

LA--10248-MS

DE85 010323

PRECO-D2: Program for Calculating Preequilibrium and Direct Reaction Double Differential Cross Sections

C. Kaibach*

DISCLAIMER

This report was prepared as an account of work sponsored by an agency of the United States Government. Neither the United States Government nor any agency thereof, nor any of 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. Reference herein to any specific commercial product, process, or service by trade name, trademark,

*Consultant at Los Alamos, Department of Physics, Duke University, Durham, NC 27706.

Los Alamos

Los Alamos National Laboratory
Los Alamos, New Mexico 87545

ANSWER *What is the name of the author of the book?*

TABLE OF CONTENTS

ABSTRACT	1
INTRODUCTION	2
I. SUMMARY OF THE EXCITON MODEL.	3
I.1 State Densities.	3
I.2 Residual Interaction Rates	6
I.3 Particle Emission Rates.	11
I.4 Fission Rates.	13
I.5 Closed Form Reaction Equations	14
I.6 Pairing Corrections.	17
II. DIRECT REACTION MODELS.	18
II.1 Nucleon Transfer	19
II.2 Knockout and Inelastic Processes with Complex Particle Degrees of Freedom	20
III. ANGULAR DISTRIBUTIONS	23
IV. ISOSPIN	24
V. DESCRIPTION OF THE PROGRAM.	25
VI. USE OF THE PROGRAM.	28
VI.1 Initial Input	28
VI.2 Initial Calculations.	29
VI.3 Initial Output.	30
VI.4 Secondary Input	32
VI.5 Secondary Calculations.	33
VI.6 Calculations with Isospin	33
VI.7 Suggested Input	34
REFERENCES	36
FIGURE 1. Flow Diagram of PRECO-D2.	38

TABLE OF CONTENTS (Cont.)

TABLE 1.	List of Variables in PRECO-D2	39
TABLE 2.	Additional Variables in Subroutines	44
APPENDIX A.	FORTRAN LISTING OF PRECO-D2	46
APPENDIX B.	SAMPLE INPUT.	71
APPENDIX C.	SAMPLE OUTPUT	72
APPENDIX D.	NEW FEATURES IN PRECO-D2.	107

PRECO-D2: PROGRAM FOR CALCULATING PREEQUILIBRIUM
AND DIRECT REACTION DOUBLE DIFFERENTIAL CROSS SECTIONS

by

C. Kalbach

ABSTRACT

The code PRECO-D2 uses the exciton model for preequilibrium nuclear reactions to describe the emission of particles with mass numbers of 1 to 4 from an equilibrating composite nucleus. A distinction is made between open and closed configurations in this system and between the multi-step direct (MSD) and multi-step compound (MSC) components of the preequilibrium cross section. Additional MSD components are calculated semi-empirically to account for direct nucleon transfer reactions and direct knockout processes involving cluster degrees of freedom. Evaporation from the equilibrated composite nucleus is included in the full MSC cross section. Output of energy differential and double differential cross sections is provided for the first particle emitted from the composite system. Multiple particle emission is not considered. This report describes the reaction models used in writing PRECO-D2 and explains the organization and utilization of the code.

INTRODUCTION

PRECO refers to a family of programs of increasing sophistication designed to calculate the energy spectra of particles emitted in nuclear reactions within a statistical model framework which includes non-equilibrium processes. The first series of programs (ending with PRECO-8 and PRECOM) ran on a PDP-6 computer with teletype input in the dialogue mode. The present series, which began with the code .RECO-A, is suitable for large computing center machines. The model used was first proposed by Griffin [GR66] and is discussed in detail in [KA77] and references therein. The formalism used here has been modified to include division of the preequilibrium cross section into multi-step direct (MSD) and multi-step compound (MSC) components [KA81]. Subroutines have been included to perform semi-empirical calculations for direct nucleon transfer and for knockout and inelastic processes involving complex particles. Additional subroutines use the total MSD (including direct) and total MSC (including evaporation) cross sections to calculate the angular distributions for

the emitted particles. This is done phenomenologically [KA81a].

I. SUMMARY OF THE EXCITON MODEL THEORY

The Griffin or exciton model is a simple statistical model which totally neglects angular momentum and shell structure. The nucleus is pictured with equally spaced single-particle states, and the interactions responsible for creating and destroying particle/hole pairs are assumed to be two-body, energy conserving and residual in nature. Particle emission rates are calculated from microscopic reversibility.

I.1 State Densities

The states of the system are enumerated by the number of excited particle, p , and hole, h , degrees of freedom which they contain. The sum $p+h$ is referred to as the exciton number and is denoted as n . The system is assumed to be formed in a unique particle/hole configuration specified by (p_o, h_o) , and the difference $p-h$ is assumed to remain constant so that $p-h = p_o-h_o$. This is not true near equilibrium but is adequate for closed-form preequilibrium calculations. Thus once p is specified for a given class of states, h and n are also specified if the initial

configuration is known. The density of equally spaced single-particle states is denoted as g_o .

For a system with excitation energy E , the overall particle/hole state density is given by

$$\omega(p, h, E) = \frac{g^n(p) [E - A(p, h)]^{n-1}}{p! h! (n-1)!} f(p) , \quad (1)$$

$$A(p, h) = E_{\text{Pauli}}(p, h) - (p^2 + h^2 + n)/4g_o , \quad (2)$$

$$E_{\text{Pauli}}(p, h) = p_m^2/g_o , \quad (3)$$

$$p_m = \text{maximum}(p, h) . \quad (4)$$

The quantity $g(p)$ contains corrections to g_o assuming that in the real nucleus the single-particle-state density varies as the square root of the energy in the well, with g_o representing the average near the Fermi surface. Thus

$$g(p) = \begin{cases} g_o \left[\frac{p}{n} \left(\frac{V+E/n}{V} \right)^{1/2} + \frac{h}{n} \left(\frac{V-E/n}{V} \right)^{1/2} \right] & \text{for } h \leq 2 \\ g_o & \text{for } h > 2 . \end{cases} \quad (5)$$

Here V is the depth of the nuclear potential well and is assumed to be 38 MeV. The variable $f(p)$ gives corrections to the state density due to the finite depth of the nuclear potential. It has the form (where alternatives of V may be read in)

$$f(p) = \begin{cases} 1 - h \left(\frac{E-V}{V} \right)^{n-1} \Theta(E-V) + \frac{h(h-1)}{2} \left(\frac{E-2V}{V} \right)^{n-1} \Theta(E-V) & \text{for } h \leq 2 \\ 1 & \text{for } h > 2 . \end{cases} \quad (6)$$

Here θ is the Heaviside function, which is zero for a negative argument and unity for a positive one.

In the MSD/MSC formalism, the density of states containing at least one unbound particle degree of freedom is also needed for each class of states. This is referred to simply as the density of unbound states. In order to specify it, we need to define the effective separation energy, S , which determines the excitation energy at which particle degrees of freedom become unbound. In the present model we assume

$$S = \min(B_n, B_p + C_p, B_\alpha + C_\alpha), \quad (7)$$

where the B 's are the binding energies and the C 's are the Coulomb barriers for the neutron, proton, and alpha particle exit channels. The Coulomb barrier for a particle of type b is given in terms of its proton number, Z_b , and the proton number, Z_B , of the residual nucleus so that

$$C_b = 0.75 \frac{Z_b Z_B}{A^{1/3}} \text{ MeV}. \quad (8)$$

The density of unbound states is written as

$$\omega^{(u)}(p, h, E) = \frac{g_p(p) g_u^{n-1}(p) [E - A_{1,0}(p, h) - S]^{n-1}}{(p-1)! h! (n-1)!} f_1(p), \quad (9)$$

and contains different effective single-particle-state densities, g_p and g_u , for the unbound particle and the remaining degrees of freedom. These are

$$g_u(p) = \begin{cases} g_o \left[\frac{p-1}{n-1} \left(\frac{V + (E-S)/n}{V} \right)^{1/2} + \frac{h}{n-1} \left(\frac{V - (E-S)/n}{V} \right)^{1/2} \right] & \text{for } h \leq 2 \\ g_o & \text{for } h > 2, \end{cases} \quad (10)$$

$$g_p(p) = \begin{cases} g_o \left(\frac{V+S+(E-S)/n}{V} \right)^{1/2} & \text{for } h \leq 2 \\ g_o \left(\frac{V+S}{V} \right)^{1/2} & \text{for } h > 2 \end{cases} \quad (11)$$

The finite well depth corrections are contained in the quantity $f_1(p)$ which, for $h \leq 2$, is given by

$$f_1(p) = \frac{1}{p} \sum_{i=1}^k \sum_{j=0}^h (-1)^{i+j+1} \binom{p}{i} \binom{h}{j} \left(\frac{E-iS-jV}{E-S} \right)^{n-1} \Theta(E-iS-jV), \quad (12)$$

while for $h > 2$ only $j=0$ terms are included. The quantity $A_{1,0}(p,h)$ is analogous to $A(p,h)$ and is defined by [KA83]

$$A_{i,0}(p,h) = \frac{p_m^2 + (p_m - i)^2}{2g_o} - \frac{(p-i)^2 + h^2 + n-i}{4g_o} . \quad (13)$$

The density of bound states is also needed and is found by a simple difference

$$\omega^{(b)}(p, h, E) = \omega(p, h, E) - \omega^{(u)}(p, h, E) . \quad (14)$$

1.2 Residual Interaction Rates

The residual interactions which are responsible for changing the exciton number of the system are assumed to be energy conserving and two-body in nature so that allowed transitions are those for which $\Delta p = \Delta h = 0, \pm 1$. The rates for

these three general categories of transitions are denoted λ_0 , λ_+ and λ_- . In addition, the superscripts u and b are used to indicate the unbound or bound character of the initial and final state in the interaction. Thus $\lambda_+^{(ub)}(p,h,E)$ is the average rate for creating a particle/hole pair starting from an unbound state in the class specified by (p,h) and going to a bound final state in the class $(p+1,h+1)$. All of the transition rates have the general form exemplified below:

$$\lambda_+^{(ub)}(p,h,E) = (2\pi/\hbar) M^2(p) \omega_+^{(ub)}(p,h,E), \quad (15)$$

where M^2 is the mean square two-body matrix element and $\omega_+^{(ub)}$ gives the density of accessible final states. The quantity M^2 is given by [KA78]

$$M^2(p) = \begin{cases} \frac{k}{A^3} \frac{n}{E} \left[\frac{E/n}{7 \text{ MeV}} \frac{E/n}{2 \text{ MeV}} \right]^{1/2} & \text{for } E/n < 2 \text{ MeV}, \\ \frac{k}{A^3} \frac{n}{E} \left[\frac{E/n}{7 \text{ MeV}} \right]^{1/2} & \text{for } 2 \text{ MeV} \leq E/n < 7 \text{ MeV}, \\ \frac{k}{A^3} \frac{n}{E} & \text{for } 7 \text{ MeV} \leq E/n < 15 \text{ MeV}, \\ \frac{k}{A^3} \frac{n}{E} \left[\frac{15 \text{ MeV}}{E/n} \right]^{1/2} & \text{for } 15 \text{ MeV} < E/n, \end{cases} \quad (16)$$

where A is the mass number of the composite nucleus and k is an empirical constant usually taken to be 135 MeV^3 .

Several quantities are useful in describing the densities of accessible final states. We therefore define

$$x_i(p) = [E - A_{i,0}(p,h) - iS] \Theta(E - A_{i,0}(p,h) - iS), \quad (17)$$

as well as the correction functions

$$f_+(p) = \begin{cases} 1 - \left(\frac{E-S-V}{E-S}\right)^{n-1} \Theta(E-S-V) - \frac{1}{2} \left(\frac{E-2S}{E-S}\right)^{n-1} \Theta(E-2S) \\ \quad + \frac{1}{2} \left(\frac{E-2S-V}{E-S}\right)^{n-1} \Theta(E-2S-V) & \text{for } h \leq 2, \\ 1 - \frac{1}{2} \left(\frac{E-2S}{E-S}\right)^{n-1} \Theta(E-2S) & \text{for } h > 2, \end{cases} \quad (18)$$

$$f_0(p) = 1 - \frac{1}{2} \left(\frac{E-2S}{E-S}\right)^{n-1} \Theta(E-2S); \quad (19)$$

and the effective single-particle-state densities

$$g_a(p+1) = \begin{cases} \frac{n-1}{n} g_u(p+1) + \frac{1}{n} g_p(p+1) & \text{for } h \leq 1, \\ g_o & \text{for } h > 1, \end{cases} \quad (20)$$

$$g_h = g_o [(V-S)/V]^{1/2}, \quad (21)$$

where we have assumed that $V > S$. The fraction of unbound states in a given class which have more than one unbound particle also enters in and is given by

$$m(p) = \frac{\sum_{i=2}^p \sum_{j=0}^h (-1)^{i+j} \binom{p}{i} \binom{h}{j} \Theta(E-iS-jV) (E-iS-jV)^{n-1}}{\sum_{i=1}^p \sum_{j=0}^h (-1)^{i+j+1} \binom{p}{i} \binom{h}{j} \Theta(E-iS-jV) (E-iS-jV)^{n-1}} \quad (22)$$

for $h \leq 2$, while for $h > 2$ only $j=0$ is considered. Finally we need the three miscellaneous quantities

$$f_u(p) = \frac{2f_1(p) m(p)}{p-1} \left(\frac{E-S}{E-2S}\right)^{n-1}, \quad (23)$$

$$G_1 = \text{maximum}(E-V, 0) , \quad (24a)$$

$$G_2 = \text{maximum}(E-V-S, 0) . \quad (24b)$$

In terms of these quantities, the necessary densities of final states for the two-body interactions are given by

$$\begin{aligned} \omega_{+}^{(uu)}(p, h, E) &= \frac{g_a(p+1) g_u^2(p+1)}{2n} \frac{x_1^{n+1}(p+1) \left[\frac{1}{n} f_+(p+1) + \frac{n-1}{n} f(p+1) \right]}{x_1^{n-1}(p)} \\ &\quad + \frac{m(p)}{1-m(p)} \omega_{+}^{(ub)}(p, h, E) , \end{aligned} \quad (25)$$

$$\begin{aligned} \omega_{+}^{(ub)}(p, h, E) &= \frac{1-m(p)}{2n(n+1)} \left\{ g^3(p+1) f(p+1) \right. \\ &\quad \left. + \frac{n^2}{2} [x_0(p+1) - x_1(p)]^2 + \frac{n}{2} [x_0^2(p+1) - x_1^2(p)] + x_0(p+1) x_1(p) \right\} \\ &\quad - 2g_p(p+1) g_u^2(p+1) f_+(p+1) \frac{x_1^{n+1}(p+1)}{x_1^{n-1}(p)} , \end{aligned} \quad (26)$$

$$\begin{aligned} \omega_{+}^{(bu)}(p, h, E) &= \frac{g_p(p+1) g_u^2(p+1) h}{x_0^{n-1}(p) f(p) - px_1^{n-1}(p) f_1(p)} \\ &\quad \left\{ \frac{x_1^{n+1}(p+1) - px_2^{n+1}(p+1)}{2n(n+1)} - \frac{1}{4n(n+1)} \right. \\ &\quad \left. \left[G_1^{n-1} [n(n-1) G_1^2 - 2(n+1)(n-1) G_1 x_1(p+1) + n(n+1) x_1^2(p+1)] \right. \right. \\ &\quad \left. \left. - p G_2^{n-1} [n(n-1) G_2^2 - 2(n+1)(n-1) G_2 x_2(p+1) + n(n+1) x_2^2(p+1)] \right] \right\} , \end{aligned} \quad (27)$$

$$\begin{aligned}
\omega_{+}^{(bb)}(p, h, E) = & - \frac{1}{2n(n+1)} \frac{1}{x_0^{n-1}(p) f(p) - p x_1^{n-1}(p) f_1(p)} \\
& \left\{ ng^3(p+1) x_0^{n+1}(p+1) f(p+1) \right. \\
& - p(n-1) g_u^3(p+1) x_1^{n+1}(p+1) f_1(p+1) \\
& - pg^3(p+1) x_1^{n-1}(p) f_1(p) \\
& \left. \left\{ \frac{n^2}{2} [x_0(p+1) - x_1(p)]^2 + \frac{n}{2} [x_0^2(p+1) - x_1^2(p)] + x_0(p+1) x_1(p) \right\} \right\} \\
& - \omega_{+}^{(bu)}(p, h, E), \tag{28}
\end{aligned}$$

$$\begin{aligned}
\omega_0^{(ub)}(p, h, E) = & \frac{[1-m(p)]g_0^2}{n} \left\{ \frac{p+2h-1}{2} n[x_0(p) - x_1(p)] f(p) \right. \\
& \left. + (p-1) x_1(p) [f(p) - 2f_0(p)] \right\}, \tag{29}
\end{aligned}$$

$$\begin{aligned}
\omega_0^{(bu)}(p, h, E) = & \frac{g_u(p) g_p(p)}{2n} \frac{1}{x_0^{n-1}(p) f(p) - p x_1^{n-1}(p) f_1(p)} \\
& \left\{ (p+2h-1) \{2x_1^n(p) + n[x_0(p) - x_1(p)]x_1^{n-1}(p)\} f(p) \right. \\
& - (p-1)(p+2h-1) \{2x_2^n(p) + n[x_1(p) - x_2(p)]x_2^{n-1}(p)\} f_1(p) \\
& \left. - 4(n-1) x_1^n(p) f_1(p) + 4(p-1)(n-2) x_2^n(p) f_2(p) \right\}, \tag{30}
\end{aligned}$$

$$\omega_{-}^{(ub)}(p, h, E) = [1-m(p)] g_h h(h-1)/2, \tag{31}$$

$$\begin{aligned}
\omega_{-}^{(bu)}(p, h, E) = & \frac{g_p(p-1) \ h p(p-1)}{4} \frac{1}{x_0^{n-1}(p) \ f(p) - px_1^{n-1}(p) \ f_1(p)} \\
& \{x_1^{n-3}(p-1)[(n-2)(n-3)x_1^2(p-1) \\
& - 2(n-1)(n-3)x_1(p-1) \ x_0(p) + (n-1)(n-2)x_0^2(p)] \ f(p) \\
& - (p-2)x_2^{n-3}(p-1)[(n-2)(n-3)x_2^2(p-1) \\
& - 2(n-1)(n-3)x_2(p-1) \ x_1(p) + (n-1)(n-2)x_1^2(p)] \ f_1(p) \\
& - 4[x_1^{n-1}(p) \ f_1(p) - (p-2)x_2^{n-1}(p) \ f_u(p)]\}. \quad (32)
\end{aligned}$$

I.3 Particle Emission Rates

The average rate for emitting a particle of type b and energy ϵ from an unbound state specified by (p, h) is derived from microscopic reversibility to be

$$w_b^{(u)}(p, h, \epsilon) d\epsilon = \frac{(2s_b+1)}{\pi^2 \ h^3} A_b \ \epsilon \ \sigma_b(\epsilon) d\epsilon \ Q_b(p) \ \frac{\omega(p-A_b, h, U)}{\omega^{(u)}(p, h, E)}, \quad (33)$$

where s_b and A_b are the spin and mass numbers of the emitted particle. The quantity σ_b is the cross section for the inverse process of composite nucleus formation, while U is the excitation energy of the residual nucleus. The quantity $Q_b(p)$ takes account of the fact that proton and neutron degrees of freedom are distinguishable. It assumes that in each pair-creation interaction protons and neutrons

are excited with the relative probabilities Z/A and N/A , and it has the form

$$Q_b(p) = \begin{cases} \frac{(p-A_b)!}{p!} \sum_{i=0}^{p-A_a} \frac{(p-A_a)!}{i! (p-A_a-i)!} \left(\frac{Z_T}{A_T}\right)^{i-Z_b} \left(\frac{N_T}{A_T}\right)^{p-A_a-i-N_b} \\ \frac{(Z_a+i)!}{(Z_a+i-Z_b)!} \frac{(p-Z_a-i)!}{(p-Z_a-i-N_b)!} & \text{for } n \leq \bar{n} \text{ and } n \leq 20 \\ 1 & \text{for } n > \bar{n} \text{ or } n > 20. \end{cases} \quad (34)$$

Here Z_b and N_b are the proton and neutron numbers of the emitted particle, and $A_a = Z_a + N_a$ is the nucleon number of the projectile. The quantities Z_T , N_T and A_T are the proton, neutron and mass numbers of the target. Finally \bar{n} is the most probable number of degrees of freedom at equilibrium (i.e., the number for which $\omega(p, h, E)$ is a maximum).

An alternative to $Q_b(p)$ has been added in this newer version of PRECO-D. It is assumed that proton and neutron particle/hole pairs are excited in proportion to the state densities of the configurations formed [GA77]. This assumption is more consistent with the assumption made in deriving the emission rates from microscopic reversibility (i.e., all configurations of a given p , h , and E are equally likely to be populated). It leads to the Q_b values of

$$Q_b^{(\pm)}(p) = \frac{\frac{(Z)^n \cdot \pi^{Z_b(N_b + h_b - N_b)}}{A} \frac{(p - A_b)!}{(p_z - Z_b)! (p_n - N_b)! h_z! h_n!}}{\sum_p \frac{(Z)^n \cdot (N)^n}{(A)^n} \frac{p!}{p_z! p_n! h_z! h_n!}}, \quad (35)$$

where the subscripts z and n refer to proton and neutron degrees of freedom, and Z , N , and A are for the composite nucleus. Either form for Q may be chosen by specifying an input parameter.

Particle emission rates for the bound states are, of course, zero.

At equilibrium the emission rates for the full system are calculated from the Weisskopf-Ewing evaporation formula and are given by

$$W_b(\varepsilon) d\varepsilon = \frac{(2s_b + 1)}{\pi^2 \hbar^3} A_b \varepsilon \sigma_b(\varepsilon) d\varepsilon \frac{\omega(U)}{\omega(E)}, \quad (36)$$

where in the present case the state densities are assumed to be those for a single Fermi gas,

$$\omega(E) \propto E^{-1} \exp[2(aE)^{1/2}], \quad (37)$$

and $a = \pi^2 g_0/6$.

I.4 Fission Rates

In order to facilitate application of the model to very heavy nuclei, fission competition in the equilibrium limit has been considered. Barrier penetration has been neglected

in order to obtain a closed-form integral over the kinetic energy degree of freedom. Thus one obtains

$$W_f(E) = \frac{1}{2\pi h} \int_0^{E-B_f} \frac{\omega(E-B_f-\varepsilon_f)}{\omega(E)} d\varepsilon_f , \quad (38)$$

where ε_f is the kinetic energy in the fission degree of freedom.

I.5 Closed-Form Reaction Equations

In the closed-form approach, the strength of the system is imagined to pass sequentially through configurations of increasing complexity until the most probable class of states at equilibrium (specified by $\bar{n}=\bar{p}+\bar{h}$) is reached. Particle emission is allowed at each stage.

In the MSD/MSC formalism it is necessary to know for each class of states how much of the reaction strength passes through unbound states and how much through bound states. These quantities are denoted $S_u(p,h)$ and $S_b(p,h)$, respectively. In addition, it is necessary to know the amount of strength which passes through unbound states of the class (p,h) and which was unbound at all previous stages as well. This is denoted $S_d(p,h)$ and is the strength responsible for the MSD cross section.

Strengths of the three types arriving at the (p,h) stage of the hierarchy through pair creation are denoted $P_u(p,h)$, $P_b(p,h)$ and $P_d(p,h)$, respectively, and differ from the

corresponding S quantities by the effect of the $\lambda_0^{(u)}$ and $\lambda_-^{(u)}$ interactions.

Initially it is assumed that the unbound states are populated in proportion to their relative state densities so that

$$P_u(p_o, h_o) = P_d(p_o, h_o) = \omega^{(u)}(p_o, h_o, E) / \omega(p_o, h_o, E), \quad (39a)$$

$$P_b(p_o, h_o) = 1 - P_u(p_o, h_o). \quad (39b)$$

At this exciton number and at each succeeding exciton number it is assumed that there will be at most one exciton-scattering interaction which takes the system from a bound to an unbound configuration or *vice versa*. Further, it is assumed that there can be at most one pair-annihilation interaction which changes the bound/unbound character of the state and that it will immediately be followed by a pair-creation interaction which does not change the bound/unbound character. Thus in effect $\lambda_0^{(ub)}$ and $\lambda_-^{(ub)}$ are treated together, and similarly for $\lambda_0^{(bu)}$ and $\lambda_-^{(bu)}$.

Because of these assumptions, strength arriving for the first time at a given hierarchy is assigned a lifetime of

$$T_u(p, E) = [\lambda_+^{(uu)}(p) + \lambda_+^{(ub)}(p) + \lambda_0^{(ub)}(p) + \lambda_-^{(ub)}(p) + \sum_b \int W_b(p, \epsilon) d\epsilon]^{-1}, \quad (40a)$$

(where the hole label has been suppressed) if it is in unbound states or

$$T_b(p, E) = [\lambda_+^{(bb)}(p) + \lambda_+^{(bu)}(p) + \lambda_0^{(bu)}(p) + \lambda_-^{(bu)}(p)]^{-1} \quad (40b)$$

if it is in bound configurations. Strength which has already

undergone an exciton-scattering or pair-annihilation interaction starting from this hierarchy is assigned a different lifetime, one of the two quantities

$$T'_u(p, E) = [\lambda_+^{(uu)}(p) + \lambda_+^{(ub)}(p) + \sum_b \int W_b(p, \varepsilon) d\varepsilon]^{-1}, \quad (41a)$$

$$T'_b(p, E) = [\lambda_+^{(bb)}(p) + \lambda_+^{(bu)}(p)]^{-1} \quad (41b)$$

The strengths passing through states of different p values are then found from a set of recursion relations:

$$S_d(p) = P_d(p), \quad (42a)$$

$$S_u(p) = P_u(p) + P_b(p) \Gamma_0^{(bu)}(p) T'_u(p)/T_u(p), \quad (42b)$$

$$S_b(p) = P_b(p) + P_u(p) \Gamma_0^{(ub)}(p) T'_b(p)/T_b(p), \quad (42c)$$

and

$$P_d(p+1) = S_d(p) \Gamma_+^{(uu)}(p) \equiv P_d(p) \Gamma_+^{(uu)}(p), \quad (43a)$$

$$P_u(p+1) = S_u(p) \Gamma_+^{(uu)}(p) + S_b(p) \Gamma_+^{(bu)}(p), \quad (43b)$$

$$P_b(p+1) = S_b(p) \Gamma_+^{(bb)}(p) + S_u(p) \Gamma_+^{(ub)}(p), \quad (43c)$$

with

$$\Gamma_+^{(uu)}(p) = \lambda_+^{(uu)}(p) T_u(p), \quad (44a)$$

$$\Gamma_+^{(ub)}(p) = \lambda_+^{(ub)}(p) T_u(p), \quad (44b)$$

$$\Gamma_0^{(ub)}(p) = [\lambda_0^{(ub)}(p) + \lambda_-^{(ub)}(p)] T_u(p), \quad (44c)$$

and similarly for $\Gamma_+^{(bb)}$, $\Gamma_+^{(bu)}$ and $\Gamma_0^{(bu)}$ but with the u and b labels interchanged.

The preequilibrium energy differential cross sections are then given in terms of the strength variables. The total MSD and MSC preequilibrium cross sections are

$$\frac{d\sigma}{d\varepsilon}(a,b)_{PRE} = \sigma_a(\varepsilon_a) \sum_{p=p_0}^{\bar{p}} S_u(p) T_u(p) W_b^{(u)}(p, \varepsilon), \quad (45a)$$

$$\frac{d\sigma}{d\varepsilon}(a,b)_{pre-MSD} = \sigma_a(\varepsilon_a) \sum_{p=p_0}^{\bar{p}} S_d(p) T_u(p) W_b^{(u)}(p, \varepsilon), \quad (45b)$$

$$\frac{d\sigma}{d\varepsilon}(a,b)_{pre-MSC} = \frac{d\sigma}{d\varepsilon}(a,b)_{PRE} - \frac{d\sigma}{d\varepsilon}(a,b)_{pre-MSD}, \quad (45c)$$

where $\sigma_a(\varepsilon_a)$ is the cross section for forming the composite nucleus when the projectile has energy ε_a . The evaporation or equilibrium components are determined by taking all of the input reaction cross section not used in the preequilibrium phase of the reaction and distributing it among the various reaction channels in proportion to their phase space. Only evaporation from the original compound nucleus is presently considered in the program. Subsequent emission from the residual nuclei is left out.

The energy differential evaporation cross section for the first emitted particle in a reaction is

$$\frac{d\sigma}{d\varepsilon}(a,b)_{EVAP} = [\sigma_a(\varepsilon_a) - \sigma_{PRE}] \frac{W_b(\varepsilon)}{\sum_b \int W_b(\varepsilon) d\varepsilon}, \quad (46)$$

where the total preequilibrium cross section, σ_{PRE} , is

$$\sigma_{PRE} = \sum_b \int \frac{d\sigma}{d\varepsilon}(a,b)_{PRE} d\varepsilon. \quad (47)$$

I.6 Pairing Corrections

While preequilibrium pairing corrections are best made and studied in the two-component exciton model, they are

sometimes included in the one-component model.

In the present code, two sets of pairing corrections are used: $\Delta_{\text{pre}}(Z,N)$ is used for states with $h \leq 2$, while the equilibrium pairing correction $\Delta_{\text{eq}}(Z,N)$ is used for $h > 2$. The quantity Δ_{pre} may be zero, the equilibrium value or a special preequilibrium pairing correction, perhaps related to the number of broken proton or neutron pairs in states with $n=n_0$ [GR73][GR74]. Likewise, $\Delta_{\text{eq}}(Z,N)$ may also be taken to be zero.

Whenever pairing energy corrections are used, the quantity E in the state density expressions should be replaced by $(E - \Delta_{\text{pre}})$ or $(E - \Delta_{\text{eq}})$ depending on the value of h , and similarly for U in the state densities for the residual nuclei.

II. DIRECT REACTION MODELS

Two classes of direct reactions which are not included in the Griffin model are nucleon transfer (stripping, pickup and exchange) and knockout or inelastic-scattering processes which involve complex particle degrees of freedom. These are treated semi-empirically in subroutines in PRECO-D. The formulae used are refined somewhat from those in [KA77].

II.1 Nucleon Transfer

The basic equation for the energy differential cross section for nucleon transfer is

$$\frac{d\sigma}{d\varepsilon}_{NU} = \frac{(2s_b+1)}{(2s_a+1)} \frac{A_b}{A_a} \frac{\varepsilon \sigma_{b-NE}(\varepsilon)}{A_a \varepsilon_a} K_{\alpha,p} \left(\frac{A_a}{E_a + V_a} \right)^{2n} \left(\frac{2860}{A_B} \right)^n$$

$$0.0127 \sum_{p_\pi} \left(\frac{2Z_T}{A_\pi} \right)^{6n_\pi} \frac{p!}{p_\pi! p_\nu!} \frac{h!}{h_\pi! h_\nu!} \frac{g_\pi^{n_\pi} g_\nu^{n_\nu}}{g^n}$$

$$\omega(p, h, \Omega) , \quad (48)$$

where the subscripts a and B refer to the projectile and residual nucleus. The inverse cross sections, $\sigma_b(\varepsilon)$, used in the main program are assumed to be the total composite nucleus formation cross section, while $\sigma_{b-NE}(\varepsilon)$ here is the total nonelastic cross section. The two are assumed to be related by

$$\sigma_b(\varepsilon) \begin{cases} = 0.95 \sigma_{b-NE}(\varepsilon) & \text{for } b = \text{nucleon} \\ = 0.85 \sigma_{b-NE}(\varepsilon) & \text{for } b = \text{complex} . \end{cases} \quad (49)$$

The quantities E_a and ε_a are the laboratory and center of mass energies of the projectile, while V_a is the average potential seen by the projectile in the direct reaction region. It is taken to be one fourth of the real central well depth or

$$V_a = \frac{50 \text{ MeV } A_a}{4} . \quad (50)$$

The constant $K_{\alpha,p}$ gives a factor of 12 enhancement to the

nucleon transfer whenever the projectile and ejectile are both tightly bound (i.e., a nucleon or an alpha particle). It is unity for all other reactions. The particle and hole numbers are given by $p=p_{\pi}+p_{\nu}$ and $h=h_{\pi}+h_{\nu}$, where p_{π} , p_{ν} , h_{π} , and h_{ν} are the numbers of stripped protons and neutrons and picked-up protons and neutrons, respectively. In general, these variables are uniquely determined by the natures of the incoming and outgoing particles. For inelastic channels of weakly bound projectiles (d , t or 3He), however, exchange of either a proton or a neutron with the target is allowed. Thus there will be two contributions to the nucleon transfer cross section, one with $p_{\pi}=h_{\pi}=1$ and $p_{\nu}=h_{\nu}=0$, and the other having $p_{\pi}=h_{\pi}=0$ and $p_{\nu}=h_{\nu}=1$. This is the only case where the sum over p_{π} is needed. Finally, g_{π} and g_{ν} are the proton and neutron single-particle-state densities, assumed to be $g_{\pi} = g_o (Z/A)$ and $g_{\nu} = g_o (N/A)$ evaluated in the residual nucleus.

The nucleon-transfer cross section as given here will be in mb/MeV. All of the energies are assumed to be in MeV and the nonelastic cross sections in mb.

II.2 Knockout and Inelastic Processes with Complex Particle Degrees of Freedom

In this type of mechanism, a complex projectile is assumed to be able to excite a proton, neutron or alpha cluster particle/hole pair in its initial interaction and

to retain its own cluster identity in the interaction.

Nucleon pair excitation by a nucleon projectile is already considered in the exciton model, so only alpha cluster excitation is considered here for incident neutrons and protons. The basic equations for knockout and inelastic scattering are

$$\frac{d\sigma}{d\varepsilon}(a,b)_{KO} = \frac{\sigma_{a-NE}}{13.5} (2s_b+1) A_b \varepsilon \sigma_{b-NE}(\varepsilon)$$

$$\frac{P_b g_a g_b [U-A(A_a, A_b)]}{\sum_{c=a,b} (2s_c+1) A_c \bar{\sigma}_c (\varepsilon_m + 2B_c) (\varepsilon_m - B_c)^2 (g_a g_b^2 / 6g_c)} . \quad (51)$$

$$\frac{d\sigma}{d\varepsilon}(a,a')_{INEL} = \frac{\sigma_{a-NE}}{13.5} (2s_a+1) A_a \varepsilon \sigma_{a-NE}(\varepsilon)$$

$$\sum_{i=n,p,\alpha} \frac{P_i g_i^2 [U-A(A_i, A_i)]}{\sum_{c=i,a} (2s_c+1) A_c \bar{\sigma}_c (\varepsilon_m + 2B_c) (\varepsilon_m - B_c)^2 (g_a g_i^2 / 6g_c)}$$

for $a = \text{complex}$, (52a)

$$\frac{d\sigma}{d\varepsilon}(a,a')_{INEL} = \frac{\sigma_{a-NE}}{13.5} (2s_a+1) A_a \varepsilon \sigma_{a-NE}(\varepsilon)$$

$$\frac{P_\alpha g_\alpha [U-A(4,4)]}{\sum_{c=a,\alpha} (2s_c+1) A_c \bar{\sigma}_c (\varepsilon_m + 2B_c) (\varepsilon_m - B_c)^2 (g_a g_\alpha^2 / 6g_c)}$$

for $a = \text{nucleon}$, (52b)

The quantity ε_m is the energy of the ground state transition for emission of a particle of type c , while σ_{a-NE} with no energy variable specified is the nonelastic cross section for the entrance channel. The cross section $\bar{\sigma}_c$ is the average for all emission energies above the Coulomb barrier

for particles of type c. The probabilities for exciting the different kinds of particle/hole pairs are

$$p_n = \frac{N_T - f Z_T}{A_T - 2f Z_T + f Z_T/2} \approx \frac{N_T}{A_T}, \quad (53a)$$

$$p_p = \frac{Z_T - f Z_T}{A_T - 2f Z_T + f Z_T/2} \approx \frac{Z_T}{A_T}, \quad (53b)$$

$$p_\alpha = \frac{f Z_T/2}{A_T - 2f Z_T + f Z_T/2} \approx \frac{f Z_T}{2 A_T}. \quad (53c)$$

Here the quantity f is defined as the fraction of the time that four nucleons in correlated orbits will "look like" an alpha cluster or, alternatively, the fraction of the possible alpha clusters that will, on the average, exist at any given time. It has been assumed that $N > Z$, so that a maximum of $Z/2$ alpha clusters is possible. In general $f \ll 1$. The exact size and systematics of f are not well known. Here the values obtained from (p, α) reactions by neglecting pickup [MIT4] are renormalized slightly and parameterized to give

$$f = \begin{cases} 0.08 & N_T \leq 116 \\ 0.02 + 0.06(126 - N_T)/10 & 116 < N_T < 126 \\ 0.02 + 0.06(N_T - 126)/3 & 126 \leq N_T < 129 \\ 0.08 & 129 \leq N_T \end{cases} \quad (54)$$

In evaluating the single-particle-state densities for the different types of emitted particles, it is assumed that all of the nucleons of a cluster are in correlated orbits, each holding a maximum of two protons and two neutrons. In addition it should be recognized that a cluster of A_b

nucleons will carry A_b times the energy of a single particle in one of these orbits. Thus we get

$$g_n = g_o N/A = (N/13) \text{ MeV}^{-1}, \quad (55a)$$

$$g_p = g_o Z/A = (Z/13) \text{ MeV}^{-1}, \quad (55b)$$

$$g_d = g_o/4 = (A/52) \text{ MeV}^{-1}, \quad (55c)$$

$$g_t = g_\tau = g_o/12 = (A/156) \text{ MeV}^{-1}, \quad (55d)$$

$$g_\alpha = g_o/16 = (A/208) \text{ MeV}^{-1}. \quad (55e)$$

In each case they are evaluated in the appropriate residual nucleus. Once again, the final cross sections are in mb/MeV.

III. ANGULAR DISTRIBUTIONS

The angular distributions are calculated phenomenologically in terms of Legendre polynomials using the formula [KA81a]

$$\begin{aligned} \frac{d^2\sigma}{d\Omega d\varepsilon}(a,b) &= a_o(\text{MSD}) \sum_{\ell=0}^6 b_\ell(e) P_\ell(\cos\theta) \\ &\quad + a_o(\text{MSC}) \sum_{\substack{\ell=0 \\ \Delta\ell=2}}^6 b_\ell(e) P_\ell(\cos\theta), \end{aligned} \quad (56)$$

where $e = \varepsilon + I_{\text{ang}} \cdot B_b$ and I_{ang} is either 0 or 1, and where

$$a_o(\text{MSD}) = \frac{1}{4\pi} \left[\frac{d\sigma}{d\varepsilon}(a,b)_{\text{pre-MSD}} + \frac{d\sigma}{d\varepsilon}(a,b)_{\text{NU}} + \frac{d\sigma}{d\varepsilon}(a,b)_{\text{KO}} \right]. \quad (57a)$$

(For the reaction (a,a') the knockout contribution is, of course, replaced by an inelastic contribution.)

$$a_o(\text{MSC}) = \frac{1}{4\pi} \left[\frac{d\sigma}{d\varepsilon}(a,b)_{\text{pre-MSC}} + \frac{d\sigma}{d\varepsilon}(a,b)_{\text{EVAP}} \right], \quad (57b)$$

$$b_\ell(\epsilon) \left\{ \begin{array}{ll} = \frac{(2\ell+1)}{1 + \exp[A_\ell(B_\ell - \epsilon)]} & \text{for } \ell \geq 1 \\ = 1 & \text{for } \ell = 1 \end{array} \right. , \quad (58)$$

$$A_\ell = 0.036 \text{ MeV}^{-1} + 0.0039 \text{ MeV}^{-1} \ell(\ell+1), \quad (59a)$$

$$B_\ell = [92 + 6(I_{\text{ang}})] \text{MeV} - 90 \text{ MeV} [\ell(\ell+1)]^{-1/2}. \quad (59b)$$

IV. ISOSPIN

For reactions induced by protons or ^3He ions with sufficient energy, states in the composite nucleus with two different total isospins can be populated: $T_<=T_z$ and $T_>=T_z+1$. Therefore, it may sometimes be desirable to perform preequilibrium calculations in which the isospin quantum number is conserved [KA74].

To perform such isospin-conserved calculations, it is necessary to know the state densities for both the $T_<$ and $T_>$ configurations in a given nucleus. At equilibrium it is usually a reasonable approximation to take

$$\omega(E, T_<) = \omega(E), \quad (60a)$$

$$\omega(E, T_>) = \omega(E - E_{\text{sym}}(Z, A)), \quad (60b)$$

where $E_{\text{sym}}(Z, A)$ is the nuclear symmetry energy and can be obtained from the empirical formula [AN65]

$$E_{\text{sym}}(Z, A) = 1.44 \text{ MeV} \frac{Z+1/2}{A^{1/3}} - 1.13 \text{ MeV} + Q_{(p,n)}(Z-1, A), \quad (61)$$

where $Q_{(p,n)}$ is the Q -value for the (p, n) reaction on the nucleus with the indicated proton and mass numbers.

For particle/hole state densities, a similar approximation has been used [KA74] giving

$$\omega(p,h,E,T_\zeta) = \omega(p,h,E), \quad (62a)$$

$$\omega(p,h,E,T_\gamma) = \omega(p,h,E-E_{\text{sym}}(Z,A)) \quad (62b)$$

Similar assumptions are assumed to apply for the T-conserving transition-state densities.

While such relations are the only ones which can currently be used in PRECO-D because isospin is not explicitly considered, they have a serious shortcoming for small values of $n=p+h$ where the state densities increase only as E^{n-1} [KA84]. Thus, isospin-dependent calculations are probably better carried out using the two-component exciton model code PRECO-E [KA84a] in which more accurate state density formulae are used.

In either code two calculations are performed, one for each isospin in the composite nucleus. The reaction cross sections are multiplied by the appropriate isospin Clebsch-Gordan coefficients and the state densities are evaluated using excitation energies measured relative to the lowest energy state of the isospin in question.

V. DESCRIPTION OF THE PROGRAM

The program PRECO-D2 is written in simple FORTRAN designed to be compatible with most computer systems. A general flow diagram is contained in Fig. 1; a list of variables is given

in Tables 1 and 2; Appendix A is a FORTRAN listing of the code; Appendix B contains sample input; Appendix C gives the corresponding output; and Appendix D lists the differences between PRECO-D2 and its predecessor, PRECO-D.

Input and output device numbers are assigned as IREA and IWRI, respectively, just following the COMMON statements, so they may be easily altered.

The Griffin, or exciton, model calculations are carried out in the main program, while direct reaction cross sections and angular distributions are calculated in subroutines. The subroutines NUTRA, KNOCK, and INEL handle, respectively, nucleon transfer, knockout involving cluster degrees of freedom, and inelastic scattering involving cluster degrees of freedom. The subroutine POLLY calculates a library of Legendre polynomials, while ANGEL generates the angular distributions for the emitted particles.

Reaction cross sections for the projectile and inverse-reaction cross sections for the emitted particles may, as an option, be calculated internally using an empirical approximation [NA79][CH81]. It has been modified here to go over to the geometrical cross section

$$\sigma_{b-\text{GEO}}(\epsilon) = \pi(1.23A^{1/3} + R_b + \lambda)^2 \quad (63)$$

at high energies, where R_b is the radius of the particle (0 for $b=n,p$; 0.8F for $b=d,t,\tau$; and 1.2F for $b=\alpha$). The parameter values are set in subroutine SIGPAR, while the entrance and exit channel cross sections are calculated in subroutine CROSS and the main program, respectively.

Initial calculations are always performed for the principal reaction channels of neutron, proton and alpha particle emission. Fission competition is included at equilibrium if a fission barrier has been specified. Spectra for other ejected light ions may be calculated at a later stage in the program. In the setup and calling of the direct reaction subroutines, it is assumed that both the incoming and outgoing particles have mass numbers no larger than 4.

The range of exciton numbers considered is determined from the input initial configuration, (p_o, h_o) , and the state density, $\omega(p, h, E)$. The difference $p-h$ is assumed to remain constant at p_o-h_o . The smallest $n=p+h$ considered is the minimum value consistent with there being at least one degree of freedom to carry the excitation energy and with both p and h being nonnegative. In most cases $(p_o, h_o) = (A_a + 1, 1)$, so the simplest configuration of the system is $(A_a, 0)$. At the other extreme, the most complex configuration considered has its exciton number given by minimum $(\bar{n}, 20)$, where \bar{n} is the exciton number corresponding to the maximum in the $\omega(p, h, E)$ vs p curve and is therefore the most probable equilibrium value of n .

For particle emission to be allowed from a state, its particle number must be greater than or equal to the nucleon number of the emitted particle (*i.e.*, $p \geq A_b$) and n must be large enough for the residual nucleus to be left with at least one exciton (*i.e.*, $n \geq A_b + 1$).

VI. USE OF THE PROGRAM

VI.1 Initial Input

The input information requested for the first phase of the calculations is given below. Each numbered item corresponds to a new card or record of input, and the appropriate formats are given in parentheses. Default values are also indicated where they exist.

- i) Angular distribution parameter, I_{ang} , and Q-factor parameter, IG, (2I3) for all problems in the data file.
- ii) Effective well depths $V(h=1)$ and $V(h=2)$ in MeV (2F10.2) for all problems in the data file.
- 1) Composite nucleus excitation energy and projectile binding energy in MeV (2F10.2).
- 2) Z and N of the target nucleus (2F10.2)
- 3) n, p, and α binding energies and fission barrier height in the composite nucleus, all in MeV (4F10.2). If no fission barrier is specified, no fission is calculated.
- 4) Pairing energies for the n, p, and α residual nuclei and for the composite nucleus to be used in the preequilibrium phase of the calculation for states with $h \leq 2$ (4F10.2).
- 5) Pairing energy corrections for states with $h > 2$ and for equilibrium calculations. These are the conventional pairing energies that would be used in an evaporation model. (4F10.2).
- 6) Entrance channel nonelastic cross section and estimated average, σ_a , for inelastic scattering (2F10.2). If either one is read in as zero, it is calculated internally using a parametric approximation to the optical model potentials indicated in Sect. VI.7 below.
- 7) Number of channel energies considered, NEPS(1), the lowest channel energy in MeV, and the increment in the channel energy also in MeV (1I3,2F10.2). If the lowest channel

energy is greater than zero, the remaining channel energies and all of the inverse reaction cross sections are calculated internally. (If the channel energy increment is zero, it defaults to 1 MeV.) If the lowest channel energy is read in as zero, input according to item 7a, below, is read in.

- 7a) NEPS(1) records, each containing (4F10.2) a channel energy in MeV and the corresponding n, p and α non-elastic cross sections in mb. The energies must be in ascending order.
- 8) Z and N of the projectile (2I3). If preferential excitation of an unpaired target nucleon is considered, then this nucleon should be counted as if it were part of the projectile.
- 9) Initial particle and hole numbers, p_o and h_o (2I3). If these are read in as zero, p_o defaults to $A_a + 1$ and h_o defaults to unity.
- 10) Single-particle-state densities, g_o , in MeV^{-1} for the composite system; the n, p and α residual nuclei; and the fission saddle point (5F10.2). If only the value for the composite nucleus is read in, the others are calculated assuming that g_o is proportional to A. If g_o for the composite nucleus is read in as zero, it defaults to $A/(13 \text{ MeV})$.
- 11) Matrix element scaling factor divided by 100 (*i.e.*, $k/100$) (1F10.2). If it is read in as zero, it defaults to 1.35 ($k=135 \text{ MeV}$).

VI.2 Initial Calculations

After initial input, the program proceeds to the calculation of state densities, transition rates and particle emission rates. The branching ratios, Γ , are generated and the closed-form calculations are executed.

At this point, the three principal particle types (n, p, and α) are considered in turn. For each, the necessary direct reaction subroutines are called to generate the

appropriate energy spectra. The results of the closed-form exciton model calculations are used to obtain MSD and MSC preequilibrium energy spectra. A first-chance evaporation spectrum is also calculated. For each emission energy the total MSD and MSC cross sections are used in the subroutine ANGEL to calculate an angular distribution. Finally, the single and double differential cross sections are printed out.

VI.3 Initial Output

The first quantities to be printed are S, the average effective separation energy, and the fraction of the total reaction cross section which goes into preequilibrium particle emission. These are part of a heading labelled OCCUPATION PROBABILITIES. Below the heading is a table with the following columns:

P	particle number of the class of states
H	hole number of the class of states
RHOU/RHO	ratio $\omega^{(u)}(p,h,E)/\omega(p,h,E)$ of unbound and total state densities
STRU/STR	ratio $P_u(p)/P(p)$ of strength arriving at states with p particle degrees of freedom which is in unbound states
STRD/STR	ratio $P_d(p)/P(p)$ of MSD strength to total strength arriving at states specified by p

Following this are two pages of output for each particle type, one containing the energy differential spectra and the other the angular distributions for the various outgoing energies. The column headings for the angle-integrated

spectra are

EPS	ϵ , the outgoing channel energy
DIRECT/NUTRA	nucleon transfer cross section from Eq. (48)
DIRECT/KNOCK	knockout or inelastic cross section from Eq. (51) or (52)
PREEQUILIBRIUM/MSD	MSD preequilibrium cross section from Eq. (45b)
PREEQUILIBRIUM/MSC	MSC preequilibrium cross section from Eq. (45c)
EQUIL/WEISS	evaporation cross section from Eq. (46)
TOTAL/MSD	"NUTRA+KNOCK"+"PREEQUILIBRIUM/MSD"
TOTAL/MSC	"PREEQUILIBRIUM/MSC"+"EQUIL/WEISS"
TOTAL/MSD+MSC	grand total

At the bottom of each column is the summed cross section for that component of the spectrum. If a fission barrier has been specified, the equilibrium fission cross section is printed below the sum of EQUIL/WEISS for the alpha spectra.

The angular distributions are printed in 10-deg increments. Angles from 0 to 90 deg are in one block of the table, while angles from 100 to 180 deg are in a second block. The first column of each block gives the channel energy; the remaining columns give the double differential cross section at the angle indicated in the column heading. The last column in the second block gives the angle-integrated cross sections and is identical to TOTAL/MSD+MSC above.

VI.4 Secondary Input

After printout of the occupation probabilities and of the neutron, proton and alpha particle spectra, the program reads the secondary input, which contains the following items:

- 12) Z and N of an additional particle type (2I3). (If $Z < 0$, input passes to item 13a below.)
- 13) spin degeneracy, $2s+1$, of the emitted particle (1I3)
- 14) binding energy of emitted particle in MeV and the single-particle-state density in its residual nucleus in MeV^{-1} (2F10.2). If the single-particle-state density is zero, it defaults to $(g_o A_{\text{res}} / A_{\text{cn}})$.
- 15) preequilibrium ($h < 2$) and equilibrium pairing corrections for the residual nucleus in MeV (2F10.2)
- 16) number of emitted particle energies considered = NEPS(2), the lowest emission channel energy, and the channel energy increment (1I3, 2F10.2). If the lowest energy is greater than zero, the remaining energies and the inverse cross sections are calculated internally. (A zero channel energy increment defaults to 1 MeV.) If the minimum energy is zero (or less), input item 16a is read in.
- 16a NEPS(2) records, each containing a particle energy in MeV and the corresponding nonelastic cross section in mb (2F10.2). Energies must be in increasing order. [This ends the secondary input.]
- 13a) NEWP (1I3)
 - If NEWP = 1, input passes to item 8 in the initial input and new parameter values are read in.
 - If NEWP = 0, input passes to item 1 in the initial input for the beginning of a new problem.
 - If NEWP = -1, the job is terminated.

VI.5 Secondary Calculations

If a new particle type was chosen in item 12) above, the program uses the stored information on the results of the closed-form calculations to obtain the preequilibrium spectra for this particle type, and the appropriate direct reaction subroutines are called. The single and double differential cross sections are then printed in the same format as for the neutron, proton and alpha particle spectra. After printout, the program returns to item 12) in the secondary input so that an additional particle type may be specified.

VI.6 Calculations with Isospin

Preequilibrium model calculations in which isospin conservation is considered can be accomplished with PRECO-D2 by actually doing two calculations, one for each isospin in the composite nucleus. The changes in the input needed for the isospin-conserved calculations relative to the general (isospin-mixed) calculations just described are indicated in the table below for the initial input. All other quantities in the initial input are as indicated in VI.1 and VI.4 above.

In the table, the second set of input for B_n and $\sigma_{n-NE}(\epsilon)$ for the $T_>$ calculations applies if the neutron channel to the $T_>$ states in the residual nucleus is open.

Neutrons decay from the $T_<$ composite nucleus states to the ground state isospin if the residual nucleus is isospin-forbidden.

Normal	$T_<$	$T_>$
E	E	$E - E_{\text{sym}}(Z, A)$
B_n	B_n	$\begin{cases} 0 \\ B_n - E_{\text{sym}}(Z, A) + E_{\text{sym}}(Z-1, A) \end{cases}$
B_p	B_p	$B_p - E_{\text{sym}}(Z, A)$
B_α	B_α	0
$\sigma_{a-\text{NE}}$	$\sigma_{a-\text{NE}}[2T_o/(2T_o+1)]$	$\sigma_{a-\text{NE}}[1/(2T_o+1)]$
$\sigma_{n-\text{NE}}(\varepsilon)$	$\sigma_{n-\text{NE}}(\varepsilon)$	$\begin{cases} 0 \\ \sigma_{n-\text{NE}}(\varepsilon)[2T_o/(2T_o+1)] \end{cases}$
$\sigma_{p-\text{NE}}(\varepsilon)$	$\sigma_{p-\text{NE}}(\varepsilon)[2T_o/(2T_o+1)]$	$\sigma_{p-\text{NE}}(\varepsilon)[1/(2T_o+1)]$
$\sigma_{\alpha-\text{NE}}(\varepsilon)$	$\sigma_{\alpha-\text{NE}}(\varepsilon)$	0
k	k	$k[E - E_{\text{sym}}(Z, A)]/E$

VI.7 Suggested Input

Appropriate values for the various input quantities are summarized below

(p_o, h_o) $(A_a+1, 1)$, where A_a is the mass number of the projectile (the default values).

T_o $A/13$ with values for the residual nuclei calculated in the program (the default).

k 135 MeV³ (input number is 1.35) (the default).

B_b from [WA77] or similar mass tables

$\sigma_{\text{e-NE}}(\epsilon)$	from [MA63]	} (The default is approximations to these optical model cross sections.)
$\sigma_{\text{t-NE}}(\epsilon)$	from [BE69]	
$\sigma_{\alpha-\text{NE}}(\epsilon)$	from [HU62]	
$\sigma_{\text{b-NE}}(\epsilon)$	from [CL72] for b=d,t, ^3He	

The choice of IANG = 0 or 1 is discussed in the second paper in [Ka81a]. In general, IANG=1 is probably to be preferred.

In addition, there is the choice of whether to use IG = 0, invoking Eq. (34), or IG = 1, invoking Eq. (35), for the Q-factors that account for proton/neutron distinguishability. For reasons of consistency in the model, IG = 1 is to be preferred.

REFERENCES

- [AN65] J. D. Anderson, C. Wong and J. W. McClure, Phys. Rev. 138 (1965) B615
- [BE69] F. D. Becchetti, Jr. and G. W. Greenlees, Phys. Rev. 182 (1969) 1190
- [CH81] A. Chatterjee, K. H. N. Murthy and S. K. Gupta, Pramāna 16 (1981) 391
- [CL72] C. K. Cline, Nucl. Phys. A193 (1972) 417
- [GA77] E. Gadioli, E. Gadioli-Erba and J. Hogan, Phys. Rev. C 16 (1977) 1404
- [GR66] J. J. Griffin, Phys. Rev. Lett. 17 (1966) 478
- [GR73] S. M. Grimes, J. D. Anderson, J. C. Davis and C. Wong, Phys. Rev. C 8 (1973) 1770
- [GR76] S. M. Grimes, J. D. Anderson, and C. Wong, Phys. Rev. C 13 (1976) 2224
- [HU62] J. R. Huizenga and G. Igo, Nucl. Phys. 29 (1962) 462
- [KA74] C. Kalbach-Cline, J. R. Huizenga and H. K. Vonach, Nucl. Phys. A222 (1974) 405
- [KA77] C. Kalbach, Z. Phys. A283 (1977) 401
- [KA78] C. Kalbach, Z. Phys. A287 (1978) 319
- [KA81] C. Kalbach, Phys. Rev. C 23 (1981) 124; and Phys. Rev. C 24 (1981) 819
- [KA81a] C. Kalbach and F. M. Mann, Phys. Rev. C 23 (1981) 112; and C. Kalbach, Phys. Rev. C 25 (1982) 3197

- [KA83] C. Kalbach, in BNL-NCS-51694 (proceedings of the IAEA advisory working group meeting on basic and applied problems of nuclear level densities, edited by M. R. Bhat) June 1983, p 113
- [KA84] C. Kalbach, "Isospin Dependence of Two-Component Particle-Hole State Densities for Nuclei," Phys. Rev. C 30, p 1310 (1984).
- [KA84a] C. Kalbach, "PRECO-E: Program for Calculating Freeequilibrium and Direct Reaction Double Differential Cross Sections Using the Two-Component Exciton Model" (to be published as a Los Alamos informal document).
- [MA63] G. S. Mani, M. A. Melkanoff and I. Iori, Centre d'Etudes Nucléaires de Saclay report CEA 2380 (1963)
- [MI74] L. Milazzo-Colli *et al.*, Nucl. Phys. A218 (1974) 274; Nuov. Cim. 30A (1975) 632; Nuov. Cim. 39A (1977) 171
- [NA79] K. H. Narasimha Murthy, A. Chatterjee and S. K. Gupta, Proc. Int'l Conf. on Nucl. Cross Sections for Technology, Knoxville, 1979, NBS Spec. Pub. 594, p 793
- [WA77] A. H. Wapstra and K. Bos, Nucl. Data Tables 19 (1977) 215

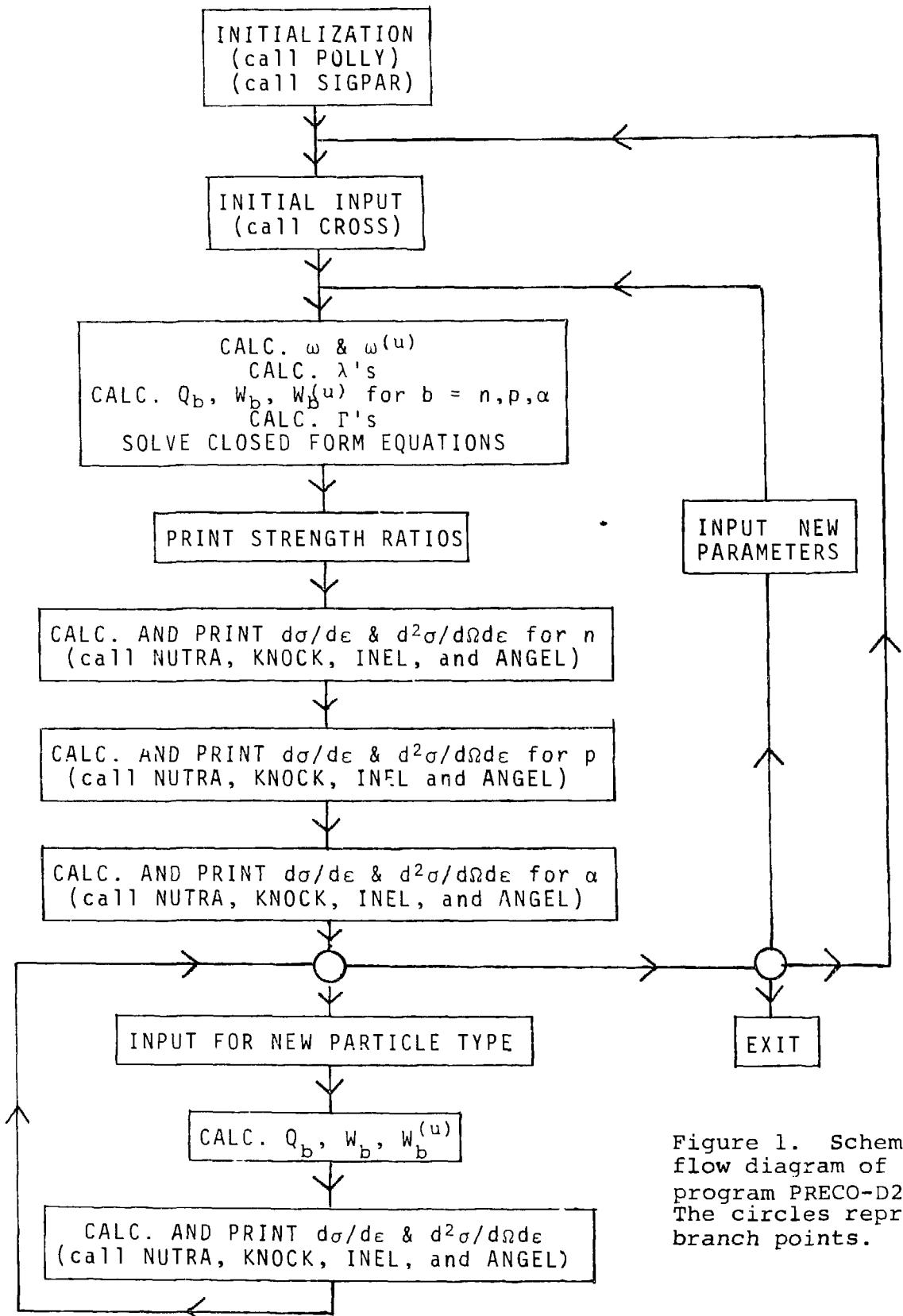


Figure 1. Schematic flow diagram of the program PRECO-D2. The circles represent branch points.

TABLE 1. LIST OF VARIABLES IN PRECO-D2

ACOM	Mass number of composite nucleus
AP	= JOUT, mass number of emitted particle
ASP,ASPR,AFISH	Level spacing parameter, a, for composite and residual nuclei and for top of fission barrier
BEA	Binding energy of projectile in composite nucleus (MeV)
BEN	Binding energy for extra emitted particle types (or temporarily of n, p or α) (MeV)
BF	Height of fission barrier (MeV)
BNEU,BPRO,BALFA	Binding energies of neutrons, protons and alpha particles in the composite nucleus
CLOSD(NP)	Time-integrated strength, $S_d(p) T_u(p)$, responsible for MSD preequilibrium cross section
CLOSU(NP)	Time-integrated strength, $S_u(p) T_u(p)$, responsible for MSD+MSC preequilibrium cross section
COUL(J)	Coulomb barriers for particles of type J, with $J=1-5$ corresponding to n, p, α , extra particle type and projectile, respectively; used in calculating direct reaction cross sections
DBL(NP)	m(p), fraction of multiple unbound states
E	Excitation energy of composite nucleus (MeV)
EC	Coulomb barrier for use in calculating total-reaction cross sections (MeV)
EMU(J)	Maximum possible emission energy for particles of type J with $J=1-3$ for n, p and α
EPS(JHALF,NE)	Particle emission energies (MeV)
ES(I)	$\Theta(E-iS)/(E-S)$ for $i=1,10$
F2	Matrix element scale factor, k

FISHW	Equilibrium cross section going into fission
FLOW,SPILL,POUR	10^{-18} , 10^{+18} , and 10^{+35} ; comparison numbers for underflow and overflow checks
FRAC(NP)	Fraction of the available strength in states with $p=NP$ which is lost due to n , p and α emission in one unit of time, $\Delta t=TAUWM$
FWD(NH)	Correction factor, $f(p)$, applied to $\omega(p,h,E)$ for $h \leq 2$ to account for the finite potential well depth
FWDHI(NP)	$f_+(p)$ and (for $h > 2$) $f_0(p)$
FWDU(NP)	$f_1(p)$
G,GNEU,GPRO, GALFA,GR,GFIS	Single-particle-state densities in the composite nucleus; in the n , p , α and extra particle residual nuclei; and at the top of the fission barrier
GF,GFU,GU,GFAV	$g(p)$, $g_u(p)$, $g_p(p)$ and $g_a(p)$ for $h \leq 2$
GFPU,GFHU	$g_p(p)$ for $h > 2$ and g_h for all h
GAMBB,GAMBU GAMUB,GAMUU	$\Gamma_+^{(bb)}(p)$, $\Gamma_+^{(bu)}(p)$, $\Gamma_+^{(ub)}(p)$, and $\Gamma_+^{(uu)}(p)$
GAMZBU,GAMZUB	$\Gamma_0^{(bu)}(p)$ and $\Gamma_0^{(ub)}(p)$
GEL(NP)	$g[E-A(p,h)]$
GELU(NP)	$g[E-A_{1,0}(p,h)-S]$ or 0, whichever is greater
GELUU(NP)	$g[E-A_{2,0}(p,h)-2S]$ or 0, whichever is greater
IANG	Angular distribution parameter. Energy parameter is $\epsilon + (IANG)(B_b)$
IG	IG=0 invokes $Q_b(p)$; IG=1 invokes $Q_b^{(G)}(p)$
IREA,IWRI	Input and output device numbers
JBAR	Most probable particle number at equilibrium
JHALF	JHALF=1 for primary calculations; JHALF=2 for secondary calculations (for additional emitted particle types)
JHI	minimum of 20 and JBAR; maximum p considered in closed-form reaction equations

JLO	Minimum p which can supply the protons and neutrons for the emitted particle
JPIN,JNIN,JIN	Z, N and A of the projectile
JRZ,JRN	Z and N of target
JPOUT,JNOUT,JOUT	Z, N, and A of emitted particle
NE	Index for emission channel energies
NEPS(1),NEPS(2)	Number of emitted particle energies for primary and secondary calculations
NEWP	Control variable for selecting new values of the model parameters or recycling to a new problem
NLOE	Particle number of simplest states which can undergo nucleon emission
NAP1	Particle number of simplest states which can undergo emission of a given type of complex particle
NP	Index for particle number, p, of states
NPART,NHOLE	Initial particle and hole numbers, p_o and h_o
NPHD,PHD	$p-h=p_o-h_o$
NDWN,NHDWN	Lower limit on particle and hole numbers
NUP	Maximum particle number accessible to system
NSD,SD	Spin degeneracy of emitted particle
PRET2	Full preequilibrium cross section for a given reaction and emission energy
PU,PB,PMSC	$P_u(p)$, $P_b(p)$ and $P_d(p)$; strengths arriving at states of given p by pair creation for the first time
PRMSC,PRMSD	MSC and MSD preequilibrium cross sections
PROB(J,NP)	$Q_b^{(G)}(p)$; J=1,2,3,4 indicates b=n,p, α and extra particle type, respectively
PTOT	Fraction of strength left at end of preequilibrium calculations
P0,P1,P2,P	Parameters P_0 , P_1 , P_2 , and p for calculating entrance channel and inverse cross sections

RHO(NP)	$\omega(p, h, E)$
RHOU(NP)	$\omega^{(u)}(p, h, E)$
RA	Radius of incident or emitted particle (used in calculating total-reaction cross sections)
RNN,RZZ	N and Z of target (same as JRN and JRZ)
RN,RZ	N/A and Z/A of target
S	Average effective separation energy
SIGCN	Entrance channel nonelastic cross section
SIGIN(J,NE)	Inverse nonelastic cross sections; J=1-4 has the same significance as for PROB
SIGBAR(J)	Average nonelastic cross section for channel J; J=1-5 has same significance as for COUL
SIZE	$g_o \cdot E$
SNOCK	Cross section for knockout or inelastic scattering involving cluster degrees of freedom
SNUTRA	Cross section for direct nucleon transfer
STRB	Strength in bound states, = S_b
STRU	Strength in unbound states, = S_u
SUMKE,SUMNE	Energy integrals of the cross sections SNOCK and SNUTRA
SUMSC,SUMSD	Energy integrals of the cross sections PRMSC and PRMSD
SUMW	Energy integral of the cross section WEISS
TAUWM	Time unit, Δt , for evaluating transition rates, = $[20 \lambda_+(p_{\min}, h_{\min})]^{-1}$
TAUWU(NP)	Lifetime of an unbound state against all two-body interactions considered and n, p and α emission; = T_u
THIBB,THIBU, THIUB,THIUU	Pair-creation rates $\lambda_+^{(bb)}$, $\lambda_+^{(bu)}$, $\lambda_+^{(uh)}$, and $\lambda_+^{(uu)}$

TLOBU,TLOUB	Pair-annihilation rates $\lambda_-^{(bu)}$ and $\lambda_-^{(ub)}$
TNOBU,TNOUB	Exciton-scattering rates $\lambda_0^{(bu)}$ and $\lambda_0^{(ub)}$
TIMHI	$\lambda_+^{(p_{\min}, h_{\min})}$
TOMSC(NE), TOMSD(NE)	Total MSC and MSD cross sections for a given emission energy
U	Residual nucleus excitation energy
WU(J,NP,NE)	$W_b^{(u)}(p, h, \epsilon)$ with J=1-4 having the same significance as for PROB
WEISS(J,NE)	Weisskopf-Ewing evaporation cross section; J=1,4 as for PROB
XL0,XL1	Parameters λ_0 and λ_1 for calculating entrance-channel and inverse-reaction cross sections
XLAMB	Parameter λ for calculating entrance-channel and inverse-reaction cross sections
XM0,XM1,XMU	Parameters μ_0 , μ_1 and μ for calculating entrance-channel and inverse-reaction cross sections
XN0,XN1,XN2,XNU	Parameters v_0 , v_1 , v_2 , and v for calculating entrance-channel and inverse-reaction cross sections

TABLE 2. ADDITIONAL VARIABLES IN SUBROUTINES

ANGEL

AL,BL	A_ℓ and B_ℓ for calculating Legendre polynomial coefficients
A0MSC,A0MAD	TOMSC and TOMSD from MAIN, each divided by 4π ; = a_0 (MSC) and a_0 (MSD)
EPSCM	EPS (JHALF,NE)
POL(J,L+1)	Legendre polynomial of order L for an angle of $10(J-i)$ degrees; (generated in subroutine POLLY)
SIGMA	Double differential cross section

CROSS

(none)

INEL

EMAX,UMAX	Maximum emission energy and maximum residual nucleus excitation energy
GBN,GBA	Nucleons and alpha particles in the residual nuclei formed by the knock-out of these particles
GII,GIR	Cluster single-particle-state density in the residual nuclei formed from emission of the cluster and from reemission of the projectile

KNOCK

EMAX,UMAX	(as in INEL)
GA,GBR.GBT	Projectile single-particle-state density in the residual nucleus and emitted-particle-state densities in the residual and target nuclei

NUTRA

GN,GP	g_ν and g_π in the residual nucleus
JPICKN,JPICKP	Number of picked-up neutrons and protons
JSTRIN,JSTRIP	Number of stripped neutrons and protons
TRANN,TRANP	Number of transferred neutrons and protons
UMAX	(as in INEL)
VAB	Effective potential, V_a , seen by the projectile in the interaction region
VELO	Projectile velocity in the interaction region

POLLY

ANG	Angle, θ , first in degrees and then in radians
POL(J,L+1)	(as in ANGEL)

SIGPAR

(none)

APPENDIX A

FORTRAN LISTING OF PRECO-D2

Los Alamos Identification No. LP-1654

```

1 C      PRECO-D
2 C      WRITTEN BY C. KALBACH
3 C      CLOSED FORM PREEQUILIBRIUM REACTION CALCULATIONS
4 C      IN GRIFFIN MODEL FORMALISM
5 C      WITH MULTI-STEP DIRECT EXTRACTED
6 C
7 C      SUBROUTINES FOR PHENOMENOLOGICAL DESCRIPTIONS OF
8 C      1. DIRECT KNOCKOUT AND INELASTIC SCATTERING INVOLVING COMPLEX
9 C          PARTICLE CLUSTERS
10 C     2. DIRECT NUCLEON TRANSFER
11 C     3. ANGULAR DISTRIBUTIONS
12 C     4. OPTION FOR APPROXIMATE REACTION CROSS SECTIONS INTERNALLY
13 C          DIRECT REACTION FORMALISM
14 C          VERIFIED FOR NUCLEON AND ALPHA INCIDENT
15 C          USE WITH CAUTION FOR D OR 3HE INCIDENT
16 C          (NO DIRECT RXN CALC FOR HEAVIER PROJECTILES)
17 C          (NO APPROX RXN CROSS SECTIONS FOR AOUT.GT.4)
18 C
19 C      PRAISE THE LORD
20 C
21      COMMON ESUM(4),RHO(31),RHOU(31),NEPS(2),ES(10),DELTA(2)
22      1,PRDB(4,30),TAUWU(31),RATIO(31,2),FRAC(31),WEISS(4,51),WU(4,31,51)
23      2,CLDSD(31),CLOSU(31),EMU(3),FWD(3),FWDU(32),DBL(31),FWDHI(32)
24      3,GEL(32),GELU(32),GELUU(32),GF(3),GU(3),GFL(3),GFAV(3),EHOLE(3)
25      4,GAMUU(31),GAMUB(31),GAMBU(31),GAMBB(31),GAMZUB(31),GAMZBU(31)
26      COMMON /ANGELS/ POL(19,7),TOMSD(51),TOMSC(51),IWRI,IANG
27      COMMON /ENERGY/ BEN,NEPS1,EPS(2,52),JHALF,JNOUT,JPOUT
28      COMMON /DIRECT/ SIGIN(4,51),JPIN,JNIN,BEA,NSD,KP,E,ACOM,G,R2Z
29      COMMON /NUTRAN/ SNUTRA(51)
30      COMMON /CLUSTR/ SNOCK(51),SIGBAR(5),SIGCN,COUL(5),BNEU,BPRO,BALFA
31      COMMON /PAR/ PO(3,3),P1(3,3),P2(3,3),XLO(3,3),XL1(3,3),XMO(3,3)
32      1,XM1(3,3),XNO(3,3),XN1(3,3),XN2(3,3)
33 C
34      CALL LINK("UNIT5=PRECOINP,UNIT6=(PRECOUT,CREATE,HC) // ")
35 C
36      IREA=5
37      IWRI=6
38      READ(IREA,5)IANG,IG
39      READ(IREA,3)EHOLE(2),EHOLE(3)
40      EHOLE(1)=EHOLE(3)
41      DO 2 NP=1,31
42      GEL(NP)=0.
43      GELU(NP)=0.
44      GELUU(NP)=0.
45      RHO(NP)=0.
46      RHOU(NP)=0.
47      GAMUU(NP)=0.
48      GAMUB(NP)=0.
49      GAMBU(NP)=0.
50      GAMBB(NP)=0.
51      GAMZUB(NP)=0.
52      GAMZBU(NP)=0.
53      FRAC(NP)=0.
54      CONTINUE
55      GEL(32)=0.
56      GELU(32)=0.
57      GELUU(32)=0.
58      DC 4 KP=1,4
59      DO 4 NP=1,50
60      4 PRDB(KP,NP)=0
61      DO 8 KP=1,4
62      DO 8 NE=1,51
63      SIGIN(KP,NE)=0.
64      8 WEISS(KP,NE)=0.
65      DO 11 I=1,19
66      DO 11 J=1,7
67      11 POL(I,J)=0.
68 C
69 C      GENERATE TABLE OF LEGENDRE POLYNOMIALS
70      CALL POLLY

```

```

71 C      GENERATE PARAMETERS FOR APPROX SIGMA INVERSE
72      CALL SIGPAR
73 C      INPUT
74 C
75 C
76      READ(IREA,3)E,BEA
77      FORMAT(5F10.2)
78      READ(IREA,3)RZZ,RNN
79      RN=RNN/(RNN+RZZ)
80      RZ=1.-RN
81      FORMAT(3I3)
82      READ(IREA,3)BNEU,BPRO,BALFA,BF
83      READ(IREA,3)PAIRN,PAIRP,PAIRA,PAIRC
84      READ(IREA,3)EPAIRN,EPAIRP,EPAIRA,EPAIRC
85      READ(IREA,3)SIGCN,SIGBAR(5)
86      READ(IREA,13)NEPS(1),EPS(1,2),DELTA(1)
87      FORMAT(1I3,2F10.2)
88      IF(NEPS(1).LE.50)GO TO 12
89      WRITE(IWR1,101)
90      101 FORMAT(1BH-TOO MANY ENERGIES)
91      GO TO 1000
92      12 IF(NEPS(1).LE.0)NEPS(1)=1
93      NEPS1=NEPS(1)+1
94      IF(EPS(1,2).GT.0.)GO TO 9
95      DELTA(1)=-1
96      DO 7 NE=2,NEPS1
97      READ(IREA,3)EPS(1,NE),SIGIN(1,NE),SIGIN(2,NE),SIGIN(3,NE)
98      CONTINUE
99      READ(IREA,5)JPIN,JNIN
100     JIN=JPIN-JNIN
101     XIN=JIN
102     ACOM=RNN+RZZ+XIN
103     IF(DELTA(1).LT.0.)GO TO 17
104     IF(DELTA(1).LE.0.)DELTA(1)=1.
105     DO 15 NE=3,NEPS1
106     DEL=NE-2
107     15 EPS(1,NE)=EPS(1,2)+DEL+DELTA(1)
108     JHALF=1
109     JNOUT=1
110     JPOUT=0
111     KP=1
112     CALL CROSS
113     JNOUT=0
114     JPOUT=1
115     KP=2
116     CALL CROSS
117     JNOUT=2
118     JPOUT=2
119     KP=3
120     CALL CROSS
121     17 EPS(1,1)=EPS(1,2)-1.
122     IF(EPS(1,1).LT.0.)EPS(1,1)=0.
123     EPS(1,NEPS1+1)=EPS(1,NEPS1)+1.
124     ATHRD=(ACOM-1.)*-0.33
125     COUL(1)=0.
126     XPIN=JPIN
127     COUL(2)=0.75*(RZZ+XPIN-1.)/ATHRD
128     ATHRD=(ACOM-4.)*-0.33
129     COUL(3)=1.50*(RZZ+XPIN-2.)/ATHRD
130     ATHRD=(RZZ+RNN)*-0.33
131     COUL(5)=0.75*XPIN-RZZ/ATHRD
132     DO 14 KP=1,3
133     ESUM(KP)=0.
134     14 SIGBAR(KP)=0.
135     EMU(1)=E-BNEU
136     EMU(2)=E-BPRO
137     EMU(3)=E-BALFA
138     DO 18 KP=1,3
139     DO 16 NE=2,NEPS1
140     IF(EPS(1,NE).LT.COUL(KP))GO TO 16
141     IF(EPS(1,NE).GT.EMU(KP))GO TO 18
142     EINT=(EPS(1,NE+1)-EPS(1,NE-1))/2.
143     ESUM(KP)=ESUM(KP)+EINT
144     SIGBAR(KP)=SIGBAR(KP)+SIGIN(KP,NE)-EINT
145     CONTINUE
146     18 IF(ESUM(KP).GT.0.)SIGBAR(KP)=SIGBAR(KP)/ESUM(KP)
147     IF(SIGBAR(5).GT.0.)GO TO 27
148     IF(JIN.NE.1)GO TO 21
149     SIGBAR(5)=SIGBAR(JPIN+1)

```

```

150      GO TO 27
151 21   IF(JIN.NE.4)GO TO 22
152      SIGBAR(5)=SIGBAR(3)
153      GO TO 27
154 22   JPOUT=JPIN
155      JNOUT=JNIN
156      JHALF=2
157      KP=4
158      NEPS1=(E-BEA)*ACOM/(2.-(ACOM-XIN))
159      DO 23 NE=2,NEPS1
160 23   EPS(2,NE)=2*(NE-1)
161      CALL CROSS
162      EPS(2,1)=1.
163      EPS(2,NEPS1+1)=EPS(2,NEPS1)+1
164      ESUM(4)=0.
165      DO 25 NE=2,NEPS1
166      IF(EPS(2,NE).LT.COUL(5))GO TO 25
167      EINT=(EPS(2,NE+1)-EPS(2,NE-1))/2.
168      ESUM(4)=ESUM(4)+EINT
169      SIGBAR(5)=SIGBAR(5)+SIGNIN(4,NE)*EINT
170 25   SIGNIN(4,NE)=0.
171      IF(ESUM(4).GT.0.)SIGBAR(5)=SIGBAR(5)/ESUM(4)
172      NEPS1=NEPS1+1
173      JHALF=1
174 27   IF(SIGCN.GT.0.)GO TO 38
175      IF(JPIN.GT.0.)GO TO 29
176      EC=2.4
177      XLAMB=XLO(1,2)/ATHRD+XL1(1,2)
178      XMU=XMO(1,2)*ATHRD+XM1(1,2)*ATHRD-ATHRD
179      XNU=XNO(1,2)*ATHRD+(RZZ+RNN)+XN1(1,2)*ATHRD-ATHRD+XN2(1,2)
180      GO TO 37
181 29   I=JPIN+1
182      J=JNIN+1
183      XLAMB=XLO(I,J)*(RZZ+RNN)+XL1(I,J)
184      XNU=(RZZ+RNN)*XM1(I,J)
185      XMU=XMO(I,J)*XNU
186      RA=1.20
187      IF(JIN.EQ.1)RA=0.
188      EC=1.44*XPIN*RZZ/(1.5*ATHRD+RA)
189      XNU=XNU+(XNO(I,J)+XN1(I,J)*EC+XN2(I,J)*EC-EC)
190      IF(JIN.EQ.2)RA=0.8
191      IF(JIN.EQ.3)RA=0.8
192      XNL=XNU/XLAMB
193      IF(XNL.GT.SPILL)XNL=0.
194      IF(XNL.LT.FLOW)GO TO 37
195      ETEST=1.2-SQRT(XNL)
196 37   ELAB=(E-BEA)*ACOM/(ACOM-XIN)
197      SIGCN=XLAMB-ELAB+XMU+XNU/ELAB
198      IF(XNL.LT.FLOW)GO TO 40
199      IF(ELAB.LT.ETEST)GO TO 40
200      GEOM=SORT(XIN*(E-BEA))
201      GEOM=1.23*ATHRD+RA+4.573/GEOM
202      GEOM=31.416*GEOM*GEOM
203      IF(SIGCN.LT.GEOM)SIGCN=GEOM
204      GO TO 38
205 40   IF(ELAB.GE.EC)GO TO 38
206      XP=PO(I,J)+P1(I,J)/EC+P2(I,J)/(EC-EC)
207      XA=-2.*XP*EC+XLAMB-XNU/(EC-EC)
208      XB=XP*EC-EC+XMU+2.*XNU/EC
209      ECUT=0.
210      CUT=XA-XA-4.*XP*XB
211      IF(CUT.GT.0.)ECUT=SQRT(CUT)
212      ECUT=(ECUT-XA)/(2.*XP)
213      SIGCN=XP*ELAB-ELAB+XA*ELAB+XB
214      IF(ELAB.LT.ECUT)SIGCN=0
215 38   READ(IREA,5)NPART,NHOLE
216      N=NPART+NHOLE
217      IF(N.GT.0)GO TO 39
218      NPART=JIN+1
219      NHOLE=1
220 39   READ(IREA,3)G,GNEU,GPRO,GALFA,GFIS
221      IF(G.LE.0.)G=ACOM/13.
222      READ(IREA,3)F2
223      IF(F2.LE.0.)F2=1.35
224      SIZE=G*(E-EPAIRC)
225      E=E-PAIRC
226      NPHD=NPART-NHOLE
227      PHD=NPHD

```

```

228      FLOW=1.E-18
229      SPILL=1.E+18
230      V=38.
231 C      EFFECTIVE G VALUES
232 C
233 C      DO 19 I=1,3
234      GF(I)=0
235      Y=I-1
236      X=NPART-NHOLE+I-1
237      IF(X.LT.0.)GO TO 19
238      PHNO=X+Y
239      IF(PHNO.LE.0.)GO TO 19
240      ENEX=E/(PHNO+38.)
241      GFP=1.+ENEX
242      GFP=SORT(GFP)
243      GFH=1.-ENEX
244      IF(GFH.LT.0.)GFH=0.
245      GFH=SORT(GFH)
246      GF(I)=G*(X*GFP+Y*GFH)/PHNO
247      CONTINUE
248 19      COULP=BPRO+Coul(2)
249      COULA=BALFA+Coul(3)
250      S=AMIN1(BNEU,COULP,COULA)
251      IF(S.LT.0.)S=0.
252      GFPU=(38.+S)/38.
253      GFPU=SORT(GFPU)*G
254      GFHU=(38.-S)/38.
255      IF(GFHU.LT.0.)GFHU=G
256      GFHU=SORT(GFHU)*G
257      DO 24 I=1,3
258      XX=1
259      GU(I)=G
260      GFU(I)=G
261      GFAV(I)=G
262      Y=I-1
263      X=NPART-NHOLE+I-1
264      IF(X.LT.0.)GO TO 24
265      PHNO=X+Y
266      IF(PHNO.LE.0.)GO TO 24
267      PHNO1=PHNO-1
268      ENEX=(E-S)/(PHNO+38.)
269      GFP=SORT(1.+ENEX)
270      GFH=1.-ENEX
271      IF(GFH.LT.0.)GFH=0
272      GFH=SORT(GFH)
273      Z=1.+ENEX+S/38.
274      GU(I)=G-SORT(Z)
275      IF(PHNO1.LE.0.)GO TO 20
276      GFU(I)=G*((X-1.)*GFP+Y*GFH)/PHNO1
277      IF(PHNO.GT.2.)GFAV(I)=((PHNO-3.)*GFU(I)+GU(I)).(PHNO-2.)
278 20      C
279 C      FWD CORRECTION FOR FULL STATE DENSITY
280 C
281 C      V=EHOLE(I)
282      IF(E.LE.V)GO TO 24
283      IF(I.EQ.1)GO TO 24
284      X=E--PHNO1
285      EV=(E-V)**PHNO1
286      XX=1.-Y=EV/X
287      EV2=E-V-V
288      IF(EV2.LE.0.)GO TO 24
289      EV2=EV2**PHNO1
290      XX=XX+Y=(Y-1.)*0.5*EV2/X
291      IF(XX.LT.0.)XX=0
292      FWD(I)=XX
293 24      E=E+PAIRC-EPAIRC
294      C
295 C      UNBOUND STATE CORRECTION FUNCTION
296 C      AND MULTIPLE UNBOUND CORRECTION. M(P)
297 C
298 C      X=E-S
299      DO 26 I=1,10
300      ES(I)=0.
301 26      DO 28 I=1,10
302      IF(X.LE.0.)GO TO 30
303      XI=I
304      Y=(E-XI*S)/X
305      IF(Y.LE.0.)GO TO 30
306

```

```

307      ES(I)=Y
308  28  CONTINUE
309      E=E+PAIRC
310  30  NTERM=I-1
311      DO 36 NP=1,30
312      NP1=NP+1
313      FT2=O.
314      FWDU(NP1)=O.
315      DBL(NP1)=O.
316      FWDHI(NP1)=O.
317      P=NP
318      NH=NP-NPHD
319      H=NH
320      IF(H.LT.O.)GO TO 36
321      PH1=P+H-1.
322      FT=ES(1)
323      IF(PH1.LE.O.)GO TO 35
324      FT=O.
325      LIM=MINO(NP,NTERM)
326      BIN=1./P
327      SI=-1.
328      IF(NH.GT.2)GO TO 32
329      EE=E-PAIRC
330      IF(EE.LE.S)GO TO 36
331      Y=EE-S
332      DO 31 I=1,LIM
333      XI=I
334      BIN=BIN=(P-XI+1.)/XI
335      SI=-SI
336      X=(EE-XI*S)/Y
337      IF(X.LT.O.)X=O.
338      IF(X.GT.O.)X=X**PH1
339      IF(X.LT.FLOW)X=O.
340      IF(X.GT.SPILL)X=O.
341      XX=SI*BIN*X
342      X=(EE-XI*S-EHOLE(NH+1))/Y
343      IF(X.LT.O.)X=O.
344      IF(X.GT.O.)X=X**PH1
345      IF(X.LT.FLOW)X=O.
346      IF(X.GT.SPILL)X=O.
347      XX=XX-SI*BIN=H-X
348      X=(EE-XI*S-EHOLE(NH+1)*2.)/Y
349      IF(X.LT.O.)X=O.
350      IF(X.GT.O.)X=X**PH1
351      IF(X.LT.FLOW)X=O.
352      IF(X.GT.SPILL)X=O.
353      XX=XX+SI*BIN=H-(H-1.)*O.5*X
354      IF(I.GT.1)FT2=FT2-XX
355  31  FT=FT+XX
356      GO TO 35
357  32  DO 34 I=1,LIM
358      XI=I
359      BIN=BIN=(P-XI+1.)/XI
360      SI=-SI
361      X=ES(I)**PH1
362      IF(X.LT.FLOW)X=O.
363      IF(X.GT.SPILL)X=O.
364      XX=SI*BIN*X
365      IF(I.GT.1)FT2=FT2-XX
366  34  FT=FT+XX
367  35  FWDU(NP1)=FT
368      IF(FT.GT.O.)FT2=FT2/FT
369      DBL(NP1)=FT2
370      X=1.
371      IF(ES(2).GT.O.)X=1.-O.5*ES(2)**PH1
372      IF(NH.GT.2)GO TO 33
373      EE=E-PAIRC-S
374      SV=S+EHOLE(NH+1)
375      X=1.
376      IF(EE.LE.S)GO TO 33
377      X=1.-O.5*((EE-S)/EE)**PH1
378      IF(EE.LE.SV)GO TO 33
379      X=X-((EE-SV)/EE)**PH1
380      SV=SV+S
381      IF(EE.LE.SV)GO TO 33
382      X=X+O.5*((EE-SV)/EE)**PH1
383  33  FWDHI(NP1)=X
384  36  CONTINUE

```

```

385      FWDU(1)=1.
386      DBL(1)=0.
387      FWDHI(1)=1.
388      FWDU(32)=1.
389      FWDHI(32)=1.
390 C      CALCULATE G*E FACTORS
391 C
392 C      POUR=1.E+35
393      IF(NPHD)54,55,56
394      NDWN=0
395 54      GO TO 58
396      NDWN=1
397 55      GO TO 58
398      NDWN=NPHD
399 56      NHDWN=NDWN-NPHD
400 58
401      NDWN1=NDWN+1
402      DO 60 NP1=NDWN1,32
403      NP=NP1-1
404      P=NP
405      NH=NP-NPHD
406      H=NH
407      Q=AMAX 1(P,H)
408      X=SIZE-Q-Q
409      IF(X.LT.0.)GO TO 62
410      IF(NH.LE.2)x=GF(NH+1)*(E-PAIRC)-Q*Q
411      GEL(NP1)=Y+(P*(P+1.)+H*(H+1.))/4.
412      IF(NP.GT.31)GO TO 60
413      GELU(NP1)=0.
414      GELUU(NP1)=0.
415      IF(P.LE.0.)GO TO 60
416      IF(E.LE.S)GO TO 60
417      P=P-1.
418      Q2=Q*Q-Q+0.5
419      X=SIZE-Q2-G*S
420      IF(NH.LE.2)x=G-(E-PAIRC)-Q2-G*S
421      IF(X.LT.0.)GO TO 60
422      GELU(NP1)=X+(P*(P+1.)+H*(H+1.))/4.
423      IF(P LE.0.)GO TO 60
424      P=P-1.
425      Q2=Q*Q-2.*Q+2.
426      X=SIZE-Q2-2.*G*S
427      IF(NH.LE.2)x=G-(E-PAIRC)-Q2-2.*G*S
428      IF(X.LT.0.)GO TO 60
429      GELUU(NP1)=X+(P*(P+1.)+H*(H+1.))/4.
430 60      CONTINUE
431 62      NUP=NP-1
432 C      CALCULATE STATE DENSITIES
433 C
434 C      FAC=0.
435      PH=NDWN+NHDWN
436      IF(NDWN.EQ.0)GO TO 61
437      DO 64 NP=1,NDWN
438      Y=NP
439
440 64      FAC=FAC+ ALOG10(Y)
441 61      IF(NHDWN.EQ.0)GO TO 63
442      DO 65 NP=1,NHDWN
443      Y=NP
444 65      FAC=FAC+ ALOG10(Y)
445 63      LIM=NDWN+NHDWN-1
446      IF(LIM.EQ.0)GO TO 67
447      DO 66 NP=1,LIM
448      Y=NP
449 66      FAC=FAC+ ALOG10(Y)
450 67      RHOL=GEL(NDWN1)
451      RHOL=(PH-1.)*ALOG10(RHOL)
452      RHOL=RHOL-FAC
453      RHO(NDWN1)=(10.-RHOL)*G
454      IF(NDWN.GT.0)GO TO 51
455      RHOL=0.
456
457 51      IF(PH.GT.1.)GO TO 57
458      RHOL=0.
459      GO TO 53
460 57      RHOL=GELU(NDWN1)
461      IF(NHDWN.LE.2)RHOL=RHOL*GFU(NHDWN+1)/G
462      IF(RHOL.EQ.0.)GO TO 59

```

```

463      RHOL=( PH- 1. )*ALOG10(RHOL)
464      RHOL=RHOL-FAC
465 53      RHOL=( 10. **RHOL )*GFP
466 59      P=NDWN
467      RHOU(NDWN1)=RHOL*FWOU(NDWN1)*P
468      DO 68 NP=NDWN1,NUP
469      NP1=NP+1
470      P=NP
471      NH=NP-NPHD
472      H=NH
473      FAC=FAC+ALOG10(P)+ALOG10(H)+ALOG10(P+H-1. )+ALOG10(P+H-2. )
474      RHOL=GEL(NP1)
475      RHOL=( P+H-1. )*ALOG10(RHOL)
476      RHOL=RHOL-FAC
477      RHO(NP1)=( 10. **RHOL )*G
478      IF(RHO(NP1).LT. 1. )GO TO 69
479      IF(NP.GT.30)GO TO 68
480      RHOL=GELU(NP1)
481      IF(NH.LE.2)RHOL=RHOL*GFU(NH+1)/G
482      IF(RHOL.EQ.0. )GO TO 74
483      RHOL=( P+H-1. )*ALOG10(RHOL)
484      RHOL=RHOL-FAC
485      RHOL=( 10. **RHOL )*GFP
486 74      RHOU(NP1)=RHOL*FWDU(NP1)*P
487 68      CONTINUE
488      GO TO 70
489 69      NUP=NP-1
490 70      DO 72 NP=NDWN1,31
491      IF(RHO(NP+1).LT.RHO(NP))GO TO 73
492 72      CONTINUE
493 73      JBAR=NP-1
494      JHI=MINO(20,JBAR)
495      JBAR1=JBAR+1
496      DO 81 I=1,3
497      NP1=I+NPHD
498      RHO(NP1)=RHO(NP1)*FWD(I)=GF(I)/G
499 81      RHOU(NP1)=RHOU(NP1)*GU(I)/GFP
500      EE=E*1.35/F2
501 C
502 C      CALCULATE TRANSITION RATES
503 C
504      DO 79 NP1=NDWN1,JBAR1
505      NP=NP-1
506      P=NP
507      NH=NP-NPHD
508      H=NH
509      PH=P+H
510      V=38
511      IF(NH.LE.2)V=EHOLE(NH+1)
512 C
513 C      PAIR ANNIHILATION
514      TLOUB=0.
515      TLOBU=0.
516      IF(NP.EQ.NDWN)GO TO 76
517      TLOUB=GFHU*H*(H-1.)*(1.-DBL(NP1))*CMAT/2.
518      TLOBU=GFP*P*(P-1.)*H*CMAT/2.
519      IF(NH.LE.3)TLOBU=TLOBU*GU(NH)/GFP
520      IF(PH.LT.4.)GO TO 71
521      TLOBU=TLOBU/2.
522      X=GEL(NP1)/G
523      IF(NH.LE.2)X=X*G/GF(NH+1)
524      XU1=GELU(NP1)/G
525      XU=GELU(NP1)/G
526      XUU1=GELUU(NP1)/G
527      XUU=GELUU(NP1)/G
528      XX=(PH-2.)*(PH-3.)*YU1-XU1-2.*((PH-1.)*(PH-3.)*X*XU1
529      XX=XX+(PH-1.)*(PH-2.)*X*X
530      IF(NH.LE.2)XY=XX+FWD(NH+1)
531      XX=XX*XU1*((PH-3.))
532      Y=X*((PH-1.))
533      IF(NH.LE.2)Y=Y*FWD(NH+1)
534      Y=Y-P*FWDU(NP1)*XU*((PH-1.))
535      X=(PH-2.)*(PH-3.)*XUU1-XUU1-2.*((PH-1.)*(PH-3.)*XU*XUU1
536      X=X+(PH-1.)*(PH-2.)*XU*XU
537      X=X*((P-2.)*FWDU(NP1)*XUU1*((PH-3.))
538      FWDUU=1.
539      DEN=0.
540      ES2=ES(2)

```

```

541      IF(NH.GT.2)GO TO 84
542      ES2=E-PAIRC-2.*S
543      IF(ES2.LE.0.)ES2=0.
544      ES2=ES2/(ES2+S)
545 84      IF(ES2.GT.0.)DEN=(P-1.)*ES2**(PH-1.)
546      IF(DEN.GT.0.)FWDUU=2.*FWDU(NP1)*DBL(NP1)/DEN
547      XT=4.*(FWDU(NP1)*XU**(PH-1.)-(P-2.)*FWDUU*XUU***(PH-1.))
548      X=XX-X-XT
549      IF(X.LT.0.)X=0.
550      TLOBU=TLOBU*X/Y
551 71      IF(RHOU(NP1).GT.RHO(NP1))RHOU(NP1)=RHO(NP1)
552      IF(RHOU(NP1).GE.RHO(NP1))TLOBU=0.
553 C
554 C      PAIR CREATION
555 76      CMAT=PH/EE
556      EPH=EE/PH
557      IF(EPH.LT.7.)CMAT=CMAT*SORT(EPH/7.)
558      IF(EPH.LT.2.)CMAT=CMAT*SORT(EPH/2.)
559      IF(EPH.GT.15.)CMAT=CMAT*SORT(15./EPH)
560      NP2=NP1+1
561      X=GEL(NP1)
562      IF(NH.LE.2)X=X-G/GF(NH+1)
563      XT=X
564      Y=GEL(NP2)
565      IF(NP.GT.NDWN)GO TO 77
566      IF(NH.LE.1)Y=Y-G/GF(NH+2)
567      A=(Y/X)**(PH-1.)
568      TIMHI =X*G-GEL(NP2)-GEL(NP2)*CMAT/(2.+(PH+1.))
569      IF(NH.LE.1)TIMHI =TIMHI *FWD(NH+2)*GF(NH+2)/G
570 77      Y=GELU(NP2)
571      X=0.
572      IF(PH.LE.1.)X=1.
573      IF(GELU(NP1).GT.0.)X=(Y/GELU(NP1))**(PH-1.)
574      FW=1.
575      IF(NH.LE.1)FW=FWD(NH+2)
576      FW=(FWDHI(NP2)+(PH-1.)*FW)/PH
577      THIUU=X*G-Y-FW*CMAT/(2.+PH)
578      IF(NH.LE.1)THIUU=THIUU*GFU(NH+2)-GFU(NH+2)-GFAV(NH+2)/(G+G+G)
579      X=GEL(NP2)
580      XX=GELU(NP1)
581      IF(NH.LE.1)X=X-G/GF(NH+2)
582      Y=PH*(X-XX)*(X-XX)+X-X-XX-XX
583      X=X-XX*0.5*PH -Y
584      IF(NH.LE.1)X=X-GF(NH+2)*GF(NH+2)*GF(NH+2)/(G+G+G)
585      THIB=X*G
586      IF(NH.LE.1)X=X-FWD(NH+2)
587      X=G-X-CMAT/(2.+PH*(PH+1.))
588      XX=THIUU-GFPU-FWDHI(NP2)/(G-FW)
589      IF(NH.LE.1)XX=THIUU*GU(NH+2)-FWDHI(NP2)/(GFAV(NH+2)-FW)
590      X=X-2.*XX/(PH+1.)
591      IF(RHOU(NP2).GE.RHO(NP2))X=0.
592      THIUU=THIUU+X=DBL(NP1)
593      THIUB=X*(1.-DBL(NP1))
594      TNOUNB=0.
595      THIBB=0.
596      THIBU=0.
597      TNODBU=0.
598      IF(RHOU(NP1).GE.RHO(NP1))GO TO 75
599      TEST=0.
600      IF(NP.GT.0)GO TO 85
601      TEST=(E-PAIRC-H-V)/P
602      IF(NH.LE.2)TEST=(E-PAIRC-V)/P
603 85      IF(TEST.GT.5)GO TO 75
604      Y=(XT/G)**(PH-1.)
605      YY=1.
606      YYY=1.
607      IF(PH.LE.1.)GO TO 82
608      YY=(GELU(NP1)/G)**(PH-1.)
609      YYY=(GELUU(NP1)/G)**(PH-1.)
610 82      Z=GEL(NP2)/G
611      I=(NH.LE.1)Z=Z-G/GF(NH+2)
612      Z=Z**(PH+1.)
613      ZZ=(GELU(NP2)/G)**(PH+1.)
614      IF(NH.LE.2)Y=Y*FWD(NH+1)
615      XX1=E-PAIRC-V
616      IF(NH.LE.2)XX1=E-PAIRC-V
617      IF(XX1.LE.0.)GO TO 78
618      XX=PH*(PH-1.)*XX1=XX1-2.* (PH+1.)*(PH-1.)*XX1=GELU(NP2)/G

```

```

619      XX=XX+PH*(PH+1.)=GELU(NP2)*GELU(NP2)/(G*G)
620      XX=XX1**=(PH-1.)=XX
621      XX1=XX1-S
622      IF(XX1.LT.0.)XX1=0.
623      X=PH*(PH-1.)*XX1*XX1-2.* (PH+1.)*(PH-1.)*XX1=GELUU(NP2)/G
624      X=X+PH*(PH+1.)*GELUU(NP2)*GELUU(NP2)/(G*G)
625      IF(XX1.GT.0.)XX=XX-P*XX1**=(PH-1.)*X
626 7B     X=(GELUU(NP2)/G)**=(PH+1.)
627     X=(GELU(NP2)/G)**=(PH+1.)-P*X-0.5*XX
628     X=X-G*G
629     Y=Y-P*YY*FWDU(NP1)
630     IF(Y.LE.0.)GO TO 75
631     THIBU=G*H*X=CMAT/(Y-2.*PH*(PH+1.))
632     IF(NH.LE.1)THIBU=THIBU=GFU(NH+2)*GU(NH+2)/(G*G*G)
633     X=PH-Z*G*G
634     IF(NH.LE.1)X=X-(GF(NH+2)**3.)*FWD(NH+2)/(G*G*G)
635     ZZ=P*(PH-1.)*ZZ=G*G*G=FWDU(NP2)
636     IF(NH.LE.1)ZZ=ZZ=(GFU(NH+2)**3.)/(G*G*G)
637     X=X-ZZ-P*THIBU=YY=FWDU(NP1)
638     X=X-CMAT/(Y-2.*PH*(PH+1.))
639     THIBB=X-THIBU
640 C
641 C      SCATTERING
642     X=GELU(NP1)
643     XT=XT-X
644     TNOUB=XT*G*(PH+H-1.)/2.
645     IF(NH.LE.2)TNOUB=TNOUB+FWD(NH+1)
646     FW=1.-2.*FWDHI(NP1)
647     IF(NH.LE.2)FW=FWD(NH+1)-2.* (1.-0.5*ES(2)**=(PH-1.))
648     TNOUB=TNOUB+(P-1.)*G*GELU(NP1)*FW/PH
649     TNOUB=TNOUB=CMAT*(1.-DBL(NP1))
650     TNOBU=(PH+H-1.)*YY*(2.*GELU(NP1)+PH*XT)
651     IF(NH.LE.2)TNOBU=TNOBU+FWD(NH+1)
652     TNOBU=TNOBU-4.* (PH-1.)*YY=GELU(NP1)*FWDU(NP1)
653     XT=GELU(NP1)-GELUU(NP1)
654     TNOBU=TNOBU-(P-1.)*(PH+H-1.)*YY*(2.*GELUU(NP1)+PH*XT)*FWDU(NP1)
655     FWDUU=1.
656     DEN=0.
657     ES2=ES(2)
658     IF(NH.GT.2)GO TO 86
659     ES2=E-PAIRC-2.*S
660     IF(ES2.LE.0.)ES2=0.
661     ES2=ES2/(ES2+S)
662 86     IF(ES2.GT.0.)DEN=(P-1.)*ES2**=(PH-1.)
663     IF(DEN.GT.0.)FWDUU=2.*FWDU(NP1)*DBL(NP1)/DEN
664     TNDBU=TNDBU+4.* (P-1.)*(PH-1.)*YY=GELUU(NP1)*FWDUU
665     TNOBU=TNOBU*P*GFPU-CMAT/(2.*PH*Y)
666     IF(NH.LE.2)TNOBU=TNOBU*GU(NH+1)*GFU(NH+1)/(G*GFPU)
667 C
668 C      TRANSITION BRANCHING RATIOS (NUMERATORS)
669 75     GAMUU(NP1)=THIUU
670     GAMUB(NP1)=THIUB
671     GAMZUB(NP1)=TNOUB+TLQUB
672     GAMBB(NP1)=THIBB
673     GAMBU(NP1)=THIBU
674     GAMZBU(NP1)=TNOBU+TLOBU
675 79     CONTINUE
676 C
677 C      NORMALIZATION
678     TAUWM=1./(TIMHI*20.)
679     DO 80 NP1=NDWN1,JBAR1
680     GAMUU(NP1)=GAMUU(NP1)-TAUWM
681     IF(GAMUU(NP1).LT.FLOW)GAMUU(NP1)=0.
682     IF(GAMUU(NP1).GT.SPILL)GAMUU(NP1)=0.
683     GAMUB(NP1)=GAMUB(NP1)-TAUWM
684     IF(GAMUB(NP1).LT.FLOW)GAMUB(NP1)=0.
685     IF(GAMUB(NP1).GT.SPILL)GAMUB(NP1)=0.
686     GAMZUB(NP1)=GAMZUB(NP1)-TAUWM
687     IF(GAMZUB(NP1).LT.FLOW)GAMZUB(NP1)=0.
688     IF(GAMZUB(NP1).GT.SPILL)GAMZUB(NP1)=0.
689     GAMBU(NP1)=GAMBU(NP1)-TAUWM
690     IF(GAMBU(NP1).LT.FLOW)GAMBU(NP1)=0.
691     IF(GAMBU(NP1).GT.SPILL)GAMBU(NP1)=0.
692     GAMBB(NP1)=GAMBB(NP1)-TAUWM
693     IF(GAMBB(NP1).LT.FLOW)GAMBB(NP1)=0.
694     IF(GAMBB(NP1).GT.SPILL)GAMBB(NP1)=0.
695     GAMZBU(NP1)=GAMZBU(NP1)-TAUWM
696     IF(GAMZBU(NP1).LT.FLOW)GAMZBU(NP1)=0.
697     IF(GAMZBU(NP1).GT.SPILL)GAMZBU(NP1)=0.

```

```

698 80  CONTINUE
699      ACUBE=ACOM+ACOM+ACOM
700      TWOP1=1.05E-12
701      TAUWM=TAUWM*TWOP1*ACUBE/1.35
702 C
703 C      CALCULATE FISSION RATES
704 C
705      IF(GFIS.EQ.0.)GFIS=G
706      FISHW=0.
707      IF(BF.EQ.0.)GO TO 136
708      E=E-EPAIRC
709      NFISH=E-BF+1.
710      AFISH=1.645*GFIS
711      ASP=1.645*G
712      REF=SQRT(ASP+E)
713      DO 134 K=1,NFISH
714      XK1=E-K-1
715      EX=(E-BF-XK1)*AFISH
716      IF(EX.LE.0.)EX=0.
717      EX=SQRT(EX)
718      EX=2.* (EX-REF)
719      EX=EX-EXP(EX)/(E-BF-XK1)
720      IF(EX.LT.FLOW)EX=0.
721      IF(EX.GT.SPILL)EX=0.
722 134  FISHW=FISHW+EX
723 136  FISHW=FISHW-2.42E+08
724      E=E+EPAIRC
725 C
726 C      CALCULATE PROBABILITY FACTORS AND EMISSION RATES
727 C
728      JNOUT=1
729      JPOUT=0
730      JOUT=1
731      JHALF=1
732      NSD=2
733      BEN=BNEU
734      GR=GNEU
735      EPAIR=EPAIRN
736      PAIR=PAIRN
737      GO TO 550
738 150  IF(JPOUT.EQ.1)GO TO 152
739 151  IF(JPOUT.EQ.2)GO TO 154
740      NLDE=NAP1
741      JNOUT=0
742      JPOUT=1
743      BEN=BPRO
744      GR=GPRO
745      EPAIR=EPAIRP
746      PAIR=PAIRP
747      GO TO 550
748 152  NLDE=MINO(NLDE,NAP1)
749      JNOUT=2
750      JPOUT=2
751      JOUT=4
752      NSD=1
753      BEN=BALFA
754      GR=GALFA
755      EPAIR=EPAIRA
756      PAIR=PAIRA
757      GO TO 550
758 C
759 C      CALCULATE FRACTIONAL STRENGTH LOSSES
760 C
761 154  DO 156 NP1=1,31
762 155  FRAC(NP1)=0.
763 156  DO 158 NP=NLDE,JBAR
764      NP1=NP+1
765 157  DO 158 NE=2,NEPS1
766      WNP=WU(1,NP,NE)+WU(2,NP,NE)+WU(3,NP,NE)
767 158  FRAC(NP1)=FRAC(NP1)+WNP*(EPS(1,NE+1)-EPS(1,NE-1))/2.
768 159  DO 186 NP1=NDWN1,JBAR1
769      YU=GAMUU(NP1)+GAMUB(NP1)+FRAC(NP1)
770      XU=YU+GAMZUB(NP1)
771      YB=GAMBB(NP1)+GAMBU(NP1)
772      XB=YB+GAMZBU(NP1)
773      IF(XU.LE.0.)GO TO 164
774      XU=1./XU

```

```

775      GAMUU(NP1)=GAMUU(NP1)*XU
776      GAMUB(NP1)=GAMUB(NP1)*XU
777      GAMZUB(NP1)=GAMZUB(NP1)*XU
778      GAMZBU(NP1)=GAMZBU(NP1)/(XU+YU)
779      FRAC(NP1)=FRAC(NP1)*XU
780 164      TAUWU(NP1)=XU
781      IF(XB.LE.0)GO TO 186
782      XB=1./XB
783      GAMBB(NP1)=GAMBB(NP1)*XB
784      GAMBU(NP1)=GAMBU(NP1)*XB
785      GAMZBU(NP1)=GAMZBU(NP1)*XB
786      GAMZUB(NP1)=GAMZUB(NP1)/(XB-YB)
787 186      CONTINUE
788      WRITE(IWRI,185)
789 185      FORMAT(1H1)
790 C
791 C      CLOSED FORM CALCULATIONS
792 C
793      L_ED=0.
794      PU=RHOU(NPART+1)/RHO(NPART+1)
795      IF(PU.GT.1.)PU=1.
796      PMSD=PU
797      PB=1.-PU
798      IF(NPART.EQ.NDWN)GO TO 202
799      DO 208 NP1=1,NPART
800      RATIO(NP1,1)=0.
801      RATIO(NP1,2)=0.
802      CLOSU(NP1)=0.
803 208      CLOSD(NP1)=0.
804 202      DO 204 NP=NPART,JB,R
805      NP1=NP+1
806      RATIO(NP1,1)=PU/(PU+PB)
807      RATIO(NP1,2)=PMSD/(PU+PB)
808      STRU=PU+PB*GAMZBU(NP1)
809      STRB=PB+PU*GAMZUB(NP1)
810      CLOSD(NP1)=PMSD*TAUWJ(NP1)
811      CLOSU(NP1)=STRU*TAUWU(NP1)
812      USED=USED+PMSD=FRAC(NP1)
813      PMSD=PMSD*GAMUU(NP1)
814      IF(PMSD.LT.0.001)PMSD=0.
815      IF(PMSD.GT.SPILL)PMSD=0.
816      PU=STRU=GAMUU(NP1)+STRB=GAMBU(NP1)
817      IF(PU.LT.FLOW)PU=0.
818      IF(PU.GT.SPILL)PU=0.
819      PB=STRB*GAMBB(NP1)+STRU=GAMUB(NP1)
820      IF(PB.LT.FLOW)PB=0.
821      IF(PB.GT.SPILL)PB=0
822 204      CONTINUE
823      PTOT=PU+PB
824      TWEIS=0.
825      DO 272 NE=2,NEPS1
826      EINT=(EPS1,NE+1)-EPS(1,NE-1))/2.
827      DO 272 KP=1,3
828 272      TWEIS=TWEIS+WEISS(KP,NE)*EINT
829      TWEIS=TWEIS+FISHW
830      FWEIS=PTOT/TWEIS
831 C
832 C      PRINT OCCUPATION PROBABILITIES
833 C
834      JRZ=RZZ+0.5
835      JRN=RNN+0.5
836      WRITE(IWRI,351)S
837 351      FORMAT(3H-S,1F7.2,4H MEV)
838      WRITE(IWRI,352)NPART,NHOLE
839 352      FORMAT(3OH-OCCUPATION PROBABILITIES PO=.1I2.5H, HO=.1I2)
840      WRITE(IWRI,353)JRZ,JRN,JPIN,JNIN
841 353      FORMAT(11H TARGET Z=.1I3.4H, N=.1I3.12H PROJ Z=.1I3.4H, N=.1I
842 13)
843      WRITE(IWRI,354)G,E
844 354      FORMAT(3H G=,1F6.3,4H, E=,1F6.3)
845      PTOT=1.-PTOT
846      WRITE(IWRI,355)F2,PTOT
847 355      FORMAT(14H SCALE FACTOR=.1FB.3,13H, FRAC PREEQ=.1F6.3)
848      WRITE(IWRI,360)
849 360      FORMAT(41H- P H RHOU/RHC STRU/STR STRD/STR)
850      WRITE(IWRI,361)
851 361      FORMAT(1H )
852      DO 364 NP1=NDWN1,JBAR1

```

```

853      NP=NP1-1
854      NH=NP-NPHD
855      X=RHOU(NP1)/RHO(NP1)
856      Y=RATIO(NP1,1)
857      Z=RATIO(NP1,2)
858      WRITE(IWRI,362)NP,NH,X,Y,Z
859 362  FORMAT(2I3,8(1PE12.3))
860 364  CONTINUE
861 C
862 C      PRINT PARTICLE SPECTRA
863 C
864      JPOUT=0
865      JNOUT=1
866      JOUT=1
867      NSD=2
868      BEN=BNEU
869      KP=1
870 450  NEPS1=NEPS(JHALF)+1
871      IF(JOUT.GT.1)GO TO 451
872      NI=NLOE
873      IF(PROB(KP.NI).EQ.0.)NI=NI+1
874 451  WRITE(IWRI,454)JPOUT,JNOUT
875 454  FORMAT(21H1PARTICLE SPECTRA Z=.1I2.4H, N=.1I2)
876      WRITE(IWRI,456)NI
877 456  FORMAT(21H FIRST EMISSION AT P=.1I2)
878      WRITE(IWRI,353)JRZ,JRN,JPIN,JNIN
879      WRITE(IWRI,458)NPART,NHOLE,G,E
880 458  FORMAT(4H PO=.1I2.5H, HO=.1I2.4H, G=.1F6.3,4H, E=.1F6.3)
881      IF(IG.EQ.0)WRITE(IWRI,457)
882 457  FORMAT(37H PAIR EXCIT. ACCORDING TO Z/A AND N/A)
883      IF(IG.GT.0)WRITE(IWRI,459)
884 459  FORMAT(47H PAIR EXCIT. ACCORDING TO 2 COMP. ST. DENSITIES)
885      WRITE(IWRI,460)SIGCN
886 460  FORMAT(25H REACTION CROSS SECTION =.1F7.1)
887      WRITE(IWRI,355)F2,PTOT
888      X=RHOU(NPART+1)/RHO(NPART+1)
889      WRITE(IWRI,462)X,USED
890 462  FORMAT(14H MSD POSSIBLE=.1F6.3,7H, USED=.1F6.3)
891      NIC=MAX0(NI,NPART)
892      WRITE(IWRI,470)NIC
893 470  FORMAT(29H CLOSED FORM SUM STARTS AT P=.1I2)
894      WRITE(IWRI,474)
895 474  FORMAT(86H- EPS -- DIRECT -- PREEQUILIBRIUM -- EQ
896 1UIL -- TOTAL (MB/MEV))
897      WRITE(IWRI,476)
898 476  FORMAT(6X,87H-- NUTRA KNOCK -- MSD MSC -- WEISS
899 1 -- MSD MSC MSD+MSC)
900      WRITE(IWRI,477)
901 477  FORMAT(1H )
902      SUMNU=0.
903      SUMKE=0.
904      SUMSD=0.
905      SUMSC=0.
906      SUMW=0.
907 C
908 C      DIRECT REACTION CALCULATIONS
909      DO 471 NE=1,51
910      SNUTRA(NE)=0.
911 471  SNOCK(NE)=0.
912      IF(JIN.GT.4) GO TO 473
913      IF(JOUT.EQ.JIN)GO TO 472
914      CALL NUTRA
915      IF(JHALF.EQ.1)CALL KNOCK
916      GO TO 473
917 472  IF(JPOUT.EQ.JPIN)CALL INEL
918      IF(JHALF.EQ.2)CALL NUTRA
919 473  SIGCN=SIGCN+0.95
920      IF(JIN.GT.1)SIGCN=SIGCN+0.89
921      DO 480 NE=2,NEPS1
922      PRMSD=0.
923      PRET2=0.
924      DO 466 NP=NIC,JBAR
925      NP1=NP+1
926      X=WU(KP,NP,NE)
927      PRMSD=PRMSD+X*CLSD(NP1)
928 466  PRET2=PRET2+X*CLDSU(NP1)
929      PRMSD=PRMSD+SIGCN
930      PRET2=PRET2+SIGCN

```

```

931      PRMSC=PRET2-PRMSD
932      WEISS(KP,NE)=WEISS(KP,NE)*SIGCN*FWEIS
933      TOMSD(NE)=PRMSD+SNUTRA(NE)+SNOCK(NE)
934      TOMSC(NE)=PRMSC+WEISS(KP,NE)
935      TOTAL=TOMSD(NE)+TOMSC(NE)
936      IF(TOTAL.LE.0.)GO TO 480
937      EINT=(EPS(JHALF,NE+1)-EPS(JHALF,NE-1))/2.
938      SUMNU=SUMNU+SNUTRA(NE)*EINT
939      SUMKE=SUMKE+SNOCK(NE)*EINT
940      SUMSD=SUMSD+PRMSD*EINT
941      SUMSC=SUMSC+PRMSC*EINT
942      SUMW=SUMW+WEISS(KP,NE)*EINT
943      WRITE(IWRI,482)EPS(JHALF,NE),SNUTRA(NE),SNOCK(NE),PRMSD,PRMSC,WEI
944      SS(KP,NE),TOMSD(NE),TOMSC(NE),TOTAL
945 480      CONTINUE
946 482      FORMAT(1F6.2,9(1PE11.3))
947      WRITE(IWRI,481)SUMNU,SUMKE,SUMSD,SUMSC,SUMW
948 481      FORMAT(GHOSUMS,6(1PE11.3))
949      IF(JQUT.NE.4)GO TO 483
950      IF(BF.EQ.0.)GO TO 483
951      FISBO=0.
952      FISSD=0.
953      FISCL=0.
954      FISMA=0.
955      FISHW=FISHW*SIGCN*FWEIS
956      WRITE(IWRI,490)FISBO,FISSD,FISCL,FISMA,FISHW
957 490      FORMAT(6H-FISS.,6(1PE11.3))
958 483      SIGCN=SIGCN/0.95
959      IF(JIN.GT.1)SIGCN=SIGCN/0.89
960 C
961 C      CALCULATE AND PRINT ANGULAR DISTRIBUTIONS
962      CALL ANGEL
963 C
964      IF(JPOUT.GT.0)GO TO 484
965      JPOUT=1
966      JNOUT=0
967      JOUT=1
968      BEN=BPRO
969      KP=2
970      JHALF=1
971      GO TO 450
972 484      IF(JNOUT.GT.0)GO TD 500
973      JPOUT=2
974      JNOUT=2
975      JOUT=4
976      NSD=1
977      BEN=BALFA
978      KP=3
979      NI=NAP1
980      JHALF=1
981      GO TO 450
982 C
983 C      RECYCLE OPTIONS
984 C
985 500      READ(IREA,5)JPOUT,JNOUT
986      JOUT=JPOUT+JNOUT
987      IF(JPOUT.GE.0)GO TO 504
988      READ(IREA,5)NEWP
989      IF(NEWP)1000,1,502
990 502      NEPS1=NEPS(1)+1
991      GO TO 9
992 504      READ(IREA,5)NSD
993      READ(IREA,3)BEN,GR
994      READ(IREA,3)PAIR,EPAIR
995      READ(IREA,13)NEPS(2),EPS(2,2),DELTA(2)
996      IF(NEPS(2).LE.50)GO TD 505
997      WRITE(IWRI,10)
998      GO TO 1000
999 505      IF(NEPS(2).LE.0)NEPS(2)=1
1000     NEPS1=NEPS(2)+1
1001     IF(EPS(2,2).LE.0.)GO TO 510
1002     IF(DELTA(2).LE.0.)DELTA(2)=1.
1003     DO 514 NE=3,NEPS1
1004     X=NE-2
1005 514     EPS(2,NE)=EPS(2,2)+X*DELTA(2)
1006     JHALF=2
1007     KP=4
1008     CALL CROSS

```

```

1009      GO TO 512
1010 510  DO 506 NE=2,NEPS1
1011      READ(IREA,3)EPS(2,NE),SIGIN(4,NE)
1012 506  CONTINUE
1013 512  EPS(2,1)=EPS(2,2)-1
1014      IF(EPS(2,1).LT.0.)EPS(2,1)=0.
1015      EPS(2,NEPS1+1)=EPS(2,NEPS1)+1.
1016      XOUT=JOUT
1017      XPUT=JPUT
1018      ATHRD=(ACDM-XOUT)**C 33
1019      COUL(4)=0.75*XPUT*(RZ+XPIN-XPUT)/ATHRD
1020      SIGBAR(4)=0.
1021      ESUM(4)=0.
1022      DO 508 NE=2,NEPS1
1023      IF(EPS(2,NE).LT.COUL(4))GO TO 508
1024      EINT=(EPS(2,NE+1)-EPS(2,NE-1))/2.
1025      ESUM(4)=ESUM(4)+EINT
1026      SIGBAR(4)=SIGBAR(4)+SIGIN(4,NE)*EINT
1027 508  CONTINUE
1028      SIGEAR(4)=SIGBAR(4)/ESUM(4)
1029      JHALF=2
1030 C      CALCULATE PROBABILITY FACTORS
1031 C
1032 C
1033 550  JTA=JPUT+JIN
1034  JTB=JNOUT+JPIN
1035  JLO=MAXO(JIN,JOUT,JTA,JTB)
1036  KP=2+JHALF
1037  IF(JOUT.EQ.1)KP=JPUT+1
1038  IF(IG.EQ.1)GO TO 551
1039 C
1040 C      1. PROB. FROM Z/A AND N/A
1041  JPOF=1
1042  IF(JPUT.LT.2)GO TO 524
1043  DO 522 J=2,JPUT
1044 522  JPOF=JPOF*J
1045 524  XPOF=JPOF
1046  JNCF=1
1047  IF(JNOUT.LT.2)GO TO 528
1048  DO 526 J=2,JNOUT
1049 526  JNOF=JNOF*J
1050 528  XNOF=JNOF
1051  APFAC=1.
1052  DO 529 N=1,JOUT
1053  X=N
1054 529  APFAC=APFAC*X
1055  XPUT=JPUT
1056  XNOUT=JNOUT
1057  CONEW=(RZ**XPUT)-(RN**XNOUT)
1058  CONEW=XPOF*XNOF/(APFAC-CONEW)
1059  DD 532 J=2,NPART
1060  JJ=J-1
1061 532  PROB(KP,JU)=1.
1062  DO 548 J=NPART,JHI
1063  XJ=J
1064  PROB(KP,J)=0.
1065  IF(J.LT.JOUT)GO TO 548
1066  XSUM=0.
1067  JMAX=J-JIN
1068  XNORM=1.
1069  DO 530 I=1,JOUT
1070  XI=I
1071 530  XNORM=XNORM*(XJ-XI+1.)/XI
1072  XMAX=JMAX
1073  JMAX=1+JMAX-1
1074  DO 546 I2=1,JMAX-1
1075  I=I2-1
1076  XI=I
1077  JP=JPIN+I
1078  JN=JIN+JMAX-I
1079  JP1=JP+1
1080  JN1=JN+1
1081  PROBI=1.
1082  IF(I.EQ.0)GO TO 536
1083  DO 534 II=1,I
1084  XII=II
1085 534  PROBI=PROBI*(XMAX-XII+1.)/XII
1086 536  PROBI=PROBI*(RZ**XI)-(RN**((XMAX-XI)))

```

```

1087      XSUM=XSUM+PROB1
1088      PROB1=PROB1/(XPOF*XNOF)
1089      IF(JP .LT. JPOUT)GO TO 546
1090      IF(JN.LT.JNOUT)/\DO TO 546
1091      IF(JPOUT.EQ.0)GO TO 540
1092      DO 538 II=1,JPOUT
1093      X=JP1-II
1094 538   PROB1=PROB1*X
1095 540   IF(JNOUT.EQ.0)GO TO 544
1096   DO 542 II=1,JNOUT
1097   X=JN1-II
1098 542   PROB1=PROB1*X
1099 544   PROB(KP,J)=PROB(KP,J)+PROB1
1100 546   CONTINUE
1101   PROB(KP,J)=PROB(KP,J)*CONEW
1102   PROB(KP,J)=PROB(KP,J)/(XNORM-XSUM)
1103 548   CONTINUE
1104   GO TO 575
1105 C
1106 C      2. PROB. FROM AVAILABLE PHASE SPACE (GADIOLI)
1107 551   JPLO=MAX0(JPIN,JPOUT)
1108   JNLO=MAX0(JNIN,JNOUT)
1109   XPOUT=JPOUT
1110   XNOUT=JNOUT
1111   XOUT=JOUT
1112   XPFAC=1.
1113   XHFAC=1
1114   DO 552 J=2,NPART
1115   JJ=J-1
1116   XJJ=JJ
1117   XPFAC=XPFAC*XJJ
1118 552   PROB(KP,JJ)=1.
1119   IF(NHOLE.LT.2)GO TO 556
1120   DO 554 J=2,NHOLE
1121   XJ=J-1
1122 554   XHFAC=XHFAC*XJ
1123 556   DO 574 JP=NPART,JHI
1124   XP=JP
1125   PROB(KP,JP)=0.
1126   IF(JP.LT.JOUT)GO TO 574
1127   XPFAC=XPFAC*XP
1128   XH=JP-NPHD
1129   XHFAC=XHFAC*XH
1130   XNORM=XPFAC*XHFAC
1131   XNUM=XHFAC
1132   IMAX=JP-JOUT
1133   IF(IMAX.LT.2) GO TO 559
1134   DO 558 I=2,IMAX
1135   XI=I
1136 558   XNUM=XNUM-XI
1137 559   SUMN=0.
1138   SUMD=0.
1139   JPMAX=JP-JNIN
1140   DO 572 JPPI=JPIN,JPMAX
1141   JPNU=JP-JPPI
1142   JHPI=JPPI-JPIN
1143   JHNU=JPNU-JNIN
1144   XPI=JPPI+JHPI
1145   XNU=JPNU+JHNU
1146   FACN=1.
1147   IF(JHPI.LT.2)GO TO 561
1148   DO 560 I=2,JHPI
1149   XI=I
1150 560   FACN=FACN-XI
1151 561   IF(JHNU.LT.2)GO TO 563
1152   DO 562 I=2,JHNU
1153   XI=I
1154 562   FACN=FACN-XI
1155 563   FACD=FACN
1156   IF(JPPI.LT.2)GO TO 565
1157   DO 564 I=2,JPPI
1158   XI=I
1159 564   FACD=FACD-XI
1160 565   IF(JPNU.LT.2)GO TO 567
1161   DO 566 I=2,JPNU
1162   XI=I
1163 566   FACD=FACD-XI
1164 567   FACD=XNORM/FACD

```

```

1165      FACD=FACD+RZ--*PI*RN--*XNU
1166      SUMD=SUMD+FACD
1167      IF(JPPI.LT.JPLD)GO TO 572
1168      IF(JPNU.LT.JNLD)GO TO 572
1169      IMAX=JPPI-JPOUT
1170      IF(IMAX.LT.2)GO TO 569
1171      DO 568 I=2,IMAX
1172      XI=I
1173 568   FACN=FACN-XI
1174 569   IMAX=JPNU-JNOUT
1175      IF(IMAX.LT.2)GO TO 571
1176      DO 570 I=2,IMA*
1177      XI=1
1178 570   FACN=FACN*XI
1179 571   FACN=XNUM/FACN
1180      FACN=FACN+RZ--*(XP1-XP0UT)*RN--*(XNU-XNOUT)
1181      SUMN=SUMN+FACN
1182 572   CONTINUE
1183      IF(SUMD.GT.0.)PROB(KP,JP)=SUMN/SUMD
1184 574   CONTINUE
1185 575   AP=JOUT
1186      IF(JHI.GE.JBAR)GO TO 620
1187      JHI=JHI+1
1188      DO 582 J=JHI,JBAR
1189 582   PROB(KP,J)=1
1190 C
1191 C      CALCULATE EMISSION RATES
1192 C
1193 620   SD=NSD
1194      CNNE=0.85
1195      IF(JOUT.EQ.1)CNNE=0.95
1196      REL=SD*AP*372000.*CNNE
1197      IF(GR.EQ.0.)GR=G*(ACOM-AP)/ACOM
1198      ASPR=1.645*GR
1199      NAP1=MAX0(JLD,NDWN)-1
1200 606   NAP1=NAP1+1
1201      NTEST=2-NAP1-NPHD
1202      IF(NTEST.LT.JOUT+1)GO TO 606
1203      REF=SQRT(1.645*SIZE)
1204      DO 612 NE=2,NEPS1
1205      U=E-EPAIR-BEN-EPS(JHALF,NE)
1206      DO 602 NP=1,30
1207 602   WU(KP,NP,NE)=0.
1208      WEISS(KF,NE)=0.
1209      IF(U.LE.0.)GO TO 612
1210      QUO=U/(E-EPAIRC)
1211      PROD=EPS(JHALF,NE)-SIGNIN(KP,NE)-REL
1212      PREX=PROD/QUO
1213      EX=SQRT(ASPR*U)
1214      EX=2.-(EX-REF)
1215      WEISS(KP,NE)=PREX-EXP(EX)
1216      IF(WEISS(KP,NE).LT.FLOW)WEISS(KP,NE)=0.
1217      IF(WEISS(KP,NE).GT.SPILL)WEISS(KP,NE)=0.
1218      FAC=1.
1219      IT=JOUT+3
1220      IF(NAP1.LT.IT)GO TO 601
1221      DO 600 I=IT,NAP1
1222      Y=I-JOUT-1
1223 600   FAC=FAC*Y
1224 601   NH=NAP1-NPHD
1225      IF(NH.LT.3)GO TO 604
1226      DO 603 I=3,N4
1227      Y=I-1
1228 603   FAC=FAC*Y
1229 604   NPH=NAP1+NH
1230      IT=JOUT+5
1231      IF(NPH.LT.IT)GO TO 607
1232      DO 605 I=IT,NPH
1233      Y=I-JOUT-3
1234 605   FAC=FAC*Y
1235 607   FAC=ALOG10(FAC)
1236      AP2=JOUT+2
1237      DO 622 NP=NAP1,JBAR
1238      NP1=NP+1
1239      P=NP
1240      IF(NP.EQ.JOUT)GO TO 608
1241      FAC=FAC+ALOG10(P-AP)
1242 608   H=NP-NPHD

```

```

1243      IF(H.EQ.0.)GO TO 609
1244      FAC=FAC+ALOG10(H)
1245 609    PH=P+H
1246      IF(PH.LE.AP2)GO TO 610
1247      FAC=FAC+ALOG10(PH-AP-1. )+ALOG10(PH-AP-2. )
1248 610    PTT=P-AP
1249      Q=AMAX1(PTT,H)
1250      RHOL=GR+U-Q-Q
1251      IF(NH.LE.2)RHOL=GR*(U+EPAIR-PAIR)-Q-Q
1252      IF(RHOL.GE.0.)GO TO 611
1253      RHOL=0.
1254      GO TO 613
1255 611    RHOL=RHOL+(PTT*(PTT+1. )+H*(H+1. ))/4.
1256      RHOL=(PH-AP-1. )+ALOG10(RHOL)
1257      RHOL=RHOL-FAC
1258      RHOL=(10.-**RHOL)*GR
1259      IF(RHOL.LE.1.)RHOL=0
1260      IF(RHOL.GT.POUR)RHOL=0.
1261      IF(H.LE.0.)GO TO 613
1262      IF(H.GT.2.)GO TO 613
1263      I=NP-NPHD+1
1264      UU=U+EPAIR-PAIR
1265      IF(UU.LE.EHOLE(I))GO TO 613
1266      X=(UU-EHOLE(I))/UU
1267      X=H*X**=(PH-AP-1. )
1268      Y=0.
1269      EH2=EHOLE(I)+EHOLE(1)
1270      IF(UU.LE.EH2)GO TO 614
1271      Y=(UU-EH2)/UU
1272      Y=0.5*H*(H-1. )+Y**=(PH-AP-1. )
1273 614    RHOL=RHOL*(1.-X+Y)
1274 613    WR=RHOL*PROD=PROB(KP,NP)*TAUWM
1275      X=0.
1276      IF(RHOU(NP1).GT.0.)X=WR/RHOU(NP1)
1277      IF(X.LT.FLOW)X=0.
1278      IF(X.GT.SPILL)X=0.
1279      WU(KP,NP,NE)=X
1280 622    CONTINUE
1281 612    CONTINUE
1282      IF(JHALF.EQ.1)GO TO 150
1283      NI=NAP1
1284      GO TO 450
1285 1000   CALL EXIT
1286 C
1287 C      HALLELUJAH
1288 C
1289      END

```

```

1290      SUBROUTINE ANGEL
1291 C
1292 C      CALCULATE ANGULAR DISTRIBUTIONS
1293 C      FROM SYSTEMATIC LEGENDRE COEFFICIENTS
1294 C      PROGRAM - MARCH 1979
1295 C      SUBROUTINE - JANUARY 1980
1296 C
1297 C      LIBRARY OF POLYNOMIALS GENERATED IN SUBROUTINE POLLY OF MAIN
1298 C
1299      DIMENSION SIGMA(10),B(7),JANG(10)
1300      COMMON /ANGELS/ POL(19,7),TOMSD(51),TOMSC(51),IWRI,IANG
1301      COMMON /ENERGY/ BEN,NEPS1,EPS(2,52),JHALF,JNOUT,JPOUT
1302      DO 10 I=1,10
1303      JANG(I)=0
1304 10     SIGMA(I)=0.
1305      DO 11 I=1,7
1306 11     B(I)=0.
1307      WRITE(IWRI,12)JPOUT,JNOUT
1308 12     FORMAT(26H1ANGULAR DISTRIBUTIONS Z=.1I2.4H N=.1I2)
1309      IF(IANG.EQ.0)WRITE(IWRI,6)
1310 6      FORMAT(2OH ENERGY PARAM. = EPS)
1311      IF(IANG.EQ.1)WRITE(IWRI,8)
1312 8      FORMAT(25H ENERGY PARAM. = EPS+B.E.)
1313      ANG=IANG
1314      WRITE(IWRI,14)
1315 14     FORMAT(29H CROSS SECTIONS IN MB/STR-MEV)
1316      DO 15 J=1,10
1317 15     JANG(J)=10.*J-10
1318      WRITE(IWRI,16)(JANG(J),J=1,10)
1319 16     FORMAT(5H- EPS,1I7,9I9)
1320      WRITE(IWRI,18)
1321 18     FORMAT(1H )
1322      B(1)=1.
1323      DO 28 NE=2,NEPS1
1324      EPSCM=EPS(JHALF,NE)+ANG=BEN
1325      AOMSD=TOMSD(NE)/12.5664
1326      AOMSC=TOMSC(NE)/12.5664
1327      TCTAL=AOMSD+AOMSC
1328      IF(TOTAL.LE.0.)GO TO 28
1329      DO 19 L=2,7
1330      XL=L-1
1331      BL=XL*(XL+1.)
1332      AL=0.036+0.0039*BL
1333      EL=92.+ANG=6.-90./SQRT(BL)
1334      X=AL=(BL-EPSCM)
1335      X=1.+EXP(X)
1336 19     B(L)=(2.-XL+1.)/X
1337      DO 24 I=1,10
1338      SIG=0.
1339      DO 20 L=1,7
1340 20     SIG=SIG+B(L)*POL(I,L)
1341      SIGMA(I)=SIG*AOMSD
1342      SIG=0.
1343      DO 22 LL=1,4
1344      L=2*LL-1
1345 22     SIG=SIG+B(L)*POL(I,L)
1346 24     SIGMA(I)=SIGMA(I)+SIG*AOMSC
1347      WRITE(IWRI,26)EPS(JHALF,NE),(SIGMA(I),I=1,10)
1348 26     FORMAT(1F6.2,1O(1PE9.2))
1349 28     CONTINUE
1350      DO 27 J=1,9
1351 27     JANG(J)=10.*J+90
1352      WRITE(IWRI,29)(JANG(J),J=1,9)
1353      WRITE(IWRI,18)
1354 29     FORMAT(5H- EPS,1I7,8I9,1OH TOTAL)
1355      DO 3B NE=2,NEPS1
1356      EPSCM=EPS(JHALF,NE)+ANG=BEN
1357      AOMSD=TOMSD(NE)/12.5664
1358      AOMSC=TOMSC(NE)/12.5664
1359      TOTAL=TOMSD(NE)+TOMSC(NE)
1360      IF(TOTAL.LE.0.)GO TO 38
1361      DO 30 L=2,7
1362      XL=L-1
1363      BL=XL*(XL+1.)
1364      AL=0.036+0.0039*BL
1365      BL=92.+ANG=6.-90./SQRT(BL)
1366      X=AL=(BL-EPSCM)
1367      X=1.+EXP(X)
1368 30     B(L)=(2.-XL+1.)/X
1369      DO 36 J=1,9

```

```

1370      I= J+10
1371      SIG=0.
1372      DO 32 L=1,7
1373 32    SIG=SIG+B(L)*POL(I,L)
1374      SIGMA(J)=SIG+AOMSD
1375      SIG=0.
1376      DO 34 LL=1,4
1377      L=2*LL-1
1378 34    SIG=SIG+B(L)*POL(I,L)
1379 36    SIGMA(J)=SIGMA(J)+SIG*AOMSC
1380    WRITE(IWRI,26)EPS(JHALF,NE),(SIGMA(I),I=1,9),TOTAL
1381 38    CONTINUE
1382    RETURN
1383    END

```

```

1384      SUBROUTINE CROSS
1385 C
1386 C      CALCULATE OPTICAL MODEL REACTION CROSS SECTIONS
1387 C      WITH EMPIRICAL PARAMETERIZATION OF
1388 C      NARASIMHA MURTHY, CHATTERJEE, AND GUPTA
1389 C
1390 C      PARAMETER VALUES SET IN SUBROUTINE SIGPAR
1391 C
1392      COMMON /ENERGY/ BEN,NEPS1,EPS(2,52),JHALF,JNOUT,JPOUT
1393      COMMON /DIRECT/ SIGIN(4,51),JPIN,JNIN,BEA,NSD,KP,E,ACOM,G,RZZ
1394      COMMON /PAR/ PO(3,3),P1(3,3),P2(3,3),XLO(3,3),XL1(3,3),XMO(3,3)
1395      1,XM1(3,3),XNO(3,3),XN1(3,3),XN2(3,3)
1396      FLOW=1.E-18
1397      SPILL=1.E+18
1398      JOUT=JPOUT+JNOUT
1399      XOLT=JOUT
1400      ATAR=ACOM-XOUT
1401      ATHRD=ATAR-0.333
1402      IF(JPOUT.GT.0.)GO TO 2
1403      XLAMB=XLO(1,2)/ATHRD+XL1(1,2)
1404      XMU=XMO(1,2)=ATHRD+XM1(1,2)=ATHRD-ATHRD
1405      XNU=XNO(1,2)*ATHRD=ATAR+XN1(1,2)=ATHRD-ATHRD+XN2(1,2)
1406      EC=2.4
1407      P=PO(1,2)
1408      NCOU=2
1409      ETEST=32.
1410      GO TO 3
1411 2     RA=1.20
1412      IF(JOUT.EQ.1)RA=0.
1413      XPOUT=JPOUT
1414      RZ=JPIN-JPOUT
1415      RZ=RZ+RZZ
1416      EC=1.44*XPOUT=RZ/(1.5*ATHRD+RA)
1417      I=JPOUT+1
1418      J=JNOUT+1
1419      P=PO(I,J)+P1(I,J)/EC+P2(I,J)/(EC-EC)
1420      XLAMB=XLO(I,J)=ATAR+XL1(I,J)
1421      A=ATAR-XM1(I,J)
1422      XMU=XMO(I,J)*A
1423      XNU=A-(XNO(I,J)+XN1(I,J)=EC+XN2(I,J)=EC*EC)
1424      IF(JOUT.EQ.2)RA=0.8
1425      IF(JOUT.EQ.3)RA=0.8
1426      XNL=XNL/XLAMB
1427      IF(XNL.GT.SPILL)XNL=0.
1428      IF(XNL.LT.FLOW)GO TO 3
1429      ETEST=1.2=SQRT(XNL)
1430 3     A=-2.*P*EC+XLAMB-XNU/(EC*EC)
1431      B=P*EC-EC+XMU+2.*XNU/EC
1432      ECUT=0.
1433      CUT=A*A-4.*P*B
1434      IF(CUT.GT.0.)ECUT=SQRT(CUT)
1435      ECUT=(ECUT-A)/(2.*P)
1436      DO 4 NE=2,NEPS1
1437      IF(EPS(JHALF,NE).GT.EC)GO TO 6
1438 4     CONTINUE
1439 6     NCOU=NE
1440      IF(NCOU.LT.3)GO TO 10
1441      NC=NCOU-1
1442      DO 8 NE=2,NC
1443      ELAB=EPS(JHALF,NE)=ACDM/ATAR
1444      SIGIN(KP,NE)=P*ELAB*ELAB+A*ELAB+B

```

```

1445 8      IF(ELAB.LT.ECUT)SIGIN(KP,NE)=0.
1446 10     DO 12 NE=NCOU,NEPS1
1447       ELAB=EPS(JHALF,NE)*ACOM/ATAR
1448       SIG=XLAMB=ELAB+XMU+XNU/ELAB
1449       GEOM=0.
1450       IF(XNL.LT.FLOW)GO TO 12
1451       IF(ELAB.LT.ETEST)GO TO 12
1452       GEOM=SORT(XOUT*EPS(JHALF,NE))
1453       GEOM=1.23*ATHRD+RA+4.573/GEOM
1454       GEOM=31.416*GEOM*GEOM
1455 12      SIGIN(KP,NE)=AMAX1(GEOM,SIG)
1456      RETURN
1457      END

1458      SUBROUTINE INEL
1459 C      PHENOMENOLOGICAL ENERGY SPECTRA
1460 C      INELASTIC SCATTERING
1461 C      (RXN MUST INVOLVE AT LEAST ONE COMPLEX PARTICLE)
1462 C      FEBRUARY 1980
1463 C
1464 C      COMMON /ENERGY/ BEN,NEPS1,EPS(2,52),JHALF,JNOUT,JPOUT
1465      COMMON /DIRECT/ SIGIN(4,51),JPIN,JNIN,BEA,NSD,KP,F,ACOM,G,RZZ
1466      COMMON /CLSTR/ SNDCK(51),SIGBAR(5),SIGCN,COUL(5),BNEU,BPRO,BALFA
1467      XPIN=JPIN
1468      XNIN=JNIN
1469      JIN=JPIN+JNIN
1470      XIN=JIN
1471      ATARG=ACOM-XIN
1472      RNN=ATARG-RZZ
1473
1474 C      CALCULATE NORMALIZATION
1475 C
1476 C      XNDR=JIN*NSD
1477      XNORM=XNDR*SIGCN/16.
1478
1479 C      1. PROJECTILE=EJECTILE DENOMINATOR
1480      XNDR=XNDR*SIGBAR(KP)/6.
1481      COUP=COUL(KP)
1482      EMAX=E-BEN
1483      XNDR=XNDR*(EMAX+2.*COUP)*(EMAX-COUP)*(EMAX-COUP)
1484      IF(JIN.GT.1.)GO TO 2
1485      GBA=(RZZ-1.)/13.
1486      IF(JNIN.EQ.1.)GBA=(RNN-1.),13.
1487      GO TO 10
1488 2      IF(JIN-3).LT.5.6
1489 4      GBN=(ACOM-1.)/52.
1490 4      GBA=(ACOM-4.)/52.
1491      GO TO 8
1492 5      GBN=(ACOM-1.)/156.
1493 5      GBA=(ACOM-4.)/156.
1494 6      GO TO 8
1495 6      GBN=(ACOM-1.)/208.
1496 6      GBA=(ACOM-4.)/208.
1497
1498 C      2. NEUTRON EXCITATION
1499 8      GII=(RNN+XNIN-1.)/13.
1500 8      GIR=RNN/13.
1501 8      P=RNN/ATARG
1502 8      XNORI=2.*SIGBAR(1)*GII*GBN/6.
1503 8      EMAX=E-BNEU
1504 8      COUP=COUL(1)
1505 8      XNORI=XNORI*(EMAX+2.*COUP)*(EMAX-COUP)*(EMAX-COUP)
1506 8      XNORB=XNDR*GIR*GIR
1507 8      XN=GIR*GIR*P/(XNORI+XNORB)
1508 8      EPN=1./GIR
1509 8      UDQN=1./GIR
1510 8
1511 C      3. PROTON EXCITATION
1512 C      GII=(RZZ+XPIN-1.)/13.
1513 8      GIR=RZZ/13.
1514 8      P=RZZ/ATARG
1515 8      XNORI=2.*SIGBAR(2)*GII*GBN/6.
1516 8      EMAX=E-BPRO
1517 8      COUP=COUL(2)
1518 8      XNORI=XNORI*(EMAX+2.*COUP)*(EMAX-COUP)*(EMAX-COUP)
1519 8      XNORB=XNDR*GIR*GIR
1520 8      XN=GIR*GIR*P/(XNORI+XNORB)
1521 8

```

```

1522      EPP=1./GIR
1523      UADP=1./GIR
1524 C
1525 C      4. ALPHA EXCITATION
1526 10      GI=(ACOM-4.)/208.
1527      GIR=ATARG/208.
1528      P=0.08*RZZ/(2.*ATARG)
1529      IF(RNN.LE.116.)GO TO 14
1530      IF(RNN.GE.129.)GO TO 14
1531      IF(RNN-126.)12,12,13
1532 12      Y=0.02+0.06*(126.-RNN)/10.
1533      P=P*Y/0.08
1534      GO TO 14
1535 13      Y=0.02+0.06*(RNN-126.)/3.
1536      P=P*Y/0.08
1537 14      XNORI=4.*SIGBAR(3)*GII*GBA/6.
1538      EMAX=E-BALFA
1539      COUP=COUL(3)
1540      XNORI=XNORI*(EMAX+2.*COUP)*(EMAX-COUP)*(EMAX-COUP)
1541      XNORB=XNORI*GIR=GIR
1542      XA=GIR*GIR=P/(XNORI+XNORB)
1543      EPA=52./RZZ+52./RNN
1544      UADA=1./GIR
1545 C
1546 C      CALCULATE ENERGY SPECTRUM
1547 C
1548      UMAX=E-BEN
1549      FLOW=1.E-18
1550      SPILL=1.E+18
1551      DO 16 NE=2,NEPS1
1552      U=UMAX-EPS(JHALF,NE)
1553      IF(U.LE.0.)GO TO 18
1554      X=0.
1555      IF(JIN.LE.1)GO TO 22
1556      UN=U-EPN
1557      IF(UN.LE.0.)GO TO 20
1558      UN=UN+UADN
1559      X=X+XN*UN
1560 20      UP=U-EPP
1561      IF(UP.LE.0.)GO TO 22
1562      UP=UP+UADP
1563      X=X+XP*UP
1564 22      UA=U-EPA
1565      IF(UA.LE.0.)GO TO 24
1566      UA=UA+UADA
1567      X=X+XA*UA
1568 24      SIG=XNORM-EPS(JHALF,NE)*SIGIN(KP,NE)-X
1569      IF(SIG.LT.FLOW)SIG=0.
1570      IF(SIG.GT.SPILL)SIG=0.
1571 16      SNOCK(NE)=SIG
1572 18      RETURN
1573      END

1574      SUBROUTINE KNOCK
1575 C
1576 C      PHENOMENOLOGICAL ENERGY SPECTRA
1577 C      KNOCKOUT OF N, P, OR ALPHA
1578 C      (RXN MUST HAVE AT LEAST ONE COMPLEX PARTICLE)
1579 C      FEBRUARY 1980
1580 C
1581      COMMON /ENERGY/ BEN,NEPS1,EPS(2,52),JHALF,JNOUT,JPOUT
1582      COMMON /DIRECT/ SIGIN(4,51),JPIN,JNIN,BEA,NSD,KP,E,ACOM,G,RZZ
1583      COMMON /CLSTR/ SNOCK(51),SIGBAR(5),SIGCN,COUL(5),BNEU,BPRO,BALFA
1584      JIN=JPIN+JNIN
1585      XIN=JIN
1586      XPIN=JPIN
1587      XNIN=JNIN
1588      JOUT=JPOUT+JNOUT
1589      XOUT=JOUT
1590      XPOUT=JPOUT
1591      XNOUT=JNOUT
1592      ATARG=ACOM-XIN
1593      RNN=ATARG-RZZ
1594      ARES=ACOM-XOUT

```

```

1595 C      CALCULATE S. P STATE DENSITIES
1596 C      AND (2S+1)*A TYPE FACTORS
1597 C
1598 C
1599 IF(JIN.GT.1)GO TO 2
1600 REL=2.
1601 GA=(RZZ+1.-XPOUT)/13.
1602 IF(JNIN.EQ.1)GA=(RNN+1.-XNOUT)/13.
1603 GO TO 8
1604 2      IF(JIN-314,5,6
1605 4      REL=6.
1606 GA=ARES/52.
1607 GO TO 8
1608 5      REL=6.
1609 GA=ARES/156.
1610 GO TO 8
1611 6      REL=4.
1612 GA=ARES/208.
1613 8      IF(JOUT.GT.1)GO TO 10
1614 GBT=RZZ/13.
1615 GBR=(RZZ+XPIN-1.)/13.
1616 IF(JNOUT.EQ.0)GO TO 12
1617 GBT=RNN/13.
1618 GBR=(RNN+XNIN-1.)/13.
1619 GO TO 12
1620 10     GBT=ATARG/208.
1621 GBR=ARES/208.
1622 C      CALCULATE NORMALIZATION
1623 C
1624 C
1625 12     XNORM=SIGCN/16.
1626 XNORB=JOUT=NSD
1627 XNORM=XNORM*XNORE
1628 XNORB=XNORB-SIGBAR(KP)*GA*GBR/6.
1629 EMAX=E-BEN
1630 COUP=COUL(KP)
1631 XNORB=XNORB*(EMAX+2.*COUP)*(EMAX-COUP)*(EMAX-COUP)
1632 XNORA=REL=SIGBAR(5)*GBT*GBT/6.
1633 EMAX=E-BEA
1634 COUP=COUL(5)
1635 XNORA=XNORA*(EMAX+2.*COUP)*(EMAX-COUP)*(EMAX-COUP)
1636 XNORM=XNORM*GA=GBR/(XNORA+XNORB)
1637 IF(JPOUT-1)14,15,16
1638 14     XNORM=XNORM*(1.-RZZ/ATARG)
1639 GO TO 20
1640 15     XNORM=XNORM=RZZ/ATARG
1641 GO TO 20
1642 16     XNORM=XNORM*0.08=RZZ/(2.*ATARG)
1643 IF(RNN.LE.116.)GO TO 20
1644 IF(RNN.GE.129.)GO TO 20
1645 IF(RNN-126.,)18,18,19
1646 18     X=0.02+0.06*(126.-RNN)/10.
1647 XNORM=XNORM*X/0.08
1648 GO TO 20
1649 19     X=0.02+0.06*(RNN-126.)/3.
1650 XNORM=XNORM*X/0.08
1651 C
1652 C      CALCULATE ENERGY SPECTRUM
1653 C
1654 20     FLOW=1.E-18
1655 S-ILL=1.E+18
1656 UMAX=E-BEN
1657 Q=JPIN+JPOUT
1658 UPAUL=Q-Q*13./ (RZZ+XPIN-XPOUT)
1659 Q=JNIN+JNOUT
1660 UPAUL=UPAUL+Q*Q*13./ (RNN+XNIN-XNOUT)
1661 UADD=0.5/GBR+0.5/GA
1662 DO 22 NE=2,NEPS1
1663 U=UMAX-EPS(1,NE)
1664 U=U-UPAUL
1665 IF(U.LT.0.)GO TO 24
1666 U=U+UADD
1667 SIG=XNORM*U-EPS(1,NE)*SIGIN(KP,NE)
1668 IF(SIG.LT.FLOW)SIG=0.
1669 IF(SIG.GT.SPI_L)SIG=0.
1670 22     SNOCK(NE)=SIG
1671 24     RETURN
1672 END

```

```

1673      SUBROUTINE NUTRA
1674 C
1675 C      PHENOMENOLOGICAL ENERGY SPECTRA
1676 C      NUCLEON TRANSFER
1677 C      PROGRAM - FEBRUARY 1979
1678 C      SUBROUTINE (SIMPLIFIED) - FEBRUARY 1980
1679 C
1680      COMMON /ENERGY/ BEN,NEPS1,EPS(2,52),JHALF,JNOUT,JPOUT
1681      COMMON /DIRECT/ SIGIN(4,51),JPIN,JNIN,BEA,NSD,KP,E,ACOM,G,RZZ
1682      COMMON /NUTRAN/ SNUTRA(51)
1683      FLOW=1.E-18
1684      SPILL=1.E+18
1685      JIN=JNIN+JPIN
1686      XIN=JIN
1687      JOUT=JNOUT+JPOUT
1688      XOUT=JCUT
1689      ATARG=ACOM-XIN
1690      ECOM=E-BEA
1691      ELABI=ECOM+ACOM/ATARG
1692      ARES=ACOM-XOUT
1693      UMAX=E-BEN
1694      VAB=0.25*50.*XIN
1695      KAB=1.
1696      IF(JHALF.EQ.2)GO TO 2
1697      IF(JIN.EQ.2)GO TO 2
1698      IF(JIN.EQ.3)GO TO 2
1699      KAB=12.
1700      2    VELO=XIN/(ELABI+VAB)
1701      X=JOUT-NSD
1702      X=X*KAB
1703      Y=6.
1704      IF(JIN.EQ.1)Y=2.
1705      IF(JIN.EQ.4)=4
1706      XHOLD=0.0127*X/(Y*ECOM-XIN)
1707      TRANP=IABS(JPOUT-JPIN)
1708      ZA=2.*RZZ/ATARG
1709      TRANN=IABS(JNOUT-JNIN)
1710      PH=TRANN+TRANP
1711      NEX=0
1712      IF(PH.GT.0.) GO TO 5
1713      NEX=1.
1714      PH=2.
1715      TRANP=2.
1716      5    X=XHOLD*(2860./ARES)**PH
1717      X=X*VELO**(2.-PH)
1718      X=X-ZA***(6.-TRANP)
1719      JP=JPIN-JPOUT
1720      Y=JP
1721      GP=(RZZ+Y)/13.
1722      JN=JNIN-JNOUT
1723      Y=JN
1724      RNN=ATARG-RZZ
1725      GN=(RNN+Y)/13.
1726      X=X*(GP**TRANP)**(GN**-TRANN)
1727      JPICKP=MAX0(JP,0).
1728      IF(JPICKP.LT.2)GO TO 6
1729      DO 4 N=2,JPICKP
1730      Y=N
1731      4    X=X/Y
1732      6    JSTRIP=JPICKP-JP
1733      IF(JSTRIP.LT.2)GO TO 10
1734      DO B N=2,JSTRIP
1735      Y=N
1736      8    X=X/Y
1737      10   JPICKN=MAX0(JN,0)
1738      IF(JPICKN.LT.2)GO TO 14
1739      DO 12 N=2,JPICKN
1740      Y=N
1741      12   X=X/Y
1742      14   JSTRIN=JPICKN-JN
1743      IF(JSTRIN.LT.2)GO TO 18
1744      DO 16 N=2,JSTRIN
1745      Y=N
1746      16   X=X/Y
1747      18   J=JPICKP+JPICKN
1748      IF(NEX.GT.0) J=1
1749      H=J
1750      JMAX=JSTRIP+JSTRIN
1751      IF(NEX.GT.0)JMAX=1
1752      P=JMAX

```

```

1753      IMAX=PH
1754      GR=ARES/13.
1755      W=O.
1756      FAC=1.
1757      DO 20 I=1,IMAX
1758      XI=I
1759 20    IF(I.GT.1)X=X/(XI-1.)
1760      PPI=USTRIP
1761      PNU=USTRIN
1762      HPI=UPICKP
1763      HNU=UPICKN
1764      Q=AMAX1(PNU,HNU)
1765      UMAX=UMAX-Q/GN
1766      Q=AMAX1(PPI,HPI)
1767      UMAX=UMAX-Q/GP
1768      UADD=(PNU-(PNU+1.))+HNU-(HNU+1.))/(4.*GN)
1769      UADD=UADD+(PPI+(PPI+1.)*HPI+(HPI+1.))/(4.*GP)
1770      DO 40 NE=2,NEPS1
1771      U=UMAX-EPS(JHALF,NE)
1772      IF(U.LT.O.)GO TO 40
1773      U=U+UADD
1774      W=X*U**((PH-1.))
1775      SIG=W-EPS(JHALF,NE)-SIGIN(KP,NE)
1776      IF(SIG.LT.FLOW)SIG=0.
1777      IF(SIG.GT.SPILL)SIG=0.
1778      SNUTRA(NE)=SIG+SNUTRA(NE)
1779 40    CONTINUE
1780      IF(NEX.NE.1) GO TO 42
1781      NEX=2
1782      TRANP=0.
1783      TRANR=2.
1784      GO TO 5
1785 42    RETURN
1786      END

```

```

1787      SUBROUTINE POLLY
1788 C
1789 C      GENERATE LIBRARY OF LEGENDRE POLYNOMIALS
1790 C
1791      COMMON /ANGELS/ POL(19,7),TOMSD(51),TOMSC(51),IWRI,IANG
1792      DO 4 J=1,19
1793      ANG=(J-1)*10
1794      ANG=ANG*3.14159/180.
1795      COSX=COS(ANG)
1796      COSX2=COSX*COSX
1797      COSX3=COSX*COSX2
1798      COSX4=COSX*COSX3
1799      COSX5=COSX*COSX4
1800      COSX6=COSX*COSX5
1801      POL(J,1)=1.
1802      POL(J,2)=COSX
1803      POL(J,3)=(3.*COSX2-1.)/2.
1804      POL(J,4)=(5.*COSX3-3.*COSX)/2.
1805      POL(J,5)=(35.*COSX4-30.*COSX2+3.)/8.
1806      POL(J,6)=(63.*COSX5-70.*COSX3+15.*COSX)/8.
1807      POL(J,7)=(231.*COSX6-315.*COSX4+105.*COSX2-5.)/16.
1808 4    CONTINUE
1809      RETURN
1810      END

```

```

1811      SUBROUTINE SIGPAR
1812 C
1813 C      STORE PARAMETERS FOR CALCULATING APPROXIMATE
1814 C      OPTICAL MODEL REACTION CROSS SECTIONS
1815 C
1816 C      NEUTRON FROM MANI ET AL
1817 C      PROTON FROM BECCHETTI-GREENLEES
1818 C      ALPHA FROM HUIZENGA AND IGO
1819 C      D FROM O.M. OF PEREY AND PEREY
1820 C      T FROM O.M. OF HAFELE, FLYNN ET AL
1821 C      3HE FROM O.M. OF GIBSON ET AL

```

```

1822 C
1823 COMMON /PAR/ PO(3,3),P1(3,3),P2(3,3),XLO(3,3),XL1(3,3),XMO(3,3)
1824 1,XM1(3,3),XNO(3,3),XN1(3,3),XN2(3,3)
1825 DO 2 I=1,3
1826 DO 2 J=1,3
1827 PO(I,J)=0.
1828 P1(I,J)=0.
1829 P2(I,J)=0.
1830 XLO(I,J)=0.
1831 XL1(I,J)=0.
1832 XMO(I,J)=0.
1833 XM1(I,J)=0.
1834 XNO(I,J)=0.
1835 XN1(I,J)=0.
1836 2 XN2(I,J)=0.
1837 PO(1,2)=-312.
1838 PO(2,1)=15.72
1839 PO(2,2)=0.798
1840 PO(2,3)=-21.45
1841 PO(3,2)=-2.88
1842 PO(3,3)=10.95
1843 P1(2,1)=9.65
1844 P1(2,2)=420.3
1845 P1(2,3)=484.7
1846 P1(3,2)=105.6
1847 P1(3,3)= 85.21
1848 P2(2,1)=-449.
1849 P2(2,2)=-1651.
1850 P2(2,3)=-1608.
1851 P2(3,2)=-1487.
1852 P2(3,3)=1146.
1853 XLO(1,2)=12.10
1854 XLO(2,1)=0.00437
1855 XLO(2,2)=0.00619
1856 XLO(2,3)=0.0186
1857 XLO(3,2)=0.00459
1858 XLO(3,3)=0.0642
1859 XL1(1,2)=-11.2/
1860 XL1(2,1)=-16.58
1861 XL1(2,2)=-7.54
1862 XL1(2,3)=-8.90
1863 XL1(3,2)=-8.93
1864 XL1(3,3)=-13.96
1865 XMO(1,2)=234.1
1866 XMO(2,1)=244.7
1867 XMO(2,2)=583.5
1868 XMO(2,3)=686.3
1869 XMO(3,2)=611.2
1870 XMO(3,3)=781.2
1871 XM1(1,2)=38.26
1872 XM1(2,1)=0.503
1873 XM1(2,2)=0.337
1874 XM1(2,3)=0.325
1875 XM1(3,2)=0.35
1876 XM1(3,3)=0.29
1877 XNO(1,2)=1.55
1878 XNO(2,1)=273.1
1879 XNO(2,2)=421.8
1880 XNO(2,3)=368.9
1881 XNO(3,2)=473.8
1882 XNO(3,3)=-304.7
1883 XN1(1,2)=-106.1
1884 XN1(2,1)=-182.4
1885 XN1(2,2)=-474.5
1886 XN1(2,3)=-522.2
1887 XN1(3,2)=-468.2
1888 XN1(3,3)=-470.
1889 XN2(1,2)=1280.8
1890 XN2(2,1)=-1.872
1891 XN2(2,2)=-3.592
1892 XN2(2,3)=-4.998
1893 XN2(3,2)=-2.225
1894 XN2(3,3)=-8.582
1895 RETURN
1896 END

```

APPENDIX B

SAMPLE INPUT

```

1   1
    38.00    38.00
    33.50     5.05
    26.00    28.00
    14.08     5.05      8.20      0.00
    0.0
    0.00
  966.00    780.

16
    1.00    740.00     0.00     0.00
    2.00   1651.00      4.00     0.00
    3.00   1484.00     36.00     0.00
    4.00   1383.00    123.00     0.00
    5.00   1372.00    258.00     2.00
    6.00   1383.00    402.00    26.00
    7.00   1388.00    517.00   140.00
    8.00   1384.00    596.00   350.00
    9.00   1374.00    655.00   570.00
   10.00   1361.00    702.00   730.00
   12.00   1340.00    779.00   980.00
   14.00   1326.00    834.00  1170.00
   16.00   1315.00    881.00  1300.00
   20.00   1282.00    932.00  1420.00
   24.00   1263.00    957.00  1500.00
   26.00   1225.00    966.00  1520.00

2   0
2   0
    4.23
    1.35
-1   0
1
1   0
2   1
    4.23
    1.35
-1   0
0
    66.00     5.05
    26.00    28.0
    14.08     5.05      8.20      0.00
    0.00
    0.00
    0.00
  30      2.00      2.00
1   0
0
    0.00
    0.00
1   1
3
    16.21
    0.00
24      2.00      2.00
1   2
2
    20.63
    0.00
22      3.00      2.00
2   1
2
    18.24
    0.00
21      6.00      2.00
-1   0
0
    20.04     4.79
    92.00   143.00
    5.69     4.79      -5.24      5.75
    0.00
    0.00
    0.00
  24      1.00
1   0
0
    16.62    16.57     17.13     17.49     17.97
-1   0
-1

```

APPENDIX C

SAMPLE OUTPUT

```
1  
-S= 10.45 MEV  
-OCCUPATION PROBABILITIES PO= 2. HO= 0  
TARGET Z= 26, N= 28 PROJ Z= 2, N= 0  
G= 4.230, E=33.500  
SCALE FACTOR= 1.350, FRAC PREEO= 0.043  
- P H RHOU/RHO STRU/STR STRD/STR  
  
2 0 9.896E-01 9.896E-01 9.896E-01  
3 1 8.373E-01 8.206E-01 8.206E-01  
4 2 5.637E-01 5.722E-01 5.240E-01  
5 3 3.264E-01 3.079E-01 2.333E-01  
6 4 1.509E-01 1.353E-01 6.615E-02  
7 5 5.744E-02 5.872E-02 1.043E-02
```

1PARTICLE SPECTRA Z= 0, N= 1
 FIRST EMISSION AT P= 3
 TARGET Z= 26, N= 28 PROJ Z= 2, N= 0
 PO= 2, HO= 0, G= 4.230, E=33.500
 PAIR EXCIT. ACCORDING TO 2 COMP. ST. DENSITIES
 REACTION CROSS SECTION = 966.0
 SCALE FACTOR= 1.350, FRAC PREEQ= 0.043
 MSD POSSIBLE= 0.990, USED= 0.006
 CLOSED FORM SUM STARTS AT P= 3
 - EPS -- DIRECT -- PREEQUILIBRIUM -- EQUIL -- TOTAL (MB/MEV)
 -- NUTRA KNOCK -- MSD MSC -- WEISS -- MSD MSC MSD+MSC
 1.00 8.383E-03 2.629E-02 0. 0. 4.379E+00 3.467E-02 4.379E+00 4.413E+00
 2.00 3.314E-02 1.103E-01 0. 0. 1.114E+01 1.435E-01 1.114E+01 1.128E+01
 3.00 3.928E-02 1.393E-01 0. 0. 8.440E+00 1.786E-01 8.440E+00 8.618E+00
 4.00 4.252E-02 1.614E-01 0. 0. 5.799E+00 2.039E-01 5.799E+00 6.003E+00
 5.00 4.547E-02 1.857E-01 0. 0. 3.908E+00 2.311E-01 3.908E+00 4.139E+00
 6.00 4.688E-02 2.070E-01 0. 0. 2.520E+00 2.539E-01 2.520E+00 2.774E+00
 7.00 4.614E-02 2.218E-01 0. 0. 1.541E+00 2.680E-01 1.541E+00 1.809E+00
 8.00 4.347E-02 2.294E-01 0. 0. 8.953E-01 2.729E-01 8.953E-01 1.168E+00
 9.00 3.934E-02 2.300E-01 0. 0. 4.967E-01 2.694E-01 4.967E-01 7.661E-01
 10.00 3.423E-02 2.244E-01 0. 0. 2.636E-01 2.586E-01 2.636E-01 5.222E-01
 12.00 2.278E-02 1.971E-01 0. 0. 6.503E-02 2.199E-01 6.503E-02 2.849E-01
 14.00 1.173E-02 1.490E-01 0. 0. 1.297E-02 1.607E-01 1.297E-02 1.737E-01
 16.00 3.353E-03 7.988E-02 0. 0. 1.907E-03 8.323E-02 1.907E-03 8.513E-02
 OSUMS 4.750E-01 2.780E+00 0. 0. 3.968E+01

UNANGULAR DISTRIBUTIONS Z = 0 N = 1
 ENERGY, PARAM = EPS+E.E.
 CROSS SECTIONS IN MEV-STR-MEV
 EDS

1PARTICLE SPECTRA Z= 1, N= 0
 FIRST EMISSION AT P= 3
 TARGET Z= 26, N= 28 PROJ Z= 2, N= 0
 PO= 2, HO= 0, G= 4.230, E=33.500
 PAIR EXCIT. ACCORDING TO 2 COMP. ST. DENSITIES
 REACTION CROSS SECTION = 966.0
 SCALE FACTOR= 1.350, FRAC PREEQ= 0.043
 MSD POSSIBLE= 0.990, USED= 0.006
 CLOSED FORM SUM STARTS AT P= 3

- EPS	-- DIRECT	-- PREEQUILIBRIUM	-- EQUIL	TOTAL (MB/MEV)				
				-- NUTRA	KNOCK	MSD	MSC	MSD+MSC
2.00	6.902E-05	3.742E-04	0.	0.	2.823E+00	4.433E-04	2.823E+00	2.823E+00
3.00	9.317E-04	4.831E-03	0.	0.	2.370E+01	5.763E-03	2.370E+01	2.371E+01
4.00	4.244E-03	2.100E-02	0.	0.	6.660E+01	2.524E-02	6.660E+01	6.663E+01
5.00	1.113E-02	5.242E-02	0.	0.	1.067E+02	6.355E-02	1.067E+02	1.068E+02
6.00	2.081E-02	9.306E-02	0.	0.	1.208E+02	1.139E-01	1.208E+02	1.209E+02
7.00	3.122E-02	1.322E-01	0.	0.	1.085E+02	1.634E-01	1.085E+02	1.087E+02
8.00	4.113E-02	1.644E-01	0.	0.	8.473E+01	2.056E-01	8.473E+01	8.493E+01
9.00	5.086E-02	1.912E-01	0.	0.	6.134E+01	2.421E-01	6.134E+01	6.158E+01
10.00	6.056E-02	2.133E-01	0.	0.	4.224E+01	2.739E-01	4.224E+01	4.251E+01
12.00	8.065E-02	2.458E-01	0.	0.	1.803E+01	3.254E-01	1.803E+01	1.836E+01
14.00	1.007E-01	2.591E-01	0.	0.	6.776E+00	3.598E-01	6.776E+00	7.136E+00
16.00	1.216E-01	2.550E-01	0.	0.	2.280E+00	3.766E-01	2.280E+00	2.657E+00
20.00	1.608E-01	1.845E-01	0.	0.	1.723E-01	3.453E-01	1.723E-01	5.176E-01
24.00	1.981E-01	3.907E-02	0.	0.	6.234E-03	2.372E-01	6.234E-03	2.434E-01
26.00	2.167E-01	0.	0.	0.	7.194E-04	2.167E-01	7.194E-04	2.174E-01
OSUMS	2.541E+00	3.610E+00	0.	0.	6.957F+02			

ANGULAR DISTRIBUTIONS Z = 1 N = 0		ENERGY PARAM. = EPS+E.E.									
CROSS SECTIONS IN MEV/STR-MEV		EPS									
		0	10	30	40	50	60	70	80	90	
2.00	2.68E-01	2.66E-01	2.65E-01	2.52E-C1	2.41E-C	2.30E-01	2.19E-01	2.11E-01	2.05E-01	2.03E-01	1.95E-01
3.00	2.28E+00	2.26E+00	2.24E+00	2.13E+00	2.03E+00	1.93E+00	1.84E+00	1.76E+00	1.71E+00	1.70E+00	1.70E+00
4.00	6.46E+00	6.41E+00	5.25E+00	6.02E+00	5.73E+00	5.43E+00	5.15E+00	4.93E+00	4.79E+00	4.74E+00	4.74E+00
5.00	1.0E+01	1.04E+01	1.0E+01	9.71E+00	9.23E+00	8.72E+00	8.25E+00	7.87E+00	7.63E+00	7.54E+00	7.54E+00
6.00	1.20E+01	1.19E+01	1.16E+01	1.11E+01	1.05E+01	9.88E+00	9.32E+00	8.87E+00	8.38E+00	8.48E+00	8.48E+00
7.00	1.09E+01	1.08E+01	1.07E+01	1.00E+01	9.48E+00	8.80E+00	8.37E+00	7.94E+00	7.66E+00	7.56E+00	7.56E+00
8.00	8.64E+00	8.55E+00	8.29E+00	7.91E+00	7.45E+00	6.97E+00	6.53E+00	6.17E+00	5.94E+00	5.86E+00	5.86E+00
9.00	6.33E+00	6.28E+00	6.09E+00	5.79E+00	5.44E+00	5.07E+00	4.72E+00	4.45E+00	4.28E+00	4.21E+00	4.21E+00
10.00	4.45E+00	4.40E+00	4.26E+00	4.04E+00	3.78E+00	3.51E+00	3.26E+00	3.06E+00	2.93E+00	2.88E+00	2.88E+00
12.00	1.99E+00	1.97E+00	1.90E+00	1.79E+00	1.66E+00	1.53E+00	1.41E+00	1.31E+00	1.22E+00	1.22E+00	1.22E+00
14.00	8.20E-01	8.08E-01	7.76E-01	7.28E-01	6.69E-01	6.09E-01	5.53E-01	5.07E-01	4.77E-01	4.63E-01	4.63E-01
16.00	3.38E-01	3.33E-01	3.18E-01	2.96E-01	2.41E-01	2.14E-01	1.92E-01	1.77E-01	1.68E-01	1.68E-01	1.68E-01
20.00	1.00E-01	9.84E-02	9.28E-02	8.44E-02	7.42E-02	6.54E-02	5.30E-02	4.24E-02	3.08E-02	2.08E-02	2.08E-02
24.00	6.13E-02	5.99E-02	5.59E-02	5.00E-02	4.30E-02	3.56E-02	2.86E-02	2.24E-02	1.73E-02	1.34E-02	1.34E-02
26.00	5.84E-02	5.70E-02	5.30E-02	4.71E-02	4.00E-02	3.27E-02	2.58E-02	2.05E-02	1.51E-02	1.14E-02	1.14E-02
-	EPS	100	110	120	130	140	150	160	170	180	TOTAL

1PARTICLE SPECTRA Z= 2, N= 2

FIRST EMISSION AT P= 4

TARGET Z= 26, N= 28 PROD Z= 2, N= 0

PO= 2, HO= 0, G= 4.230, E=33.500

PAIR EXCIT. ACCORDING TO 2 COMP. ST. DENS TIES

REACTION CROSS SECTION = 966.0

SCALE FACTOR= 350, FRAC PREEQ= 0.043

MSD POSSIBLE= 0.990, USED= 0.006

CLOSED FORM SUM STARTS AT P= 4

	EPS	DIRECT	PREEQUILIBRIUM	EQUIL	TOTAL	(MB/MEV)
	--	--	--	--	MSC	MSD+MSC
	--	NUTRA	KNOCK	--	WEISS	--
5.00	3.109E-04	4.894E-05	4.307E-03	6.932E-02	1.531E-01	4.667E-03
6.00	4.596E-03	7.038E-04	4.898E-02	7.152E-01	1.419E+00	5.428E-02
7.00	2.727E-02	4.046E-03	2.225E-01	2.903E+00	5.230E+00	2.539E-01
8.00	1.335E-02	1.049E-02	4.563E-01	5.222E+00	8.652E+00	5.402E-01
9.00	1.260E-01	1.726E-02	5.957E-01	5.857E+00	9.049E+00	7.390E-01
10.00	1.674E-01	2.177E-02	5.999E-01	4.948E+00	7.237E+00	7.891E-01
12.00	2.313E-01	2.608E-02	4.754E-01	2.543E+00	3.493E+00	7.328E-01
14.00	2.687E-01	2.380E-02	3.191E-01	9.870E-01	1.342E+00	6.116E-01
16.00	2.733E-01	0.	1.915E-01	3.163E-01	4.241E-01	4.648E-01
20.00	1.877E-01	0.	5.414E-02	2.721E-02	2.273E-02	2.419E-01
24.00	0.	0.	6.450E-03	1.521E-03	3.195E-04	6.450E-03
OSUMS	3.053E+00	1.649E-01	4.627E+00	3.031E+01	4.639E+01	

:ANGULAR DISTRIBUTIONS Z= 2 N= 2

ENERGY PARAM. = EPS+B.E.

CROSS SECTIONS IN MB/STR-MEV

- EPS	0	10	20	30	40	50	60	70	80	90	
5.00	2.34E-02	2.32E-02	2.25E-02	2.14E-02	2.01E-02	1.88E-02	1.76E-02	1.66E-02	1.59E-02	1.56E-02	
6.00	2.29E-01	2.27E-01	2.20E-01	2.09E-01	1.96E-01	1.82E-01	1.69E-01	1.59E-01	1.52E-01	1.49E-01	
7.00	8.95E-01	8.85E-01	8.55E-01	8.11E-01	7.57E-01	7.01E-01	6.49E-01	6.07E-01	5.79E-01	5.68E-01	
8.00	1.57E+00	1.55E+00	1.50E+00	1.41E+00	1.32E+00	1.21E+00	1.12E+00	1.04E+00	9.88E-01	9.66E-01	
9.00	1.74E+00	1.72E+00	1.66E+00	1.56E+00	1.45E+00	1.33E+00	1.22E+00	1.13E+00	1.06E+00	1.04E+00	
10.00	1.48E+00	1.46E+00	1.41E+00	1.32E+00	1.22E+00	1.11E+00	1.01E+00	9.31E-01	8.76E-01	8.51E-01	
12.00	8.28E-01	8.16E-01	7.81E-01	7.29E-01	6.65E-01	5.99E-01	5.38E-01	4.87E-01	4.51E-01	4.33E-01	
14.00	4.02E-01	3.95E-01	3.76E-01	3.48E-01	3.14E-01	2.78E-01	2.45E-01	2.16E-01	1.95E-01	1.83E-01	
16.00	1.94E-01	1.90E-01	1.80E-01	1.65E-01	1.47E-01	1.27E-01	1.09E-01	9.32E-02	8.10E-02	7.27E-02	
20.00	6.68E-02	6.53E-02	6.12E-02	5.49E-02	4.75E-02	3.96E-02	3.22E-C2	2.57E-02	2.04E-02	1.63E-02	
24.00	2.06E-03	2.00E-03	1.86E-03	1.65E-03	1.40E-03	1.14E-03	8.99E-04	6.97E-04	5.38E-04	4.23E-04	
- EPS	100	110	120	130	140	150	160	170	180	TOTAL	
5.00	1.58E-02	1.64E-02	1.73E-02	1.84E-02	1.97E-02	2.09E-02	2.19E-02	2.25E-02	2.28E-02	2.27E-01	
6.00	1.51E-01	1.57E-01	1.66E-01	1.77E-01	1.90E-01	2.02E-01	2.12E-01	2.19E-01	2.21E-01	2.19E+00	
7.00	5.73E-01	5.96E-01	6.32E-01	6.78E-C1	7.29E-01	7.78E-01	8.19E-01	8.46E-01	8.50E-01	8.39E+00	
8.00	9.75E-01	1.01E+00	1.08E+00	1.16E+00	1.25E+00	1.34E+00	1.42E+00	1.47E+00	1.48E+00	1.44E+01	
9.00	1.05E+00	1.09E+00	1.16E+00	1.26E+00	1.36E+00	1.46E+00	1.54E+00	1.60E+00	1.62E+00	1.56E+01	
10.00	8.57E-01	8.93E-01	9.54E-01	1.03E+00	1.12E+00	1.21E+00	1.28E+00	1.33E+C0	1.35E+00	1.30E+01	
12.00	4.33E-01	4.50E-01	4.81E-01	5.24E-01	5.72E-01	6.19E-01	6.59E-01	6.85E-01	6.95E-01	6.77E+00	
14.00	1.79E-01	1.84E-01	1.95E-01	2.12E-01	2.31E-01	2.51E-01	2.68E-01	2.79E-01	2.83E-01	2.94E+00	
16.00	6.23E-02	6.76E-02	6.98E-02	7.44E-02	8.04E-02	8.67E-02	9.22E-02	9.60E-02	9.73E-02	1.21E+00	
20.00	1.34E-02	1.15E-02	1.02E-02	9.62E-03	9.39E-03	9.40E-03	9.52E-03	9.64E-03	9.69E-03	2.92E-01	
24.00	3.45E-04	2.99E-04	2.77E-04	2.73E-04	2.81E-04	2.96E-04	3.11E-04	3.22E-04	3.26E-04	8.29E-03	

1
-S= 10.28 MEV
-OCCUPATION PROBABILITIES PO= 2, HO= 1
TARGET Z= 26, N= 28 PROJ Z= 1, N= 0
G= 4.230, E=33.500
SCALE FACTOR= 1.350, FRAC PREEQ= 0.537
- P H RHOU/RHO STRU/STR STRD/STR

1 0 1.000E+00 0.
2 1 8.680E-01 8.680E-01 8.680E-01
3 2 6.725E-01 6.594E-01 5.956E-01
4 3 4.489E-01 4.075E-01 3.082E-01
5 4 2.498E-01 2.030E-01 1.112E-01
6 5 1.201E-01 9.209E-02 2.593E-02
7 6 4.876E-02 4.558E-02 3.496E-03

80

PARTICLE SPECTRA Z= 0, N= 1
 FIRST EMISSION AT P= 2
 TARGET Z= 26, N= 28 PROJ Z= 1, N= 0
 PO= 2, HO= 1, G= 4.230, E=33.500
 PAIR EXCIT. ACCORDING TO 2 COMP. ST. DENSITIES
 REACTION CROSS SECTION = 966.0
 SCALE FACTOR= 1.350, FRAC PREEQ= 0.537
 MSD POSSIBLE= 0.868, USED= 0.405
 CLOSED FORM SUM STARTS AT P= 2

			PREEQUILIBRIUM			EQUIL			TOTAL (MB/MEV)		
- EPS --	DIRECT	--	--	MSD	MSC	--	WEISS	--	MSD	MSC	MSD+MSC
1.00	0.	0.	1.134E+00	6.178E-01	2.381E+00	1.134E+00	2.999E+00	4.133E+00			
2.00	0.	0.	4.321E+00	1.756E+00	6.059E+00	4.321E+00	7.815E+00	1.214E+01			
3.00	0.	0.	4.993E+00	1.527E+00	4.591E+00	4.993E+00	6.117E+00	1.111E+01			
4.00	0.	0.	5.335E+00	1.240E+00	3.155E+00	5.335E+00	4.395E+00	9.730E+00			
5.00	0.	0.	5.704E+00	1.020E+00	2.126E+00	5.704E+00	3.146E+00	8.850E+00			
6.00	0.	0.	5.958E+00	8.293E-01	1.371E+00	5.958E+00	2.201E+00	8.159E+00			
7.00	0.	0.	6.030E+00	6.594E-01	8.384E-01	6.030E+00	1.498E+00	7.528E+00			
8.00	0.	0.	5.938E+00	5.138E-01	4.873E-01	5.938E+00	1.001E+00	6.939E+00			
9.00	0.	0.	5.722E+00	3.935E-01	2.704E-01	5.722E+00	6.639E-01	6.385E+00			
10.00	0.	0.	5.415E+00	2.967E-01	1.435E-01	5.415E+00	4.402E-01	5.855E+00			
12.00	0.	0.	4.637E+00	1.613E-01	3.541E-02	4.637E+00	1.968E-01	4.833E+00			
14.00	0.	0.	3.677E+00	8.137E-02	7.067E-03	3.677E+00	8.844E-02	3.765E+00			
16.00	0.	0.	2.524E+00	3.672E-02	1.039E-03	2.524E+00	3.776E-02	2.562E+00			
OSUMS	0.	0.	7.746E+01	9.598E+00	2.158E+01						

1 ANGULAR DISTRIBUTIONS Z= 0 N= 1

ENERGY PARAM. = EPS+B.E

CROSS SECTIONS IN MB/STR-MEV

- EPS	0	10	20	30	40	50	60	70	80	90
1.00	5.18E-01	5.11E-01	4.92E-01	4.63E-01	4.28E-01	3.90E-01	3.54E-01	3.22E-01	2.97E-01	2.80E-01
2.00	1.63E+00	1.60E+00	1.54E+00	1.44E+00	1.33E+00	1.20E+00	1.07E+00	9.65E-01	8.77E-01	8.14E-01
3.00	1.60E+00	1.58E+00	1.51E+00	1.41E+00	1.29E+00	1.15E+00	1.02E+00	9.05E-01	8.09E-01	7.38E-01
4.00	1.51E+00	1.49E+00	1.42E+00	1.32E+00	1.20E+00	1.06E+00	9.32E-01	8.14E-01	7.15E-01	6.39E-01
5.00	1.48E+00	1.45E+00	1.39E+00	1.28E+00	1.15E+00	1.02E+00	8.82E-01	7.60E-01	6.56E-01	5.74E-01
6.00	1.46E+00	1.43E+00	1.36E+00	1.25E+00	1.12E+00	9.82E-01	8.43E-01	7.17E-01	6.09E-01	5.23E-01
7.00	1.42E+00	1.40E+00	1.33E+00	1.22E+00	1.09E+00	9.43E-01	8.02E-01	6.74E-01	5.64E-01	4.76E-01
8.00	1.38E+00	1.36E+00	1.28E+00	1.18E+00	1.04E+00	8.98E-01	7.58E-01	6.30E-01	5.21E-01	4.33E-01
9.00	1.33E+00	1.30E+00	1.23E+00	1.12E+00	9.91E-01	8.49E-01	7.11E-01	5.86E-01	4.79E-01	3.92E-01
10.00	1.27E+00	1.24E+00	1.17E+00	1.06E+00	9.35E-01	7.96E-01	6.62E-01	5.40E-01	4.37E-01	3.54E-01
12.00	1.12E+00	1.09E+00	1.03E+00	9.29E-01	8.08E-01	6.81E-01	5.58E-01	4.48E-01	3.56E-01	2.82E-01
14.00	9.24E-01	9.04E-01	8.46E-01	7.59E-01	6.55E-01	5.45E-01	4.41E-01	3.48E-01	2.72E-01	2.11E-01
16.00	6.66E-01	6.51E-01	6.07E-01	5.41E-01	4.62E-01	3.80E-01	3.03E-01	2.35E-01	1.80E-01	1.38E-01

- EPS	100	110	120	130	140	150	160	170	180	TOTAL
1.00	2.72E-01	2.71E-01	2.78E-01	2.89E-01	3.04E-01	3.19E-01	3.32E-01	3.41E-01	3.44E-01	4.13E+00
2.00	7.78E-01	7.66E-01	7.75E-01	8.01E-01	8.37E-01	8.75E-01	9.08E-01	9.31E-01	9.39E-01	1.21E+01
3.00	6.91E-01	6.69E-01	6.67E-01	6.81E-01	7.04E-01	7.32E-01	7.57E-01	7.75E-01	7.81E-01	1.11E+01
4.00	5.86E-01	5.55E-01	5.43E-01	5.45E-01	5.57E-01	5.74E-01	5.90E-01	6.01E-01	6.06E-01	9.73E+00
5.00	5.15E-01	4.76E-01	4.55E-01	4.48E-01	4.51E-01	4.59E-01	4.68E-01	4.75E-01	4.77E-01	8.85E+00
6.00	4.59E-01	4.14E-01	3.86E-01	3.72E-01	3.67E-01	3.68E-01	3.71E-01	3.74E-01	3.76E-01	8.15E+00
7.00	4.09E-01	3.61E-01	3.28E-01	3.09E-01	2.98E-01	2.94E-01	2.93E-01	2.94E-01	2.94E-01	7.53E+00
8.00	3.65E-01	3.15E-01	2.80E-01	2.58E-01	2.44E-01	2.36E-01	2.33E-01	2.31E-01	2.31E-01	6.94E+00
9.00	3.26E-01	2.76E-01	2.41E-01	2.17E-01	2.01E-01	1.92E-01	1.86E-01	1.83E-01	1.83E-01	5.39E+00
10.00	2.90E-01	2.42E-01	2.07E-01	1.83E-01	1.67E-01	1.56E-01	1.50E-01	1.47E-01	1.46E-01	5.85E+00
12.00	2.26E-01	1.84E-01	1.53E-01	1.31E-01	1.16E-01	1.06E-01	9.93E-02	9.57E-02	9.45E-02	4.83E+00
14.00	1.66E-01	1.32E-01	1.08E-01	9.01E-02	7.78E-02	6.94E-02	6.39E-02	6.08E-02	5.98E-02	3.77E+00
16.00	1.06E-01	8.28E-02	6.62E-02	5.44E-02	4.61E-02	4.04E-02	3.66E-02	3.44E-02	3.38E-02	2.56E+00

8
N

```

1PARTICLE SPECTRA Z= 1. N= 0
FIRST EMISSION AT P= 2
TARGET Z= 26, N= 28      PROJ Z= 1, N= 0
PO= 2., HO= 1, G= 4.230, E=33.500
PAIR EXCIT. ACCORDING TO 2 COMP. ST. DENSITIES
REACTION CROSS SECTION = 966.0
SCALE FACTOR= 1.350, FRAC PREEQ= 0.537
MSD POSSIBLE= 0.868, USED= 0.405
CLOSED FORM SUM STARTS AT P= 2

```

- EPS --		DIRECT		PREEQUILIBRIUM		EQUIL		TOTAL		(MB/MEV)	
--	NUTRA	KNOCK	--	MSD	MSC	--	WEISS	--	MSD	MSC	MSD+MSC
2.00	O.	7.104E-06	7.376E-02	3.037E-01	1.534E+00	7.377E-02	1.838E+00	1.911E+00			
3.00	O.	9.228E-05	8.229E-01	2.677E+00	1.288E+01	6.330E-01	1.556E+01	1.639E+01			
4.00	O.	4.039E-04	3.191E+00	7.890E+00	3.620E+01	3.191E+00	4.409E+01	4.728E+01			
5.00	O.	1.016E-03	7.073E+00	1.327E+01	5.801E+01	7.074E+00	7.128E+01	7.836E+01			
6.00	O.	1.818E-03	1.124E+01	1.582E+01	6.565E+01	1.124E+01	8.148E+01	9.272E+01			
7.00	O.	2.606E-03	1.441E+01	1.508E+01	5.901E+01	1.441E+01	7.409E+01	8.850E+01			
8.00	O.	3.273E-03	1.630E+01	1.261E+01	4.607E+01	1.631E+01	5.867E+01	7.498E+01			
9.00	O.	3.849E-03	1.739E+01	9.917E+00	3.335E+01	1.739E+01	4.327E+01	6.066E+01			
10.00	O.	4.348E-03	1.794E+01	7.566E+00	2.297E+01	1.794E+01	3.054E+01	4.848E+01			
12.00	O.	5.162E-03	1.813E+01	4.282E+00	9.810E+00	1.813E+01	1.409E+01	3.222E+01			
14.00	O.	5.664E-03	1.736E+01	2.411E+00	3.687E+00	1.736E+01	6.098E+00	2.346E+01			
16.00	O.	5.891E-03	1.612E+01	1.384E+00	1.241E+00	1.613E+01	2.624E+00	1.875E+01			
20.00	O.	5.286E-03	1.227E+01	4.382E-01	9.379E-02	1.228E+01	5.320E-01	1.281E+01			
24.00	O.	3.428E-03	7.248E+00	1.163E-01	3.397E-03	7.252E+00	1.197E-01	7.372E+00			
26.00	O.	O.	4.263E+00	5.051E-02	3.921E-04	4.263E+00	5.091E-02	4.314E+00			
OSUMS	O.	9.034E-02	2.940E+02	1.086E+02	3.783E+02						

1ANGULAR DISTRIBUTIONS Z= 1 N= 0

ENERGY PARAM = EPS+P.E.

CROSS SECTIONS IN MC/STR-MEV

- EPS	0	10	20	30	40	50	60	70	80	90	.
2.00	1.86E-01	1.85E-01	1.80E-01	1.74E-01	1.66E-01	1.58E-01	1.50E-01	1.44E-01	1.40E-01	1.38E-01	
3.00	1.62E+00	1.61E+00	1.57E+00	1.51E+00	1.44E+00	1.36E+00	1.29E+00	1.23E+00	1.19E+00	1.17E+00	
4.00	4.78E+00	4.74E+00	4.62E+00	4.44E+00	4.21E+00	3.97E+00	3.75E+00	3.56E+00	3.43E+00	3.36E+00	
5.00	8.14E+00	8.06E+00	7.84E+00	7.51E+00	7.10E+00	6.67E+00	6.26E+00	5.91E+00	5.67E+00	5.53E+00	
6.00	9.94E+00	9.84E+00	9.56E+00	9.12E+00	8.59E+00	8.02E+00	7.48E+00	7.03E+00	6.69E+00	6.50E+00	
7.00	9.87E+00	9.77E+00	9.46E+00	9.00E+00	8.43E+00	7.83E+00	7.25E+00	6.76E+00	6.38E+00	6.16E+00	
8.00	8.78E+00	8.68E+00	8.39E+00	7.95E+00	7.41E+00	6.83E+00	6.28E+00	5.79E+00	5.42E+00	5.17E+00	
9.00	7.52E+00	7.43E+00	7.17E+00	6.76E+00	6.27E+00	5.73E+00	5.22E+00	4.76E+00	4.40E+00	4.15E+00	
10.00	6.43E+00	6.34E+00	6.10E+00	5.73E+00	5.28E+00	4.79E+00	4.31E+00	3.88E+00	3.54E+00	3.29E+00	
12.00	4.94E+00	4.87E+00	4.66E+00	4.34E+00	3.94E+00	3.51E+00	3.09E+00	2.71E+00	2.39E+00	2.14E+00	
14.00	4.11E+00	4.05E+00	3.85E+00	3.56E+00	3.20E+00	2.81E+00	2.42E+00	2.07E+00	1.77E+00	1.52E+00	
16.00	3.65E+00	3.59E+00	3.40E+00	3.12E+00	2.78E+00	2.41E+00	2.04E+00	1.71E+00	1.42E+00	1.19E+00	
20.00	2.88E+00	2.82E+00	2.66E+00	2.41E+00	2.11E+00	1.78E+00	1.47E+00	1.19E+00	9.84E-01	7.62E-01	
24.00	1.86E+00	1.82E+00	1.70E+00	1.52E+00	1.31E+00	1.08E+00	8.68E-01	6.80E-01	5.26E-01	4.06E-01	
26.00	1.15E+00	1.13E+00	1.05E+00	9.30E-01	7.91E-01	6.46E-01	5.11E-01	3.94E-01	2.89E-01	2.27E-01	
- EPS	100	110	120	130	140	150	160	170	180	TOTAL	

2.00	1.38E-01	1.41E-01	1.46E-01	1.53E-01	1.60E-01	1.67E-01	1.72E-01	1.76E-01	1.77E-01	1.91E+00	
3.00	1.18E+00	1.20E+00	1.25E+00	1.30E+00	1.37E+00	1.43E+00	1.48E+00	1.51E+00	1.52E+00	1.64E+01	
4.00	3.37E+00	3.44E+00	3.57E+00	3.73E+00	3.92E+00	4.10E+00	4.25E+00	4.35E+00	4.29E+00	4.73E+01	
5.00	5.53E+00	5.64E+00	5.84E+00	6.12E+00	6.44E+00	6.74E+00	7.00E+00	7.17E+00	7.12E+00	7.84E+01	
6.00	6.46E+00	6.57E+00	6.81E+00	7.13E+00	7.50E+00	7.86E+00	8.17E+00	8.37E+00	8.44E+00	9.27E+01	
7.00	6.09E+00	6.16E+00	6.36E+00	6.65E+00	7.00E+00	7.33E+00	7.62E+00	7.81E+00	7.88E+00	8.85E+01	
8.00	5.07E+00	5.10E+00	5.24E+00	5.46E+00	5.73E+00	6.00E+00	6.23E+00	6.39E+00	6.44E+00	7.50E+01	
9.00	4.02E+00	4.06E+00	4.08E+00	4.23E+00	4.42E+00	4.52E+00	4.79E+00	4.91E+00	4.95E+00	6.07E+01	
10.00	3.13E+00	3.03E+00	3.10E+00	3.19E+00	3.21E+00	3.44E+00	3.56E+00	3.64E+00	3.67E+00	4.85E+01	
12.00	1.96E+00	1.85E+00	1.81E+00	1.80E+00	1.83E+00	1.88E+00	1.92E+00	1.95E+00	1.96E+00	3.22E+01	
14.00	1.34E+00	1.21E+00	1.12E+00	1.08E+00	1.06E+00	1.05E+00	1.06E+00	1.06E+00	1.07E+00	1.35E+01	
16.00	1.00E+00	8.70E-01	7.75E-01	7.11E-01	6.71E-01	6.48E-01	6.36E-01	6.31E-01	6.29E-01	1.88E+01	
20.00	6.14E-01	5.03E-01	4.21E-01	3.63E-01	3.22E-01	2.94E-01	2.76E-01	2.66E-01	2.63E-01	1.28E+01	
24.00	3.15E-01	2.48E-01	2.00E-01	1.65E-01	1.41E-01	1.24E-01	1.13E-01	1.07E-01	1.05E-01	7.37E+00	
26.00	1.73E-01	1.34E-01	1.06E-01	8.65E-02	7.26E-02	6.30E-02	5.67E-02	5.31E-02	5.19E-02	4.31E+00	

1PARTICLE SPECTRA Z= 2, N= 2

FIRST EMISSION AT P= 4

TARGET Z= 26, N= 2E PROJ Z= 1, N= 0

PO= 2, HQ= 1, G= 4 230, E=33.500

PAIR EXCIT. ACCORDING TO 2 COMP. ST. DENSITIES

REACTION CROSS SECTION = 966.0

SCALE FACTOR= 1.350, FRAC PREEQ= 0.537

MSD POSSIBLE= 0.868, USED= 0.405

CLOSED FORM SUM STARTS AT P= 4

EPS	DIRECT		PREEQUILIBRIUM		EQUIL		TOTAL (ME/MEV)		
	NUTRA	KNOCK	MSD	MSC	WEISE	MSD	MSC	MSD+MSC	
5.00	2.009E-03	7.507E-05	4.508E-04	8.793E-03	8.224E-02	2.545E-03	9.103E-01	9.357E-02	
6.00	2.809E-02	1.098E-03	5.232E-03	8.289E-02	7.624E-01	3.442E-02	8.453E-01	8.797E-01	
7.00	1.572E-01	6.433E-03	2.383E-02	3.071E-01	2.812E+00	1.575E-01	3.119E+00	3.306E+00	
8.00	3.972E-01	1.706E-02	4.920E-02	5.051E-01	4.654E+00	4.635E-01	5.159E+00	5.622E+00	
9.00	6.385E-01	2.883E-02	6.490E-02	5.208E-01	4.869E+00	7.323E-01	5.390E+00	6.122E+00	
10.00	7.899E-01	2.759E-02	6.624E-02	4.090E-01	3.895E+00	8.938E-01	4.304E+00	5.198E+00	
12.00	9.302E-01	1.946E-02	5.420E-02	1.934E-01	1.882E+00	1.034E+00	2.076E+00	3.110E+00	
14.00	8.933E-01	5.345E-02	3.745E-02	7.912E-02	7.240E-01	9.842E-01	8.031E-01	1.787E+00	
16.00	7.183E-01	4.825E-02	2.268E-02	3.051E-02	2.290E-01	7.892E-01	2.595E-01	1.049E+00	
20.00	2.321E-01	0.	5.071E-03	3.595E-03	1.231E-02	2.372E-01	1.590E-02	2.531E-01	
24.00	0.	0.	0.	0.	1.737E-04	0.	1.737E-04	1.737E-04	
OSUMS	9.138E+00	4.605E-01	5.146E-01	2.689E+00	2.497E+01				

ANGULAR DISTRIBUTIONS Z = 2 N= 2		ENERGY PARAM. = EPS+E/E CROSS SECTIONS IN MEV/SIR-MEV									
-EPS	0	10	20	30	40	50	60	70	80	90	100
5.00	9.70E-02	9.60E-03	9.31E-03	8.87E-03	8.34E-03	7.78E-03	7.26E-03	6.84E-03	6.46E-03	6.16E-03	6.45E-03
6.00	9.32E-02	9.21E-02	8.92E-02	8.47E-02	7.94E-02	7.37E-02	6.85E-02	6.42E-02	6.01E-02	5.60E-02	6.01E-02
7.00	3.60E-01	3.55E-01	3.43E-01	3.25E-01	3.03E-01	2.80E-01	2.59E-01	2.41E-01	2.29E-01	2.24E-01	2.24E-01
8.00	6.32E-01	6.24E-01	6.02E-01	5.68E-01	5.27E-01	4.85E-01	4.44E-01	4.14E-01	3.88E-01	3.77E-01	3.77E-01
9.00	7.18E-01	7.08E-01	6.81E-01	6.41E-01	5.92E-01	5.40E-01	4.91E-01	4.51E-01	4.22E-01	4.06E-01	4.06E-01
10.00	6.43E-01	6.34E-01	6.08E-01	5.70E-01	5.23E-01	4.73E-01	4.27E-01	3.87E-01	3.58E-01	3.41E-01	3.41E-01
12.00	4.44E-01	4.37E-01	4.17E-01	3.87E-01	3.50E-01	3.11E-01	2.74E-01	2.42E-01	2.16E-01	1.99E-01	1.99E-01
14.00	3.04E-01	2.99E-01	2.83E-01	2.60E-01	2.32E-01	2.02E-01	1.73E-01	1.48E-01	1.27E-01	1.11E-01	1.11E-01
16.00	2.09E-01	1.93E-01	1.76E-01	1.55E-01	1.33E-01	1.11E-01	9.18E-02	7.58E-02	6.33E-02	5.33E-02	5.33E-02
20.00	6.12E-02	5.98E-02	5.60E-02	5.2C2E-02	4.33E-02	3.61E-02	2.92E-02	2.31E-02	1.81E-02	1.42E-02	1.42E-02
24.00	2.49E-05	2.44E-05	2.28E-05	2.04E-05	1.76E-05	1.48E-05	1.23E-05	1.04E-05	9.25E-06	8.85E-06	8.85E-06
-EPS	100	110	120	130	140	150	160	170	180	TOTAL	

98

1
-S= 10.28 MEV
-OCCUPATION PROBABILITIES PO= 2. HQ= 1
TARGET Z= 26, N= 28 PROJ Z= 1, N= 0
G= 4.231, E=66.000
SCALE FACTOR= 1 350. FRAC PREEQ= 0.916
- P H RHOu/RHO STRU/STR STRD/STR

P	H	RHOu/RHO	STRU/STR	STRD/STR
1	0	1.000E+00	0.	0.
2	1	1.000E+00	1.000E+00	1.000E+00
3	2	1.000E+00	1.000E+00	1.000E+00
4	3	1.000E+00	1.000E+00	1.000E+00
5	4	9.581E-01	9.387E-01	9.387E-01
6	5	8.680E-01	8.558E-01	8.151E-01
7	6	7.513E-01	7.499E-01	6.399E-01
8	7	6.175E-01	6.154E-01	4.353E-01
9	8	4.794E-01	4.751E-01	2.442E-01

1PARTICLE SPECTRA Z= 0, N= 1
 FIRST EMISSION AT P= 2
 TARGET Z= 26, N= 28 PRDJ Z= 1, N= 0
 PO= 2, HO= 1, G= 4.231, E=66.000
 PAIR EXCIT. ACCORDING TO 2 COMP. ST. DENSITIES
 REACTION CROSS SECTION = 840.8
 SCALE FACTOR= 1.350, FRAC PREEQ= 0.916
 MSD POSSIBLE= 1.000, USED= 0.882
 CLOSED FORM SUM STARTS AT P= 2

		DIRECT		PREEQUILIBRIUM		EQUIL		TOTAL (MB/MEV)	
		-- NUTRA	KNDCK	-- MSD	MSC	-- WEISS	-- MSD	MSC	MSD+MSC
2.00	O.	O.		2.657E+00	3.926E-01	1.026E+00	2.657E+00	1.419E+00	4.076E+00
4.00	O.	O.		4.255E+00	4.440E-01	1.016E+00	4.255E+00	1.460E+00	5.714E+00
6.00	O.	O.		5.112E+00	3.653E-01	7.289E-01	5.112E+00	1.094E+00	6.206E+00
8.00	O.	O.		5.580E+00	2.653E-01	4.591E-01	5.580E+00	7.244E-01	6.304E+00
10.00	O.	O.		5.821E+00	1.793E-01	2.570E-01	5.821E+00	4.463E-01	6.267E+00
12.00	O.	O.		5.933E+00	1.155E-01	1.466E-01	5.933E+00	2.621E-01	6.195E+00
14.00	O.	O.		5.971E+00	7.187E-02	7.689E-02	5.971E+00	1.488E-01	6.120E+00
16.00	O.	O.		5.811E+00	4.351E-02	3.873E-02	5.811E+00	8.224E-02	5.893E+00
18.00	O.	O.		5.605E+00	2.573E-02	1.880E-02	5.605E+00	4.453E-02	5.649E+00
20.00	O.	O.		5.366E+00	1.489E-02	8.804E-03	5.366E+00	2.369E-02	5.390E+00
22.00	O.	O.		5.105E+00	8.414E-03	3.981E-03	5.105E+00	1.240E-02	5.118E+00
24.00	O.	O.		4.828E+00	4.634E-03	1.737E-03	4.828E+00	6.371E-03	4.834E+00
26.00	O.	O.		4.538E+00	2.475E-03	7.298E-04	4.538E+00	3.205E-03	4.541E+00
28.00	O.	O.		4.239E+00	1.274E-03	2.947E-04	4.239E+00	1.568E-03	4.240E+00
30.00	O.	O.		3.931E+00	6.260E-04	1.140E-04	3.931E+00	7.400E-04	3.932E+00
32.00	O.	O.		3.616E+00	2.905E-04	4.205E-05	3.616E+00	3.325E-04	3.616E+00
34.00	O.	O.		3.294E+00	1.253E-04	1.471E-05	3.294E+00	1.400E-04	3.294E+00
36.00	O.	O.		2.965E+00	4.917E-05	4.845E-06	2.965E+00	5.401E-05	2.965E+00
38.00	O.	O.		2.629E+00	1.702E-05	1.488E-06	2.629E+00	1.851E-05	2.629E+00
40.00	O.	O.		2.284E+00	4.956E-06	4.204E-07	2.284E+00	5.376E-06	2.284E+00
42.00	O.	O.		1.931E+00	1.123E-06	1.072E-07	1.931E+00	1.230E-06	1.931E+00
44.00	O.	O.		1.568E+00	1.720E-07	2.399E-08	1.568E+00	1.960E-07	1.568E+00
46.00	O.	O.		1.193E+00	1.313E-08	4.505E-09	1.193E+00	1.703E-08	1.193E+00
48.00	O.	O.		8.051E-01	2.136E-10	6.536E-10	8.051E-01	8.672E-10	8.051E-01
50.00	O.	O.		4.025E-01	0.	6.121E-11	4.025E-01	6.121E-11	4.025E-01
OSUMS	O.	O.		1.895E+02	3.676E+00	7.074E+00			

1ANGULAR DISTRIBUTIONS Z= 0 N= 1

ENERGY PARAM. = EPS+E.E.

CROSS SECTIONS IN MB/STR-MEV

- EPS	0	10	20	30	40	50	60	70	80	90	
2.00	6.42E-01	6.32E-01	6.05E-01	5.64E-01	5.13E-01	4.58E-01	4.03E-01	3.52E-01	3.09E-01	2.73E-01	
4.00	9.83E-01	9.67E-01	9.23E-01	8.54E-01	7.70E-01	6.79E-01	5.86E-01	5.05E-01	4.34E-01	3.75E-01	
6.00	1.16E+00	1.14E+00	1.08E+00	9.98E-01	8.92E-01	7.77E-01	6.64E-01	5.60E-01	4.71E-01	3.98E-01	
8.00	1.27E+00	1.25E+00	1.18E+00	1.08E+00	9.59E-01	8.26E-01	6.96E-01	5.77E-01	4.76E-01	3.93E-01	
10.00	1.36E+00	1.33E+00	1.26E+00	1.14E+00	1.00E+00	8.54E-01	7.09E-01	5.79E-01	4.68E-01	3.79E-01	
12.00	1.43E+00	1.40E+00	1.32E+00	1.19E+00	1.04E+00	8.72E-01	7.15E-01	5.74E-01	4.56E-01	3.62E-01	
14.00	1.50E+00	1.47E+00	1.37E+00	1.23E+00	1.06E+00	8.26E-01	7.16E-01	5.66E-01	4.42E-01	3.44E-01	
16.00	1.53E+00	1.50E+00	1.40E+00	1.24E+00	1.07E+00	8.74E-01	6.96E-01	5.41E-01	4.15E-01	3.17E-01	
18.00	1.56E+00	1.52E+00	1.41E+00	1.25E+00	1.03E+00	8.56E-01	6.71E-01	5.13E-01	3.86E-01	2.89E-01	
20.00	1.57E+00	1.53E+00	1.42E+00	1.24E+00	1.04E+00	8.32E-01	6.42E-01	4.81E-01	3.55E-01	2.61E-01	
22.00	1.59E+00	1.54E+00	1.42E+00	1.23E+00	1.02E+00	8.03E-01	6.08E-01	4.47E-01	3.23E-01	2.32E-01	
24.00	1.59E+00	1.54E+00	1.41E+00	1.22E+00	9.92E-01	7.69E-01	5.71E-01	4.11E-01	2.90E-01	2.05E-01	
26.00	1.59E+00	1.54E+00	1.40E+00	1.19E+00	9.59E-01	7.30E-01	5.31E-01	3.73E-01	2.57E-01	1.77E-01	
28.00	1.53E+00	1.53E+00	1.38E+00	1.16E+00	9.21E-01	6.88E-01	4.88E-01	3.34E-01	2.25E-01	1.52E-01	
30.00	1.56E+00	1.51E+00	1.35E+00	1.13E+00	8.77E-01	6.41E-01	4.43E-01	2.95E-01	1.93E-01	1.27E-01	
32.00	1.54E+00	1.48E+00	1.31E+00	1.08E+00	8.26E-01	5.89E-01	3.96E-01	2.55E-01	1.62E-01	1.05E-01	
34.00	1.50E+00	1.44E+00	1.27E+00	1.03E+00	7.70E-01	5.34E-01	3.48E-01	2.17E-01	1.33E-01	8.44E-02	
36.00	1.45E+00	1.38E+00	1.21E+00	9.66E-01	7.06E-01	4.76E-01	2.99E-01	1.79E-01	1.07E-01	6.66E-02	
38.00	1.38E+00	1.32E+00	1.14E+00	8.93E-01	6.36E-01	4.14E-01	2.49E-01	1.43E-01	8.25E-02	5.13E-02	
40.00	1.30E+00	1.23E+00	1.05E+00	8.08E-01	5.60E-01	3.50E-01	2.01E-01	1.10E-01	6.16E-02	3.86E-02	
42.00	1.19E+00	1.12E+00	9.46E-01	7.11E-01	4.76E-01	2.85E-01	1.55E-01	8.02E-02	4.40E-02	2.85E-02	
44.00	1.04E+00	9.83E-01	8.18E-01	6.00E-01	3.86E-01	2.19E-01	1.12E-01	5.47E-02	2.99E-02	2.06E-02	
46.00	8.66E-01	8.11E-01	6.64E-01	4.73E-01	2.92E-01	1.55E-01	7.29E-02	3.38E-02	1.92E-02	1.45E-02	
48.00	6.39E-01	5.95E-01	4.78E-01	3.30E-01	1.93E-01	9.49E-02	4.05E-02	1.79E-02	1.13E-02	9.53E-03	
50.00	3.51E-01	3.24E-01	2.56E-01	1.70E-01	9.32E-02	4.15E-02	1.56E-02	6.92E-03	5.32E-03	4.89E-03	
- EPS	100	110	120	130	140	150	160	170	180	TOTAL	
2.00	2.47E-01	2.29E-01	2.19E-01	2.14E-01	2.13E-01	2.14E-01	2.17E-01	2.19E-01	2.20E-01	4.08E+00	
4.00	3.31E-01	2.99E-01	2.78E-01	2.66E-01	2.60E-01	2.58E-01	2.59E-01	2.60E-01	2.60E-01	5.71E+00	
6.00	3.41E-01	3.00E-01	2.71E-01	2.53E-01	2.42E-01	2.36E-01	2.34E-01	2.33E-01	2.33E-01	6.21E+00	
8.00	3.29E-01	2.81E-01	2.47E-01	2.24E-01	2.09E-01	2.00E-01	1.95E-01	1.93E-01	1.92E-01	6.30E+00	
10.00	3.10E-01	2.58E-01	2.21E-01	1.95E-01	1.77E-01	1.65E-01	1.58E-01	1.55E-01	1.51E-01	6.27E+00	
12.00	2.89E-01	2.36E-01	1.96E-01	1.69E-01	1.49E-01	1.36E-01	1.28E-01	1.24E-01	1.22E-01	6.20E+00	
14.00	2.69E-01	2.15E-01	1.75E-01	1.47E-01	1.27E-01	1.13E-01	1.04E-01	9.94E-02	9.78E-02	6.12E+00	
16.00	2.44E-01	1.90E-01	1.52E-01	1.25E-01	1.06E-01	9.24E-02	8.36E-02	7.87E-02	7.71E-02	5.89E+00	
18.00	2.18E-01	1.67E-01	1.31E-01	1.06E-01	8.79E-02	7.53E-02	6.69E-02	6.21E-02	6.06E-02	5.65E+00	
20.00	1.93E-01	1.46E-01	1.13E-01	8.93E-02	7.27E-02	6.11E-02	5.34E-02	4.89E-02	4.74E-02	5.39E+00	
22.00	1.69E-01	1.26E-01	9.57E-02	7.48E-02	5.99E-02	4.94E-02	4.23E-02	3.82E-02	3.68E-02	5.12E+00	
24.00	1.46E-01	1.07E-01	8.06E-02	6.22E-02	4.90E-02	3.96E-02	3.32E-02	2.95E-02	2.83E-02	4.83E+00	
26.00	1.25E-01	9.02E-02	6.73E-02	5.14E-02	3.99E-02	3.16E-02	2.59E-02	2.25E-02	2.14E-02	4.54E+00	
28.00	1.05E-01	7.52E-02	5.58E-02	4.03E-02	2.24E-02	2.50E-02	1.99E-02	1.69E-02	1.59E-02	4.24E+00	
30.00	8.69E-02	6.21E-02	4.60E-02	3.47E-02	2.62E-02	1.97E-02	1.51E-02	1.23E-02	1.14E-02	3.93E+00	
32.00	7.09E-02	5.08E-02	3.79E-02	2.85E-02	2.11E-02	1.53E-02	1.12E-02	8.74E-03	7.93E-03	3.62E+00	
34.00	5.71E-02	4.14E-02	3.12E-02	2.34E-02	1.70E-02	1.18E-02	8.05E-03	5.88E-03	5.17E-03	3.29E+00	
36.00	4.54E-02	3.36E-02	2.58E-02	1.93E-02	1.36E-02	8.87E-03	5.53E-03	3.65E-03	3.06E-03	2.97E+00	
38.00	3.58E-02	2.74E-02	2.14E-02	1.59E-02	1.08E-02	6.45E-03	3.51E-03	1.98E-03	1.53E-03	2.63E+00	
40.00	2.81E-02	2.23E-02	1.77E-02	1.30E-02	8.32E-03	4.42E-03	1.93E-03	8.11E-04	5.37E-04	2.28E+00	
42.00	2.19E-02	1.81E-02	1.45E-02	1.04E-02	6.18E-03	2.71E-03	1.55E-04	1.31E-04	6.57E-05	1.93E+00	
44.00	1.69E-02	1.44E-02	1.16E-02	8.10E-03	4.29E-03	1.32E-03	3.81E-05	9.93E-05	7.20E-05	1.57E+00	
46.00	1.25E-02	1.09E-02	8.78E-03	5.90E-03	2.65E-03	2.49E-04	4.68E-04	3.68E-05	4.42E-04	1.19E+00	
48.00	8.42E-03	7.29E-03	5.91E-03	3.82E-03	1.32E-03	4.23E-04	5.67E-04	3.60E-04	9.25E-04	8.05E-01	
50.00	4.19E-03	3.57E-03	2.95E-03	1.86E-03	3.97E-04	5.75E-04	3.89E-04	5.30E-04	1.04E-03	4.02E-01	

1PARTICLE SPECTRA Z= 1, N= 0

FIRST EMISSION AT P= 2

TARGET Z= 26, N= 28 PROJ Z= 1, N= 0

POF= 2, HO= 1, G= 4.231, E=66.000

PAIR EXCIT. ACCORDING TO 2 COMP. ST. DENSITIES

REACTION CROSS SECTION = 840.8

SCALE FACTOR= 1.350, FRAC PREFQ= 0.916

MSD POSSIBLE= 1.000, USED= 0.882

CLOSED FORM SUM STARTS AT P= 2

- EPS --	DIRECT	PREEQUILIBRIUM		EQUIL		TOTAL (ME/MEV)		
		NUTRA	KNGOK	MSD	MSC	WEISS	MSD	MSC
2.00	O.	1.957E-09	8.526E-05	3.661E-05	1.522E-04	8.526E-05	1.888E-04	2.741E-04
4.00	O.	8.199E-05	2.764E+00	9.473E-01	3.438E+00	2.764E+00	4.386E+00	7.150E+00
6.00	O.	2.685E-04	7.151E+00	1.894E+00	6.013E+00	7.151E+00	7.906E+00	1.505E+01
8.00	O.	5.313E-04	1.143E+01	2.262E+00	6.288E+00	1.143E+01	8.550E+00	1.998E+01
10.00	O.	7.667E-04	1.364E+01	1.951E+00	4.745E+00	1.364E+01	6.696E+00	2.034E+01
12.00	O.	9.713E-04	1.463E+01	1.463E+00	3.108E+00	1.463E+01	4.570E+00	1.920E+01
14.00	O.	1.146E-03	1.496E+01	1.012E+00	1.874E+00	1.496E+01	2.887E+00	1.785E+01
16.00	O.	1.293E-03	1.494E+01	6.649E-01	1.067E+00	1.494E+01	1.732E+00	1.667E+01
18.00	O.	1.414E-03	1.475E+01	4.203E-01	5.802E-01	1.475E+01	1.001E+00	1.575E+01
20.00	O.	1.508E-03	1.449E+01	2.580E-01	3.038E-01	1.449E+01	5.618E-01	1.505E+01
22.00	O.	1.578E-03	1.421E+01	1.548E-01	1.537E-01	1.422E+01	3.084E-01	1.452E+01
24.00	O.	1.626E-03	1.372E+01	9.102E-02	7.527E-02	1.372E+01	1.663E-01	1.389E+01
26.00	O.	1.652E-03	1.301E+01	5.258E-02	3.574E-02	1.301E+01	8.832E-02	1.310E+01
28.00	O.	1.657E-03	1.227E+01	2.985E-02	1.645E-02	1.227E+01	4.629E-02	1.232E+01
30.00	O.	1.644E-03	1.152E+01	1.662E-02	7.331E-03	1.152E+01	2.395E-02	1.154E+01
32.00	O.	1.613E-03	1.076E+01	9.055E-03	3.162E-03	1.076E+01	1.222E-02	1.077E+01
34.00	O.	1.565E-03	9.994E+00	4.804E-03	1.317E-03	9.996E+00	6.122E-03	1.000E+01
36.00	O.	1.503E-03	9.228E+00	2.467E-03	5.288E-04	9.229E+00	2.996E-03	9.232E+00
38.00	O.	1.427E-03	8.461E+00	1.216E-03	2.038E-04	8.462E+00	1.420E-03	8.464E+00
40.00	O.	1.339E-03	7.695E+00	5.697E-04	7.515E-05	7.696E+00	6.448E-04	7.697E+00
42.00	O.	1.239E-03	6.932E+00	2.502E-04	2.637E-05	6.923E+00	2.765E-04	6.932E+00
44.00	O.	1.130E-03	6.172E+00	1.011E-04	8.747E-06	6.173E+00	1.099E-04	6.173E+00
46.00	O.	1.013E-03	5.416E+00	3.667E-05	2.719E-06	5.417E+00	3.939E-05	5.417E+00
48.00	O.	9.055E-04	4.758E+00	1.171E-05	7.986E-07	4.758E+00	1.251E-05	4.758E+00
50.00	O.	7.935E-04	4.107E+00	3.065E-06	2.155E-07	4.108E+00	3.280E-06	4.108E+00
52.00	O.	6.713E-04	3.432E+00	5.902E-07	5.190E-08	3.433E+00	6.421E-07	3.433E+00
54.00	O.	5.388E-04	2.729E+00	6.854E-08	1.078E-08	2.730E+00	7.932E-08	2.730E+00
56.00	O.	3.961E-04	1.993E+00	2.978E-09	1.819E-09	1.993E+00	4.797E-09	1.993E+00
58.00	O.	0.	1.220E+00	0.	2.219E-10	1.220E+00	2.219E-10	1.220E+00
60.00	O.	0.	4.039E-01	0.	1.461E-11	4.039E-01	1.461E-11	4.039E-01
OSUMS	O.	6.055E-02	5.134E+02	2.247E+01	5.542E+01			

ANGULAR DISTRIBUTIONS Z = 1 N = 0		ENERGY PARAM. = EPS+E E									
CROSS SECTIONS IN ME/STR-MEV		EPS									
	0	10	20	30	40	50	60	70	80	90	
2.00	3.10E-05	3.07E-05	2.99E-05	2.86E-05	2.70E-05	2.53E-05	2.35E-05	2.10E-05	2.07E-05	1.97E-05	
4.00	8.66E-01	8.57E-01	8.30E-01	7.90E-01	7.40E-01	6.86E-01	6.31E-01	5.82E-01	5.40E-01	5.08E-01	
6.00	1.97E+00	1.95E+00	1.88E+00	1.75E+00	1.65E+00	1.51E+00	1.38E+00	1.25E+00	1.14E+00	1.06E+00	
8.00	2.85E+00	2.81E+00	2.70E+00	2.54E+00	2.34E+00	2.12E+00	1.90E+00	1.69E+00	1.52E+00	1.38E+00	
10.00	3.16E+00	3.12E+00	2.99E+00	2.79E+00	2.55E+00	2.28E+00	2.01E+00	1.76E+00	1.55E+00	1.38E+00	
12.00	3.26E+00	3.21E+00	3.08E+00	2.84E+00	2.57E+00	2.27E+00	1.98E+00	1.71E+00	1.47E+00	1.28E+00	
14.00	3.28E+00	3.23E+00	3.07E+00	2.83E+00	2.54E+00	2.22E+00	1.91E+00	1.62E+00	1.37E+00	1.16E+00	
16.00	3.30E+00	3.24E+00	3.08E+00	2.82E+00	2.51E+00	2.17E+00	1.84E+00	1.53E+00	1.27E+00	1.05E+00	
18.00	3.34E+00	3.27E+00	3.09E+00	2.82E+00	2.49E+00	2.13E+00	1.78E+00	1.46E+00	1.19E+00	9.68E-01	
20.00	3.39E+00	3.33E+00	3.13E+00	2.84E+00	2.48E+00	2.10E+00	1.73E+00	1.40E+00	1.12E+00	8.96E-01	
22.00	3.47E+00	3.40E+00	3.19E+00	2.87E+00	2.49E+00	2.08E+00	1.70E+00	1.35E+00	1.06E+00	8.33E-01	
24.00	3.52E+00	3.44E+00	3.21E+00	2.87E+00	2.47E+00	2.04E+00	1.64E+00	1.28E+00	9.93E-01	7.64E-01	
26.00	3.51E+00	3.45E+00	3.19E+00	2.83E+00	2.41E+00	1.97E+00	1.55E+00	1.27E+00	9.10E-01	6.88E-01	
28.00	3.50E+00	3.41E+00	3.15E+00	2.78E+00	2.34E+00	1.88E+00	1.47E+00	1.11E+00	8.27E-01	6.14E-01	
30.00	3.47E+00	3.38E+00	3.11E+00	2.72E+00	2.26E+00	1.80E+00	1.37E+00	1.02E+00	7.45E-01	5.42E-01	
32.00	3.44E+00	3.34E+00	3.06E+00	2.65E+00	2.18E+00	1.70E+00	1.28E+00	9.29E-01	6.64E-01	4.73E-01	
34.00	3.39E+00	3.29E+00	2.99E+00	2.57E+00	2.08E+00	1.60E+00	1.18E+00	8.37E-01	5.84E-01	4.08E-01	
36.00	3.33E+00	3.25E+00	2.92E+00	2.48E+00	2.04E+00	1.49E+00	1.07E+00	7.44E-01	5.07E-01	3.46E-01	
38.00	3.25E+00	3.14E+00	2.82E+00	2.37E+00	1.86E+00	1.38E+00	9.66E-01	6.52E-01	4.32E-01	2.89E-01	
40.00	3.16E+00	3.04E+00	2.72E+00	2.25E+00	1.74E+00	1.25E+00	8.57E-01	5.61E-01	3.62E-01	2.36E-01	
42.00	3.05E+00	2.92E+00	2.59E+00	2.12E+00	1.60E+00	1.13E+00	7.47E-01	4.74E-01	2.96E-01	1.89E-01	
44.00	2.91E+00	2.44E+00	1.97E+00	1.46E+00	1.46E-01	9.97E-01	6.38E-01	3.90E-01	2.36E-01	1.49E-01	
46.00	2.74E+00	2.62E+00	2.27E+00	1.80E+00	1.30E+00	8.63E-01	5.31E-01	3.12E-01	1.83E-01	1.14E-01	
48.00	2.60E+00	2.47E+00	2.12E+00	1.65E+00	1.16E+00	7.41E-01	4.36E-01	2.45E-01	1.39E-01	8.66E-02	
50.00	2.42E+00	2.29E+00	1.95E+00	1.48E+00	1.01E+00	6.19E-01	3.46E-01	1.61E-01	1.02E-01	6.49E-02	
52.00	2.07E+00	1.73E+00	1.29E+00	8.47E-01	4.94E-01	2.61E-01	1.32E-01	7.19E-02	4.77E-02		
54.00	1.90E+00	1.78E+00	1.37E+00	1.06E+00	6.11E-01	3.69E-01	1.81E-01	8.64E-02	4.79E-02	3.44E-02	
56.00	1.51E+00	1.41E+00	1.15E+00	8.04E-01	4.93E-01	2.48E-01	1.12E-01	5.05E-02	2.98E-02	2.38E-02	
58.00	1.01E+00	9.39E-01	7.49E-01	5.07E-01	2.88E-01	1.35E-01	5.46E-02	2.40E-02	1.64E-02	1.45E-02	
60.00	3.69E-01	3.40E-01	2.65E-01	1.73E-01	9.14E-02	3.83E-02	5.39E-02	5.04E-03			
62.00	1.10E-01	9.51E-02	5.71E-02	2.81E-02	1.04E-02	3.01E-02	4.48E-02	4.81E-02			
64.00	3.10E-02	3.10E-02	1.20E-02	4.00E-03	1.10E-03	2.00E-03	3.00E-03	3.00E-03			
66.00	9.30E-03	9.30E-03	2.00E-03	4.00E-04	8.00E-04	1.00E-04	1.00E-04	1.00E-04			
68.00	2.60E-03	2.60E-03	5.00E-04	1.00E-04	2.00E-04	2.00E-04	2.00E-04	2.00E-04			
70.00	7.00E-04	7.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
72.00	1.80E-04	1.80E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
74.00	4.50E-05	4.50E-05	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
76.00	1.10E-05	1.10E-05	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
78.00	2.80E-06	2.80E-06	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
80.00	7.00E-07	7.00E-07	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
82.00	1.80E-07	1.80E-07	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
84.00	4.50E-08	4.50E-08	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
86.00	1.10E-08	1.10E-08	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
88.00	2.80E-09	2.80E-09	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
90.00	7.00E-10	7.00E-10	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
92.00	1.80E-10	1.80E-10	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
94.00	4.50E-11	4.50E-11	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
96.00	1.10E-11	1.10E-11	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
98.00	2.80E-12	2.80E-12	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
100.00	7.00E-13	7.00E-13	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
110.00	1.80E-13	1.80E-13	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
120.00	4.50E-14	4.50E-14	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
130.00	1.10E-14	1.10E-14	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
140.00	2.80E-15	2.80E-15	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
150.00	7.00E-16	7.00E-16	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
160.00	1.80E-16	1.80E-16	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
170.00	4.50E-17	4.50E-17	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
180.00	1.10E-17	1.10E-17	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
190.00	2.80E-18	2.80E-18	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
200.00	7.00E-19	7.00E-19	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
210.00	1.80E-19	1.80E-19	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
220.00	4.50E-20	4.50E-20	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
230.00	1.10E-20	1.10E-20	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
240.00	2.80E-21	2.80E-21	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
250.00	7.00E-22	7.00E-22	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
260.00	1.80E-22	1.80E-22	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
270.00	4.50E-23	4.50E-23	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
280.00	1.10E-23	1.10E-23	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
290.00	2.80E-24	2.80E-24	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
300.00	7.00E-25	7.00E-25	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
310.00	1.80E-25	1.80E-25	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
320.00	4.50E-26	4.50E-26	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
330.00	1.10E-26	1.10E-26	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
340.00	2.80E-27	2.80E-27	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
350.00	7.00E-28	7.00E-28	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
360.00	1.80E-28	1.80E-28	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
370.00	4.50E-29	4.50E-29	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
380.00	1.10E-29	1.10E-29	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
390.00	2.80E-30	2.80E-30	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
400.00	7.00E-31	7.00E-31	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
410.00	1.80E-31	1.80E-31	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
420.00	4.50E-32	4.50E-32	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
430.00	1.10E-32	1.10E-32	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04	1.00E-04			
440.00	2.80E-33	2.80E-33	1.00E-04</								

2.00	1.92E-05	1.89E-05	1.90E-05	1.93E-05	1.98E-05	2.03E-05	2.07E-05	2.10E-05	2.11E-05	2.14E-04
4.00	4.87E-01	4.77E-01	4.75E-01	4.80E-01	4.90E-01	5.11E-01	5.19E-01	5.21E-01	5.24E+00	7.15E+20
6.00	9.96E-01	9.60E-01	9.46E-01	9.48E-01	9.61E-01	9.79E-01	9.91E-01	1.01E+00	1.01E+00	1.51E+21
8.00	1.28E+00	1.21E+00	1.17E+00	1.16E+00	1.16E+00	1.18E+00	1.19E+00	1.21E+00	1.21E+00	2.00E+21
10.00	1.25E+00	1.15E+00	1.09E+00	1.06E+00	1.05E+00	1.06E+00	1.06E+00	1.07E+00	1.07E+00	2.03E+21
12.00	1.12E+00	1.01E+00	9.40E-01	8.94E-01	8.69E-01	8.58E-01	8.56E-01	8.56E-01	8.56E-01	1.92E+01
14.00	9.96E-01	8.75E-01	7.89E-01	7.32E-01	6.96E-01	6.75E-01	6.60E-01	6.59E-01	6.59E-01	7.8E+01
16.00	8.86E-01	7.58E-01	6.66E-01	6.01E-01	5.58E-01	5.20E-01	5.14E-01	5.08E-01	5.03E-01	1.67E+01
18.00	7.95E-01	6.64E-01	5.68E-01	5.00E-01	4.53E-01	4.21E-01	4.01E-01	3.91E-01	3.87E-01	5.8E+01
20.00	7.20E-01	5.89E-01	4.92E-01	4.23E-01	3.73E-01	3.40E-01	3.18E-01	3.08E-01	3.03E-01	5.1E+01
22.00	6.57E-01	5.26E-01	4.31E-01	3.62E-01	3.13E-01	2.79E-01	2.57E-01	2.44E-01	2.40E-01	4.45E+01
24.00	5.92E-01	4.65E-01	3.74E-01	3.08E-01	2.61E-01	2.28E-01	2.07E-01	1.94E-01	1.90E-01	3.9E+01
26.00	5.23E-01	4.04E-01	3.19E-01	2.58E-01	2.15E-01	1.85E-01	1.64E-01	1.53E-01	1.49E-01	3.1E+01
28.00	4.59E-01	3.48E-01	2.71E-01	2.16E-01	1.76E-01	1.49E-01	1.30E-01	1.20E-01	1.16E-01	2.3E+01
30.00	3.98E-01	2.98E-01	2.28E-01	1.79E-01	1.44E-01	1.19E-01	1.05E-01	9.33E-02	9.01E-02	1.5E+01
32.00	3.41E-01	2.51E-01	1.90E-01	1.47E-01	1.17E-01	9.52E-02	8.65E-02	7.98E-02	7.69E-02	9.0E+01
34.00	2.89E-01	2.10E-01	1.57E-01	1.24E-01	9.43E-02	7.53E-02	6.125E-02	5.91E-02	5.71E-02	7.0E+01
36.00	2.41E-01	1.74E-01	1.29E-01	9.81E-02	7.56E-02	5.91E-02	4.71E-02	4.15E-02	3.98E-02	5.38E+01
38.00	1.98E-01	1.42E-01	1.05E-01	7.95E-02	6.04E-02	4.60E-02	3.59E-02	3.29E-02	3.09E-02	4.27E+01
40.00	1.64E-01	1.15E-01	8.53E-02	6.42E-02	4.80E-02	3.55E-02	2.66E-02	2.36E-02	2.16E-02	3.0E+00
42.00	1.28E-01	9.22E-02	6.90E-02	5.19E-02	3.81E-02	2.45E-02	1.92E-02	1.49E-02	1.30E-02	2.03E+00
44.00	1.01E-01	7.37E-02	5.59E-02	4.20E-02	2.30E-02	1.73E-02	1.33E-02	9.29E-03	8.01E-03	1.71E+00
46.00	7.83E-02	5.88E-02	4.54E-02	3.40E-02	2.35E-02	1.48E-02	1.08E-02	7.69E-03	6.35E-03	4.34E+00
48.00	6.15E-02	4.79E-02	3.77E-02	2.79E-02	1.84E-02	1.051E-02	7.66E-03	5.07E-03	3.60E-03	4.24E+00
50.00	4.84E-02	3.91E-02	3.13E-02	2.28E-02	1.41E-02	6.90E-03	2.135E-03	1.20E-03	4.75E-04	4.11E+00
52.00	3.78E-02	3.18E-02	2.56E-02	1.82E-02	1.02E-02	3.89E-03	1.35E-03	6.53E-05	2.70E-05	3.43E+00
54.00	2.89E-02	2.49F-02	2.02E-02	1.33E-02	6.82E-03	4.48E-03	1.35E-04	6.29E-04	4.27E-04	2.73E+00
56.00	2.09E-02	1.81E-02	1.44E-02	9.66E-03	3.89E-03	2.61E-03	1.06E-03	1.06E-04	1.36E-05	4.99E+00
58.00	1.27E-02	1.10E-02	8.94E-03	5.70E-03	1.63E-03	1.44E-03	4.02E-04	1.96E-04	2.13E-05	2.22E+00
60.00	4.16E-03	3.50E-03	2.98E-03	1.67E-03	2.55E-04	5.52E-04	1.36E-04	4.36E-04	4.64E-05	4.64E-01

1PARTICLE SPECTRA Z= 2, N= 2

FIRST EMISSION AT P= 4

TARGET Z= 26, N= 28 PROJ Z= 1, N= 0

PO= 2, HO= 1, G= 4.221, E=66.000

PAIR EXCIT. ACCORDING TO 2 COMP. ST. DENSITIES

REACTION CROSS SECTION = 840.8

SCALE FACTOR= 1.350, FRAC PREEQ= 0.916

MSD POSSIBLE= 1.000, USED= 0.842

CLOSED FORM SUM STARTS AT P= 4

EPS	DIRECT		PREEQUILIBRIUM		EQUIL		TOTAL MSC	(ME/MEV) MSD+MSC
	NUTRA	KNOCK	MSD	MSC	WEISS	MSD		
6.00	3.860E-03	2.125E-04	9.239E-03	1.035E-02	8.395E-02	1.331E-02	9.430E-02	1.076E-01
8.00	3.047E-02	1.743E-03	5.062E-02	5.212E-02	3.665E-01	8.283E-02	4.187E-01	5.015E-01
10.00	8.499E-02	5.057E-03	9.770E-02	9.127E-02	5.601E-01	1.878E-01	6.513E-01	8.391E-01
12.00	1.453E-01	9.007E-03	1.154E-01	9.639E-02	5.192E-01	2.697E-01	6.156E-01	8.853E-01
14.00	1.927E-01	1.247E-02	1.059E-01	7.779E-02	3.698E-01	3.111E-01	4.476E-01	7.587E-01
16.00	2.284E-01	1.546E-02	8.714E-02	5.525E-02	2.323E-01	3.310E-01	2.880E-01	6.190E-01
18.00	2.536E-01	1.799E-02	6.747E-02	3.624E-02	1.356E-01	3.391E-01	1.719E-01	5.109E-01
20.00	2.693E-01	2.007E-02	5.025E-02	2.247E-02	7.468E-02	2.297E-01	9.714E-02	4.369E-01
22.00	2.767E-01	2.171E-02	3.670E-02	1.335E-02	3.926E-02	3.351E-01	5.261E-02	3.877E-01
24.00	2.767E-01	2.293E-02	2.634E-02	7.672E-03	1.982E-02	3.260E-01	2.749E-02	3.534E-01
26.00	2.704E-01	2.373E-02	1.873E-02	4.299E-03	9.634E-03	3.129E-01	1.393E-02	3.268E-01
28.00	2.588E-01	2.414E-02	1.323E-02	2.363E-03	4.515E-03	2.961E-01	6.877E-03	3.030E-01
30.00	2.428E-01	2.416E-02	9.309E-03	1.279E-03	2.040E-03	2.762E-01	3.318E-03	2.796E-01
32.00	2.233E-01	2.380E-02	6.530E-03	6.825E-04	8.871E-04	2.536E-01	1.570E-03	2.552E-01
34.00	2.012E-01	2.309E-02	4.566E-03	3.592E-04	3.709E-04	2.288E-01	7.301E-04	2.295E-01
36.00	1.773E-01	2.203E-02	3.182E-03	1.858E-04	1.485E-04	2.025E-01	3.344E-04	2.028E-01
38.00	1.524E-01	2.063E-02	2.205E-03	9.391E-05	5.679E-05	1.753E-01	1.507E-04	1.754E-01
40.00	1.273E-01	1.891E-02	1.515E-03	4.601E-05	2.060E-05	1.478E-01	6.662E-05	1.478E-01
42.00	1.028E-01	1.688E-02	1.028E-03	2.157E-05	7.046E-06	1.207E-01	2.962E-03	1.207E-01
44.00	7.949E-02	1.456E-02	6.830E-04	9.488E-06	2.250E-06	9.473E-02	1.174E-05	9.473E-02
46.00	5.807E-02	1.195E-02	4.390E-04	3.790E-06	6.618E-07	7.045E-02	4.452E-06	7.045E-02
48.00	3.916E-02	9.068E-03	2.671E-04	1.295E-06	1.761E-07	4.649E-02	1.471E-06	4.649E-02
50.00	2.336E-02	5.929E-03	1.475E-04	1.331E-07	4.125E-08	2.941E-02	3.743E-07	2.944E-02
52.00	1.123E-02	0.	6.745E-05	4.612E-08	8.1295E-09	1.130E-02	5.425E-08	1.130E-02
54.00	3.302E-03	0.	1.926E-05	0.	1.243E-09	3.321E-03	1.243E-09	3.321E-03
56.00	0.	0.	0.	0.	1.232E-10	0.	1.232E-10	1.232E-10

OSUMS 7.466E+00 7.311E-01 1.418E+00 9.445E-01 4.839E+00

		ANGULAR DISTRIBUTIONS Z= 2 N= 2									
		ENERGY PARAM. = EPS+B.E.									
		CROSS SECTIONS IN MB/STR-MEV									
- EPS		0 10 20 30 40 50 60 70 80 90									
		30 40 50 60 70 80 90									
6.00		1.21E-02	1.19E-02	1.15E-02	1.09E-02	1.02E-02	9.41E-03	8.68E-03	8.06E-03	7.60E-03	7.35E-03
8.00		5.97E-02	5.89E-02	5.67E-02	5.34E-02	4.94E-02	4.51E-02	4.11E-02	3.77E-02	3.51E-02	3.36E-02
10.00		1.07E-01	1.06E-01	1.02E-01	9.50E-02	8.70E-02	7.85E-02	7.05E-02	6.36E-02	5.83E-02	5.50E-02
12.00		1.24E-01	1.22E-01	1.17E-01	1.11E-01	1.08E-01	9.81E-02	8.73E-02	7.70E-02	6.81E-02	6.12E-02
14.00		1.19E-01	1.17E-01	1.11E-01	1.02E-01	9.13E-02	8.00E-02	7.93E-02	7.70E-02	6.99E-02	5.24E-02
16.00		1.09E-01	1.07E-01	1.01E-01	9.25E-02	8.18E-02	7.06E-02	5.98E-02	5.04E-02	4.29E-02	3.74E-02
18.00		1.02E-01	9.97E-02	9.38E-02	8.49E-02	7.43E-02	6.31E-02	5.24E-02	4.30E-02	3.55E-02	2.98E-02
20.00		9.73E-02	9.52E-02	8.92E-02	8.01E-02	6.93E-02	5.79E-02	4.72E-02	3.78E-02	3.02E-02	2.45E-02
22.00		9.53E-02	9.30E-02	8.68E-02	7.73E-02	6.61E-02	5.45E-02	4.36E-02	3.42E-02	2.65E-02	2.08E-02
24.00		9.45E-02	9.21E-02	8.55E-02	7.56E-02	6.39E-02	5.19E-02	4.08E-02	3.13E-02	2.37E-02	1.80E-02
26.00		9.40E-02	9.15E-02	8.45E-02	7.42E-02	6.20E-02	5.02E-02	3.83E-02	2.87E-02	2.13E-02	1.57E-02
28.00		9.33E-02	9.07E-02	8.33E-02	7.24E-02	5.98E-02	4.71E-02	3.50E-02	2.46E-02	1.90E-02	1.37E-02
30.00		9.18E-02	8.91E-02	8.14E-02	7.01E-02	5.71E-02	4.33E-02	3.29E-02	2.36E-02	1.67E-02	1.18E-02
32.00		8.94E-02	8.65E-02	7.86E-02	6.70E-02	5.38E-02	4.10E-02	2.97E-02	2.09E-02	1.44E-02	9.92E-03
34.00		8.57E-02	8.28E-02	7.47E-02	6.30E-02	4.99E-02	3.72E-02	2.64E-02	1.80E-02	1.21E-02	8.16E-03
36.00		8.09E-02	7.80E-02	6.98E-02	5.82E-02	4.53E-02	3.30E-02	2.28E-02	1.52E-02	9.89E-03	6.52E-03
38.00		7.48E-02	7.19E-02	6.39E-02	5.26E-02	4.01E-02	2.86E-02	1.92E-02	1.23E-02	7.92E-03	5.04E-03
40.00		6.76E-02	6.48E-02	5.71E-02	4.63E-02	3.46E-02	2.40E-02	1.56E-02	9.67E-03	5.94E-03	3.76E-03
42.00		5.93E-02	5.66E-02	4.94E-02	3.94E-02	2.88E-02	1.94E-02	1.21E-02	7.24E-03	4.30E-03	2.69E-03
44.00		5.01E-02	4.77E-02	4.09E-02	3.23E-02	2.30E-02	1.49E-02	8.59E-03	5.12E-03	2.95E-03	1.83E-03
46.00		4.03E-02	3.81E-02	3.25E-02	2.50E-02	1.73E-02	1.08E-02	6.17E-03	3.36E-03	1.88E-03	1.18E-03
48.00		2.99E-02	2.83E-02	2.38E-02	1.79E-02	1.20E-02	7.13E-03	3.86E-03	2.00E-03	1.09E-03	7.10E-04
50.00		1.97E-02	1.86E-02	1.54E-02	1.13E-02	7.26E-03	4.10E-03	2.03E-03	1.01E-03	5.56E-04	3.85E-04
52.00		8.25E-03	7.72E-03	6.31E-03	4.14E-03	2.76E-03	1.46E-03	6.84E-04	3.16E-04	1.80E-04	1.37E-04
54.00		2.65E-03	2.47E-03	1.98E-03	1.36E-03	7.95E-04	3.89E-04	1.65E-04	7.29E-05	4.63E-05	3.93E-05
56.00		5.41E-11	5.00E-11	3.92E-11	2.59E-11	1.43E-11	6.57E-12	2.80E-12	1.59E-12	1.45E-12	1.50E-12
100		110	120	130	140	150	160	170	180	TOTAL	


```

1 PARTICLE SPECTRA Z = 1. N = 2
FIRST EMISSION AT P = 3
TARGET Z = 26. N = 28 PROJ Z =
PO = 2, HO = 1, G = 4.231, E = 66.0000
PAIR EXCIT. ACCORDING TO 2 COMP.
SCALE FACTOR = 840.0
SECTION = 840.0
FRAC PREE = 1.350
MSD POSSIBLE = 1.000, USED = 0.882
CLOSED FORM SUM STARTS AT P = 3
EPS = DIRECT

```

```

TARGET Z = 26, N = 28 PROJ Z = 1, N = 0
PO = 2., HO = 1., G = 4.231, E = 66.000
PAIR EXCIT. ACCORDING TO 2 COMP. ST. DENSITIES
REACTION CROSS SECTION = 840.8
SCAT. FACTOR = 1.750, ENERG. = 0.916

```

ANGULAR DISTRIBUTIONS Z= 1 N= 2

ENERGY PARAM. = EPS+B.E.		CROSS SECTIONS IN MB/STR-MEV									
- EPS	O	10	20	30	40	50	60	70	80	90	
3.00	1.25E-03	1.23E-03	1.16E-03	1.07E-03	9.53E-04	8.31E-04	7.16E-04	6.17E-04	5.41E-04	4.90E-04	
5.00	1.89E-02	1.85E-02	1.75E-02	1.59E-02	1.02E-02	1.20E-02	8.56E-03	7.29E-03	6.42E-03	6.42E-03	
7.00	3.17E-02	3.17E-02	9.98E-02	2.69E-02	2.34E-02	1.98E-02	1.63E-02	1.11E-02	9.41E-03	9.41E-03	
9.00	3.77E-02	3.69E-02	3.44E-02	3.08E-02	2.65E-02	2.20E-02	1.78E-02	1.42E-02	1.14E-02	9.31E-03	
11.00	4.03E-02	3.93E-02	3.65E-02	3.24E-02	2.75E-02	2.25E-02	1.79E-02	1.39E-02	1.08E-02	8.47E-03	
13.00	4.22E-02	4.11E-02	3.80E-02	3.34E-02	2.81E-02	2.26E-02	1.76E-02	1.33E-02	1.00E-02	7.67E-03	
15.00	4.42E-02	4.29E-02	3.94E-02	3.44E-02	2.85E-02	2.26E-02	1.72E-02	1.28E-02	9.35E-03	6.88E-03	
17.00	4.61E-02	4.48E-02	4.10E-02	3.54E-02	2.90E-02	2.26E-02	1.69E-02	1.22E-02	8.71E-03	6.23E-03	
19.00	4.83E-02	4.68E-02	4.25E-02	3.64E-02	2.94E-02	2.25E-02	1.64E-02	1.16E-02	8.07E-03	5.62E-03	
21.00	5.03E-02	4.86E-02	4.40E-02	3.72E-02	2.96E-02	2.26E-02	1.62E-02	1.13E-02	7.39E-03	5.02E-03	
23.00	5.20E-02	5.03E-02	4.51E-02	3.78E-02	2.95E-02	2.17E-02	1.51E-02	1.01E-02	6.66E-03	4.42E-03	
25.00	5.36E-02	5.16E-02	4.59E-02	3.79E-02	2.91E-02	2.09E-02	1.41E-02	9.18E-03	5.87E-03	3.81E-03	
27.00	5.45E-02	5.23E-02	4.62E-02	3.76E-02	2.83E-02	1.98E-02	1.30E-02	8.14E-03	5.04E-03	3.21E-03	
29.00	5.47E-02	5.24E-02	4.58E-02	3.67E-02	2.70E-02	1.83E-02	1.16E-02	7.01E-03	4.21E-03	2.64E-03	
31.00	5.49E-02	5.15E-02	4.51E-02	3.55E-02	2.52E-02	1.55E-02	1.00E-02	5.82E-03	3.38E-03	2.10E-03	
33.00	5.20E-02	4.94E-02	4.23E-02	3.27E-02	2.28E-02	1.44E-02	8.36E-03	4.62E-03	2.61E-03	1.63E-03	
35.00	4.85E-02	4.59E-02	3.89E-02	2.94E-02	1.98E-02	1.20E-02	6.60E-03	3.47E-03	1.91E-03	1.22E-03	
37.00	4.07E-02	3.97E-02	3.39E-02	2.50E-02	1.63E-02	9.35E-03	4.83E-03	2.40E-03	1.31E-03	8.88E-04	
39.00	3.55E-02	3.32E-02	2.73E-02	1.96E-02	1.22E-02	6.59E-03	3.16E-03	1.48E-03	8.31E-04	6.14E-04	
41.00	2.48E-02	2.31E-02	1.86E-02	1.30E-02	7.67E-03	3.85E-03	1.68E-03	7.51E-04	4.58E-04	3.78E-04	
43.00	1.04E-02	9.59E-03	7.60E-03	5.09E-03	2.84E-03	1.29E-03	5.03E-04	2.21E-04	1.61E-04	1.46E-04	
45.00	1.09E-11	1.00E-11	7.73E-12	4.29E-12	2.58E-12	1.09E-12	4.29E-13	2.63E-13	2.75E-13	2.94E-13	
- EPS	100	110	120	130	140	150	160	170	180	TOTAL	

1PARTICLE SPECTRA Z= 2. N= 1
 FIRST EMISSION AT P= 3
 TARGET Z= 26. N= 2B PROJ Z= 1. N= 0
 PO= 2. HO= 1. G= 4.231. E=66.000
 PAIR EXCIT. ACCORDING TO 2 COMP. ST. DENSITIES
 REACTION CROSS SECTION = 840.8
 SCALE FACTOR= 1.350. FRAC PREiQ= 0.916
 MSC POSSIBLE= 1.000. USED= 0.882
 CLOSED FORM SUM STARTS AT P= 3
 EPS -- DIRECT -- PREEQUILIBRIUM -- EQL -- TOTAL (MB/MEV)
 -- NUTRA KNOCK -- MSD MSC -- WEISS -- MSD MSC
 -- MSD MSC

	8.00	1.442E-02	2.340E-02	6.324E-03	2.349E-02	3.722E-02	2.981E-02	6.763E-C2
10.00	4.048E-02	4.772E-02	9.930E-03	3.208E-02	8.820E-02	4.201E-02	1.302E-01	
12.00	6.898E-02	0.0	5.978E-02	9.407E-03	2.677E-02	1.288E-01	3.558E-02	1.643E-01
14.00	9.352E-02	0.0	6.055E-02	7.066E-03	1.669E-02	1.541E-01	2.375E-02	1.778E-01
16.00	1.142E-01	0.0	5.616E-02	4.771E-03	9.404E-03	1.704E-01	1.418E-02	1.845E-01
18.00	1.311E-01	0.0	4.979E-02	3.027E-03	4.880E-03	1.809E-01	7.907E-03	1.888E-01
20.00	1.443E-01	0.0	4.303E-02	1.839E-03	2.374E-03	4.873E-01	4.213E-03	1.915E-01
22.00	1.539E-01	0.0	3.664E-02	1.092E-03	1.092E-03	1.905E-01	2.171E-03	1.927E-01
24.00	1.599E-01	0.0	3.091E-02	6.132E-04	4.764E-04	1.908E-01	1.090E-03	1.919E-01
26.00	1.625E-01	0.0	2.592E-02	3.362E-04	1.973E-04	1.884E-01	5.335E-04	1.890E-01
28.00	1.618E-01	0.0	2.164E-02	1.768E-04	7.741E-05	1.834E-01	2.542E-04	1.837E-01
30.00	1.578E-01	0.0	1.799E-02	8.827E-05	2.866E-05	1.758E-01	1.699E-04	1.759E-01
32.00	1.506E-01	0.0	1.488E-02	4.112E-05	9.951E-06	1.655E-01	5.107E-05	1.655E-01
34.00	1.403E-01	0.0	1.221E-02	1.743E-05	3.212E-06	1.525E-01	2.054E-05	1.525E-01
36.00	1.270E-01	0.0	9.906E-03	6.441E-06	9.518E-07	1.369E-01	7.393E-06	1.369E-01
38.00	1.108E-01	0.0	7.878E-03	1.927E-05	2.542E-07	1.186E-01	2.182E-06	1.186E-01
40.00	9.170E-02	0.0	6.050E-03	4.029E-07	5.947E-08	9.775E-02	4.623E-07	9.775E-02
42.00	6.988E-02	0.0	4.349E-03	4.083E-08	1.165E-08	7.423E-02	5.249E-08	7.423E-02
44.00	4.539E-02	0.0	2.696E-03	0.0	1.758E-09	4.809E-02	4.809E-02	
46.00	1.833E-02	0.0	1.008E-03	0.0	1.696E-10	1.934E-02	1.696E-10	1.934E-02
OSUMS	4.304E+00	0.	1.055E+00	8.945E-02	2.340E-01			

1
-S= 5.69 MEV
-OCCUPATION PROBABILITIES PO= 2. HO= 1
TARGET Z= 92, N=143 PROJ Z= 1, N= 0
G=16.620, E=20.040
SCALE FACTOR= 1.350, FRAC PREQ= 0.255
- P H RHOU/RHO STRU/STR STRD/STR

	P	H	RHOU/RHO	STRU/STR	STRD/STR
1	0	1.000E+00	0.	0.	
2	1	8.741E-01	8.741E-01	8.741E-01	
3	2	7.182E-01	6.998E-01	6.662E-01	
4	3	5.258E-01	5.045E-01	4.229E-01	
5	4	3.387E-01	3.060E-01	2.183E-01	
6	5	1.997E-01	1.611E-01	9.085E-02	
7	6	1.091E-01	7.821E-02	3.004E-02	
8	7	5.519E-02	3.832E-02	7.700E-03	
9	8	2.570E-02	2.048E-02	1.477E-03	
10	9	1.087E-02	1.202E-02	0.	

1PARTICLE SPECTRA Z= 0, N= 1
 FIRST EMISSION AT P= 2
 TARGET Z= 92, N=143 PROJ Z= 1, N= 0
 PD= 2, HO= 1, G=16.620, E=20.040
 PAIR EXCIT. ACCORDING TO 2 COMP. ST. DENSITIES
 REACTION CROSS SECTION = 743.2
 SCALE FACTOR= 1.350, FRAC_PREQ= 0.255
 MSD POSSIBLE= 0.874, USED= 0.214
 CLOSED FORM SUM STARTS AT P= 2
 EPS -- NUTRA DIRECT -- PREEQUILIBRIUM -- EQUIL -- TOTAL (MB/MEV) MSD+MSC
 -- KNOCK -- MSD MSC -- WEIS -- MSD MSC
 1.00 0. 8.505E+00 7.638E+00 8.488E+01 8.506E+00 9.252E+01 1.010E+02
 2.00 0. 0. 1.579E+01 7.858E+00 5.678E+01 1.579E+01 6.463E+01 8.043E+01
 3.00 0. 0. 1.768E+01 5.205E+00 2.123E+01 1.768E+01 2.644E+01 4.412E+01
 4.00 0. 0. 1.736E+C1 3.241E+00 6.436E+00 1.736E+01 9.677E+00 2.704E+01
 5.00 0. 0. 1.614E+01 2.010E+00 1.596E+00 1.614E+01 3.707E+00 1.985E+01
 6.00 0. 0. 1.454E+01 1.243E+00 3.949E-01 1.454E+01 1.638E+00 1.618E+01
 7.00 0. 0. 1.281E+C1 7.590E-01 8.119E-02 1.281E+01 8.402E-01 1.366E+01
 8.00 0. 0. 1.108E+01 4.533E-01 1.459E-02 1.108E+01 4.678E-01 1.155E+01
 9.00 0. 0. 9.388E+00 2.616E-01 2.243E-03 9.388E+00 2.638E-01 9.651E+00
 10.00 0. 0. 7.729E+00 1.439E-01 2.851E-04 7.729E+00 1.442E-01 7.873E+00
 11.00 0. 0. 6.083E+00 7.401E-02 2.825E-05 6.083E+00 7.403E-02 6.157E+00
 12.00 0. 0. 4.408E+00 3.471E-02 1.966E-06 4.408E+00 3.474E-02 4.442E+00
 13.00 0. 0. 2.647E+00 1.407E-02 7.680E-08 2.647E+00 1.407E-02 2.661E+00
 14.00 0. 0. 7.266E-01 3.087E-03 8.253E-10 7.266E-01 3.087E-03 7.297E-01
 OSUMS 0. 0. 1.449E+02 2.894E+01 1.715E+02

ANGULAR DISTRIBUTIONS Z=0 N=1

CROSS SECTIONS IN MB/STR-MEV										TOTAL	
-	EPS	0	10	20	30	40	50	60	70	80	90
1.00	1.01E+01	9.98E+00	9.75E+00	9.40E+00	8.96E+00	8.50E+00	8.07E+00	7.70E+00	7.43E+00	7.29E+00	
2.00	8.63E+00	8.55E+00	8.33E+00	8.00E+00	7.56E+00	7.11E+00	6.66E+00	6.27E+00	5.97E+00	5.77E+00	
3.00	5.29E+00	5.13E+00	4.89E+00	4.58E+00	4.24E+00	3.91E+00	3.60E+00	3.34E+00	3.14E+00	3.14E+00	
4.00	3.75E+00	3.70E+00	3.58E+00	3.38E+00	3.14E+00	2.88E+00	2.59E+00	2.33E+00	2.10E+00	1.91E+00	
5.00	3.02E+00	2.99E+00	2.88E+00	2.70E+00	2.49E+00	2.25E+00	2.01E+00	1.78E+00	1.57E+00	1.40E+00	
6.00	2.61E+00	2.57E+00	2.47E+00	2.32E+00	2.13E+00	1.91E+00	1.69E+00	1.48E+00	1.29E+00	1.13E+00	
7.00	2.29E+00	2.25E+00	2.16E+00	2.02E+00	1.85E+00	1.65E+00	1.45E+00	1.26E+00	1.09E+00	9.45E-01	
8.00	1.99E+00	1.96E+00	1.88E+00	1.76E+00	1.60E+00	1.42E+00	1.25E+00	1.08E+00	9.24E-01	7.93E-01	
9.00	1.71E+00	1.69E+00	1.61E+00	1.53E+00	1.42E+00	1.32E+00	1.21E+00	1.05E+00	9.04E-01	7.56E-01	
10.00	1.43E+00	1.41E+00	1.35E+00	1.28E+00	1.18E+00	1.09E+00	9.99E-01	8.66E-01	7.40E-01	5.30E-01	
11.00	1.15E+00	1.13E+00	1.08E+00	1.00E+00	9.00E-01	7.92E-01	6.83E-01	5.80E-01	4.89E-01	4.11E-01	
12.00	8.50E-01	8.36E-01	7.97E-01	7.36E-01	6.61E-01	5.79E-01	4.96E-01	4.19E-01	3.51E-01	2.93E-01	
13.00	5.22E-01	5.13E-01	4.88E-01	4.50E-01	4.02E-01	3.51E-01	2.99E-01	2.52E-01	2.09E-01	1.73E-01	
14.00	1.46E-01	1.44E-01	1.37E-01	1.26E-01	1.17E-01	9.73E-02	8.26E-02	6.70E-02	5.70E-02	4.70E-02	
	EPS	100	110	120	130	140	150	160	170	180	

!PARTICLE SPECTRA Z= 1, N= 0
 FIRST EMISSION AT P= 2
 TARGET Z= 92, N=143 PROJ Z= 1, N= 0
 PO= 2, HO= 1, G=16.620, E=20.040
 PAIR EXCIT. ACCORDING TO 2 COMP. ST. DENSITIES
 REACTION CROSS SECTION = 743.2
 SCALE FACTOR= 1.350, FRAC PREEQ= 0.255
 MSD POSSIBLE= 0.874, USED= 0.214
 CLOSED FORM SUM STARTS AT P= 2

		DIRECT		PREEQUILIBRIUM		EQUIL		TOTAL (MB/MEV)		
- EPS --	-- NUTRA	KNOCK	-- MSD	MSC	-- WEISS	-- MSD	MSC	MSD+MSC		
9.00	0.		1.570E-03	1.182E-01	3.811E-03	9.428E-05	1.197E-01	3.905E-03	1.237E-01	
10.00	0.		7.964E-03	5.501E-01	1.206E-02	7.402E-05	5.581E-01	1.214E-02	5.702E-01	
11.00	0.		1.586E-02	1.024E+00	1.513E-02	1.967E-05	1.039E+00	1.515E-02	1.055E+00	
12.00	0.		2.300E-02	1.410E+00	1.399E-02	3.136E-06	1.433E+00	1.399E-02	1.447E+00	
13.00	0.		2.632E-02	1.562E+00	1.055E-02	3.026E-07	1.588E+00	1.055E-02	1.599E+00	
14.00	0.		2.200E-02	1.295E+00	6.371E-03	1.443E-08	1.317E+00	6.371E-03	1.323E+00	
15.00	0.		0.	3.628E-01	1.529E-03	1.443E-10	3.628E-01	1.529E-03	3.643E-01	
OSUMS	0.		9.671E-02	6.322E+00	6.345E-02	1.914E-04				

ANGULAR DISTRIBUTIONS Z= 1 N= 0

ENERGY PARAM. = EPS+B/E.

CROSS SECTIONS IN MB/STR-MEV

	10	20	30	40	50	60	70	80	90		
- EPS 0	9.00	2.15E-02	2.12E-02	2.03E-02	1.89E-02	1.72E-02	1.53E-02	1.34E-02	1.15E-02	9.90E-03	8.48E-03
10.00	1.02E-01	1.00E-01	9.58E-02	8.91E-02	8.08E-02	7.16E-02	6.23E-02	5.35E-02	4.56E-02	3.87E-02	3.10E-02
11.00	1.93E-01	1.90E-01	1.81E-01	1.68E-01	1.52E-01	1.34E-01	1.16E-01	9.92E-02	8.40E-02	7.10E-02	
12.00	2.71E-01	2.67E-01	2.54E-01	2.36E-01	2.12E-01	1.87E-01	1.61E-01	1.36E-01	1.15E-01	9.64E-02	
13.00	3.07E-01	3.02E-01	2.87E-01	2.65E-01	2.38E-01	2.09E-01	1.79E-01	1.51E-01	1.26E-01	1.05E-01	
14.00	2.60E-01	2.56E-01	2.43E-01	2.24E-01	2.00E-01	1.75E-01	1.49E-01	1.25E-01	1.04E-01	8.62E-02	
15.00	7.33E-02	7.21E-02	6.84E-02	6.29E-02	5.61E-02	4.87E-02	4.13E-02	3.45E-02	2.85E-02	2.34E-02	
- EPS 100	1.10	1.20	1.30	1.40	1.50	1.60	1.70	1.80	1.90	TOTAL	
9.00	7.29E-03	6.34E-03	5.58E-03	5.00E-03	4.56E-03	4.25E-03	4.04E-03	3.91E-03	3.87E-03	3.91E-03	1.24E-01
10.00	3.31E-02	2.85E-02	2.49E-02	2.22E-02	2.01E-02	1.86E-02	1.76E-02	1.70E-02	1.68E-02	5.70E-02	
11.00	6.02E-02	5.15E-02	4.47E-02	3.95E-02	3.55E-02	3.27E-02	3.07E-02	2.96E-02	2.92E-02	1.05E+00	
12.00	8.12E-02	6.90E-02	5.95E-02	5.22E-02	4.67E-02	4.27E-02	4.00E-02	3.84E-02	3.79E-02	1.45E+00	
13.00	8.82E-02	7.45E-02	6.38E-02	5.56E-02	4.94E-02	4.50E-02	4.19E-02	4.02E-02	3.96E-02	1.60E+00	
14.00	7.17E-02	6.01E-02	5.12E-02	4.43E-02	3.92E-02	3.55E-02	3.30E-02	3.15E-02	3.11E-02	1.32E+00	
15.00	1.94E-02	1.61E-02	1.37E-02	1.18E-02	1.04E-02	9.33E-03	8.64E-03	8.24E-03	8.11E-03	3.64E-01	

1PARTICLE SPECTRA Z= 2, N= 2
 FIRST EMISSION AT P= 4
 TARGET Z= 92, N=143 PROJ Z= 1, N= 0
 PO= 2, HO= 1, G=16.620, E=20.040
 PAIR EXCIT. ACCORDING TO 2 COMP. ST. DENSITIES
 REACTION CROSS SECTION = 743.2
 SCALE FACTOR= 1.350, FRAC PREEQ= 0.255
 MSD POSSIBLE= 0.874, USED= 0.214
 CLOSED FORM SUM STARTS AT P= 4

	DIRECT		PREEQUILIBRIUM		EQUIL		TOTAL (MB/MEV)		
	--	--	--	--	--	--	MSD	MSC	MSD+MSC
	NUTRA	KNOCK	MSD	MSC	WEISS				
21.00	5.665E-02	8.147E-02	1.319E-04	6.297E-05	2.667E-05	1.383E-01	8.964E-05	1.383E-01	
22.00	8.886E-02	1.555E-01	1.570E-04	5.829E-05	6.133E-06	2.445E-01	6.442E-05	2.446E-01	
23.00	7.333E-02	1.553E-01	1.055E-04	3.221E-05	6.527E-07	2.287E-01	3.286E-05	2.288E-01	
24.00	2.867E-02	0.	3.577E-05	9.743E-06	3.202E-08	2.870E-02	9.775E-06	2.871E-02	
OSUMS	2.475E-01	3.923E-01	4.302E-04	1.632E-04	3.348E-05				
-FISS.	0.	0.	0.	0.	3.543E+02				

ANGULAR DISTRIBUTIONS Z= 2 N= 2
 ENERGY PARAM. = EPS+E_c
 CROSS SECTIONS IN MB/STR-MEV

- EPS	0	10	20	30	40	50	60	70	80	90
21.00	2.54E-02	2.50E-02	2.39E-02	2.22E-02	2.00E-02	1.77E-02	1.53E-02	1.31E-02	1.10E-02	9.31E-03
22.00	4.60E-02	4.52E-02	4.32E-02	3.99E-02	3.60E-02	3.16E-02	2.72E-02	2.31E-02	1.94E-02	1.63E-02
23.00	4.40E-02	4.33E-02	4.12E-02	3.81E-02	3.42E-02	2.99E-02	2.56E-02	2.16E-02	1.81E-02	1.51E-02
24.00	5.65E-03	5.55E-03	5.28E-03	4.87E-03	4.36E-03	3.80E-03	3.24E-03	2.72E-03	2.26E-03	1.87E-03
- EPS	100	110	120	130	140	150	160	170	180	TOTAL
21.00	7.88E-03	6.72E-03	5.81E-03	5.10E-03	4.56E-03	4.17E-03	3.90E-03	3.74E-03	3.69E-03	3.38E-01
22.00	1.37E-02	1.16E-02	9.99E-03	8.72E-03	7.77E-03	7.07E-03	6.60E-03	6.32E-03	6.23E-03	2.45E-01
23.00	1.26E-02	1.06E-02	9.08E-03	7.89E-03	6.99E-03	6.34E-03	5.89E-03	5.64E-03	5.55E-03	2.29E-01
24.00	1.55E-03	1.30E-03	1.11E-03	9.56E-04	8.44E-04	7.62E-04	7.07E-04	6.74E-04	6.64E-04	2.87E-02

APPENDIX D
NEW FEATURES IN PRECO-D2

1. Pairing energies are read in separately for the preequilibrium and equilibrium phases of the reaction. (No pairing energies were read in in PRECO-D.)
2. Different effective well depths for states with $h=1$ and $h=2$ may be read in. (The well depth was always 38 MeV in PRECO-D.)
3. Default values are supplied for p_o , h_o , g_o , and K .
4. The composite nucleus-formation cross section and the inverse-reaction cross sections may be generated internally using a parametric approximation to the optical-model cross sections. See Section VI.7 for the optical potentials which are approximated.
5. For nucleon transfer reactions, the residual configurations in which transfer produces passive particles and/or holes are not explicitly considered, because they are already counted in the general-state density formula.
6. For knockout and inelastic-scattering reactions involving cluster degrees of freedom, the single-cluster-state densities have been reduced. In PRECO-D they were derived by considering the number of ways in which the constituent nucleons could be put into correlated single-nucleon states. The desired quantity is, instead, given as [<# cluster states]/[MeV carried by cluster], which adds

a factor of $1/A_{\text{cluster}}$. In addition, the quantity of interest is really the number of clusters that can be accommodated per MeV of cluster energy, not the number of ways of putting the constituent nucleons in a given cluster into the correlated single-nucleon states. This adds a factor of $1/2$ for deuterons, tritons and ^3He ions. See Section II.2 for the current single-cluster-state densities.

7. The closed-form reaction equations have been corrected to guarantee conservation of strength.

8. Either ϵ or $\epsilon+B_b$ may be used as the energy parameter for calculating angular distributions. (In PRECO-D the energy parameter was always taken to be ϵ .)

9. Either $Q_b(p)$ or $Q_b^{(G)}(p)$ may be used for handling proton/neutron distinguishability in calculating particle emission rates. (In PRECO-D only $Q_b(p)$ was available.)

10. For each particle spectrum, output for a given ϵ is suppressed if all of the component cross sections are zero.