



**AUSTRALIAN ATOMIC ENERGY COMMISSION
RESEARCH ESTABLISHMENT
LUCAS HEIGHTS**

**A METHOD FOR INVERTING THE CHANNEL MATRIX IN
NUCLEAR RESONANCE THEORY**

by

W.K. BERTRAM

November 1971

APPROVED FOR PUBLICATION

ISBN 0 642 99445 5

AUSTRALIAN ATOMIC ENERGY COMMISSION

RESEARCH ESTABLISHMENT

LUCAS HEIGHTS

A METHOD FOR INVERTING THE CHANNEL MATRIX
IN NUCLEAR RESONANCE THEORY

by

W. K. BERTRAM

ABSTRACT

A technique is presented for inverting the channel matrix which occurs in R-matrix theory. The collision matrix can be expressed in a form which does not involve matrix inversions; it can be evaluated by the application of a recurrence relation. The connection between this method and the level matrix formalism is examined.

Note: This work has been submitted to a journal. Further details can be obtained from the author or the Director of the Research Establishment.

National Library of Australia card number and ISBN 0 642 99445 5

The following descriptors have been selected from the INIS Thesaurus to describe the subject content of this report for information retrieval purposes. For further details please refer to IAEA-INIS-12 (INIS: Manual for Indexing) and IAEA-INIS-13 (INIS: Thesaurus) published in Vienna by the International Atomic Energy Agency.

ANALYTICAL SOLUTION; COLLISION MATRIX; CONFIGURATION; CROSS SECTIONS;
ENERGY LEVELS; R-MATRIX; RECURSION RELATIONS; RESONANCE

1. INTRODUCTION

The resonance structure of nuclear reaction cross sections is usually studied by the R-matrix theory developed by Wigner and Eisenbud (1947) (Lane and Thomas 1958). In this theory there occurs the channel matrix, the inverse of which is required for the calculation of the collision matrix. When the number of reaction channels is large the inversion of this matrix can be carried out analytically only in the special case where the R-matrix is of rank one (Teichmann 1950, Newton 1952) as is the case when it contains one single level. The cross section is then given by the familiar Breit-Wigner single level formula.

The single level theory has enjoyed a great deal of success in its description of resonances for many nuclei, but there are nuclei, notably the fissionable, for which the single level formula has proved inadequate. Several multilevel theories have been developed which take into account the interference between resonances (Thomas 1955, Vogt 1958, Reich and Moore 1958). All these theories are based on a method for expressing the collision matrix in a form in which the channel matrix has been eliminated and the problem of inverting the channel matrix is replaced by one of inverting the level matrix.

This paper presents an alternative approach to the problem of inverting the channel matrix. The collision matrix can be expressed in a form which does not involve matrix inversions; it can be obtained by the application of a relatively simple recurrence relation. The relationship between this method and the level matrix expansion is discussed and approximations for the collision matrix are derived.

2. INVERSION OF THE CHANNEL MATRIX

The collision matrix U is connected with the R-matrix through the relation (Lane and Thomas 1958)

$$U = \Omega \left[I + 2i P^{\frac{1}{2}} (I - R L^0)^{-1} R P^{\frac{1}{2}} \right] \Omega . \quad \dots(1)$$

R is a symmetric matrix with elements

$$R_{cc'} = \sum_{\lambda} \frac{\gamma_{\lambda c} \gamma_{\lambda c'}}{E_{\lambda} - E} \quad \dots(2)$$

where $\gamma_{\lambda c}$ and E_{λ} are real constants.

Ω , L^0 and P are diagonal matrices with elements

$$\Omega_c = e^{i\phi_c} \quad \dots(3)$$

$$L_c^0 = S_c - B_c + iP_c \quad \dots(4)$$

where, for channel c , ϕ_c is the hard sphere scattering phase, S_c the level shift, P_c the penetration factor and B_c is a constant arising from the boundary conditions.

The central problem in calculating U is the determination of the matrix

$$S = (I - RL^0)^{-1} R \quad \dots(5)$$

Our method for solving this problem is based on the well-known result (Lane and Thomas 1958) that if

$$R = R_1 + R_1' \quad \dots(6)$$

S can be expressed as a sum of two terms, one of them involving only R_1 and the other involving both R_1 and R_1' . That is

$$S = (I - RL^0)^{-1} R = (I - R_1 L^0)^{-1} R_1 + (I - R_1 L^0)^{-1} (I - R_1' L^{(1)})^{-1} R_1' \times \\ \times (I - L^0 R_1)^{-1} \quad \dots(7)$$

where

$$L^{(1)} = L^0 (I - R_1 L^0)^{-1} \quad \dots(8)$$

We note from equation (2) that the R matrix is a sum over single level R -matrices,

$$R = \sum_{\lambda} R_{\lambda} \quad \dots(9)$$

Each R_{λ} has the property that

$$(R_{\lambda})^2 = R_{\lambda} \text{Trace} (R_{\lambda}) \quad \dots(10)$$

and therefore

$$(I - R_{\lambda} L^0)^{-1} = I + \frac{R_{\lambda} L^0}{1 - \text{Trace} R_{\lambda} L^0} \quad \dots(11)$$

Thus writing

$$R = R_1 + \sum_{\lambda=2} R_\lambda \quad \dots(12)$$

and substituting this in equation (7) we obtain

$$S = \frac{R_1 L^0}{1 - \text{Trace } R_1 L^0} + \left(I + \frac{R_1 L^0}{1 - \text{Trace } R_1 L^0} \right) (I - R_1' L^{(1)})^{-1} R_1' \times \\ \times \left(I + \frac{L^0 R_1}{1 - \text{Trace } R_1 L^0} \right) \quad \dots(13)$$

$$\text{and } L^{(1)} = L^0 + \frac{L^0 R_1 L^0}{1 - \text{Trace } R_1 L^0} \quad \dots(14)$$

We may now write

$$R_1' = R_2 + R_2' \quad \dots(15)$$

and use equation (7) again to expand the matrix S_1 where

$$S_1 = \left(I - R_1' L^{(1)} \right)^{-1} R_1' \quad \dots(16)$$

This procedure can be repeated as many times as there are levels, or at least until we have included enough levels to allow us to treat the remainder as a constant background, R_∞ .

Thus after eliminating k levels from the R -matrix the elimination of the next level is achieved by using

$$(I - R_k' L^{(k)})^{-1} R_k' = \frac{R_{k+1}}{1 - T_{k+1 k}} + \left(I + \frac{R_{k+1} L^{(k)}}{1 - T_{k+1 k}} \right) \\ (I - R_{k+1}' L^{(k+1)})^{-1} R_{k+1}' \times \left(I + \frac{L^{(k)} R_{k+1}}{1 - T_{k+1 k}} \right) \quad \dots(17)$$

so that equation (20) becomes

$$S = (L^0)^{-1} \sum_{k=0}^{N-1} \frac{L^{(k)} R_{k+1} L^{(k)}}{1 - T_{k+1 k}} (L^0)^{-1} \quad \dots(22)$$

This can be simplified even further by noting that if we carry out a summation from $k = 0$ to $N - 1$ on both sides of equation (18) we get

$$L^{(N)} = L^0 + \sum_{k=0}^{N-1} \frac{L^{(k)} R_{k+1} L^{(k)}}{1 - T_{k+1 k}} \quad \dots(23)$$

Thus we obtain the following simple form for S ,

$$S = (I - RL^0)^{-1} R = (L^0)^{-1} (L^{(N)} - L^0) (L^0)^{-1} \quad \dots(24)$$

The matrix $L^{(N)}$ can be calculated from the recurrence relation (equation (18)). Therefore the calculation of S does not involve any matrix inversions.

3. CONNECTION WITH THE LEVEL MATRIX FORMALISM

In the level matrix formalism (Lane and Thomas 1958) a particular element $S_{cc'}$ of S is expressed in the form

$$S_{cc'} = \left[(I - RL^0)^{-1} R \right]_{cc'} = \sum_{\lambda\mu} \gamma_{\lambda c} \gamma_{\mu c'} A_{\lambda\mu} \quad \dots(25)$$

where the matrix A is the inverse of the level matrix C which has elements

$$C_{\lambda\mu} = (E_\lambda - E) \delta_{\lambda\mu} - \xi_{\lambda\mu} \quad \dots(26)$$

where

$$\xi_{\lambda\mu} = \sum_c \gamma_{\lambda c} \gamma_{\mu c} L^0_c \quad \dots(27)$$

The element $S_{cc'}$ can be evaluated if the values of $E_\lambda, \gamma_{\lambda c}, \gamma_{\lambda c'}$ and $\xi_{\lambda\mu}$ are known for all λ, μ . If there are N levels, the number of parameters required to determine any one particular element of S is $\frac{1}{2}N(N+1)+3N$ which is independent of the number of channels.

On the other hand if we wish to use the recurrence relation (equation (18)) to determine a given element $S_{cc'}$, then it is necessary to know the values of E_λ and $\gamma_{\lambda c}$ for all values of λ and c . In this case we require $N(M+1)$ parameters, where M is number of channels. However, since the two methods are equivalent, there must be a means of calculating $L_{cc'}^{(N)}$ using only those parameters which occur in the level matrix formalism. A method is found easily by introducing quantities $\beta_{\lambda c}^{(k)}$ and $\xi_{\lambda\mu}^{(k)}$ which are defined as

$$\beta_{\lambda c}^{(k)} = \sum_{c'} \gamma_{\lambda c'} L_{cc'}^{(k)} \quad \dots(28)$$

$$\text{and} \quad \xi_{\lambda\mu}^{(k)} = \sum_c \gamma_{\lambda c} \beta_{\mu c}^{(k)} = \sum_{cc'} \gamma_{\lambda c} \gamma_{\mu c'} L_{cc'}^{(k)} \quad \dots(29)$$

For $k = 0$ these quantities are just the parameters which occur in the level matrix.

From equation (18) we have

$$L_{cc'}^{(n)} = L_{cc'}^{(n-1)} + \frac{\left(L_{cc'}^{(n-1)} R_n L_{cc'}^{(n-1)} \right)}{1 - \text{Trace} (R_n L_{cc'}^{(n-1)})} \quad \dots(30)$$

$$\text{but} \quad \left(L_{cc'}^{(n-1)} R_n L_{cc'}^{(n-1)} \right)_{cc'} = \frac{\beta_{nc}^{(n-1)} \beta_{nc'}^{(n-1)}}{E_n - E} \quad \dots(31)$$

$$\text{and} \quad \text{Trace} (R_n L_{cc'}^{(n-1)}) = \frac{\xi_{nn}^{(n-1)}}{E_n - E} \quad \dots(32)$$

Therefore

$$L_{cc'}^{(n)} = L_{cc'}^{(n-1)} + \frac{\beta_{nc}^{(n-1)} \beta_{nc'}^{(n-1)}}{E_n - E - \xi_{nn}^{(n-1)}} \quad \dots(33)$$

If we multiply both sides of equation (33) by $\gamma_{n+1, c'}$ and carry out the summation over c' we find

$$\beta_{n+1, c}^{(n)} = \beta_{n+1, c}^{(n-1)} + \frac{\beta_{nc}^{(n-1)} \xi_{n+1, n}^{(n-1)}}{E_n - E - \xi_{nn}^{(n-1)}} \quad \dots(34)$$

By repeating this procedure, this time multiplying throughout by γ_{kc} we obtain

$$\xi_{n+1, k}^{(n)} = \xi_{n+1, k}^{(n-1)} + \frac{\xi_{nk}^{(n-1)} \xi_{n+1, n}^{(n-1)}}{E_n - E - \xi_{nn}^{(n-1)}} \quad \dots(35)$$

We see then that starting only with the values of $E_\lambda, \beta_{\lambda c}^{(0)}, \beta_{\lambda c'}^{(0)}$ and $\xi_{\lambda\mu}^{(0)}$ we can use the recurrence relations (equations (34) and (35)) to determine $\beta_{nc}^{(n-1)}, \beta_{nc'}^{(n-1)}$ and $\xi_{nn}^{(n-1)}$ and thus $S_{cc'}$.

4. APPROXIMATIONS

In the level matrix formalism, approximations for S are usually obtained by writing the level matrix C in the form (Schmidt 1966)

$$C = D + \xi' \quad \dots(36)$$

where D is the diagonal part of C and ξ' the off-diagonal part.

Approximations for S are then found by the expansion,

$$A = D^{-1} + D^{-1} \xi' D^{-1} + \dots \quad \dots(37)$$

Instead of this procedure we can use a perturbation method based on equation (23) for obtaining S . As a first approximation we replace $L^{(k)}$ by L^0 on the right of equation (23) and obtain

$$L^{(N)} \approx L^{\circ} + \sum_{k=0}^{N-1} \frac{L^{\circ} R_{k+1} L^{\circ}}{1 - T_{k+1}^{\circ}} \quad \dots(38)$$

and hence

$$S_{cc'} \approx \sum_{\lambda} \frac{\gamma_{\lambda c} \gamma_{\lambda c'}}{E_{\lambda} - E - \xi_{\lambda\lambda}^{\circ}} \quad \dots(39)$$

Exactly the same result is obtained by using only the first term on the right of equation (37) to calculate S. A better approximation is obtained by substituting

$$L^{(k)} \approx L^{\circ} + \sum_{i=0}^{k-1} \frac{L^{\circ} R_{i+1} L^{\circ}}{1 - T_{i+1}^{\circ}} \quad \dots(40)$$

into equation (23) where

$$S = \sum_{k=0}^{N-1} \frac{L^{\circ} R_{k+1} L^{\circ}}{1 - T_{k+1}^{\circ}} + \sum_{k=0}^{N-1} \sum_{i=0}^{k-1} \frac{L^{\circ} R_{k+1} L^{\circ} R_{i+1} L^{\circ} + L^{\circ} R_{i+1} L^{\circ} R_{k+1} L^{\circ}}{(1 - T_{i+1}^{\circ})(1 - T_{k+1}^{\circ})} +$$

$$+ \sum_{k=0}^{N-1} \sum_{i,j=0}^{k-1} \frac{L^{\circ} R_{i+1} L^{\circ} R_{k+1} L^{\circ} R_{j+1} L^{\circ}}{(1 - T_{i+1}^{\circ})(1 - T_{k+1}^{\circ})(1 - T_{j+1}^{\circ})} \quad \dots(41)$$

If we put $L^{(k)} = L^{\circ}$ in the evaluation of the traces in equation (41) we obtain the expression

$$S_{cc'} \approx \sum_{\lambda} \frac{\gamma_{\lambda c} \gamma_{\lambda c'}}{E_{\lambda} - E - \xi_{\lambda\lambda}^{\circ}} + \sum_{\lambda} \sum_{\mu \neq \lambda} \frac{\gamma_{\lambda c} \gamma_{\mu c'} \xi_{\lambda\mu}^{\circ}}{(E_{\lambda} - E - \xi_{\lambda\lambda}^{\circ})(E_{\mu} - E - \xi_{\mu\mu}^{\circ})} \quad \dots(42)$$

where we have omitted the third term in equation (41). This expression for S_{cc} is the same as that obtained from the level matrix formulation using the first two terms of equation (37). Note that the analysis of resonances using a theory based on the expansion (equation (37)) will never give the correct position and total width of the resonances no matter how many terms of equation (37) are taken. This is because for a finite number of terms, the poles of A , and therefore also those in S , are always at $E = E_\lambda - \xi_{\lambda\lambda}$ whereas their true positions are given by the eigenvalues of the level matrix.

Our perturbation method does not have this problem since the traces in equation (41) are calculated using $L^{(k)}$ given by equation (40).

5. CONCLUSION

We have shown how the channel matrix can be inverted so that the resulting expression for the collision matrix no longer involves matrix inversions. By examining the relation between this method and the more conventional method involving the level matrix, we found that parametrization of the collision matrix in terms of level matrix parameters is still possible. It also appears that our method has some advantages over the level matrix formalism when it comes to finding approximations for the collision matrix.

6. ACKNOWLEDGEMENT

I wish to thank Mr. J. L. Cook for stimulating discussions on many aspects of R-matrix theory.

7. REFERENCES

- Lane, A. M. and Thomas, R. G. (1958) - Rev. Mod. Phys. 30, 257
 Newton, T. D. (1952) - Can. J. Phys. 30, 53.
 Reich, C. W. and Moore, M. S. (1958) - Phys. Rev. 111, 929.
 Schmidt, J. J. (1966) - 'Neutron Cross Sections for Fast Reactor Materials'
 Vol. 1. KFK120 (EANDC-E-35U)
 Teichmann, T. (1950) - Phys. Rev. 77, 506.
 Thomas, R. G. (1955) - Phys. Rev. 97, 224.
 Vogt, E. (1958) - Phys. Rev. 112, 203.
 Vogt, E. (1960) - Phys. Rev. 118, 724.
 Wigner, E. P. and Eisenbud, L. (1947) - Phys. Rev. 71, 29.

