# Lawrence Berkeley National Laboratory 

 Recent WorkTitle
THE CLASSICAL S-MATRIX FOR COULOMB EXCITATION

## Permalink

https://escholarship.org/uc/item/5d73q101
Author
Donangelo, R.
Publication Date
1976-02-05

LBL-4347
Preprint
. 1

FEB 5 1Y/b
LIBRARY AND
DOCUMENTS SECTION

THE CLASSICAL S-MATRIX FOR COULOMB EXCITATION
R. Donangelo, M. W. Guidry, and J. O. Rasmussen

December 1975

Prepared for the U. S. Energy Research and
Development Administration under Contract W-7405-ENG-48

## For Reference

Not to be taken from this room


## DISCLAIMER

This document was prepared as an account of work sponsored by the United States Government. While this document is believed to contain correct information, neither the United States Government nor any agency thereof, nor the Regents of the University of California, nor any of their employees, makes any warranty, express or implied, or assumes any legal responsibility for the accuracy, completeness, or usefulness of any information, apparatus, product, or process disclosed, or represents that its use would not infringe privately owned rights. Reference herein to any specific commercial product, process, or service by its trade name, trademark, manufacturer, or otherwise, does not necessarily constitute or imply its endorsement, recommendation, or favoring by the United States Government or any agency thereof, or the Regents of the University of California. The views and opinions of authors expressed herein do not necessarily state or reflect those of the United States Government or any agency thereof or the Regents of the University of California.

# THE CLASSICAL S-MATRIX FOR COULOMB EXCITATION* 

R. Donangelo, M. W. Guidry and J. O. Rasmussen

Lawrence Berkeley Laboratory
University of California
Berkeley, California 94720

ABSTRACT
An expression of the classical S-matrix for Coulomb excitation is derived and directly evaluated without resorting to stationary phase integration methods. The results obtained are in quantitative agreement with other quantum mechanical and semi-classical predictions. This and the simplicity of the method suggest the feasibility of extending it to cases where other methods are not easily applied.

[^0]The electromagnetic interaction that appears in the collision of a deformed target nucleus with a nuclear projectile excites rotational states of the target. Until recently only light projectile ions were available and partial wave, coupled channel, quantum mechanical calculations of this Coulomb excitation process were feasible.

With the advent of heavier projectiles exact quantum mechanical calculations have become impractical. However, the shorter de Broglie wavelength of heavy ions brings the problem closer to the realm of Classical Mechanics, and in recent years the Uniform Semi-classical Approximation (USCA), developed mainly by W. Miller, ${ }^{1-3)}$ was applied to this problem ${ }^{4,5)}$ with results that compare very well with the predictions of the semi-classical method of Winther and de Boer. ${ }^{6}$ )

Miller derives an integral expression for the classical S-matrix which he refers to as the initial value representation of the S-matrix. ${ }^{2}$ ) The USCA method consists in evaluating this integral expression by the stationary phase approximation.

In this initial value representation, the Coulomb excitation of a two-dimensional rotor by a backscattered projectile has an S-matrix given by:

$$
\begin{equation*}
S_{I_{2} \leftarrow 0}=\frac{1}{2 \pi} \int_{0}^{2 \pi} d \beta_{1}\left[\frac{\partial \beta_{2}\left(\beta_{1}\right)}{\partial \beta_{1}}\right]^{\frac{1}{2}} \exp \left(i\left\{\phi_{2}\left(\beta_{1}\right)+\bar{\beta}_{2}\left(\beta_{1}\right)\left[I_{2}\left(\beta_{1}\right)-I_{2}\right]\right\}\right) \tag{1}
\end{equation*}
$$

where the final spins $I_{2}$ are $I_{2}=0,2,4, \ldots$ (in units of $h$ ). Here $\beta_{1}$ is the initial orientation angle between the rotor axis and the beam axis, and $\bar{\beta}_{2}\left(\beta_{1}\right)$ is the final angle shift which is defined as

$$
\begin{equation*}
\bar{\beta}_{2}\left(\beta_{1}\right)=\beta_{2}\left(\beta_{1}\right)-\frac{m R_{2}\left(\beta_{1}\right)}{P_{2}\left(\beta_{1}\right)} \frac{\hbar I_{2}\left(\beta_{1}\right)}{\mathscr{q}} \tag{2}
\end{equation*}
$$

where $\beta_{2}\left(\beta_{1}\right)$ is the orientation angle, $I_{2}\left(\beta_{1}\right)$ is the angular momentum of the rotor (in units of $\hbar$ ), $R_{2}\left(\beta_{1}\right)$ is the projectile-target distance, and $P_{2}\left(\beta_{1}\right)$ is the relative momentum of the projectile in the center-of-mass system, all evaluated at a point along the trajectory such that the interaction has already taken place (the spin I has taken a constant value). The reduced mass of the system and the moment-ofinertia of the rotor are indicated by $m$ and $\ell$ respectively.

Finally, $\phi_{2}\left(\beta_{1}\right)$ in Eq. (1) is the classical phase in units of $h$ evaluated along the trajectory that has an initial oricntation $\beta_{1}$ and a final one $\beta_{2}$.

$$
\begin{equation*}
\phi_{2}\left(\beta_{1}\right)=-\frac{1}{\hbar} \int_{t\left(\beta_{1}\right)}^{t\left(\beta_{2}\right)}[R(t) \dot{P}(t)+\beta(t) \hbar \dot{I}(t)] d t \tag{3}
\end{equation*}
$$

where all variables were defined above.
The reason for taking the angle shift $\bar{\beta}_{2}$ and not the angle variable $\beta_{2}$ in the expression for the $S$-matrix is that Eq. (1) must be independent of the point in the asymptotic region where we terminate the trajectory. This requirement is satisfied by an expression such as Eq. (2). Wong and Marcus ${ }^{7}$ ) give a more general relationship which can also be used in this case.

A real deformed nucleus must be considered as a three-dimensional rotor. The final $\operatorname{spin}$ takes the values $I_{2}=0.5,2.5,4.5, \ldots$ since to the quantum mechanical spin $I$ there corresponds a classical spin $I+\frac{1}{2} .4$ ) From geometrical considerations ${ }^{4,9)}$ the different orientations $\beta_{1}$ must be weighted by the factor $\left(8 \pi \sin \beta_{1}\right)^{\frac{1}{2}}$ with $0 \leqslant \beta_{1} \leqslant \pi$.

From Eq. (1) and preceding considerations we obtain the following expression for the Coulomb excitation S-matrix of a deformed nucleus by
a backscattered projectile

$$
\begin{equation*}
S_{I_{2} \leftarrow 0}=\frac{1}{\sqrt{\pi}} \int_{0}^{\pi} \mathrm{d} \beta_{1}\left[\sin \beta_{1} \frac{\partial \bar{\beta}_{2}\left(\beta_{1}\right)}{\partial \beta_{2}}\right]^{\frac{1}{2}} \exp \left(\mathrm{i}\left\{\phi_{2}\left(\beta_{1}\right)+\bar{\beta}_{2}\left(\beta_{1}\right)\left[\mathrm{I}_{2}\left(\beta_{1}\right)-\mathrm{I}_{2}\right]\right\}\right) \tag{4}
\end{equation*}
$$

where all magnitudes retain their previous definition and $I_{2}=0.5$, $2.5,4.5, \ldots$

This integral is not difficult to integrate even for massive targots amd projectiles. In the case of ${ }^{40} \mathrm{Ar}$ on ${ }^{238} \mathrm{U}$ at 170 MeV laboratory energy shown in Fig. 1, it was sufficient to take a partition of the interval $[0, \pi]$ into 100 points to evaluate the excitation probabilities defined by

$$
\begin{equation*}
P_{(I)}=\left|S_{I+0}\right|^{2} \tag{5}
\end{equation*}
$$

with good accuracy. We compared our results with the Winther-de Boer ${ }^{6)}$ predictions since no quantum mechanical codes are feasible in this case. The energy levels were taken from the rotational model with constant moment of inertia but it would be very simple to allow for variable moment of inertia by taking the appropriate expression for the classical Hamiltonian used to calculate the trajectories. Equation (4) is useful even at energies near the Coulomb barrier where the onset of the nuclear potential makes the Winther-de Boer method no longer valid. The nuclear potential can be easily included in the classical equations of motion used to obtain the variables in the integrand of Eq. (4).

The fact that the stationary phase approximation is not used makes expression (4) more accurate where that approximation breaks down. Figure 2 and Table I show such a case; in this example the exponent in Eq. (4) does not change very rapidly and the stationary
phase approximation is not so good. The direct evaluation of the classical S-matrix gives much better agreement with the exact quantummechanical results than the USCA for this case.

There are also cases where more than one internal degree of freedom must be considered, such as in rotational-vibrational excitation. For these cases the uniform approximation to the classical S-matrix is more complicated ${ }^{11)}$ and the multi-dimensional search for the stationary points may be as difficult from a numerical point of view ${ }^{10}$ ) as the direct evaluation of the classical S-matrix.

Acknowledgments: We thank Dr. Herbert Massmann for stimulating discussions, and Drs. Robert Anholt, Jean Paul Boisson and Prof. Hiroyasu Ejiri for useful comments on our manuscript. R. Donangelo acknowledges support from an International Atomic Energy Agency fellowship.

$$
-6-
$$

## REFERENCES

${ }^{1)}$ W. H. Miller, J. Chem. Phys. 53 (1970) 1949.
${ }^{2)}$ W. H. Miller, J. Chem. Phys. 53 (1.970) 3578.
${ }^{3}$ )W. H. Miller, J. Chem. Phys. 54 (1971) 5386.
${ }^{4)}$ H. Massmann and J. O. Rasmussen, Nuc1. Phys. A243 (1975) 155.
${ }^{5)}$ S. I.evit, II. Smilansky and D. Pelte, Phys. Lett. 53 B (1974) 39.
${ }^{6)}$ A. Winther and J. de Boer, in "Coulomb Excitation", K. Alder and A. Winther eds. (Academic Press, New York, 1965) p. 303.
${ }^{7}$ W. H. Wong and R. A. Marcus, J. Chem. Phys. 55 (1971) 5663.
${ }^{8)}$ F. Roesel, J. X. Saladin and K. Alder, "Quantum Mechanical Coupled Channel Code for Coulomb Excitation', University of Pittsburgh Progress Report (1972).
$\left.{ }^{9}\right)_{\text {H }}$. Massmann, Ph.D. Thesis, University of California; Berkeley (1975), unpublished.
${ }^{10)}$ M. W. Guidry, H. Massmann and J. O. Rasmussen (unpublished).
${ }^{11)}$ R. A. Marcus, J. Chem. Phys. 57 (1972) 4903.

TABLE I. Comparison of the ratios of the normalized excitation probabilities predicted by the Integral Expression (4) and by the USCA method to the Quantum-Mechanical coupled channel results of the AROSA code ${ }^{8)}$ in the case depicted in Fig. 2.

| SPIN | INTEGRAL/Q.M. | USCA/Q.M. |
| :---: | :---: | :---: |
| 0 | 1.06 | 0.97 |
| 2 | 0.90 | 1.06 |
| 4 | 0.99 | 0.90 |
| 6 | 1.13 | 0.61 |
|  |  | 1.15 |

## FIGURE CAPTIONS

Fig. 1. Calculation of Coulomb excitation probabilities to excite members of the rotational ground band in ${ }^{238} \mathrm{U}$ for backscattering of ${ }^{40}$ Ar at $E_{1 a b}=170 \mathrm{MeV}$. The quadrupole moment of ${ }^{238} \mathrm{U}$ is taken to be $11.12 \mathrm{eb}^{2}$. The results given by the Winther-de Boer code are indicated by dark circles joined by full lines. The classical S-matrix values are indicated by open circles joined by dashed lines. The energy levels of ${ }^{238} \mathrm{U}$ are taken from the rotational model.

Fig. 2. Calculation of Coulomb excitation probabilities to excite the rotational band in ${ }^{152} \mathrm{Sm}$ for backscattering of $8^{8} \mathrm{Be}$ at $E_{1 a b}=30 \mathrm{MeV}$. The Quantum Mechanical calculations were performed using the coupled channel code AROSA. ${ }^{8)}$ The quadrupole moment of ${ }^{152} \mathrm{Sm}$ is taken to be $5.85 \mathrm{eb}^{2}$. Quantum Mechanical results are shown as dark circles joined by full lines, USCA results by open squares joined by dotted lines and integral expression (4) values are indicated by open circles joined by dashed lines. The energy levels of ${ }^{152}$ Sm are taken from the rotational model for all calculations.

$$
00104404174
$$



Fig. 1


Fig. 2

## LEGAL NOTICE

This report was prepared as an account of work sponsored by the United States Government. Neither the United States nor the United States Energy Research and Development Administration, nor any of their employees, nor any of their contractors, subcontractors, or their employees, makes any warranty, express or implied, or assumes any legal liability or responsibility for the accuracy, completeness or usefulness of any information, apparatus, product or process disclosed, or represents that its use would not infringe privately owned rights.

TECHNICAL INFORMATION DIVISION
LAWRENCE BERKELEY LABORATORY
UNIVERSITY OF CALIFORNIA
BERKELEY, CALIFORNIA 94720


[^0]:    *This work was supported by the U. S. Energy Research and Development Administration.

