C THIS SUBROUTINE SOLVES A LINEAR EQUATION OF THE
C FORM C1*S+CO = 0

C CLIN=-CO/C1
RETURN
END

```
SUBROUTINE POLY 2 (CO.C1.C2.CQUAC)
      IMPLICIT COMPLEX (C)
      DIMENSION CQUAD(2)
      THIS SUBROUTINE SOLVES A QUACRATIC EQUATION OF THE
C
      FGRM C2*(S**2)+C1*S+CG = 0
C
      CC=(C1 **2)-(4.*C2*C0)
      CQ=CSQRT(CQ)
      CDR=(2.*C2)
      CQUAD(1)=(-C1+CC)/CDR
      CQUAD(2) = (-C1-CQ) /CDR
      RETURN
      END
      SUBROUTINE POLY 3 (CO, C1, C2, C3, CUBE)
       IMPLICIT COMPLEX (C)
       DIMENSION CUBE(3) .CAUX1(3) .CAUX2(3)
       THIS SUBREUTINE SOLVES A CUBIC EQUATION OF THE
C
       FORM: C3*(S**31+C2*(S**2)+C1*S+CC = 0
C
       REFERENCE......EQN. 3.8.2. OF AMS 55. PAGE 17.
C
       CA2=C2/C3
       CA1=C1/C3
       CAC=CO/C3
       CQ=(CA1/3.1-((CA2**2)/9.1
       CR=((CA1*CA2-3.*CA0)/6.)-((CA2**3)/27.)
       CRQ=CSQRT((CR**2) + (CQ**3))
       TPI=2.*3.14159265
       CRP=CR+CRQ
       CRM=CR-CRG
       CS 1=CEXP(CLOG(CRP)/3.)
       CS2=CEXP(CLGG(CRM)/3.)
       CQTY =-CQ
       C I= (0., 1.)
       EPS=1.E-5
       DO 20 IC=1,3
       TIC=FLOAT(IC)
       CAUX1(IC)=CEXP(CI*TP[*TIC/3.1+CS1
       DO 10 JC=1.3
        TJC=FLOAT(JC)
       CAUX2(JC)=CEXP(CI*TPI*TJC/3.1*CS2
       CP=CAUX1(IC) * CAUX2(JC)-CQTY
        DELR=ABS (REAL (CP))
        DELI=ABS (AIMAG (CP))
        IF (DELR.LE.EPS. AND. DELI.LE.EPS) GO TO 30
        CONTINUE
  10
        CONTINUE
  20
        CS1=CAUX1(IC)
   30
        CS2=CAUX2(JC)
        CSP=(CS1+CS2)/2.
        CSM={CS1-C S2)/2.
        CR3=CI*SQRT(3.)
        CA23=CA2/3.
        CUBE(1)=CS1+CS2-CA23
        CUBE(2) = -CSP-CA23+(CR3+CSM)
        CUBE(3) =-CSP-CA23-(CR3*CSM)
        RETURN
```

END

```
COMPLEX FUNCTION OFCTS (CS, CSHIFT)
     IMPLICIT COMPLEX (C)
     DIMENSION COR(IG), CFD(10), CNR(16), CFN(16)
     CS=CS+CSHIFT
     NP=10
     CDR(1)=(3.5,1.5)
     CDR(2)=(1.75, 2.25)
     CDR(3) = (1.5, 2.75)
     CDR (4)=(2.25, 3.25)
     CDR(5)=(2.75,3.5)
     CDR (6)=(2.25.3.75)
     CDR(7)=(1.25.0.75)
     CCR(8)=(2.75,1.75)
     CDR(9) = (1.5,0.5)
     CDR(10)=(2.25,1.25)
     DC 10 I=1.NP
     CFD(I)=CS-CDR(I)
10
     CONTINUE
     CDENOM= (1 .. 0.)
     90 20 I=1.NP
     CDENCM=CDENOM*CFD(I)
20
     CCNTINUE
     NO=16
     CNR(1)=(4.5,.5)
     CNR(2)=(4.25,4.75)
     CNR(3)=(3.75,1.25)
     CNR(4) = (3.75, 1.75)
     CNR (5) = (4.5,2.5)
     CNR(6)=(4.25,2.75)
     CNR(7)=(2.75,3.75)
     CNR(8)=(1.25.2.25)
     CNR(9)=(1.75,2.75)
     CNR(10)=(4.5,4.5)
     CNR(11)=(.75,1.5)
     CNR(12)=(.5,1.25)
     CNR(13)=(.25,1.75)
     CNR(14)=(.5,4.25)
     CNR(15)=(.75,4.5)
      CNR(16)=(.25.4.75)
      DO 30 I=1.NO
      CFN(I)=CS-CNR(I)
30
      CONT INUE
     C NUM= (1.,0.)
      00 40 I=1.NC
      C NUM = C NUM + CFN(I)
      CONTINUE
40
      CFCT S=CNUM/CDENUM
      A=0 .
      CFCTS=CFCTS*CEXP(A*CS)
      RETURN
      END
```

APPENDIX B

A List of Relevant Definitions and Some Interesting but not Very Useful Results

Analytic function:

If A(s) has a derivative at a point s_0 and also at each point in some neighborhood of s_0 , then A(s) is said to be analytic at s_0 . The terms holomorphic, monogenic and regular are also sometimes used [9].

Entire function:

An entire function is one which is analytic everywhere in the plane and may have its singularities only at infinity [9].

Meromorphic function:

A meromorphic function is one whose only singularities, except at infinity, are poles [4].

Argument number:

The argument number N_a is the number of excess zeros over poles (N_O-N_p) of a meromorphic function, inside a simple closed contour. Note that N_a is the order of the pole at infinity when $N_a>0$ and it is the order of the zero at infinity when $N_a<0$.

Exceptional point:

For a given function, certain values may be exceptional, in the sense that the function can not take these values [4]. For instance, the points $(0 \pm j1)$ are exceptional points of the meromorphic function $\tan(s)$.

A major problem in the procedure developed in this note, of finding the poles and zeros of a given complex meromorphic function lies in obtaining the number of poles N_p , deterministically in a given contour. If N_p is known unambigiously, there is no difficulty in accurately determining the pole locations. Since the excess number of zeros over poles, i.e., (N_0-N_p) can be easily determined from the principle of the argument, attempt was made to obtain another expression involving N_0 and N_p so that one can solve a set of simultaneous equations for N_0 and N_p . In this unsuccessful attempt, the following results were obtained. The following results are not useful in determining N_0 and N_p , when used along with the known argument number. When M(s) has all its poles simple,

$$\frac{1}{2\pi j} \oint_{C} s \left\{ \frac{M'(s)}{M(s)} \right\}^{2} ds = (N_{o} - N_{p})^{2}$$
(B.1)

$$\frac{1}{2\pi j} \oint_{C} s \left\{ \frac{M''(s)}{M(s)} \right\} ds = (N_{o} - N_{p})^{2} - (N_{o} - N_{p})$$
 (B.2)

$$\frac{1}{2\pi j} \oint_{C} \frac{M''(s)}{M'(s)} ds = (N'_{o} - N'_{p}) = (N'_{o} - 2N_{p})$$
 (B.3)

with $N_0' = \text{number of zeros of } M'(s) \text{ in } C$,

 N_{p}' = number of poles of M'(s) in C

= twice the number of poles of M(s) in $C = 2N_p$.

$$\frac{1}{2\pi j} \oint_{C} \frac{M'''(s)}{M''(s)} ds = (N'_{1} - 3N_{p})$$
 (B.4)

with $N_1' = \text{number of zeros of } (M'(s) - 1)$ in C.

In this context, it may be noted that N_{p} can be obtained deterministically, if

i) we develop a method of determining the number of essential singularities of M(s) in C. or ii) we know an exceptional value ϵ of M(s) in C as illustrated below.

i) Essential singularities

Observe that the function $\exp(M(s))$ has no zeros in C, but all the poles of M(s) become essential singularities of $\exp(M(s))$ within C. So, if one can determine the number of essential singularities of a given function in a given contour, this concept is useful in determining the number of poles of a function inside the contour.

ii) Exceptional value

Let ϵ be a known exceptional value of M(s) inside the contour C. The existence of ϵ is improbable, if not impossible, since the affinity of a function for every value is the same. If ϵ does exist, one may define a new function a(s)

$$a(s) = \frac{1}{M(s) - \varepsilon}$$
 (B.5)

then the argument number of a(s) is given by

$$\frac{1}{2\pi j} \oint_C \frac{a'(s)}{a(s)} ds = N_p$$
 (b.6)

and thus N_p , the number of poles of M(s) in C is easily determined. Note that a(s) is an analytic function in C since M(s) \dagger ϵ and the poles of M(s) in C become the zeros of a(s).