

NOTES ON ECIS94

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Summary :

This report includes all the “Notes” written on different versions of the Code **ECIS**. The Chapter I relates the evolution of these codes since the first study of how to solve coupled-channels equation by the method of “Equations Couplées en Itérations Séquentielles”. The Chapter II indicate how these codes have been used on various computers, because such indications can help to solve future problems.

Chapter III present the “generalised optical model” in the Dirac and the Schrödinger formalism and the numerical methods. Chapter IV deals with the special treatment of long range interaction. Possibilities to describe excitation of particle and target are the subject of Chapter V and zero-range transfer reactions are presented in Chapter VI. Miscellaneous topics are gathered in Chapter VII and Chapter VIII is a description of the last version of the code, called **ECIS94**.

Sommaire :

Ce rapport reprend toutes les “Notes” écrites sur les différentes version du programme **ECIS**. Le Chapitre I retrace l'évolution de ces programmes depuis la première étude de la résolution des équations de voies couplées par la méthode des “Equations Couplées en Itérations Séquentielles”. Le Chapitre II montre comment ces programmes ont été utilisés sur différents ordinateurs car ces indications peuvent être précieuses pour l'avenir.

Le Chapitre III présente le “modèle optique généralisé” dans le formalisme de Dirac et de Schrödinger ainsi que les méthodes numériques. Le sujet du Chapitre IV est le traitement des interactions à longue portée. La possibilité de décrire l'excitation de la particule et de la cible est étudiée dans le Chapitre V et celle de réaction de transfert avec portée nulle dans le Chapitre VII. Divers sujets sont rassemblés dans le Chapitre VIII et le Chapitre IX est la description de la dernière version du programme, appelée **ECIS94**.

Chapter I

Introduction - History

The name “ECIS” is made of the first letters of “Equations Couplées en Itérations Séquentielles”, by reference to the method of solution used in the codes, although the usual method of solution is also present and has been written with as much care.

A Numerical methods

The numerical methods are a generalisation of those studied for the optical model in : [1] MELKANOFF, M.A., SAWADA, T. and RAYNAL, J., “*Nuclear Optical Model Calculations*”, published in “Methods in Computational Physics.6: Nuclear Physics” ALDER, B., FERNBACH, S. and ROTENBERG, M., eds. (Academic Press, New York, 1966) page 1. The application to coupled equations has been presented in : [2] RAYNAL, J., “*Optical-Model and Coupled-Channel Calculations in Nuclear Physics*”, published in “Computing as a Language of Physics”, ICTP International Seminar Course, Trieste, Italy, Aug.2-10, 1971 (IAEA, 1972), page 281.

A.1 Integration methods

The programme uses integration methods related to the Numerov method described in : [3] NOUMEROV, B. V., “*A Method of Extrapolation of Perturbations*”, Monthly Notices Roy. Astr. Soc. **84** (1924) page 592. The Numerov method do not deal with the function but with a combination of the function and its second derivative depending on the step size.

Unhappily, the name of Numerov method is used in many Numerical Analysis books, in many articles of Journal of Computational Physics and of Computational Physics Communications for another method which deals only with the functions and which is named Cowell method in Ref. [1] and [2] as it can been found in : [4] COWELL, P. H. and CROMMELIN, A. C. D., Appendix to Greenwich Observations for 1909, Edinburgh (1910) page 84.

The text of Ref. [3] is unambiguous on this point. NOUMEROV presents his method for an inhomogeneous equation (see Equation (20) of Ref. [2]) and begins by :

“I have pointed out in my previous work* the advantage which is to be got in the numerical integration of the equations of perturbed motion by the introduction of special rectangular co-ordinates. In fact, if we denote by \bar{x} , \bar{y} , \bar{z} the special co-ordinates connected with ordinary heliocentric rectangular co-ordinates by the equations

$$\bar{x} = x \left(1 + \frac{k^2 \omega^2}{12r^3} \right); \quad \bar{y} = y \left(1 + \frac{k^2 \omega^2}{12r^3} \right); \quad \bar{z} = z \left(1 + \frac{k^2 \omega^2}{12r^3} \right) \quad (I.1)$$

*B. Noumerov, “Méthode nouvelle de la détermination des orbites et le calcul des éphémérides en tenant compte des perturbations”, *Publications de l'Observatoire Central de Russie*, vol. ii., Moscou, 1923.

where k is the Gaussian constant, ω the interval of integration, and r the radius vector,...

The reference [5] given above by NOUMEROV himself is quite difficult to find, but Ref. [3] can be found in any Astronomy library.

In astronomical problems, NOUMEROV had to solve a third order equation at each step. In coupled channel calculations, the use of his method leads to solve a system of linear equations at each step. In Ref. [1] was presented a method which replaces this resolution by a matrix multiplication. This “Modified Numerov method” has a truncation error of the same order, but of reversed sign for a constant potential. Details, comparison with methods used by other authors and discussions can be found in Ref. [2].

The programme **ECIS** uses the “Modified Numerov method” of integration with the usual methods of coupled channels calculations (**LO(21)=.TRUE.**), but only in the Schrödinger formalism. With iterations (**LO(21)=.FALSE.**), which are compulsory with deformed spin-orbit and in Dirac formalism, it uses the “Modified Numerov method” if **LO(27)=.FALSE.** or the Numerov method if **LO(27)=.TRUE.** The option **LO(26)=.TRUE.** (**LO(30)=.TRUE.** in **ECIS79**) given as “integration stabilised for long range potentials” means the inclusion of the truncation error as described in Ref. [2], assuming a constant potential; for usual coupled equations (**LO(21)=.TRUE.**), only the diagonal potential is taken into account; no improvement of the inelastic scattering has to be expected from that and it effects mainly the elastic result.

More details will be given in section (III.C).

A.2 Iteration methods

The iteration methods are described Ref.[2]. The first description was in : [6] RAYNAL, J., “*Equations couplées et D.W.B.A.*”, published in “Sur quelques Méthodes de Physique Nucléaire Théorique” Aussois (France) Feb 26 - March 2 1968 (Institut de Physique Nucléaire de Lyon) (LYCEN-6804) pages 179 and another one can be found in : [7] RAYNAL, J., “*Recurrence relations for distorted-wave Born approximation Coulomb excitation integrals and their use in coupled channel calculations*”, Phys. Rev. **C23** (1981) page 2571.

A.2.a Scheme of iterations

The principle of the iterations is to write all the non-diagonal terms as inhomogeneous terms of single equations. The procedure is obtained by considering an expansion of the solution in powers of the strength of these inhomogeneous terms and by setting a recurrence relation between different powers. However, this notion of power expansion is not respected because the last known solution is used to compute the inhomogeneous term with the hope to obtain a quicker result. The computation will depend upon the order of equations: it is why these iterations are called “sequential”.

In **ECIS79**, if **LO(26)=.FALSE.**, the inhomogeneous differential equations are solved: (“Differential method”). If **LO(26)=.TRUE.**, an irregular solution of the single inhomogeneous equations is obtained, a Green’s function is built with it and the “optical” solution and its integral with the inhomogeneous term computed (“Integral method”). There is no noticeable difference of time between these two methods. However, the solution by the “Differential method” can involve differences of large numbers and fail to give the good result: it is the case when there is a closed channel with a large Q value, for which the “Integral method” gives good results.

However, since the code **ECIS88** as described by [8] RAYNAL, J., “*Coupled Channel Calculations and Computer Code ECIS*”, published in “Workshop on Applied Nuclear Theory and Nuclear Model Calculations for Nuclear Technology Applications”, Trieste, Italy, Feb. 15-March 18, 1988,(Trieste, 1988), MEHTA, M., K. and SCHMIDT, J., J., eds. (World Scientific, 1989) page 506, the generalisation of the Optical Model in the Dirac phenomenology introduced by [9] ARNOLDS, L. C., CLARK, B. C., MERCER, R. L. and SCWANDT, F., “*Dirac potential model analysis of \bar{p} - ^{40}Ca elastic scattering at 180 MeV and the wine-bottle-bottom shape*”, Phys. Rev. **C23** (1981) page 1949 can be used. The equations to solve have been described in : [10] RAYNAL, J., “*Coupled Channels Description of Inelastic Scattering*

with a Dirac Equation”, published in “Proceedings of the 6th International Symposium on Polarization Phenomena in Nuclear Physics”, Osaka, Japan, Aug. 26-30, 1985, KONDO, M., KOBAYASHI, S. and TANIFUJI, M., eds., J. Phys. Soc. Japan **55** (Suppl.) (1986) page 922, in : [11] RAYNAL, J., “*Formalisme des Voies Couplées et Programmes ECIS*”, published in “Réunion des Spécialistes sur l’Utilisation du Modèle Optique pour le Calcul des Sections Efficaces Neutroniques au-dessous de 20MeV”, NEANDC-222 “U”, Paris, France, Nov. 13-15, 1985 (OCDE, 1986) page 63 and in : [12] RAYNAL, J., “*Ambiguity on the imaginary potentials in the Dirac formalism for the elastic and the inelastic scattering of nucleons*”, Phys. Lett. **B196** (1987) page 7. If these equations are correctly solved and not approximated as described in : [13] RAYNAL, J. and SHERIF, H., S., “*Comparison between Dirac Equation and its Equivalent Schrödinger Equation for Inelastic Scattering*”, published in “Proceedings of the 6th International Symposium on Polarization Phenomena in Nuclear Physics”, Osaka, Japan, Aug. 26-30, 1985, KONDO, M., KOBAYASHI, S. and TANIFUJI, M., eds., J. Phys. Soc. Japan **55** (Suppl.) (1986) page 924, it is difficult to include the reorientation terms in the left side (uncoupled equations) of the iteration procedure. So, since **ECIS88**, the reorientation terms are not used in the calculation of the uncoupled solutions, except if **LO(29)=.TRUE.**, which can be used only in the Schrödinger formalism.

Because the “Differential method” of iterations has no practical equivalent for Dirac equations and no clear advantage in the Schrödinger formalism, there is only the “Integral method” (Green’s function) since the code **ECIS88**. (The option **LO(29)** of **ECIS79**, “matching with derivatives” has also been discarded).

A.2.b Pade approximants

This iteration procedure is not converging in all cases. Problems arise chiefly for neutron scattering at very low energy or for heavy ions scattering for the “grazing” J -value. After four iterations, the code looks after the convergence of Pade approximants (see in : [14] PADE, H., “*Sur la représentation approchée d’une fonction par des fractions rationnelles*”, Ann. Sci. Ec. Norm. Sup. Paris **9** (1892) page 1 and **16** (1899) page 395 and in : [15] WALL, H. S., “*Continued Fractions*”, Van Nostrand, New York 1948) built with the results already obtained.

To explain the role of the Pade approximants let us multiply the inhomogeneous terms by some factor x . There are some values (complex) of x for which there are solutions of the system which are purely outgoing for all the equations: there are the **Weinberg states**, and the x ’s are the Weinberg eigenvalues. The solution of the system of coupled equations can be expressed with Weinberg eigenfunctions and eigenvalues. If one does not use the “sequential” method but does an exact power expansion, this expansion converge only when the smallest $|x|$ is smaller than unity. The Pade approximant generates poles which are the Weinberg eigenvalues and can be smaller than unity. It can be verified, in case of non-convergence that the poles of the Pade approximants with small absolute value are the same for all the equations. The use of the sequential approach do not change this situation.

The ability of Pade approximants to construct the poles from the power series depends drastically of the precision of the intermediate calculations. The limit is about $|x| > 1/6$ and is better with the **CDC** version of the programme than with the **IBM** version. It can be useful to use the **CDC** version on an **IBM** after doubling its precision. However, for low energy neutrons, the computation with the usual methods (**LO(21)=.TRUE.**) can be quite quick and can be used if one is not interested by spin-orbit deformation.

If **LO(22)=.TRUE.**, Pade approximants are not used. If **LO(22)=.FALSE.** Pade approximants are computed after the fourth iteration if there is no natural convergence and if convergence has been obtained for all the precedent equations in this iteration. There is convergence when two Pade approximants differ by less than the precision required, even if there are not the ones evaluated with the maximum number of parameters. The iteration procedure stops when the maximum number of iterations is reached or when a phase-shift in the last iteration was larger than 10^{10} because there is no more hope to find a precise result.

When convergence has not been obtained with Pade approximants, if **LO(23)=.TRUE.**, the equations for this J -value and parity are solved by usual methods as if **LO(21)=.TRUE.** and a warning printed. On the contrary, if **LO(23)=.FALSE.**, in the last iteration, the Pade approximants are computed even if the precedent equations did not converge and the mean value of the two nearest results is kept as the solution. The shift to usual methods of integration is forbidden in Dirac formalism.

See also section (III.D.4).

A.2.c Increase of the imaginary potential

In low energy neutron scattering on a target with $0^+ - 2^+ - 4^+$ rotational band, there are usually convergence problems which come from a large difference between coupled and uncoupled solutions of the Schrödinger equations: the imaginary potential needed in coupled channel equations is weaker than the one needed to describe elastic scattering. An attempt to avoid this problem is to use a stronger imaginary potential and to introduce the difference as a reorientation term. For that, the reorientation terms must be in the second member (`LO(29)=.FALSE.`). The ratio of increase is a data of the input. Some examples are given in Ref. [8] and Ref. [11].

Advantages of this process in heavy-ion inelastic scattering have not been studied.

B History of different versions

The Buck and Hill's code **INCH** for coupled channels was the starting point of **ECIS**. Some features remains, chiefly in the input of data. The names given to the different versions are those with which they have been kept on a tape for the oldest one, and those which have been explicitly introduced into the code for the latest ones.

B.1 Earlier codes

B.1.a Code ECIS68

In **ECIS68** the integration method of the code **INCH** was changed into the Numerov Method with matching at two points as described in Ref. [2] in order to get more precise results. The geometrical coefficients were replaced by those written at the Département de Calcul Electronique of Saclay by : [16] LAFON, R., "*Sous-Programmes DFCG-DFC3J, DFR6J et DFR9J (IBM 360)*", Report DCE-Saclay No. 326 (1967). There was only the "Differential **ECIS** method", including Pade approximants. The use of a large working space cut into arrays by calls to subroutines to avoid fixed dimensions, was already introduced. This code was used to compare the **ECIS** method to the usual methods. There was no comparison with experimental data. Results can be found in Ref. [6] and in : [17] RAYNAL, J., "*An iterative procedure for coupled channel calculations*", Communication IV-129 in "Dubna publication D-3893" International Symposium on Nuclear Structure, Dubna, URSS (1968).

B.1.b Code ECIS69

In **ECIS69**, the Coulomb functions for the open channels were replaced by those written at the Department de Calcul Electronique Saclay by : [18] BARDIN, C., DANDEU, Y., GAUTIER, L., GUILLERMIN, J., LENA, T. and PERNET, J.M., Note CEA-N-906 (1968) and [19] BARDIN, C., DANDEU, Y., GAUTHIER, C., GUILLERMIN, J., LENA, T., PERNET, J.-M., WOLTER, H. H. and TAMURA, T., "*Coulomb functions in entire (η, ρ) -plane*", Comp. Phys. Comm. **3** (1972) page 72. The code was largely rewritten to take into account the full Thomas form for the spin-orbit deformation. In peculiar, the helicity formalism as in : [20] JACOB, M. and WICK, G. C., "*On the General Theory of Collisions for Particles with Spin*", Annals of Physics **7** (1959) page 404 or in : [21] RAYNAL, J., "*Aspects Géométriques des Réactions*", Note CEA-N-1529 (1972) (see also in : [22] RAYNAL, J., "*Utilisation de Faisceaux de Deutons Polarisés et Détermination des Paramètres du Modèle Optique*", Thesis, Faculté d'Orsay, June 24 1964 (Rapport CEA-N-2511-1964) and translation ANL-TRANS-258, the Chapter II.) was introduced for the amplitudes. A 20 points Gauss-Legendre integration was used for the form factors of the rotational model. Results obtained were reported in : [23] RAYNAL, J., "*Couplage LS dans les descriptions macroscopiques et microscopiques des réactions nucléaires*", published in "Symposium sur les

Mécanismes de Réaction Nucléaire et Phénomènes de Polarisation” Québec, Canada, Sept. 1-2, 1969 (Les Presses de l’Université Laval, 1970) page 75 and in : [24] RAYNAL, J., “*Potentiel spin-orbite déformé en équations couplées (abstract)*”, published in “International Conference on Properties of Nuclear States”, Montreal, Canada, Aout 25-30, 1969 HARVEY, M., CUSON, R. Y., GEIGER, J., S. and PEARSON, J., M. eds. (Les Presses de l’Université de Montréal, 1969) page 771.

B.1.c Code ECIS70

The version **ECIS70** is shortly described in Ref. [2]. The asymmetric rotational model is added: the form factors are calculated by integration on the sphere with 36 points, the weights of which were obtained once for all by the inversion of the matrix of rotation matrix elements at these points (this procedure means that the potential is supposed to be expanded only with 36 rotation matrix elements, the coefficients of which are obtained by solving a set of 36 linear equations). The number of multipoles is limited to 15 (id est $\lambda = 8$, whereas 36 is $\lambda = 14$). In the usual method to solve coupled equations, a Schmidt’s orthogonalisation procedure is introduced every n points to avoid a loss of independence between the solutions. There was also an attempt to write on a scratch tape the couplings between equations. These couplings are sums of form factors multiplied by a geometrical coefficient: if they are computed before for the iteration method, the time can be divided by two in the rotational model. The use of a scratch tape turned out to be worse than the computation at each iteration. Results obtained at that time were presented in : [25] RAYNAL, J., “*Effets de l’interaction L.S nucléon-nucléon sur l’asymétrie de la diffusion inélastique de protons*”, Colloque sur les Mécanismes des Réactions Nucléaires, Grenoble, France, March 16-18, 1970, GUGENBERGER, P., eds. J. Physique **31C2** (1970) page 92, in : [26] LOMBARD, R., M., MAYER, B. and RAYNAL, J., “*Calculs d’asymétrie de protons polarisés de 20,3 MeV sur ^{24}Mg , ^{26}Mg et ^{28}Si* ”, Colloque sur les Mécanismes des Réactions Nucléaires, Grenoble, France, March 16-18, 1970, GUGENBERGER, P., eds. J. Physique **31C2** (1970) page 90 and in : [27] RAYNAL, J., “*Nuclear structure effects on asymmetry of proton scattering*”, published in “Polarization Phenomena in Nuclear Reactions: Proceedings of the 3rd International Symposium”, Madison, USA, Aug.31 - Sept.4, 1970 BARSCHALL, H., H. and HAEBERLI W., eds. (The University of Wisconsin Press, 1971) page 798.

ECIS70 was given to Karlsruhe where Dr. G. SCHWEIMER adapted it for automatic search on alpha inelastic scattering. In this code the spin orbit deformation was limited to spin one half, only cross-section, polarisation, analysing power and spin-flip were computed.

B.1.d Code ECIS71

The **ECIS71** code was an attempt to introduce automatic search in **ECIS70**. The integral version of the **ECIS** method was added and turned out to be equivalent to the differential version at the point of view of time. The spin orbit deformation is extended to any spin. The different attempts of automatic search are described in Ref. [2].

With **ECIS71** began the rewriting of the subroutines for integration, geometrical coefficients and reduced rotation matrix element in IBM Assembler language. A subroutine written by RENARDY defines as working array all the region left free by the programme and allows the use of a **LOAD MODULE** with different sizes. At that time were written Ref. [2] and [28] RAYNAL, J., “*Spin-Orbit Interaction in Inelastic Nucleon Scattering*”, published in “The Structure of Nuclei”, International Course on Nuclear Theory, Trieste, Jan. 13 - March 12, 1971, (IAEA,1972) page 75. Results can be found in : [29] RAYNAL, J., “*Interaction spin-orbite dans la diffusion inélastique de nucléons*”, published in “Sur Certains Aspects Microscopiques des Réactions Nucléaires”, La Toussuire, France, Feb. 15-19, 1971 (IPN, Université Claude Bernard, Lyon I) LYCEN 7104, page C7.1 and in : [30] De SWINIARSKI, R. and RAYNAL, J., “*Déformations dans la couche s-d et diffusion inélastique de protons polarisés*”, published in “Sur Certains Aspects Microscopiques des Réactions Nucléaires”, La Toussuire, France, Feb. 15-19, 1971 (IPN, Université Claude Bernard, Lyon I) LYCEN 7104, page S9.1.

B.1.e Code ECIS72

In **ECIS72**, there were further attempts of automatic search in the same approach as in **ECIS71**. The vibrational model is generalised to mixtures of one phonon and two phonons states. Publications of 1972 are in : [31] LOMBARD, R., KAMITSUBO, H., RAYNAL, J. and GOSSET, J., “*Etude des états collectifs de ^{32}S et de ^{34}S par diffusion inélastique de protons polarisés*”, C. R. Acad. Sci. **B274** (1972) page 761, in : [32] De SWINIARSKI, R., BACHER, A., D., RESMINI, F., G., PLATTNER, G., R., HENDRIE, D., L. and RAYNAL, J., “*Determination of deformation parameters of ^{20}Ne and ^{22}Ne by inelastic scattering of polarized protons*”, Phys. Rev. Lett. **28** (1972) page 1139 and in : [33] RAYNAL, J., “*Sur l’influence de l’interaction spin-orbite à haute énergie*”, published in “Compte-Rendu de la Conférence Européenne de Physique Nucléaire. Vol 2”, Aix-en-Provence, France, June 26 - July 1, 1972, page 149.

B.1.f Code ECIS73

In **ECIS73**, the observables which can be taken into account for automatic search are generalised to any one, in the notations of Ref. [21]. They can be defined in the laboratory system or with an axis of quantification perpendicular to the reaction plane. Publications of 1973 are in : [34] De SWINIARSKI, R., GENOUX-LUBAIN, A., BAGIEU, G., CAVAGNAC, J., F., WORLEDGE, D., H. and RAYNAL, J., “*A coupled-channels analysis of 30 MeV proton scattering from low-lying positive-parity states in ^{19}F , ^{20}Ne , ^{21}Ne* ”, Phys. Lett. **43B** (1973) page 27, in : [35] KUREPIN, A., B., LOMBARD, R., M. and RAYNAL, J., “*Method for identification of the nuclear collective modes*”, Phys. Lett. **45B** (1973) page 184 and in : [36] LOMBARD, R., M. and RAYNAL, J., “*Polarized-proton inelastic scattering on ^{32}S and possible evidence for an hexadecapole phonon state*”, Phys. Rev. Lett. **31** (1973) page 1015.

In the same time, Dr G. SCHWEIMER used **ECIS** in Karlsruhe and introduced an automatic search on parameters, using methods very different of those of Ref. [1] and [2]. He introduced also :

- the folding model,
- angular distribution which are sums of levels

but he was not interested in polarisation effects and could deal only with one cross-section for each level. He used different subroutines with the same name to treat different problems: for example, the difference between rotational and vibrational model is obtained by loading different subroutines for nuclear matrix elements and for form factors.

B.1.g Code ECIS75

In June 1973, in collaboration with Dr. G. SCHWEIMER, best points of the the Karlsruhe version were included to the code, giving **ECIS75**. Each subroutine includes many comment cards, and in peculiar explanation of inputs and outputs at their beginning. This was already done by Dr. G. SCHWEIMER in his KARLSRUHE version of **ECIS**. All the cards are identified by the four first characters of the subroutine and a sequence number in columns 73-80.

All personal attempts of automatic searches were stopped: the subroutines developed by Dr. G. SCHWEIMER in KARLSRUHE were introduced in the code. However, the use of different modules instead of loading different models at the same time with overlay, was not adopted. The folding model was introduced, after generalisation to the spin-orbit potential. Its meaning is the folding of a nucleon-nucleus potential with the intrinsic wave function of the incoming particle (the meaning of the KARLSRUHE version was the folding of the density distribution of the target with an alpha-nucleon interaction). The folding distribution can be Gaussian, the sum of two Yukawa or a Woods-Saxon distribution.

Other features from the KARLSRUHE version intruded in **ECIS75** are :

- unresolved angular distributions,

- constrained asymmetric rotational model,
- anharmonic vibrational model,
- use of the mean value of the two nearest Pade results.

Some other details were added independently:

- J-dependence of imaginary potentials (dropped since **ECIS78**),
- symmetrised χ^2 for the cross-sections,
- factorisation of $(1 - x \cos \theta)$ in the amplitudes,
- possibility to save a search on a tape, if limited by the time.

A description of the **INPUT** was written on cards (681 cards). All these cards images were inserted in **FORMATs** in a programme, which could be used to reproduce the listing of these cards. However, this description was not inserted in the programme itself.

Publications of 1975 are in : [37] RAYNAL, J., “*Application des équations couplées*”, published in “3ème Session d’Etudes Biennale de Physique Nucléaire”, La Toussuire, France, Feb. 10-15, 1975 (IPN, Université Claude Bernard, Lyon I) LYCEN 7502, page C7.1 and in : [38] RAYNAL, J., “*Inelastic Scattering (Coupled Channels)*”, published in “Proceedings of the 4th International Symposium on Polarisation Phenomena in Nuclear Reactions”, Zurich, Switzerland, Aug. 25-29 1975, GRUEBLER, W. and KONIG, V., eds., (Birkhauser, 1976) page 677.

B.1.h Code ECIS76

The code **ECIS** was adapted on a **CDC** computer by Dr. Marek SIEMASKO, from Katowice. At the same time, Dr. M. SIEMASKO introduced Hauser-Feshbach corrections for compound nucleus. To do that in the **IBM** versions, eight subroutines had to be modified, one of which was also translated into Assembler language. The resulting programme is **ECIS76**, with his two mean versions, one for low precision computers as the **IBM** and the other for large precision computers as **CDC**. However, due to the use of **LOGICAL*1** and **INTEGER*2**, the **IBM** version cannot be run on other computers as **UNIVAC**, **IRIS** or Japanese **FACCOM** : some modifications had to be done on an **IRIS** computer at Bordeaux, a **SIEMENS** in Berlin, a **FACCOM** in Japan and an **UNIVAC** in Copenhagen, leading to a third version, the **UNIVAC** one.

B.2 Codes with Coulomb corrections

B.2.a Code ECIS78

With the introduction of Coulomb corrections for heavy ions, as described in Ref. [7], the code has been enlarged. In peculiar, in collaboration with Prof. H. V. von GERAMB, the possibility to use external form factors has been added, but these form factors can be read only by points. Publication of 1978 in : [39] RAYNAL, J., “*Déformation de l’interaction spin-orbite: relations avec la structure nucléaire et effets possibles sur la section efficace de diffusion inélastique*”, published in “La Physique Neutronique et les Données Nucléaires: Compte-Rendu”, Harwell, GB, Sept. 1978 (OCDE, 1978) page 372.

B.2.b Code ECIS79

In **ECIS79**, the management of memories on **CDC** or **UNIVAC** computers has been changed. In the previous versions, it was possible to use a **LOAD MODULE** with different sizes with respect to the problem, but this size had to be known beforehand. In **ECIS79**, the core is requested when it is needed, as long as it is available. However, such procedure cannot be used on an **IBM** computer on which this management of memories must stay dummy.

The possibility of external potentials has been extended by the use of standard form factors as Woods-Saxon and its derivatives, rotational form factors, Laguerre polynomials and bound states in a potential.

The code **ECIS79** was given to the Nuclear Data Bank of the OCDE in 1982.

Publications obtained with it are in : [40] RAYNAL, J., “*Coupled channel calculations of heavy ion inelastic scattering with DWBA approximation for Coulomb interaction (communicated paper)*”, International Conference on Nuclear Physics, Berkeley, USA, Aug. 24-30 1980, in : [41] RAYNAL, J., “*Coulomb effects in proton inelastic scattering on heavy target (communicated paper)*”, International Conference on Nuclear Physics, Berkeley, USA, Aug. 24-30 1980 and in : [42] RAYNAL, J., “*Strong channel coupling method for cross-section calculations (lecture notes)*”, Workshop on Nuclear Model Computer Codes, Trieste, Italy, Jan. 16 - Feb. 3, 1984. For heavy ions problems, they are in : [43] BILWES, B., BILWES, R., BAEZA, A., DIAZ, J., FERRERO, J., L. and RAYNAL, J., “*Inelastic scattering of ^{32}S on ^{28}Si* ”, 12th Winter Meeting on Nuclear Physics, Bormio, Italy, Jan. 23-27 1984, *Ricerca Scientifica ed Educazione Permanente* **35** (Suppl.) (1984) page 422, in : [44] BAEZA, A., DIAZ, J., FERRERO, J., L., BILWES, B., BILWES, R. and RAYNAL, J., “*Mutual excitations of $^{32}\text{S} + ^{28}\text{Si}$ at 90 and 97.09 MeV*”, *Phys. Lett.* **149B** (1984) page 73 and in : [45] BAEZA, A., BILWES, B., BILWES, R., DIAZ, J., FERRERO, J., L. and RAYNAL, J., “*Inelastic scattering of ^{32}S on ^{28}Si* ”, *Nucl. Phys.* **A437** (1985) page 93.

B.3 Codes with Dirac formalism

B.3.a Codes ECIS85 to ECIS87

In the Spring 1984 began some attempt to extend the Dirac phenomenology to inelastic scattering with the collaboration of Pr. H. SHERIF who, for instance, provided comparison with other codes for the elastic scattering. One of the points was to avoid the use of relativistic Coulomb functions by using Coulomb corrections as described in Ref. [7]; anyway, relativistic Coulomb functions do not take into account the effect of the anomalous magnetic moment. There was no difficulty to compute these corrections as long as they were used only for elastic scattering. The formalism was presented in Ref. [11] and earliest results were reported in Ref. [10] and Ref. [13]. Difference between **ECIS86** and **ECIS87** is essentially external form-factors generalised in Hamburg with Pr. H. V. von GERAMB in November 1986. During the Spring of 1987, **ECIS87** was left in various places in Canada and United-States. Results are presented in Ref. [12], in : [46] De SWINIARSKI, R., PHAM, D., L. and RAYNAL, J., “*Analyse of 800 MeV inelastic polarized proton scattering from ^{16}O and ^{54}Fe through the coupled Schrödinger or Dirac equations formalism (abstract)*”, published in “Conference on Nuclear Physics and Particle Physics”, Birmingham, GB, April 6-8 1987 (The Institute of Physics) page 26 and in : [47] SHIM, S., CLARK, B., C., COOPER, E., D., HAMA, S., MERCER, R., L., RAY, L., RAYNAL, J. and SHERIF, H., S., “*Comparison of relativistic and nonrelativistic approaches to the collective model treatment of $p+^{40}\text{Ca}$ inelastic scattering*”, *Phys. Rev.* **C42** (1990) 1592.

B.3.b Code ECIS88

After completion of the relativistic Coulomb corrections, this code was described in Ref. [8] and given to the Nuclear Data Bank of the OCDE. However, some points were still missing, as the Bessel expansion of the form factors.

Results obtained in Dirac formalism or their comparison to results of the Schrödinger formalism can be found in : [48] RAYNAL, J., SHERIF, H., S., KOBOS, A., M., COOPER, E., D. and JOHANSSON, J., L., “*Dirac coupled channel calculations and nucleon scattering at large momentum transfer*”, *Phys. Lett.* **B218** (1989) page 403, in : [49] De SWINIARSKI, R., PHAM, D., L. and RAYNAL, J., “*Dirac coupled-channels analysis of inelastic scattering of 800 MeV polarized protons from ^{16}O , ^{24}Mg and ^{26}Mg* ”, *Phys. Lett.* **B213** page 247, in : [50] RAYNAL, J., “*Which potentials have to be surface peaked to reproduce large angle proton scattering at high energy*”, *Australian J. Phys.* **43** (1990) page 9, in : [51] RAYNAL, J., “*Inelastic scattering of protons at 800 MeV on ^{40}Ca at large angles*”, 7th International Conference on Polarization Phenomena in Nuclear Physics, Paris, France, July 9-13 1990, Abstracts of contributed papers, page 32F and in : [52] De SWINIARSKI, R., PHAM, D., L. and RAYNAL, J., “*Comparison*

of higher order deformations in several *s-d* shell nuclei obtained through Schrödinger and Dirac coupled-channel analysis of 900 MeV polarized protons inelastic scattering experiments”, Z. Physik **A343** (1992) page 179.

Results related to heavy ions are given in : [53] KIENER, J., GSOTTSCHNEIDER, G., GILS, H., J., REBEL, H., CORCIALCIUC, V., BASU, S., K., BAUR, G. and RAYNAL, J., “Investigation of the sequential break-up ${}^6\text{Li} \rightarrow {}^6\text{Li}^*(3_1+) \rightarrow \alpha + d$ of 156 MeV ${}^6\text{Li}$ projectiles on ${}^{208}\text{Pb}$ in the very forward angle hemisphere”, Z. Phys. **A339** (1991) page 489 and in : [54] CORCIALCIUC, V., REBEL, H., KIENER, J., GSOTTSCHNEIDER, G., GILS, H., J., RAYNAL, J. and BAUR, G., “Analysis of the sequential break-up ${}^6\text{Li} \rightarrow {}^6\text{Li}^*(3_1+) \rightarrow \alpha + d$ of 156 MeV ${}^6\text{Li}$ projectiles on ${}^{208}\text{Pb}$ observed in the very forward angle hemisphere (abstract)”, Frühjahrstagung Darmstadt 1991 - Physik der Hadronen und Kerne, Darmstadt, Germany, March 11-15 1991, Verh. Dtsch. Phys. Ges. **6** (1991) page 6. But they still used primarily ECIS79.

B.3.c Codes ECIS89 to ECIS90

At the end of 1988, Dr A. Ventura of the Centro di Calcolo ENEA at Bologna, Italy, was interested by the Bessel expansion of form factors. This part of the code has been completed with him. Some other changes have been done :

- use of the deformation lengths instead of the deformations when fixed relative deformations lengths are wanted,
- use of “symmetrised” Woods-Saxon form factors.

These codes have been given to some laboratories.

B.3.d Code ECIS94

This version will be the third given to the Nuclear Data Bank of the OCDE. Besides the use of Bessel expansion for form factors, the use of deformation lengths and the use of “symmetrised” Woods-Saxon potentials, it includes :

- two bound states transitions for particle hole excitations, with the possibility of the particle in the continuum,
- expansion of cross-sections in terms of Legendre polynomials,
- possibility of angular distribution for uncoupled states without giving explicitly all the reduced nuclear matrix elements,
- for Coulomb excitation, use of a magnetic multipole.

Chapter II

Use on various computers

The code **ECIS** involves two versions with respect to the internal precision of the computer on which it is used :

- I If this precision is smaller than 10^{-8} , potentials and wave functions are stored in single precision but many quantities are in double precision: this happens for **IBM**, **UNIVAC**, **VAX**, **PRIME** computers.
- II If this precision is larger than 10^{-8} , as it is for the **CDC** and **CRAY** computers, the code is completely in single precision.

Some subroutines are identical between the two versions; for some others, the only difference is the change of the card **DOUBLE PRECISION** into comment card by adding C in column 1.

Note that it is possible to run the two versions on a **CONVEX** computer, using the “-cfc” option for the **CDC** version. In fact, it is the **CDC** version used this way which turns out to be the faster on a **CONVEX** computer.

The code **ECIS79** involved about 14250 **FORTRAN** cards, including 2300 **COMMENT** cards. It was intended to be run with an **OVERLAY** structure. The code **ECIS88** involved about 16700 **FORTRAN** cards, including 2500 **COMMENT** cards and the code **ECIS94** involves about 18500 **FORTRAN** cards, including 3340 **COMMENT** cards. Most of the topics related in this Chapter are related to the code **ECIS79**. **As their knowledge can help to solve future problem, these topics are not obsolete.**

Even if the use of **OVERLAY** is no more of interest, there are two topics very machine dependent in the code **ECIS**: the management of the working array and the control of the time limit of the job. How to by-pass these two points is explained on comment cards in the **MAIN** subroutine. For example, the **IBM** version of the **MAIN** subroutine of **ECIS79**, **ECIS88** and **ECIS94** is :

```
|C 24/09/79  IBM                                MAIN-000
|C THE ASSEMBLER SUBROUTINE ECIS(I) DEFINES THE DOUBLE PRECISION ARRAY W MAIN-001
|C AND CALLS CALC(W,W,W,IDMX) WHERE IDMX IS THE LENGTH OF W          MAIN-002
|C W IS ALL THE REGION LEFT FREE BY THE PROGRAM EXCEPT 2*I K FOR BUFFERS MAIN-003
|C THE SUBROUTINE ECIS CAN BE ELIMINATED WITH THE FOLLOWING MAIN:      MAIN-004
|C     REAL*8 W(10000)                                                MAIN-005
|C     CALL CALC(W,W,W,10000)                                         MAIN-006
|C     STOP                                                            MAIN-007
|C     END                                                            MAIN-008
|C THE ASSEMBLER SUBROUTINE STIM(I) HAS NO FORTRAN EQUIVALENT         MAIN-009
|C IT GIVES THE REMAINING TIME FOR THE JOB IN UNITS OF 26 MICROSECONDS MAIN-010
|C IT CAN BE REPLACED BY:      (ANY USE OF TIME WILL BE SENSELESS)    MAIN-011
|C     SUBROUTINE STIM(I)                                             MAIN-012
|C     DATA K /0/                                                  MAIN-013
```

C	K=K-1	MAIN-014
C	I=K	MAIN-015
C	RETURN	MAIN-016
C	END	MAIN-017
C*****		MAIN-018
	CALL ECIS(3)	MAIN-019
	STOP	MAIN-020
	END	MAIN-021

As shown above, each subroutine begins by a comment card giving the date of last revision and on which computer it can be used. All the cards are identified by the name of the subroutine in columns 73-76, a “—” in column 77 and by a number in columns 78-80, starting by 000.

A The IBM version

The dynamic allocation of memory is not possible on **IBM**.

A.1 Use on IBM360/370 - AMDHAL- SIEMENS-FUJITSU

The **FORTRAN** subroutines of **ECIS79** for integration and for geometrical coefficients has been translated into **ASSEMBLER** language. These subroutines involved around 7000 cards and only the 1200 cards of geometrical coefficients and reduced matrix element can still be used in **ECIS88** and **ECIS94**. The translation into **ASSEMBLER** language was done for an **IBM360/91** or a similar computer, on which **DO LOOP** smaller than 16 machine instructions are very quick. These **ASSEMBLER** subroutines can be used on other computers: their output statements do not work on a **SIEMENS-FUJITSU**. However, the **ASSEMBLER** subroutines **ECIS** and **STIM** are very machine dependent and cannot be replaced everywhere.

A.1.a Use with JCL

The subroutine **ECIS** defines as working array all the space available for the **JOB** in its region, except for twice its argument 'K' for **BUFFERS**. It is :

ECIS	TITLE '15/06/79-ECIS-CREATION DE ZONE DE TRAVAIL-ECIS79'	ECIS-000
ECIS	CSECT	ECIS-001
*	APPEL PAR CALL ECIS(N) OU N EST LA TAILLE EN DOUBLES K RESERVEE	ECIS-002
*	AUX BUFFERS.	ECIS-003
*	CE SOUS PROGRAMME CREE UNE ZONE DE TRAVAIL DE TAILLE MAXIMUM	ECIS-004
*	ET APPELE CALC EN LUI PASSANT TROIS FOIS LA ZONE CREEE ET SA TAILLE	ECIS-005
*	EN DOUBLE MOTS.	ECIS-006
	SAVE (14,12),T,*	ECIS-007
	USING ECIS,12	ECIS-008
	LR 12,15	ECIS-009
	SR 15,15	ECIS-010
	SPM 15	ECIS-011
	LA 4,SAVE	ECIS-012
	ST 13,4(4)	ECIS-013
	ST 4,8(13)	ECIS-014
	LR 13,4	ECIS-015
	L 1,0(1)	ECIS-016
	L 1,0(1)	ECIS-017
	SLL 1,11	ECIS-018
	ST 1,MIN	ECIS-019

	GETMAIN VU,A=ADGET,LA=MIN,SP=2	ECIS-020
FREE	L 0,MIN	ECIS-021
	O 0,=X'02000000' W IN SUBPOOL 2	ECIS-022
	FREEMAIN R,LV=(0),A=ADGET	ECIS-023
	L 0,ADGET+4	ECIS-024
	S 0,MIN	ECIS-025
	SRL 0,3	ECIS-026
	ST 0,LONG	ECIS-027
	L 1,ADGET	ECIS-028
	A 1,MIN	ECIS-029
	ST 1,ARG	ECIS-030
	ST 1,ARG+4	ECIS-031
	ST 1,ARG+8	ECIS-032
	LA 1,ARG	ECIS-033
	L 15,=V(CALC)	ECIS-034
	BALR 14,15 CALL CALC(W,W,W,INT)	ECIS-035
	L 13,4(13)	ECIS-036
	RETURN (14,12),T	ECIS-037
ADGET	DS D	ECIS-038
MIN	DC A(0,MAX)	ECIS-039
ARG	DC A(0,0,0,LONG)	ECIS-040
LONG	DS A	ECIS-041
SAVE	DS 18A	ECIS-042
MAX	EQU X'400000' 4000K	ECIS-043
	END	ECIS-044

Use the indication at the end of a **JOB** to know the best **REGION**. If data are read with a large **BLOCKSIZE**, from a tape for example, this argument has to be changed into :

	CALL ECIS(10)	MAIN-019
--	---------------	----------

instead of :

	CALL ECIS(3)	MAIN-019
--	--------------	----------

The subroutine **STIM** gives the remaining time for the **JOB** in units of 1/38500 seconds. It is :

STIM	TITLE '15/06/79-STIM-TEMPS D UNITE CENTRALE RESTANT-ECIS79'	STIM-000
** CALL	STIM(N) N TEMPS RESTANT EN UNITES DE 26 MICROSECONDES	STIM-001
**	N/384 TEMPS RESTANT EN CENTIEMES DE SECONDES	STIM-002
**	VERSION COMPATIBLE MVT/VS2/TSO	STIM-003
STIM	CSECT	STIM-004
	USING *,15	STIM-005
	STM 0,12,20(13)	STIM-006
	L RCVT,16 ->CVT	STIM-007
	L RSHPC,88(,RCVT) ->SHPC(MVT)	STIM-008
	L RW,0(,RCVT) ->TCBWORDS	STIM-009
	TM X'74'(RCVT),X'12' VS2?	STIM-010
	BO VS2	STIM-011
	L RTCB,4(,RW) ->CURRENT TCB	STIM-012
	L RTCB,X'7C'(,RTCB) ->JOB STEP TCB	STIM-013
	L RTCB,X'84'(,RTCB) ->TCB DE L'INIT	STIM-014
	L RTME,120(,RTCB) TCBTME	STIM-015
	SR RN,RN	STIM-016
LTIM	L RTOX,12(,RTME) TOX	STIM-017

	LR	RN,RTOX		STIM-018
	S	RN,0(,RSHPC)	TOX-SHPC	STIM-019
	SLL	RN,1	UNITES 13,... MICROSECONDES	STIM-020
	A	RN,80	TOX-SHPC+TIMER,EN 13 MICROSECONDES	STIM-021
	C	RTOX,12(,RTME)	SHPC MODIFIEE?	STIM-022
	BNE	LTIM		STIM-023
	SRL	RN,1	26,... MICROSECONDES	STIM-024
	B	STORE		STIM-025
	SPACE	2		STIM-026
VS2	DS	OH		STIM-027
	L	RA,12(,RW)	->ASCB	STIM-028
	L	RW,X'50'(,RA)	JOB STEP TIME LIMIT,SECONDES	STIM-029
	MH	RW,=H'100'	CENTIEMES DE S	STIM-030
	LM	RP,RI,X'40'(RA)	ELAPSED STEP TIME EN STCK	STIM-031
	SRDL	RP,12	BIN(63) MICROSECONDES	STIM-032
	D	RP,=F'10000'	=>RP=MICROSECONDES,RI=1/100 DE S	STIM-033
	SR	RI,RW	-TEMPS RESTANT A L'ETAPE EN 1/100 DE S	STIM-034
	M	RP,=F'384'	TIMER UNITS	STIM-035
	LPR	RN,RI		STIM-036
STORE	L	2,0(1)		STIM-037
	ST	RN,0(,2)		STIM-038
RETOUR	LM	0,12,20(13)		STIM-039
	BR	14		STIM-040
	LTORG			STIM-041
RN	EQU	0		STIM-042
RTOX	EQU	2		STIM-043
RCVT	EQU	3		STIM-044
RTCB	EQU	3		STIM-045
RTME	EQU	3		STIM-046
RSHPC	EQU	4		STIM-047
RW	EQU	5		STIM-048
RA	EQU	6	->ASCB	STIM-049
RP	EQU	8	REGISTRE PAIR	STIM-050
RI	EQU	RP+1		STIM-051
	END			STIM-052

It can have to be changed by a local subroutine.

The OVERLAY cards are not the same for the different codes. Those of ECIS79 were :

OVERLAY A	OVLY-001
INSERT CALX	OVLY-002
OVERLAY B	OVLY-003
INSERT LECT,LECD,DEPH,OBSE	OVLY-004
OVERLAY B	OVLY-005
INSERT INPA	OVLY-006
OVERLAY B	OVLY-007
INSERT INPB	OVLY-008
OVERLAY A	OVLY-009
INSERT COLF,FCOU,FCZO,PSI,YFRI,YFCL,YFAS,YFIR,COCL,SIGM,CORI,CORH	OVLY-010
OVERLAY A	OVLY-011
INSERT VARI,FIT1,FIT2,FITE,REST,EVAL	OVLY-012
OVERLAY A	OVLY-013
INSERT DJ6J	OVLY-014
OVERLAY C	OVLY-015
INSERT DJCG,REDM,VIBM,ROAM,ROTM	OVLY-016

OVERLAY C	OVLY-017
INSERT CAL1	OVLY-018
OVERLAY D	OVLY-019
INSERT POTE,DERI,COPO	OVLY-020
OVERLAY E	OVLY-021
INSERT VIBP,ROTD,ROTP	OVLY-022
OVERLAY E	OVLY-023
INSERT FOLD,HULT,FINT	OVLY-024
OVERLAY E	OVLY-025
INSERT EXTP,INTP,STDP	OVLY-026
OVERLAY D	OVLY-027
INSERT QUAN,DJ9J,DCGS,SCAM,CORA,CORB	OVLY-028
OVERLAY F	OVLY-029
INSERT INTI,INSI,INSH,COUP,SECM,PADE	OVLY-030
OVERLAY F	OVLY-031
INSERT INCH,CPCC,LINS	OVLY-032
OVERLAY D	OVLY-033
INSERT SCHE,RESU,SCAT,GRAL,EMRO,RESC,CPSF	OVLY-034

Those of ECIS88 are :

OVERLAY A	OV88-000
INSERT CALX	OV88-001
OVERLAY B	OV88-002
INSERT LECL,LECT,LECD,DEPH,OBSE	OV88-003
OVERLAY B	OV88-004
INSERT INPA,INPB,INPC	OV88-005
OVERLAY A	OV88-006
INSERT COLF,FCOU,FCZO,PSI,YFRI,YFCL,YFAS,YFIR,COCL,SIGM,CORI,CORO,CORZ	OV88-007
OVERLAY A	OV88-008
INSERT VARI,FIT1,FIT2,FITE,REST,EVAL,EXTP	OV88-009
OVERLAY A	OV88-010
INSERT DJ6J	OV88-011
OVERLAY C	OV88-012
INSERT DJCG,REDM,VIBM,ROAM,ROTM	OV88-013
OVERLAY C	OV88-014
INSERT CAL1	OV88-015
OVERLAY D	OV88-016
INSERT POTE,DERI,COPO,ROTZ,ROTD,ROTP,FOLD,INTP,STDP,STBF	OV88-017
OVERLAY D	OV88-018
INSERT QUAN,DJ9J,DCGS,SCHE,CORA,MTCH,CONU,PADE,DIAG	OV88-019
OVERLAY E	OV88-020
INSERT INTI,INSI,INSH	OV88-021
OVERLAY E	OV88-022
INSERT INTR,INRI,INRH	OV88-023
OVERLAY E	OV88-024
INSERT INCH,LINS	OV88-025
OVERLAY D	OV88-026
INSERT SCIN,RESU,SCAT,GRAL,EMRO,RESC	OV88-027

and none have been written for ECIS94. The difference is SCAM instead of SCHE in OV88-17 and SCHE,LCSP instead of SCIN in OV88-27.

A.1.b Use with MTS

This system do not allow **OVERLAY**, control of working space and of time. Follow indications in **COMMENT** cards of the **MAIN** routine.

A.2 Use on UNIVAC

At the first attempt to run the code **ECIS79** on an **UNIVAC** computer, the following indications are inserted at the top of the **MAIN** subroutine

C 30/01/80	MAIN FOR UNIVAC	ECIS79	ECIS-000
C THIS PROGRAM IS LIMITED TO 126K PLUS 5 TIMES THE PARAMETER OF			ECIS-001
C @XQT,X WHERE X IS TRANSFORMED TO 1 FOR A TO 26 FOR Z BY IOPTF.			ECIS-002
C THE SUBROUTINE PMARZ AVOIDS THE AUTOMATIC CHANGES OF PAGE.			ECIS-003
C***** TO BE ABLE TO USE A BLANK FOR .FALSE. THE USER OF UNIVAC MUS			ECIS-004
C INTRODUCE:			ECIS-005
C DO 29 I=1,100	AND		ECIS-006
C 29 LO(I)=.FALSE.	BEFORE CALX-123		ECIS-007
C LX=.FALSE.	BEFORE LECD-033		ECIS-008
C LT1=.FALSE.	AND		ECIS-009
C LT2=.FALSE.	BEFORE OBSE-097		ECIS-010
C LO(37)=.FALSE.	BEFORE EVAL-033		ECIS-011
C LO(35)=.FALSE.	BEFORE REST-038		ECIS-012
C*****			ECIS-013
REAL*8 W(2048)			ECIS-014
COMMON /LARGE/ W			ECIS-015
CALL PMARZ(66,0,66)			ECIS-016
CALL MEMINI(126+IOPTF(0)*5)			ECIS-017
CALL CALC(W,W,W,2048)			ECIS-018
STOP			ECIS-019
END			ECIS-020

Without these modifications, a **LOGICAL** read previously as **.TRUE.** cannot be read with a blank but must be read with a **F**. After **ECIS79**, the logicals are set **.FALSE.** before being read.

The use of files is done with :

N\$TAB 21,1,1,1,1,0 6 7 5 0	.	NTAB-000
END	.	NTAB-001

The subroutine **IOPTF**, written in Copenhagen, allows to run jobs with a larger core in special cases, for instance overnight.

	AXR\$.	IOPT-000
\$(1), IOPTF* ER	OPT\$.	IOPT-001
	JZ A0,2,X11	.	IOPT-002
	LSC A1,A0	.	IOPT-003
	ANA,U A2,8	.	IOPT-004
	LA A0,A2	.	IOPT-005
	J 2,X11	.	IOPT-006
	END	.	IOPT-007

There is a dynamic allocation of core by the **FORTRAN** subroutine **MEMO** which is more similar to the one used on **CDC** than to the one used on **IBM**. This allocation is obtained by :

```

| C 14/05/80 UNIVAC                                     ECIS79  MEMO-000
|       SUBROUTINE MEMO(IDMT,NPLACE,NQ,IX,W,LO)          MEMO-001
| C THIS SUBROUTINE GETS THE WORKING SPACE.             MEMO-002
| C IDMT      PREVIOUS SPACE                             MEMO-003
| C NPLACE    REQUESTED SPACE                           MEMO-004
| C NQ        SPACE TO BE UPDATED                       MEMO-005
| C IX: CONTROL NUMBER IX=1      FIRST CALL   STORE MINIMUM SIZE IN ID MEMO-006
| C              IX=2      DIMINUTION OF SPACE MEMO-007
| C              IX=3      REQUEST NPLACE+ISTART, AT LEAST NPLACE MEMO-008
| C              IX=4      REQUEST NPLACE MEMO-009
| C              IX=5      ABSOLUTE REQUEST MEMO-010
| C W:  WORKING SPACE MEMO-011
| C LO: LOGICAL CONTROLS  IF LO(94)=.TRUE. PRINT THE SIZE OF W MEMO-012
| C OUTPUT VARIABLES: IDMT  SIZE OF WORKING SPACE MEMO-013
| C              LO(216)=.TRUE. IF COMPUTATION CANNOT CONTINUE MEMO-014
| C*****MEMO-015
|       DIMENSION W(1) MEMO-016
|       LOGICAL LO(250) MEMO-017
|       DATA ISTART,IMOD /512,64/ MEMO-018
|       IF (IX-2) 1 , 2 , 3 MEMO-019
| 1 ID=IDMT MEMO-020
|   NT=MAX0(-IMOD,ISTART-IDMT) MEMO-021
|   NPLACE=0 MEMO-022
|   GO TO 4 MEMO-023
| 2 NT=MAX0(NPLACE,ID)-IDMT MEMO-024
|   GO TO 5 MEMO-025
| 3 NT=NPLACE-IDMT MEMO-026
|   IF (IX.NE.3) GO TO 4 MEMO-027
|   NT=NT+ISTART MEMO-028
| 4 NT=IMOD*(NT/IMOD+1) MEMO-029
| 5 CALL MEMORY(NT,NX,IR) MEMO-030
|   IDMT=IDMT+NT+IR MEMO-031
|   LO(216)=(IDMT.LT.NPLACE).AND.((IX-5)*(IX-3).EQ.0) MEMO-032
|   NQ=NQ+NT+IR MEMO-033
|   IF (LO(94)) WRITE (6,1000) IX,NT,IR,NPLACE,IDMT,NX MEMO-034
|   IF (.NOT.LO(216)) RETURN MEMO-035
|   WRITE (6,1001) IX,IDMT,NPLACE MEMO-036
|   WRITE (6,1002) MEMO-037
|   RETURN MEMO-038
| 1000 FORMAT (/8H REQUEST,I2,4H FOR,I9,9H MEMORIES,I6,25H ARE MISSING. MEMO-039
| 1 NPLACE =,I9,7H SIZE =,I9,15H TOTAL LENGTH ,I10/) MEMO-040
| 1001 FORMAT (33H NOT ENOUGH PLACE ..... REQUEST,I2,I10,8H ALLOWED,I10MEMO-041
| 1,10H REQUESTED/) MEMO-042
| 1002 FORMAT (/34H ...STOP... NEXT TIME, USE @XQT,Z/) MEMO-043
|   END MEMO-044

```

with the machine language subroutine :

```

|       AXR$ . SUBROUTINE WRITTEN IN STRASBOURGMEMI-000
| $ (1) . ON THE 23/01/80 MEMI-001
| MEMORY* . MEMI-002
|       L A1,*0,X11 . CALL MEMORY(I,J,K) MEMI-003
|       SZ *2,X11 . I NUMBER OF MEMORIES REQUESTED MEMI-004
|       L A0,LASTD . J RETURNS THE SIZE OF THE MEMI-005
|       JZ A1,RETOUR . PROGRAM MEMI-006
|       A A0,A1 . K RETURNS DIFFERENCE BETWEEN MEMI-007

```

	JP	A1,MORE	.	NUMBER OF MEMORIES GIVEN	MEMI-008
	ER	LCORE\$.	AND NUMBER OF MEMORIES	MEMI-009
	J	LMCORE	.	REQUESTED	MEMI-010
	MORE	TLE	A0,MAXCOR	.	MEMI-011
	J	ERMCORE	.		MEMI-012
	ANU	A0,MAXCOR	.		MEMI-013
	SN	A1,*2,X11	.		MEMI-014
	L	A0,MAXCOR	.		MEMI-015
	ERMCORE	ER	MCORE\$.	MEMI-016
	LMCORE	S	A0,LASTD	.	MEMI-017
	RETOUR	S	A0,*1,X11	.	MEMI-018
	J	4,X11	.		MEMI-019
	MEMINI*	L	A0,*0,X11	.	ENTRY USED TO STORE THE MAXIMUM
		LSSL	A0,10	.	SIZE FOR THE RUN
		AN,U	A0,1	.	MEMI-022
		S	A0,MAXCOR	.	CALL MEMINI(IMAX)
		J	2,X11	.	IMAX MAXIMUM SIZE IN K
	\$ (0)			.	MEMI-025
	MAXCOR	+	0	.	MEMI-026
	LASTD	+	LASTD\$.	MEMI-027
	END			.	MEMI-028

The subroutine STIM has been also written in Copenhagen :

	AXR\$.	STIM-000
	\$ (1),STIM*	LXI,U	A1,PCTBD\$.
		LA,U	A2,RPCTA\$.
		LDJ	A1,\$+1	.
		LA	A0,03,A2	.
		ANA	A0,013,A2	.
		LDJ	A1,\$+1	.
		MI,U	A0,77	.
		DI,U	A0,10	.
		SA	A0,*0,X11	.
		J	2,X11	.
	END			.

The programme ECIS79 had to be used with following OVERLAY :

```

|SEG S1
|IN ECISR.ECIS,.CALC,.HORA,.STIM,.NTAB$,.MEMO,.MEMORY
|SEG A1*
|IN ECISR.CALX
|SEG B1*
|IN ECISR.LECT,.LECD,.IOPTF
|SEG B2*,B1
|IN ECISR.DEPH,.OBSE
|SEG A2*,A1
|IN ECISR.COLF,.FCOU,.FCZO,.PSI,.YFRI,.YFCL,.YFAS,.YFIR,.COCL,.SIGM
|IN ECISR.CORI,.CORH
|SEG A3*,A1
|IN ECISR.VARI,.FIT1,.FIT2,.FITE,.REST,.EVAL
|SEG A4*,A1
|IN ECISR.DJ6J
|SEG C1*

```

```

|IN ECISR.DJCG,.REDM,.VIBM,.ROAM,.ROTM
|SEG C2*,C1
|IN ECISR.CAL1
|SEG D1*
|IN ECISR.POTE
|SEG E1*
|IN ECISR.ROTP,.ROTX
|SEG E2*,E1
|IN ECISR.VIBP
|SEG E3*,E1
|IN ECISR.DERI,.COPO
|SEG G1*
|IN ECISR.FOLD,.HULT,.FINT
|SEG G2*,G1
|IN ECISR.EXTP,.INTP,.STDP
|SEG D2*,D1
|IN ECISR.QUAN,.DJ9J,.DCGS,.SCAM,.CORA,.CORB
|SEG F1*
|IN ECISR.INTI,.INSI,.INSH,.COUP,.SECM,.PADE
|SEG F2*,F1
|IN ECISR.INCH,.CPCC,.LINS
|SEG D3*,D1
|IN ECISR.SCHE,.RESU,.SCAT,.GRAL,.EMRO,.CPSF
|SEG L1*,()
|IN LARGE
|SEG B3*,B1
|IN ECISR.INPA
|SEG B4*,B1
|IN ECISR.INPB

```

A.3 Use on VAX

Around 50 FORTRAN cards have to be changed in ECIS79, to deal with the larger floating value which is 10^{38} . The IBM versions of ECIS88 and ECIS94 need no change. There is no use of OVERLAY on this computer and the working array can be taken the largest possible. At the Washington State University, it has been possible to use as MAIN routine:

```

|C 26/11/81  VAX                                MAIN-000
|C THE ASSEMBLER SUBROUTINE ECIS(I) DEFINES THE DOUBLE PRECISION ARRAY W MAIN-001
|C AND CALLS CALC(W,W,W,IDMX) WHERE IDMX IS THE LENGTH OF W MAIN-002
|C W IS ALL THE REGION LEFT FREE BY THE PROGRAM EXCEPT 2*I K FOR BUFFERS MAIN-003
|C THE SUBROUTINE ECIS CAN BE ELIMINATED WITH THE FOLLOWING MAIN: MAIN-004
|    REAL*8 W(400000)                            MAIN-005
|    CALL CALC(W,W,W,400000)                      MAIN-006
|    STOP                                          MAIN-007
|    END                                          MAIN-008

```

There is no control of the time allowed for the JOB. The CPU elapsed time is given by following subroutine STIM :

```

|C 12/10/81  VAX                                STIM-000
|    SUBROUTINE STIM(I)                          STIM-001
|C RETURNS INTEGER VALUE OF CURRENT CPU TIME IN HUNDREDTH OF SECONDS STIM-002
|C          ACCUMULATED BY PROCESS                STIM-003
|C          RJS          12-OCT-81    (WASHINGTON STATE UNIVERSITY) STIM-004

```

```

|      IMPLICIT INTEGER*4 (A-Z)                      STIM-005
|      PARAMETER JPI$_CPUTIM='0407'X                STIM-006
|      INTEGER*2 LIST(8)                             STIM-007
|      DATA LIST/4,JPI$_CPUTIM,6*0/                STIM-008
|      EQUIVALENCE (TIMELOC,LIST(3))                 STIM-009
|      TIMELOC=%LOC(TIME)                             STIM-010
|      STAT=SYS$GETJPI(,,,LIST,,, )                 STIM-011
|      I=TIME                                          STIM-012
|      RETURN                                          STIM-013
|      END                                            STIM-014

```

The subroutine **HORA** has been modified to suppress the transformation from **IBM** units.

A.4 Use on PR1ME

The code **ECIS79** could not be used with **F77**, due to compilation errors found in the function **SIGM** and the strange behaviour of equivalences. With **FTN**, in which there is no **DASIN** and no **DSINH**, the two following subroutines must be added to the coulomb functions:

```

|      FUNCTION DASIN(X)
|      IMPLICIT REAL*8 (A-H,O-Z)
|      DASIN=DATAN2(X,DSQRT(1-X*X))
|      RETURN
|      END
|      FUNCTION DSINH(X)
|      IMPLICIT REAL*8 (A-H,O-Z)
|      DSINH=0.5D0*(DEXP(X)-DEXP(-X))
|      RETURN
|      END}

```

Further modifications had to be done : **DFLOAT** has to be replaced by **DBLE(FLOAT())** and a statement had to be modified in subroutine **FITE**. The **OVERLAY** is not needed. The **MAIN** routine is :

```

|C 27/01/82  PRIME                                     MAIN-000
|C THE ASSEMBLER SUBROUTINE ECIS(I) DEFINES THE DOUBLE PRECISION ARRAY W MAIN-001
|C AND CALLS CALC(W,W,W,IDMX) WHERE IDMX IS THE LENGTH OF W MAIN-002
|C W IS ALL THE REGION LEFT FREE BY THE PROGRAM EXCEPT 2*I K FOR BUFFERS MAIN-003
|C THE SUBROUTINE ECIS CAN BE ELIMINATED WITH THE FOLLOWING MAIN: MAIN-004
|      REAL*8 W(160000)                                MAIN-005
|      COMMON W                                         MAIN-006
|      CALL CALC(W,W,W,160000)                          MAIN-007
|      STOP                                           MAIN-008
|      END                                           MAIN-009

```

The **STIM** subroutine is :

```

|C 17/01/82  PRIME                                     STIM-000
|      SUBROUTINE STIM(I)                               STIM-001
|C RETURNS INTEGER VALUE OF CURRENT CPU TIME IN HUNDREDTH OF SECONDS STIM-002
|      CALL CTIM$(I)                                    STIM-003
|      RETURN                                          STIM-004
|      END                                            STIM-005

```

The compilation is :

```

| FTN  XX  -SA -L -X -INTL -BIG -64V

```

A.5 Use on BURROUGH

The code **ECIS79** has been run on the **BURROUGH 7700** of the Technische Hogeschool in Eindhoven, Netherlands.

Some errors without effect on other computers have been found:

- in subroutine **HORA**, **M0** and **M2** must be put on data,
- in subroutine **CORH** the array **SP** was not declared as single precision: add **REAL*4 SP** after **IMPLICIT** or change **SP** into **OP** and use **IMPLICIT REAL*8 (A-H,P-Z)**,
- in the subroutine **CAL1**, in calls to **CORB**, an argument was of wrong type; use :

```

1,DW(NXC1),DW(NX1),NC,ISM,KAB,NW(NNIV),NW(NT),NW(NIVQ),DW(NFG),LMAXCAL1-075
21,LMAX3,INC,NW(NWR),LO)                                CAL1-076
1DW(NXC1),DW(NX),NCXN,ISM,LMAX2,NW(NNIV),NW(NT),NW(NIVQ),DW(NFG),LMCAL1-137
2AX1.LMAX3.INC.NW(NML).LO)                                CAL1-138

```

- in subroutine **SCAT**, the subroutine **EMRO** was not called with an array as argument; use :

	1),X2,X3,SQX,MX,B(1),1)	SCAT-073
	1),X2,X3,SQX,MX,B(2),1)	SCAT-075
	W1=B(1)*B(2)	SCAT-076

- an error was found in subroutine **HULT**.

These changes have been done to the programme on the 10/05/82. Furthermore, the parity of the levels cannot be read on this computer unless one uses :

```
IF (SPI.IS.SIGM) IPI(IV,1)=1
```

A.6 Use on SUN

This version of **ECIS88** was run on a **SUN** computer without difficulty. As the system was **UNIX**, the subroutine **STIM** described for a **CONVEX** computer was used.

B The CDC version

The object deck of **ECIS79** was conveniently handled as a **SOURCE** file for **UPDATE**. Using **UPDATE,D.**, the identification of the cards appears in the compilation. The **UPDATE** identification is the name of the subroutine and the sequence number is the number of the card plus 2.

Due to the relative small size of **CDC** computers, the code must be used with segmentation (**SEGLOAD**). These segmentation cards for **ECIS79** are :

ECIS	GLOBAL	AZ,DBLP,RECOU,TITR,FCL.C.	OVCD-001
RESU	GLOBAL	RESC	OVCD-002
ECIS	INCLUDE	CALC	OVCD-003
LECT	INCLUDE	LECD,DEPH,OBSE	OVCD-004
FITE	INCLUDE	VARI,REST,EVAL	OVCD-005
QUAN	INCLUDE	CORB,SCAM	OVCD-006
ROTP	INCLUDE	VIBP,EXTP,COPO	OVCD-007
RESU	INCLUDE	SCHE,CPSF	OVCD-008
A	TREE	ECIS-(CALX-(INPA,INPB,LECT),COLF,DJ6J-(REDM,B),FITE)	OVCD-009
B	TREE	CAL1-(POTE-(ROTP,FOLD),QUAN-(INTI,INCH),RESU)	OVCD-010
	END		OVCD-011

The time control is obtained by the CDC subroutine **SECOND** and a call to the CERN subroutine **TIMING** at the beginning of the **JOB**.

The CDC version uses the dynamic allocation of memories, using the **MACRO** instruction **MEMORY**. For that, the working array must be extended without overlapping the programme. The CERN subroutines **MEMORY** and **INCLCM**, used previously, include the size of the local computer. The maximum core can depend on the user, the hour of computation, or on the **JOB** card. A **COMPASS** subroutine **INCM** calling the **MACRO MEMORY** has been written by LE FUR, in order to get the maximum core allowed by the installation :

	IDENT	INCM	INCM-000
	ENTRY	INCM	INCM-001
	VFD	36/4HINCM,24/INCM	INCM-002
	SVA0	DATA 0	INCM-003
	SVX1	DATA 0	INCM-004
	TAIL	DATA 0	INCM-005
	INCM	DATA 0	INCM-006
	SB7	1	INCM-007
	SX6	A0	INCM-008
	BX7	X1	INCM-009
	SA6	SVA0	INCM-010
	SA7	A6+B7	INCM-011
	SA4	X1	INCM-012
	LX4	30	INCM-013
	BX6	X4	INCM-014
	SA6	TAIL	INCM-015
	MEMORY	SCM,TAIL,RECALL,,NABORT	INCM-016
	SB7	1	INCM-017
	SA1	SVX1	INCM-018
	SA2	TAIL	INCM-019
	AX2	30	INCM-020
	BX6	X2	INCM-021
	SA6	X1	INCM-022
	SA2	A1-B7	INCM-023
	SA0	X2	INCM-024
	EQ	INCM	INCM-025
	END		INCM-026

called by the **FORTRAN** subroutine:

	SUBROUTINE MEMORY(N,NTOT,IER)	INCM-027
	DATA KX /0/	INCM-028
	IF (KX.EQ.0) CALL INCM(KX)	INCM-029
	N2=N	INCM-030
	KZ=KX+N2	INCM-031
	DO 1 I=1,1000	INCM-032
	KX=KX+N2	INCM-033
	KY=KX	INCM-034
	CALL INCM(KX)	INCM-035
	IF (KX.GE.KY.AND.N2.EQ.N) GO TO 2	INCM-036
	N2=(2*N2)/3	INCM-037
	IF (N2.EQ.0) GO TO 2	INCM-038
	1 CONTINUE	INCM-039
	2 NTOT=KX	INCM-040
	IER=KX-KZ	INCM-041
	RETURN	INCM-042
	END	INCM-043

B.1 Use on CDC 7600

The working array can be in the **SCM** or in the **LCM**.

B.1.a Using SCM

The **OVERLAY** structure on a CDC 7600 puts the **BLANK COMMON** at the top of the root. In these circumstances, if the working array is in it, it cannot be extended without overlapping the programme.

The solution chosen (for **ECIS79**) is to add two cards before the **END** of the **OVERLAY** cards to create a second level:

B	TREE CAL1-(POTE-(ROTP,FOLD),QUAN-(INTI,INCH),RESU)	OVCD-010
	LEVEL	OVCD-011
	TREE CALS	OVCD-012
	END	OVCD-013

and the subroutine **CALS** is :

CDC 15/06/79	ECIS79	CALS-000
SUBROUTINE CALS		CALS-001
C WITHOUT THE SUBROUTINE MEMORY TO GET THE SPACE, THIS SUBROUTINE MUST		CALS-002
C BE REPLACED BY		CALS-003
C DIMENSION W(10000)		CALS-004
C IDMX=10000		CALS-005
C CALL CALC(W,W,IDMX)		CALS-006
C*****		CALS-007
DIMENSION W(10)		CALS-008
CALL CALC(W,W,10)		CALS-009
RETURN		CALS-010
END		CALS-011

The **MAIN** programme is:

CALL CALS	ECIS-008
STOP	ECIS-009
END	ECIS-010

The subroutine **CALS** and the second level are overwritten by the working array but the computation never go out of the subroutine **CALC**. The **UPDATE** instruction is :

```
|*COMPILE ECIS.CALS
```

B.1.b Using LCM

The code **ECIS** in its CDC version has been written in such a way that a quantity in the working array never appears in the call of a subroutine except as an array. All the working array can be shifted into the **LCM** by inserting :

```
| LEVEL,A1,A2,A3.....}
```

in each subroutine, where **A1,A2,A3...** are the parts of the working array which are used. The **MAIN** programme is then :

	COMMON W(512)	LEV2-002
	LEVEL 2,W	LEV2-003
	CALL CALC(W,W,512)	LEV2-004
	STOP	ECIS-009
	END	ECIS-010

The UPDATE (still for ECIS79) instruction is :

	*ID LEV2	-ECIS79-
	*D ECIS.10	LEV2-001
	COMMON W(512)	LEV2-002
	LEVEL 2,W	LEV2-003
	CALL CALC(W,W,512)	LEV2-004
	
	*D INCM.018	LEV2-013
	MEMORY LCM,TAIL,RECALL,,NABORT	LEV2-014
	
	*I FIT1.39	LEV2-125
	LEVEL 2,I,W	LEV2-126
	*I FIT2.55	LEV2-127
	LEVEL 2,A,D,IP	LEV2-128
	*C ECIS.FIT2	LEV2-129

B.2 Use on CDC 6600

On a CDC 6600, it is convenient to reduce the size of the **BUFFER** in the first card of the **MAIN** routine.

In some places, **ECIS** can be run as on a CDC 7600 with **SCM**. In some other places, the use of the **MACRO MEMORY** is inhibited unless the compilation is **STATIC**. The working array can be in the **BLANK COMMON** and the use of the subroutine **CALS** and of the second level of **OVERLAY** avoided.

The size of the working array can be controlled by the **CM** parameter of the **JOB** card with the following **MAIN** routine :

	CDC 16/10/78	ECIS-000
	PROGRAM ECIS(INPUT,OUTPUT,PUNCH,TAPE5=INPUT,TAPE6=OUTPUT,TAPE7=PUNECIS-001	
	1CH,TAPE1)	ECIS-002
	C	ECIS-003
	C THE MAIN SUBROUTINE DEFINES THE WORKING SPACE W AND CALL THE	ECIS-004
	C SUBROUTINE CALC WITH THE LENGTH OF W AS LAST ARGUMENT.	ECIS-005
	C	ECIS-006
	C*****ECIS-007	
	COMMON // W(1)	ECIS-008
	CALL CALC(W,W,LENGTHB(W))	ECIS-009
	STOP	ECIS-010
	END	ECIS-011

and the **COMPASS** subroutine written in Amsterdam :

	IDENT LENGTHB	LENGTHB2
	ENTRY LENGTHB	LENGTHB3
	MEMREPLY BSSZ 1	LENGTHB4
	LENGTHB BSS 1	LENGTHB5
	SX7 X1	LENGTHB6
	SYSTEM MEM,R,MEMREPLY	LENGTHB7

	SA1	MEMREPLY	LENGTHB8
	AX1	30	LENGTHB9
	IX6	X1-X7	LENGTH10
	EQ	LENGTHB	LENGTH11
	END		LENGTH12

The indication at the end of a JOB can be used to choose the best CM in a similar computation.

B.3 Use on CRAY

The code ECIS79 has been run on a CRAY computer. Due to the large size, the OVERLAY structure is not needed. No information on time and memory management.

The code ECIS88 has been used with memory management. The MAIN subroutine is then :

	CRAY 21/03/88	ECIS88	ECIS-000
	PROGRAM ECIS(INPUT,OUTPUT,PUNCH,TAPE5=INPUT,TAPE6=OUTPUT,TAPE7=PUNECIS-001		
	1CH,TAPE8)		ECIS-002
	COMMON W(10)		ECIS-003
	CALL CALC(W,W,10)		ECIS-004
	STOP		ECIS-005
	END		ECIS-006

and the subroutine MEMO is :

	CRAY 07/06/86	ECIS88	MEMO-000
	SUBROUTINE MEMO(IDMT,NPLACE,NQ,IX,LO)		MEMO-001
	C THIS SUBROUTINE GETS THE WORKING SPACE.		MEMO-002
	C IDMT PREVIOUS SPACE		MEMO-003
	C NPLACE REQUESTED SPACE		MEMO-004
	C NQ SPACE TO BE UPDATED		MEMO-005
	C IX: CONTROL NUMBER IX=1 FIRST CALL STORE MINIMUM SIZE IN ID		MEMO-006
	C IX=2 DIMINUTION OF SPACE		MEMO-007
	C IX=3 REQUEST NPLACE+ISTART, AT LEAST NPLACE		MEMO-008
	C IX=4 REQUEST NPLACE		MEMO-009
	C IX=5 ABSOLUTE REQUEST		MEMO-010
	C LO: LOGICAL CONTROLS IF LO(94)=.TRUE. PRINT THE SIZE OF W		MEMO-011
	C OUTPUT VARIABLES: IDMT SIZE OF WORKING SPACE		MEMO-012
	C LO(216)=.TRUE. IF COMPUTATION CANNOT CONTINUE		MEMO-013
	C*****MEMO-014		
	LOGICAL LO(250)		MEMO-015
	DATA ISTART,IMOD /4096,1024/		MEMO-016
	IF (IX-2) 1 , 2 , 3		MEMO-017
	1 ID=IDMT		MEMO-018
	NT=MAX0(IDMT,ISTART)		MEMO-019
	NPLACE=0		MEMO-020
	CALL MEMORY('MAXFL',IDMAX)		MEMO-021
	CALL MEMORY('CURFL',NX)		MEMO-022
	IDMAX=IDMAX-NX+IDMT		MEMO-023
	GO TO 4		MEMO-024
	2 NT=MAX0(NPLACE,ID)-IDMT		MEMO-025
	CALL MEMORY('UC',NT)		MEMO-026
	IR=0		MEMO-027
	GO TO 5		MEMO-028
	3 NT=NPLACE-IDMT		MEMO-029

	IF (IX.NE.3) GO TO 4	MEMO-030
	NT=NT+ISTART	MEMO-031
	4 NT=IMOD*(NT/IMOD+1)	MEMO-032
	NV=MINO(NT,IDMAX-IDMT)	MEMO-033
	IR=NV-NT	MEMO-034
	CALL MEMORY('UC',NV)	MEMO-035
	5 CALL MEMORY('CURFL',NX)	MEMO-036
	IDMT=IDMT+NT+IR	MEMO-037
	LO(216)=(IDMT.LT.NPLACE).AND.((IX-5)*(IX-3).EQ.0)	MEMO-038
	NQ=NQ+NT+IR	MEMO-039
	IF (LO(94)) WRITE (6,1000) IX,NT,IR,NPLACE,IDMT,NX,NX	MEMO-040
	IF (.NOT.LO(216)) RETURN	MEMO-041
	WRITE (6,1001) IX,IDMT,NPLACE	MEMO-042
	WRITE (6,1002)	MEMO-043
	RETURN	MEMO-044
	1000 FORMAT (/8H REQUEST,I2,4H FOR,I9,9H MEMORIES,I6,25H ARE MISSING.	MEMO-045
	1 NPLACE =,I9,7H SIZE =,I9,15H TOTAL LENGTH ,I8,4X,08,1HB/)	MEMO-046
	1001 FORMAT (33H NOT ENOUGH PLACE REQUEST,I2,I10,8H ALLOWED,I10	MEMO-047
	1,10H REQUESTED/)	MEMO-048
	1002 FORMAT (/48H ...STOP... NEXT TIME, USE LARGER MFL PARAMETER/)	MEMO-049
	END	MEMO-050

The subroutine STIM is :

	CRAY 09/01/88	ECIS88	STIM-000
	SUBROUTINE STIM(K)		STIM-001
	CALL TREMAIN(T)		STIM-002
	K=38500.*T		STIM-003
	RETURN		STIM-004
	END		STIM-005

without changing the IBM subroutine HORA.

C Use on a CONVEX

On a CONVEX computer, the two versions can be used. As there can be only control of the elapsed time, a subroutine STIM, identical for the two versions, was written to stop the JOB just before an hour (the allowed time of the QUEUE for long JOB). This subroutine is :

	C 15/12/87 CONVEX IBM AND CDC VERSIONS	ECIS88	STIM-000
	SUBROUTINE STIM(I)		STIM-001
	C RETURNS INTEGER VALUE OF CPU ELAPSED TIME IN HUNDREDTHS OF SECONDS		STIM-002
	DIMENSION A(2)		STIM-003
	B=ETIME(A)		STIM-004
	I=38500*(3300.-B)		STIM-005
	RETURN		STIM-006
	END		STIM-007

It can be used on any computer with the UNIX system.

The IBM version has to be compile with "fc -72" and the CDC version with "fc -cfc -72". Even if all the subroutines has been compiled beforehand and "fc" is used only to create a LOAD MODULE, one must use "fc -cfc" to get the right answer from the subroutine STIM.

Chapter III

Coupled channels

The inelastic scattering of nucleons exciting low lying collective states of the nuclei is usually described by coupled channel calculations using a collective model. A Dirac phenomenology using a scalar and a vector potential has been introduced in Ref. [9] to describe elastic scattering at quite large energies. Strictly speaking, these calculations are only valid for infinite mass targets because the center-of-mass is separable in non relativistic Schrödinger equation and not in Dirac equation. In the same approximation, deformation or vibration can be introduced in the scalar and in the vector potential of the Dirac formalism to study the inelastic scattering in the collective model.

For elastic scattering, there is a fully equivalent Schrödinger equation for the Dirac equation. The presence of some $1/r^2$ terms for charged particles do not affect strongly the results. For inelastic scattering, this transformation has to be done on the tri-dimensional Dirac equation. Then, the spin orbit potential appears as the cross product of the gradient of the potential with a gradient acting on the wave function. Such expression for the spin-orbit is used since a long time in the description of nucleon inelastic scattering by : [55] SHERIF, H., BLAIR, J. S., “*Inelastic proton scattering and the deformed spin dependent optical potential*”, Physics Letters **26B** (1968) 489. and : [56] SHERIF, H., “*Spin-dependent effects in proton inelastic scattering*”, Thesis University of Washington (1968). However, to write the Schrödinger equation, the wave function has been multiplied by the square root of the potential. As long as the potential is a given function of the radius, it is the only difference between Dirac and Schrödinger formalism. The potential used has usually Woods-Saxon form-factors in the two formalisms; the Schrödinger potential equivalent to the Dirac one gets a *wine-bottle-bottom* shape as said in the title of Ref. [9].

Coupled channel calculations are necessary even at large energies as 500 or 800 MeV as shown in Ref. [48] and [50]. At these energies, the elastic scattering of protons to backwards angles is well described when the strong low-lying collective states are taken into account. The difference between Schrödinger and Dirac approaches seems limited to the influence of the shape of the potential as shown in : [57] COOPER, E., D., “*Are the Dirac imaginary potentials surface peaked?*”, Nucl. Phys. **A495** (1989) page 483 and in Ref. [50].

A Dirac and Schrödinger formalisms

In order to describe inelastic nucleon scattering, we must start from Dirac equation because spin-orbit effects can not be understood without it.

A.1 The Dirac equation

To describe the scattering of a nucleon with “magnetic anomalous moment” in a Coulomb field $V_C(\vec{r})$, the Dirac equation is :

$$\left(\frac{\hbar}{i} \vec{\alpha} \cdot \vec{\nabla} + \beta m + V_C(\vec{r}) + \frac{i\mu\hbar}{2m} \beta \vec{\alpha} \cdot \{ \vec{\nabla} V_C(\vec{r}) \} \right) \psi(\vec{r}) = E \psi(\vec{r}). \quad (\text{III.1})$$

More generally, this equation can be generalised to describe the scattering of this nucleon by a nucleus into the equation :

$$\left(\frac{\hbar}{i} \vec{\alpha} \cdot \vec{\nabla} + \beta \{m + V_S(\vec{r})\} + V_V(\vec{r}) + \frac{i\hbar}{2m} \beta \vec{\alpha} \cdot \{\vec{\nabla} V_T(\vec{r})\} \right) \psi(\vec{r}) = E \psi(\vec{r}) \quad (\text{III.2})$$

where $V_S(\vec{r})$, $V_V(\vec{r})$, $V_T(\vec{r})$ are three finite range complex potentials of which the real and the imaginary part are approximated by a Woods-Saxon form factor but can be replaced by any other form factor given by a theory.

The Equ. (III.1) is valid in the laboratory system and the Equ. (III.2) can be used only for a target with an infinite mass with respect to the nucleon, so that the laboratory system is also the center of mass system and m is the rest mass of the nucleon.

We consider the mass m as a parameter which can be :

- the rest mass if `LO(98)=.TRUE.`,
- the reduced mass if `LO(98)=.FALSE.`, which allows a correct limit at low energy.

In this equation, the vector potential $V_V(\vec{r})$ and the tensor potential $V_T(\vec{r})$ include the Coulomb potential $V_C(\vec{r})$; in fact, for a nucleon with an “anomalous magnetic moment” μ_n , the tensor potential is reduced here to the Coulomb potential multiplied (at the low energy limit) by $\mu_n + \frac{1}{2}z$ where z is the charge of the nucleon.

The $\vec{\alpha}$ and β are four-dimensional Dirac matrices. They can be expressed in terms of two-dimensional unit matrix and Pauli matrices $\vec{\sigma}$ as follows and allow to write the wave function :

$$\vec{\alpha} = \begin{vmatrix} 0 & \vec{\sigma} \\ \vec{\sigma} & 0 \end{vmatrix}, \quad \beta = \begin{vmatrix} 1 & 0 \\ 0 & -1 \end{vmatrix}, \quad \psi(\vec{r}) = \begin{pmatrix} F(\vec{r}) \\ iG(\vec{r}) \end{pmatrix} \quad (\text{III.3})$$

in terms of the **large component** $F(\vec{r})$ and the **small component** $G(\vec{r})$. They allow to write a set of two linear coupled equations with the Pauli matrices :

$$\begin{aligned} \hbar \vec{\sigma} \cdot \vec{\nabla} G(\vec{r}) &= [E - m - V_V(\vec{r}) - V_S(\vec{r})] F(\vec{r}) + \frac{\hbar}{2m} \vec{\sigma} \cdot [\vec{\nabla} V_T(\vec{r})] G(\vec{r}) \\ -\hbar \vec{\sigma} \cdot \vec{\nabla} F(\vec{r}) &= [E + m - V_V(\vec{r}) + V_S(\vec{r})] G(\vec{r}) + \frac{\hbar}{2m} \vec{\sigma} \cdot [\vec{\nabla} V_T(\vec{r})] F(\vec{r}) \end{aligned} \quad (\text{III.4})$$

A.2 Reduction to Schrödinger equation

The terms $[\vec{\nabla} V_T(\vec{r})]$ are almost eliminated by the substitution :

$$G(\vec{r}) = \exp\left(\frac{V_T(\vec{r})}{2m}\right) \hat{G}(\vec{r}), \quad F(\vec{r}) = \exp\left(-\frac{V_T(\vec{r})}{2m}\right) \hat{F}(\vec{r}), \quad (\text{III.5})$$

respectively, to get :

$$\begin{aligned} \hbar \vec{\sigma} \cdot \vec{\nabla} \hat{G}(\vec{r}) &= \exp\left(-\frac{V_T(\vec{r})}{m}\right) [E - m - V_V(\vec{r}) - V_S(\vec{r})] \hat{F}(\vec{r}) \\ -\hbar \vec{\sigma} \cdot \vec{\nabla} \hat{F}(\vec{r}) &= \exp\left(\frac{V_T(\vec{r})}{m}\right) [E + m - V_V(\vec{r}) + V_S(\vec{r})] \hat{G}(\vec{r}) \end{aligned} \quad (\text{III.6})$$

From the second equation, using :

$$D(\vec{r}) = E + m - V_V(\vec{r}) + V_S(\vec{r}), \quad D_T(\vec{r}) = \exp\left(\frac{V_T(\vec{r})}{m}\right) D(\vec{r}) \quad (\text{III.7})$$

we get :

$$\hat{G}(\vec{r}) = -\frac{1}{D_T(\vec{r})} \hbar \vec{\sigma} \cdot \vec{\nabla} \hat{F}(\vec{r}) \quad (\text{III.8})$$

which gives in the first equation, using :

$$(\vec{\sigma} \cdot \vec{A})(\vec{\sigma} \cdot \vec{B}) = (\vec{A} \cdot \vec{B}) + i\vec{\sigma} \cdot (\vec{A} \times \vec{B}) \quad (\text{III.9})$$

and after multiplication by $D_T(\vec{r})$:

$$-\hbar^2 \left\{ \Delta - \frac{1}{D_T(\vec{r})} [\vec{\nabla} D_T(\vec{r})] \cdot \vec{\nabla} - i\vec{\sigma} \cdot [\vec{\nabla} D_T(\vec{r})] \times \vec{\nabla} \right\} \hat{F}(\vec{r}) = D(\vec{r}) [E - m - V_V(\vec{r}) - V_S(\vec{r})] \hat{F}(\vec{r}) \quad (\text{III.10})$$

This equation include first derivatives of $\hat{F}(\vec{r})$ which can be eliminated by using :

$$f(\vec{r}) = D_T(\vec{r})^{-\frac{1}{2}} \hat{F}(\vec{r}) = D(\vec{r})^{-\frac{1}{2}} F(\vec{r}) \quad (\text{III.11})$$

to obtain :

$$\begin{aligned} \hbar^2 \left\{ \Delta - \frac{3}{4} \left(\frac{\vec{\nabla} D_T(\vec{r})}{D_T(\vec{r})} \right)^2 + \frac{1}{2} \frac{\Delta D_T(\vec{r})}{D_T(\vec{r})} - i\vec{\sigma} \cdot [\vec{\nabla} \ln D_T(\vec{r})] \times \vec{\nabla} \right\} f(\vec{r}) \\ + \left\{ -2EV_V(\vec{r}) - 2mV_S(\vec{r}) + V_V^2(\vec{r}) - V_S^2(\vec{r}) + \hbar^2 k^2 \right\} f(\vec{r}) = 0 \end{aligned} \quad (\text{III.12})$$

with $\hbar^2 k^2 = E^2 - m^2$.

This equation can be written :

$$\left\{ -\frac{\hbar^2}{2m} [\Delta + k^2] + V_1(\vec{r}) - \vec{\sigma} \cdot [\nabla \ln V_2(\vec{r})] \times \frac{\vec{\nabla}}{i} \right\} f(\vec{r}) = 0 \quad (\text{III.13})$$

with :

$$\begin{aligned} V_1(\vec{r}) &= \frac{\hbar^2}{2m} \left\{ \frac{3}{4} \left(\frac{\vec{\nabla} D_T(\vec{r})}{D_T(\vec{r})} \right)^2 - \frac{1}{2} \frac{\Delta D_T(\vec{r})}{D_T(\vec{r})} \right\} + \frac{1}{2m} \left\{ 2EV_V(\vec{r}) + 2mV_S(\vec{r}) - V_V(\vec{r})^2 + V_S(\vec{r})^2 \right\} \\ V_2(\vec{r}) &= D_T(\vec{r}) \end{aligned} \quad (\text{III.14})$$

A.3 The Schrödinger formalism

As long as there is only elastic scattering, the potentials have no angular dependence. Using :

$$\vec{\nabla} = \frac{\vec{r}}{r} \frac{d}{dr} - i \frac{1}{r^2} \vec{r} \times \vec{L} \quad (\text{III.15})$$

the spin term of the Dirac equation for $D_T(\vec{r})$ independent of \hat{r} reduces to :

$$-i\vec{\sigma} \cdot \left[\frac{\vec{r}}{r} \frac{d}{dr} \ln D_T(r) \right] \times \vec{\nabla} = \frac{1}{r} \frac{d}{dr} \ln D_T(r) (\vec{\sigma} \cdot \vec{L}) \quad (\text{III.16})$$

and Equ. (III.13) can be written :

$$\left\{ -\frac{\hbar^2}{2m} [\Delta + k^2] + V_1(r) + \frac{1}{r} \left[\frac{d}{dr} V_2(r) \right] (\vec{\sigma} \cdot \vec{L}) \right\} f(\vec{r}) = 0 \quad (\text{III.17})$$

The Schrödinger formalism is the use of Equ. (III.17).

An important difference is the use of the Coulomb potential $V_C(r)$ directly as a part of $V_1(r)$; it does not contribute to the spin-orbit potential. This allows the use of nonrelativistic coulomb functions. In the equivalent Schrödinger equation, the Coulomb potential is multiplied by the ratio E/m which is larger than 1 and its square appears also in $V_1(r)$ and $V_2(r)$ at large distances. This terms in $1/r^2$ are taken into account with the relativistic Coulomb functions. However, taking into account an “anomalous magnetic moment”, terms in $1/r^2$ appears into the spin-orbit potential in both formalisms. The elastic scattering in the Dirac formalism cannot allow to determine a tensor potential, as shown in : [58] CLARK, B., C., HAMA, S., KALBERMANN, S., G., COOPER, E., D. and MERCER, R., L., “*Equivalent local Dirac potentials*”, Phys. Rev. **C31** (1985) page 694, because it depends only upon the two potentials $V_1(r)$ and $V_2(r)$ defined in Equ. (III.14). Nevertheless, it does not means that a known tensor potential has not to be used.

A.4 Radial dependence of potentials

The potentials $V_S(\vec{r})$, $V_V(\vec{r})$ and $V_V(\vec{r})$ in Dirac equation, $V_1(\vec{r})$ and $V_2(\vec{r})$ in Schrödinger equation are usually complex with a real and an imaginary parts parametrised by a depth V multiplying a Woods Saxon form factor :

$$f(r, a, R) = \frac{1}{1 + \exp\left(\frac{r-R'}{a}\right)} \quad \text{where} \quad R' = RA^{1/3} \quad (\text{III.18})$$

where A is the mass of the target. **These are called volume potentials.**

The imaginary part of V_1 can include also **a surface potential** with the form factor :

$$f'(r, a, R) = \frac{1}{4a} \frac{d}{dR'} \left\{ \frac{1}{1 + \exp\left(\frac{r-R'}{a}\right)} \right\} \quad \text{where} \quad R' = RA^{1/3} \quad (\text{III.19})$$

and we generalise that to its real part. So, the two formalism deals with the same number of potentials.

Instead of the Woods-Saxon potential defined by Equ. (III.18), one can use a “symmetrised” Woods-Saxon potential which is :

$$f_s(r, a, R) = \left[\frac{1}{1 + \exp\left(\frac{r-R'}{a}\right)} \right] \left[\frac{1}{1 + \exp\left(\frac{r+R'}{a}\right)} \right] \quad (\text{III.20})$$

The potentials $V_V(\vec{r})$ and $V_1(\vec{r})$ include the Coulomb potential $V_C(\vec{r})$ which is :

$$V_C(\vec{r}) = \begin{cases} \frac{1}{2} Z e^2 (3 - (r/R'_c)^2) / R'_c, & \text{for } r < R'_c \\ Z e^2 / r, & \text{for } r > R'_c \end{cases} \quad (\text{III.21})$$

where $R'_c = R_c A^{1/3}$ and Z is the product of charges of the particle and the target. $V_C(\vec{r})$ can also be computed from a diffuse charge with a density distribution given by a Woods-Saxon form factor, eventually multiplied by a factor $(1 + cr^2)$ where c is a “third Coulomb parameter”.

The potential can also be obtain from a microscopic description or replaced by a Fourier expansion on Bessel functions.

B The generalised optical model

In the generalised optical model, the potential describes also the different states of the target nucleus. See for example in : [59] TAMURA, T., “*Analyse of the Scattering of Nuclear Particles by Collective Nuclei in Terms of the Coupled-Channel Calculation*”, Rev. Mod. Phys. **37** (1965) 679 and in : [60] TAMURA, T., “*Coupled-channel approach to nuclear reactions*”, Ann. Rev. Nucl. Sci. **19** (1969) 99.

B.1 The macroscopic models

We give here the physical meaning of the macroscopic models used in **ECIS**.

B.1.a The symmetric rotational model

In the symmetric rotational model, the potentials are some functions $V(\vec{r}, \hat{r}')$, where \hat{r}' is the intrinsic axis of the nucleus. These potentials are parametrised by quadrupole and hexadecupole deformations β_2 and β_4, \dots using a radius $R(\theta)$ such that :

$$R(\theta) = R_0 [1 + \beta_2 Y_2^0(\theta) + \beta_4 Y_4^0(\theta) \dots] \quad (\text{III.22})$$

where θ is the angle between \vec{r} and \hat{r}' . There are only even deformations. This radius is used instead of R in the usual expressions of the optical model. The potential can be expanded into multipoles :

$$V(\vec{r}, \hat{r}') = 4\pi \sum_{\lambda, \mu} V_\lambda(r) Y_\lambda^\mu(\vec{r}) Y_\lambda^{\mu*}(\hat{r}'), \quad V_\lambda(r) = \frac{1}{2} \int_0^\pi V(\vec{r}, \hat{r}') P_\lambda(\cos \theta) \sin \theta d\theta, \quad (\text{III.23})$$

where there are only even values of λ . This is obtained by a symmetric ten-points Legendre integral over θ of the potential multiplied by an adequate Legendre polynomial $P_\lambda(\cos \theta)$.

In this model, the target states are :

$$|\Phi_M^I\rangle = \sqrt{\frac{2I+1}{16\pi^2}} \left[\xi_K(r') R_{M,K}^{(J)}(\Omega) + (-)^{I-j} \xi_{-K}(r') R_{M,-K}^{(J)}(\Omega) \right] \quad (\text{III.24})$$

where $\xi_K(r')$ is the intrinsic function, Ω the rotation between the laboratory system and the intrinsic frame and j appears in the rotation of π around an axis perpendicular to the axis of symmetry for each component of the intrinsic state.

B.1.b The vibrational model

In the vibrational model, the radius is replaced by :

$$R = R_0 \left[1 + \sum_{\lambda, \mu} \alpha_\lambda^\mu Y_\lambda^\mu(\theta, \phi) \right] \quad (\text{III.25})$$

where :

$$\alpha_\lambda^\mu = \frac{\beta_\lambda^\mu}{\sqrt{2\lambda+1}} \left(b_{\lambda, \mu}^* + (-)^{\mu} b_{\lambda, -\mu} \right) \quad (\text{III.26})$$

in which $b_{\lambda, \mu}^*$ are phonon creation operators and $b_{\lambda, \mu}$ are phonon annihilation operators. The potential is usually expanded in powers of the α 's :

$$V(r, R_0) + \frac{d}{dR_0} V(r, R_0) \sum_{\lambda, \mu} \alpha_\lambda^\mu Y_\lambda^\mu(\theta, \phi) + \frac{1}{2} \frac{d^2}{dR_0^2} V(r, R_0) \left(\sum_{\lambda, \mu} \alpha_\lambda^\mu Y_\lambda^\mu(\theta, \phi) \right)^2 \quad (\text{III.27})$$

with only the first derivative for the **first order vibrational model**, the first and the second derivatives for the **second order vibrational model**, but also the third derivative with nuclear matrix elements to be given in the **anharmonic vibrational model**.

In the harmonic vibrational model, the target states are :

$$|\Phi_M^I\rangle = b_{IM}^* |0\rangle, \quad |\Phi_M^I\rangle = |(L_1 L_2) IM\rangle = \frac{1}{\sqrt{1 + \delta_{L_1, L_2}}} \left[b_{L_1}^* b_{L_2}^* \right]_M^I |0\rangle, \quad (\text{III.28})$$

for the 1-phonon and the 2-phonons states respectively.

B.1.c The vibration-rotational model

In the vibration-rotational model there is a static deformation like in the rotational model and a dynamical vibration α_L^M of the intrinsic state. Like in the vibrational model R_0 of Equ. (III.22) has to be replaced by R of the Equ. (III.25) but the expansion given by Equ. (III.27) is limited to the first derivative. Here, the value of M is important. The form factors are :

$$V_l^M(r) \propto \beta_L^M \int \frac{d}{dR} V(\vec{r}, \hat{r}') [Y_L^M(\theta) - x Y_l^M(\theta)] Y_l^{M*}(\theta) d\theta \quad (\text{III.29})$$

where the term $x Y_l^M(\theta)$ is there only for the form factors with $l = 0$ and $l = 1$ to correct them for center of mass motion or translation and is defined by the condition that $\int V_l^M(r) r^{l+2} dr = 0$.

In the literature, the derivation of this model is presented with the summation of the two spherical harmonics, giving a sum with Clebsch-Gordan coefficients.

The nuclear states are those given by Equ. (III.24) with or without α_L^M acting on $\xi_K(r')$.

B.1.d The asymmetric rotational model

In the asymmetric rotational model, The radius is given by :

$$R(\theta) = R_0 \left(1 + \beta_2 [\cos \gamma Y_2^0(\theta) + \frac{1}{\sqrt{2}} \sin \gamma \{Y_2^2(\theta) + Y_2^{-2}(\theta)\}] + \beta_4 \cos \gamma' Y_4^0(\theta) + \dots \right) \quad (\text{III.30})$$

and the potential are :

$$V(\vec{r}, \hat{r}') = V_0(\vec{r}, \hat{r}') + V_2^0(\vec{r}, \hat{r}') Y_2^0(\theta) + V_2^2(\vec{r}, \hat{r}') \frac{1}{\sqrt{2}} [Y_2^2(\theta) + Y_2^{-2}(\theta)] + \dots \quad (\text{III.31})$$

They are obtained by integration on the sphere with 36 points, the weights of which were obtained once for all by the inversion of the matrix of spherical harmonics at these points. This means that the potential is supposed to be expanded only with 36 terms, the coefficients of which are obtained by solving a set of 36 equations. The number of multipoles is limited to 15 (i. e. $L = 8$, whereas 36 is $L = 14$).

The nuclear states, as described in : [61] DAVYDOV, A., S. and FILIPPOV, G. F., “*Rotational states in even atomic nuclei*”, Nucl. Phys. **8** (1958) page 237, can be written :

$$\begin{aligned} |\Phi_M^{I_n} > &= \sqrt{\frac{2I+1}{8\pi^2}} \xi(r') \left[\cos \beta_n^1 R_{M,0}^{(J)}(\Omega) + \sin \beta_n^1 \cos \beta_n^2 \frac{R_{M,2}^{(J)} + R_{M,-2}^{(J)}}{\sqrt{2}} \right. \\ &\quad \left. + \sin \beta_n^1 \sin \beta_n^2 \cos \beta_n^3 \frac{R_{M,4}^{(J)} + R_{M,-4}^{(J)}}{\sqrt{2}} + \dots \right] \end{aligned} \quad (\text{III.32})$$

for $I = 2m$ even with the possibility of $1 + m$ different states and :

$$|\Phi_M^{I_n} > = \sqrt{\frac{2I+1}{8\pi^2}} \xi(r') \left[\cos \beta_n^2 \frac{R_{M,2}^{(J)} - R_{M,-2}^{(J)}}{\sqrt{2}} + \sin \beta_n^2 \cos \beta_n^3 \frac{R_{M,4}^{(J)} - R_{M,-4}^{(J)}}{\sqrt{2}} + \dots \right] \quad (\text{III.33})$$

for $I = 2m + 1$ odd with the possibility of only m different states.

If the asymmetric rotational model is “constrained”, the first state must be the 0^+ and the second state must be the first 2^+ . If there is a third state, it must be the second 2^+ . There is no protection in the code against using only elastic scattering, but no γ will be taken into account. The value read as β_1^1 for the first 2^+ is taken as the value of the γ in Equ. (III.30). The result given in Ref. [61] can be written :

$$\tan \beta_1^1 = - \frac{3 \sin \gamma \cos 3\gamma - \cos \gamma \sin 3\gamma}{\sin \gamma \sin 3\gamma + 3 \cos \gamma \cos 3\gamma + \sqrt{9 - 8 \sin^2 3\gamma}}, \quad \beta_2^1 = \beta_1^1 + \frac{\pi}{2} \quad (\text{III.34})$$

because a sign does not matter.

B.2 The coupled equations

We shall take as example the rotational model. The equation to solve is :

$$[T + V - E] \Psi = 0 \quad (\text{III.35})$$

where T is the kinetic energy and V the generalised optical potential.

B.2.a Schrödinger formalism

The total wave function is written :

$$\Psi(\vec{r}) = \frac{1}{r} \sum_{l,j,I,J} f_{ljIJ}(r) Y_{lsjIJM} \quad (\text{III.36})$$

where :

$$Y_{lsjIJM} = i^l \sum_{\mu, \sigma, m, m'} \langle j, I, m, m' | J, M \rangle \left(\langle l, s, \mu, \sigma | j, m \rangle Y_l^\mu(\hat{r}) | s \sigma \rangle \right) \Phi_I^{m'} \quad (\text{III.37})$$

are the target-spin-angular functions involving the target state ϕ_I .

Projecting the total equation on all the functions Y_{lsjIJM} , we obtain a system of coupled second order differential equations :

$$\frac{\hbar^2}{2m_i} \left[\frac{d^2}{dr^2} - \frac{2\eta_i k_i}{r} - \frac{l_i(l_i + 1)}{r^2} + k_i^2 \right] f_i(r) = \sum_{i'} \left\{ \sum_{\lambda} G_{ii'}^\lambda V^\lambda(r) \right\} f_{i'}(r) \quad (\text{III.38})$$

where i and i' stands for any set of quantum numbers (l, j, I, J) , k_i and η_i are the wave number and the coulomb parameter for equation i and $G_{ii'}^\lambda$ is a geometrical coefficient for the transition form factor $V^\lambda(r)$. For $\lambda = 0$ and $i = i'$, the second member of Equ. (III.38) includes the optical model. In general, $G_{ii'}^\lambda$ is the product of a “nuclear” part, which is the “reduced matrix element” and a geometric coefficient of the partial waves. For example, in the rotational model, a state of the target, member of a rotational band starting with 0^+ is described by Equ. (III.24) with $K = 0$:

$$|\Phi_{I,M}\rangle = \sqrt{\frac{2I+1}{8\pi^2}} R_{M0}^{(I)*}(\Omega) \chi(r') \quad (\text{III.39})$$

where $\chi(r')$ is the intrinsic wave function. The result of integration on the nuclear state is :

$$\begin{aligned} \langle \Phi_{I_f, M_f} | V(\vec{r}, \hat{r}') | \Psi_{I_i, M_i} \rangle &= \sqrt{4\pi} \sum_{\lambda} V_{\lambda}(r) Y_{\lambda}^{\mu}(\hat{r}) \times (-)^{M_i} \sqrt{(2I_i + 1)(2I_f + 1)(2\lambda + 1)} \\ &\quad \begin{pmatrix} I_i & I_f & \lambda \\ 0 & 0 & 0 \end{pmatrix} \begin{pmatrix} I_i & I_f & \lambda \\ M_f & -M_i & \mu \end{pmatrix} \end{aligned} \quad (\text{III.40})$$

and the total result is :

$$\begin{aligned} &\sum_{\lambda} V_{\lambda}(r) (-)^{J+\lambda+\frac{1}{2}} \sqrt{(2I_i + 1)(2I_f + 1)(2j_i + 1)(2j_f + 1)} \\ &\quad \begin{pmatrix} I_i & I_f & \lambda \\ 0 & 0 & 0 \end{pmatrix} \begin{pmatrix} j_f & \lambda & j_i \\ -\frac{1}{2} & 0 & \frac{1}{2} \end{pmatrix} \begin{Bmatrix} I_i & I_f & \lambda \\ j_f & j_i & J \end{Bmatrix} \end{aligned} \quad (\text{III.41})$$

where there is a part which depends only upon the target and a part which depends only upon the nucleon, the two of them related by a $6j$ coefficient.

If there are n channels and if the spin of the particle is s_i , the spin of the target I_i and the product of the intrinsic parities η_i for the channel i , the total number of coupled equations for a total spin J sufficiently large and a parity η is :

$$N = \sum_{i=1}^n N_i = \sum_{n=1}^n \frac{1}{2} [(2I_i + 1)(2s_i + 1) + \epsilon'_i (-)^{J+\eta}] \quad (\text{III.42})$$

where $\epsilon'_i = 0$ if I_i or s_i is half integer and $\epsilon'_i = \epsilon_i(-)^{I_i+s_i}$ if I_i and s_i are integers.

For the scattering of protons on a $0^+ - 2^+ - 4^+ - 6^+$ rotational band, this number is 28. For the scattering of α particles on the same levels, there are 16 equations or 12 equations according to the parity.

In the most general problem, the interaction V can be written as the scalar product of a tensor operator Q^{I^t} acting on the target, a tensor operator Q^S acting on the spin of the particle and the tensor $i^L Y_L^M(\hat{r})$ multiplying a radial form factor $V_{LSI^t}(r)$. Taking into account a factor $1/\sqrt{4\pi}$ introduced in the form factor, the geometrical coefficient is :

$$\frac{1}{\sqrt{4\pi}} \langle [(l_f \ s_f)^{j_f} I_f]^J | [i^L Y_L^M(\hat{r}) Q^S]^I \cdot Q^{I^t} | [(l_i \ s_i)^{j_i} I_i]^J \rangle = \langle s_f || Q^S || s_i \rangle \langle I_f || Q^{I^t} || I_i \rangle \quad (\text{III.43})$$

$$\times (-)^{J+I_f+j_i+\frac{1}{2}(l_f+l_i+L)} \sqrt{(2I^t+1)(2L+1)(2l_f+1)(2l_i+1)(2j_f+1)(2j_i+1)} \\ \begin{pmatrix} l_f & l_i & L \\ 0 & 0 & 0 \end{pmatrix} \begin{Bmatrix} j_f & j_i & I^t \\ I_i & I_f & J \end{Bmatrix} \begin{Bmatrix} l_f & l_i & L \\ s_f & s_i & S \\ j_f & j_i & I^t \end{Bmatrix} \quad (\text{III.44})$$

where the part of the second member written in Equ. (III.43) is the “**reduced nuclear matrix element**” and the one written in Equ. (III.44) is the geometrical factor computed by the code. This expression has the advantage to reduce the geometrical factor to the usual one of macroscopic excitations when $S = 0$ but is not symmetric when particle and target are exchanged : the reduced matrix element for a given L and $S = L$, $I^t = 0$ is $\sqrt{2S+1}$ the one for the same L and $S = 0$, $I^t = L$. Nevertheless, Equ. (III.44) was used up to **ECIS88**. To reduce this disadvantage, a factor $\sqrt{2S+1}$ has been introduced in the geometrical coefficient (III.44) in the code **ECIS94**. This correspond to use operators coupled to a scalar instead of the scalar product because :

$$(-)^{I^t} < [(l_f s_f)^{j_f} I_f]^J | [i^L Y_L^M(\hat{r}), Q^S, Q^{I^t}]_0^0 | [(l_i s_i)^{j_i} I_i]^J > \\ = \sqrt{2I^t+1} < [(l_f s_f)^{j_f} I_f]^J | [i^L Y_L^M(\hat{r}) Q^S]^{I^t} . Q^{I^t} | [(l_i s_i)^{j_i} I_i]^J > \quad (\text{III.45})$$

is symmetric for the exchange of Q^S and Q^{I^t} . In fact, the separation between reduced matrix element and geometrical coefficient correspond to the use of the expression given in Equ. (III.45) divided by $\sqrt{(2S+1)(2I^t+1)}$ in order to coincide with the usual notations for the macroscopic models.

Deformed spin-orbit interaction and magnetic multipole Coulomb interaction need a different approach. In the macroscopic models, there is no Q^S and this coefficient can be simplified.

B.2.b Spin-orbit deformation

It has been shown by **DWBA** calculations in Ref. [55] and [56] that the spin-orbit interaction should be the “full Thomas form” which is :

$$\vec{\nabla} V(\vec{r}) \times \frac{\vec{\nabla}}{i} \vec{\sigma} \quad (\text{III.46})$$

derived from Dirac’s equation by elimination of small components, in order to fit experimental data in proton elastic scattering.

B.2.b.i Spin-orbit deformation for spin one-half

For a multipole $V_\lambda(r) Y_\lambda^\mu(\hat{r})$ of the interaction $V(\vec{r})$, it can be shown by elementary manipulation of Pauli matrices as given by Equ. (III.9) and with the expression of the gradient given by Equ. (III.15) that the expression above is :

$$\vec{\nabla} [V_\lambda(r) Y_\lambda^\mu(\hat{r})] \times \frac{\vec{\nabla}}{i} \vec{\sigma} = \frac{1}{r} \left[\frac{d}{dr} V_\lambda(r) \right] Y_\lambda^\mu(\hat{r}) (\vec{\sigma} \cdot \vec{L}) - \frac{V_\lambda(r)}{r} \left[(\vec{\sigma} \cdot \vec{L}) Y_\lambda^\mu(\hat{r}) \right] \frac{d}{dr} \\ + \frac{V_\lambda(r)}{r^2} \left[(\vec{\sigma} \cdot \vec{L}) Y_\lambda^\mu(\hat{r}) \right] (\vec{\sigma} \cdot \vec{L}) - \frac{V_\lambda(r)}{r^2} \left[\vec{L} Y_\lambda^\mu(\hat{r}) \right] \cdot \vec{L} \quad (\text{III.47})$$

(see in Ref. [23] and [28]).

For any spin, the eigenvalue of $2(\vec{L} \cdot \vec{\sigma})$ is :

$$2(\vec{L} \cdot \vec{\sigma}) = \frac{1}{2} [j(j+1) - l(l+1) - s(s+1)] \quad (\text{III.48})$$

that is for spin $\frac{1}{2}$:

$$\bullet \quad \vec{L} \cdot \vec{\sigma} = l \quad \text{if} \quad j = l + \frac{1}{2},$$

- $\vec{L} \cdot \vec{\sigma} = -l - 1$ if $j = l - \frac{1}{2}$.

When taken between partial waves $|l_f j_f\rangle$ and $|l_i j_i\rangle$, the operators $\vec{L} \cdot \vec{\sigma}$ can be replaced by their eigenvalues γ_f on $|l_f j_f\rangle$ or γ_i on $|l_i j_i\rangle$. The same holds for $[\vec{L} Y_\lambda^\mu(\hat{r})] \cdot \vec{L}$. Taking carefully into account on what acts the operator, one gets the same geometrical factor G_{if}^λ as for a central term, multiplied by :

$$\frac{1}{r} \left[\frac{d}{dr} V_\lambda(r) \right] \gamma_i + \frac{V_\lambda(r)}{r} (\gamma_i - \gamma_f) \frac{d}{dr} + \frac{V_\lambda(r)}{r^2} \left[\lambda(\lambda + 1) - (\gamma_f - \gamma_i)(\gamma_f - \gamma_i \pm 1) \right] \quad (\text{III.49})$$

where ± 1 holds if the wave function is/is not multiplied by r . For the optical model, $\lambda = 0$ and $\gamma_i = \gamma_f$ and the expression (III.49) reduces to the usual one. This expression as been shown to be equivalent to the zero-range limit of a two-body spin-orbit interaction in Ref.[23] and [28] using the helicity formalism as defined in : [62] RAYNAL, J., “*Multipole expansion of a two-body interaction in helicity formalism and its application to nuclear structure and nuclear reaction calculations*”, Nucl. Phys. **A97** (1967) 572.

To compare with some earlier works using :

$$\frac{1}{r} \left[\frac{d}{dr} V_\lambda(r) \right] \frac{\gamma_i + \gamma_f}{2} \quad (\text{III.50})$$

and to be able to study the effects of each terms in the above expressions, six parameters z_1, z_2, z_3, z_4, z_5 and z_6 have been introduced to get :

$$\begin{aligned} & \frac{1}{r} \left[\frac{d}{dr} V_\lambda(r) \right] (z_1 + z_3 \gamma_i + z_4 \gamma_f) + \frac{V_\lambda(r)}{r} z_6 (\gamma_i - \gamma_f) \frac{d}{dr} \\ & + z_5 \frac{V_\lambda(r)}{r^2} \left[z_2 \lambda(\lambda + 1) - r(\gamma_f - \gamma_i)(\gamma_f - \gamma_i \pm 1) \right] \end{aligned} \quad (\text{III.51})$$

These parameters allow also to increase the spin-orbit transition without changing anything else.

The deformed spin-orbit in the “full Thomas form” introduces a first derivative of the wave function in the Schrödinger equation. The quick integration methods cannot deal with first derivatives (one has to use some Runge-Kutta method). However, the derivative terms are non diagonal because they have a factor $\gamma_i - \gamma_f$. **The deformed spin-orbit can be used only with the iteration method.**

B.2.b.ii Behaviour of the spin-orbit deformation

The behaviour of the deformed spin orbit interaction is a vector behaviour: the difference $\gamma_i - \gamma_f$ is also found to be the ratio of the geometries for a transfer of spin and a scalar interaction :

$$\left\{ \begin{array}{ccc} l_f & l_i & L \\ \frac{1}{2} & \frac{1}{2} & 1 \\ j_f & j_i & J \end{array} \right\} = \frac{\gamma_f - \gamma_i}{\sqrt{3J(J+1)}} \left\{ \begin{array}{ccc} l_f & l_i & L \\ \frac{1}{2} & \frac{1}{2} & 0 \\ j_f & j_i & J \end{array} \right\} \quad (\text{III.52})$$

To compare deformed spin-orbit and other interactions, let us write consider a $0^+ \rightarrow 2^+$ reaction. There is one equation for the 0^+ state and 5 equations for the 2^+ state. Let us consider their behaviour when the angular momenta increase. Among the five geometrical coefficients, there are :

1. those between $l_i = j_i \pm \frac{1}{2}$ and $l_f = j_f \pm \frac{1}{2}$:
 - for a central interaction, we shall take them as reference,
 - for a $\Delta S = 1$ interaction, Equ. (III.52) shows that there are of the same order,
 - for a spin-orbit interaction, the factor γ_i increases linearly with l and the other are constant.
2. those between $l_i = j_i \pm \frac{1}{2}$ and $l_f = j_f \mp \frac{1}{2}$ which :
 - are a factor l smaller than the first ones for a central term,

- for a $\Delta S = 1$ interaction, the Equ. (III.52) shows that there are multiplied by a factor l and become of the same order as the central interaction (1);
- for a spin-orbit interaction, the factors γ_i and $\gamma_i - \gamma_f$ increase linearly with l and the factor $6 - (\gamma_i - \gamma_f)(\gamma_i - \gamma_f + 1)$ increases quadratically.

This behaviour shows that spin-orbit interaction should become predominant at high energy.

B.2.b.iii Spin-orbit deformation for any spin

The “full Thomas form” of the spin-orbit interaction is conserved when the nucleon-nucleus potential is folded with the intrinsic wave function of the particle if this wave function involves only relative S-states. This is well known for the optical model of deuteron (see for example Ref. [22] Chapter IV). If the nucleon-nucleus spin-orbit form factor is $\frac{1}{r} \frac{d}{dr} f(r)$, the deuteron spin-orbit form factor is $\frac{1}{r} \frac{d}{dr} F(r)$ where $F(r)$ is obtained from $f(r)$ by folding, only if the deuteron D-wave is neglected.

This property of the folding for a structureless incident particle allows to extend the “full Thomas form” of the spin-orbit to any spin as done in Ref. [38]. For a spin \vec{s} , $2\vec{s}$ can be replaced by the sum of $2|s|$ Pauli matrices σ_i in Equ. (III.9) and Equ. (III.47) derived for each Pauli matrix. In this result, the term without Pauli matrices is multiplied by $2|s|$, the terms with one Pauli matrix are summed up to $2\vec{s}$ and the terms with two Pauli matrices need some recoupling. The total result is :

$$\frac{1}{r} \frac{dV_\lambda}{dr} \gamma_i + (\gamma_i - \gamma_f) \frac{V_\lambda}{dr} \frac{d}{dr} + \frac{V_\lambda}{2r^2} \left\{ |s| [\lambda(\lambda + 1) - l_i(l_i + 1) - l_f(l_f + 1)] + \gamma_f + \frac{\gamma_i \gamma_f}{2|s|} + F \right\} \quad (\text{III.53})$$

with :

$$F = -2|s| f_{l_i s j_i} f_{l_f s j_f} \begin{Bmatrix} l_i & j_i & s-1 \\ j_f & l_f & \lambda \end{Bmatrix} \begin{Bmatrix} l_i & j_i & s \\ j_f & l_f & \lambda \end{Bmatrix}^{-1} \quad (\text{III.54})$$

and :

$$\begin{aligned} f_{l s j} &= \langle l s j | (\vec{\sigma} \cdot \vec{L}) | l s - 1 j \rangle \\ &= - \frac{\sqrt{[l(l+1) - (j-s)(j-s+1)][(j+s)(j+s+1) - l(l+1)]}}{2|s|} \end{aligned} \quad (\text{III.55})$$

The parametrisation introduced for spin one-half has been extended to :

$$\begin{aligned} &\frac{1}{r} \frac{dV_\lambda}{dr} (z_1 + z_3 \gamma_i + z_4 \gamma_f) + z_6 (\gamma_i - \gamma_f) \frac{V_\lambda}{r} \frac{d}{dr} \\ &+ z_5 \frac{V_\lambda}{2r^2} \left\{ |s| [z_2 \lambda(\lambda + 1) - l_i(l_i + 1) - l_f(l_f + 1)] + \gamma_f + \frac{\gamma_i \gamma_f}{2|s|} + F \right\} \end{aligned} \quad (\text{III.56})$$

B.2.c Dirac formalism

The large and the small component have a different angular dependence. This point is taken into account by writing instead of Equ. (III.36) :

$$|\Psi(\vec{r})\rangle = \frac{1}{r} \sum_{l, j, I, J} \begin{pmatrix} F_{l, j, I, J}(r) \\ -i G_{l, j, I, J}(r) (\vec{\sigma} \cdot \hat{r}) \end{pmatrix} Y_{l s j I J M} \quad (\text{III.57})$$

where $Y_{l s j I J M}$ is the same as for Schrödinger equation and is given by Equ. (III.37). Here, l stands instead of the parity quantum number.

The operator $-(\vec{\sigma} \cdot \hat{r})$ interchanges the spin angular functions of the large and the small components.

Projecting the total equation on all the function Y_{lsjIJM} , we obtain the set of linear coupled equations :

$$\begin{aligned} \hbar \left[\frac{d}{dr} - \frac{K_i}{r} - \frac{1}{2m} \frac{dV_T^0(r)}{dr} \right] G_i(r) + [E_i - m - V_S^0 - V_V^0] F_i(r) &= S_i(r) \\ -\hbar \left[\frac{d}{dr} + \frac{K_i}{r} + \frac{1}{2m} \frac{dV_T^0(r)}{dr} \right] F_i(r) + [E_i + m + V_S^0 - V_V^0] G_i(r) &= T_i(r) \end{aligned} \quad (\text{III.58})$$

where the suffix i stands for (l, s, j, I, J) , V^0 indicates the monopole parts of the interactions and :

$$K_i = (-)^{l-j-\frac{1}{2}}(j + \frac{1}{2}) \quad (\text{III.59})$$

The second members are :

$$\begin{aligned} S_i(r) &= \sum_{\lambda,j} G_{i,j}^\lambda \left[\{V_S^\lambda + V_V^\lambda\} F_j(r) + \frac{\hbar}{2m} \left\{ \left(\frac{d}{dr} + \frac{K_i - K_j}{r} \right) V_T^\lambda \right\} G_j(r) \right] \\ T_i(r) &= \sum_{\lambda,j} G_{i,j}^\lambda \left[\{V_S^\lambda - V_V^\lambda\} G_j(r) + \frac{\hbar}{2m} \left\{ \left(\frac{d}{dr} + \frac{K_j - K_i}{r} \right) V_T^\lambda \right\} F_j(r) \right] \end{aligned} \quad (\text{III.60})$$

where V^λ are the multipoles of the interactions and the geometrical coefficients are the same as for the Schrödinger equation :

- between F_i and F_j , expressions are identical,
- between G_i and G_j , the operators $(\vec{\sigma} \cdot \hat{r})$ commute with the expression and their product being unity, the result is the same as between F_i and F_j ,
- between F_i and G_j , the operators $(\vec{\sigma} \cdot \vec{\nabla})$ and $(\vec{\sigma} \cdot \hat{r})$ are treated like the spin-orbit deformation in the Schrödinger equation.

There is no reason to limit the coupling to macroscopic ones ($\Delta S = 0$). For $\Delta S \neq 0$, the rules given above are no more valid and the coefficient of the central term in the second equation (III.60) is no more the same as the coefficient of the first one and must be computed separately.

The codes **ECIS88** and **ECIS94** offer two possibilities :

- **to solve exactly the Dirac equation** as described above with `LO(100)=.TRUE.`,
- **to write the equivalent Schrödinger equation and to solve it, neglecting the presence of $D(\vec{r})^{\frac{1}{2}}$ in the definition given by Equ. (III.11) of the function** if `LO(99)=.TRUE.`

The result is exactly the same for the elastic scattering.

B.3 Solutions and angular distributions

We assume that there is no long range interactions. They shall be explained in Chapter (IV).

B.3.a Solution of the equation

Beyond a matching point for which all the potentials except the Coulomb one vanish, the solution $f_i(r)$ of the Schrödinger equation is a superposition of the regular Coulomb function $F_{l_i}(\eta_i; k_i r)$ and the irregular Coulomb function $G_{l_i}(\eta_i; k_i r)$. For the equivalent Schrödinger equation of the Dirac equation, there are long range terms r^{-2} and r^{-3} which implies the use of *relativistic Coulomb functions* instead of the usual, non-relativistic ones.

Coulomb corrections described in Ref. [7] allow the use of the non-relativistic Coulomb functions in the Dirac phenomenology for nucleon scattering. They were derived for heavy-ion

scattering to reduce the matching radius to the point where nuclear interactions vanish and to correct the non relativistic Coulomb functions at this point for the effect of the long range Coulomb interaction which decrease only as r^{-l-1} . We consider here that the Coulomb functions are corrected for these long range effects.

The solution must vanish at the origin, due to the factor $1/r$ in the definition of the radial functions. They must have a plane wave incoming part for the initial channel i and an outgoing wave for all the channels, normalised to the same flux (that is divided by \sqrt{k}). Their asymptotic form is :

$$\begin{aligned} f_f(r)_{r \rightarrow \infty} &\Rightarrow \delta_{i,f} F_{l_f}(\eta_f; k_f r) + C_f^i \left(\frac{k_i m_f}{k_f m_i} \right)^{\frac{1}{2}} [G_{l_f}(\eta_f; k_f r) + i F_{l_f}(\eta_f; k_f r)] \\ &\Rightarrow \delta_{i,f} F_{l_f}(\eta_f; k_f r) + \overline{C}_f^i [G_{l_f}(\eta_f; k_f r) + i F_{l_f}(\eta_f; k_f r)] \end{aligned} \quad (\text{III.61})$$

and numerical integration gives the coefficients \overline{C}_f^i by linear combination of the n solutions which vanish at the origin. The \overline{C}_f^i are multiplied by $\sqrt{k_f m_i / (k_i m_f)}$ to get the C_f^i with are symmetric for the exchange of i and f as will be shown by Equ. (III.86). The same relation holds in Dirac formalism for the solution of the equivalent Schrödinger equation. An equivalent definition of the C_f^i will be given by Equ. (III.104) and (III.105).

By identification we obtain :

$$\begin{aligned} f_{\mu_f \sigma_f \mu_i \sigma_i}(\theta) &= \frac{4\pi}{k_i} \sum \exp(i\sigma_{l_i} + i\sigma_{l_f}) C_f^i < l_i, s, \nu_i, \sigma_i | j_i, m_i > < j_i, I_i, m_i, \mu_i | J, M > \\ &\times < l_f, s, \nu_f, \sigma_f | j_f, m_f > < j_f, I_f, m_f, \mu_f | J, M > Y_{l_i}^{\nu_i*}(\hat{k}_i) Y_{l_f}^{\nu_f}(\hat{k}_f) \end{aligned} \quad (\text{III.62})$$

where σ and μ are the projections of the spins of the particle and the target, σ_{l_i} and σ_{l_f} are the Coulomb phase-shifts.

B.3.b Helicity formalism and cross-section

These amplitude are simplified by the helicity formalism (defined in Ref. [20]; see also Ref. [21] and [22]) in which the spin of the particle is projected on its momentum and the spin of the target is projected on the inverse direction.

The axis of quantisation is along \vec{k}_i for the initial state and along \vec{k}_f for the final state and the helicity of the target is opposite of the projection of its spin.

The helicity amplitude involve only a reduced rotation matrix element $r_{m,m'}^{(J)}(\theta)$:

$$f_{\mu_f \sigma_f \mu_i \sigma_i}^{hel.}(\theta) = f_{\mu_f \sigma_f \mu_i \sigma_i}^{(Coul) hel.}(\theta) \delta_{if} + \sum_J f_{\mu_f \sigma_f \mu_i \sigma_i}^{(J) hel.} r_{\sigma_f - \mu_f, \sigma_i - \mu_i}^{(J)}(\theta) \quad (\text{III.63})$$

with :

$$\begin{aligned} f_{\mu_f \sigma_f \mu_i \sigma_i}^{(J) hel.}(\theta) &= \frac{1}{k_i} \sum \exp(i\sigma_{l_i} + i\sigma_{l_f}) C_f^i \sqrt{(2l_i + 1)(2l_f + 1)} < l_i, s_i, 0, \sigma_i | j_i, \sigma_i > \\ &\times < j_i, I_i, \sigma_i, -\mu_i | J, M_i > < l_f, s_f, 0, \sigma_f | j_f, \sigma_f > < j_f, I_f, \sigma_f, -\mu_f | J, M_f > \end{aligned} \quad (\text{III.64})$$

The coulomb helicity amplitude, which appears only in the incident channel, is the usual coulomb amplitude multiplied by the reduced rotation matrix elements for the spin s_i of the particle and the spin I_i of the target :

$$f_{\mu_f \sigma_f \mu_i \sigma_i}^{(Coul) hel.}(\theta) = -\frac{\eta}{2k \sin^2 \frac{\theta}{2}} \exp \left(-2i\eta \ln \sin \frac{\theta}{2} + 2i\sigma_0 \right) r_{\sigma_f, \sigma_i}^{(s_i)}(\theta) r_{-\mu_f, -\mu_i}^{(I_i)}(\theta) \quad (\text{III.65})$$

Note that in the codes **ECIS**, σ_0 is used only to compute Coulomb corrections and that the Coulomb phase-shift σ_l is replaced everywhere else by $\sigma_l - \sigma_0$, independently for each level.

The direct interaction differential cross-section is :

$$\frac{d\sigma(\theta)}{d\Omega} = \frac{1}{(2s_i + 1)(2I_i + 1)} \sum_{\sigma_i \sigma_f \mu_i \mu_f} |f_{\mu_f \sigma_f \mu_i \sigma_i}^{hel.}(\theta)|^2 \Phi \quad (\text{III.66})$$

and the total differential cross-section is obtained by adding the **differential compound cross-section** which is computed from the coefficients C_f^i .

In the code ECIS, the compound differential cross-section is expressed as a sum of terms, each of them being a sum of squares of reduced rotation matrix elements; usually, these squares are expanded into sums of Legendre polynomials.

B.3.c Observables

The formalism used to express the observables as been described in Ref. [21]. All the observables P can be expresses as :

$$P(\theta) \quad \frac{d\sigma(\theta)}{d\Omega} = \sum_{\lambda_i \mu_i} x_{\lambda_1 \mu_1 \lambda_2 \mu_2}^{\lambda_3 \mu_3 \lambda_4 \mu_4} \quad A_{\lambda_1 \mu_1 \lambda_2 \mu_2}^{\lambda_3 \mu_3 \lambda_4 \mu_4} \quad (\text{III.67})$$

where :

$$A_{\lambda_1 \mu_1 \lambda_2 \mu_2}^{\lambda_3 \mu_3 \lambda_4 \mu_4} = \text{Trace} \left\{ f^{hel.}(\theta) [\tau_{\lambda_1 \mu_1}^{s_i} \otimes \tau_{\lambda_2 \mu_2}^{I_i}] f^{hel.}(\theta)^+ [\tau_{\lambda_3 \mu_3}^{s_f} \otimes \tau_{\lambda_4 \mu_4}^{I_f}] \right\} \quad (\text{III.68})$$

and, in peculiar, A_{0000}^{0000} is the cross-section. Here, the τ are tensor operators in spin space defined for $0 \leq \lambda \leq 2s$ and $-\lambda \leq \mu \leq \lambda$ by :

$$\langle s \, q | \tau_{\lambda \mu}^s | s \, q' \rangle = (-)^{s-q'} \sqrt{2s+1} \langle s, s, q - q', \lambda | \mu, ec69 \rangle \quad (\text{III.69})$$

Due to the helicity formalism, the description of the polarisation of the outgoing particles is along the outgoing direction in the center of mass system. The description can be shortened, using the relations :

$$\begin{array}{ll} \text{hermiticity} & \left(A_{\lambda_1 \mu_1 \lambda_2 \mu_2}^{\lambda_3 \mu_3 \lambda_4 \mu_4} \right)^* = (-)^{\sum \mu} A_{\lambda_1 - \mu_1 \lambda_2 - \mu_2}^{\lambda_3 - \mu_3 \lambda_4 - \mu_4} \\ \text{parity} & A_{\lambda_1 \mu_1 \lambda_2 \mu_2}^{\lambda_3 \mu_3 \lambda_4 \mu_4} = (-)^{\sum \lambda + \mu} A_{\lambda_1 - \mu_1 \lambda_2 - \mu_2}^{\lambda_3 - \mu_3 \lambda_4 - \mu_4} \end{array} \quad (\text{III.70})$$

As consequences :

1. if the sum on λ is even, A is real,
2. if the sum on λ is odd, A is pure imaginary,
3. for the same λ 's and opposite μ 's, the A differs by $(-)^{\sum \lambda + \sum \mu}$.

For example, the spin-flip SP is :

$$SP(\theta) \quad \frac{d\sigma(\theta)}{d\Omega} = \frac{1}{2} (A_{0000}^{0000} + A_{1100}^{1100} + A_{1-100}^{1100}) \quad (\text{III.71})$$

Experiments are in the laboratory system; if some axis in the reaction plane is requested to describe the final polarisation, a small rotation of the spin is needed for the scattering matrix. These rotations will be explained in section (VIIA).

Some observables as Q are defined with only the incident direction :

$$Q(\theta) \quad \frac{d\sigma(\theta)}{d\Omega} = \sqrt{2} A_{1000}^{1100} \quad (\text{III.72})$$

and need a rotation of angle θ on the scattering matrix.

There are many ways to define an observable for the codes **ECIS** (with an axis perpendicular to the scattering plane, using magnetic quantum numbers, usual tensors operators). The codes transform informations into A 's, store information for necessity of rotation of the scattering matrix and compute do-loops to use Equ. (III.68).

C Integration methods

More details on these topics can be found in Ref. [1] for single equations, in Ref. [2] for coupled equations.

C.1 Single equation

The second order linear differential equation without first derivative :

$$f''(r) = V(r) f(r) \quad (\text{III.73})$$

is easily solved, using the relation between the function and its second derivative at three equidistant points :

$$\xi(r+h) = f(r+h) - \frac{h^2}{12} f''(r+h) = 2f(r) + \frac{5h^2}{6} f''(r) - f(r-h) + \frac{h^2}{12} f''(r-h) \quad (\text{III.74})$$

with the truncation error :

$$\Delta = -\frac{h^6 f^{(VI)}(r)}{240} \quad (\text{III.75})$$

(it is called Cowell method in Ref. [1]). If there is no need of the function $f(r)$, the integration can be performed using :

$$\xi(r+h) = 2\xi(r) - \xi(r-h) + u(r), \quad u(r) = \frac{h^2 V(r)}{1 - \frac{h^2}{12} V(r)} \xi(r) \quad (\text{III.76})$$

which is the Numerov's method and has the same truncation error.

For coupled equations, $V(r)$ is a matrix and the computation of $u(r)$ involves the resolution of a linear system of equations. It is easier to use the **Modified Numerov method** by expanding the fraction up to the terms in h^4 :

$$u(r) = \left\{ h^2 V(r) + \frac{h^4}{12} V^2(r) \right\} \xi(r) \quad (\text{III.77})$$

with a truncation error :

$$\Delta = \frac{h^6 V^3(r) f(r)}{144} - \frac{h^6 f^{(VI)}(r)}{240} \quad (\text{III.78})$$

For a constant potential, this truncation error is $-\frac{2}{3}$ the previous one. Such an expansion in the Cowell method has been used in : [63] BEURTEY, R., GUILLOU and RAYNAL J., "*Etude de la diffusion élastique des particules chargées à l'aide du modèle optique*", Le Journal de Physique et le Radium **21** (1960) page 402. If needed, the wave function can be obtained by one of the two expressions :

$$\begin{aligned} f(r) &= \xi(r) + \frac{1}{12} u(r) \\ f(r) &= \frac{1}{12} [\xi(r-h) + 10\xi(r) + \xi(r+h)] \end{aligned} \quad (\text{III.79})$$

with an error of the same order.

Note that (see Ref. [1]) :

- the Schrödinger equation is easily solved numerically by these methods, starting from 0 at the origin and a small value ϵ at the point h ,
- to have a precise result, the starting values have to be modified for the angular momentum $l = 1$,
- for high angular momenta, the first few points do not matter for the result,
- as the solution increases quickly, it has to be divided by a large number as soon as it becomes too large,

- for coupled equations (see Ref. [2]), the starting values are ϵ for only one component, but a Schmidt's orthogonalisation procedure can be needed from time to time to avoid that solutions align between them,
- the Schrödinger equivalent of the Dirac equation for the angular momentum $l = 0$ with Coulomb potential down to the origin can need a power expansion as starting values, but it is not used in ECIS.

Matching with the asymptotic expressions can be done by writing the matching conditions at the points $R \pm h$, using :

$$\mathcal{F}_{\uparrow}(\eta; \nabla \pm \langle \rangle) = \mathcal{F}_{\uparrow}(\eta; \nabla \pm \langle \rangle) - \frac{\langle \epsilon}{\infty \epsilon} \mathcal{F}_{\uparrow}''(\eta; \nabla \pm \langle \rangle), \quad \mathcal{G}_{\uparrow}(\eta; \nabla \pm \langle \rangle) = \mathcal{G}_{\uparrow}(\eta; \nabla \pm \langle \rangle) - \frac{\langle \epsilon}{\infty \epsilon} \mathcal{G}_{\uparrow}''(\eta; \nabla \pm \langle \rangle) \quad (\text{III.80})$$

where the values of $\mathcal{F}_{\uparrow}(\eta; \nabla \pm \langle \rangle)$ are obtained by writing three Numerov integration between $r - h$ and $r + h$ with steps $h/2$ and :

$$7[f(r+h) - f(r-h)] + 16[f(r+h/2) - f(r-h/2)] - 3h^2[f''(r+h/2) - f''(r-h/2)] = 30hf'(r) \quad (\text{III.81})$$

and eliminating the functions at $R \pm h/2$ between these four equations.

The linear system of matching conditions :

$$\begin{aligned} \xi_i^k(R+h) &= \sum_j \alpha_j^k \left\{ \mathcal{F}_j(\eta; \mathcal{R} + \langle \rangle) \delta_{ij} + \overline{\mathcal{C}}_j^{\dagger} [\mathcal{G}_j(\eta; \mathcal{R} + \langle \rangle) + \mathcal{F}_j(\eta; \mathcal{R} + \langle \rangle)] \right\} \\ \xi_i^k(R-h) &= \sum_j \alpha_j^k \left\{ \mathcal{F}_j(\eta; \mathcal{R} - \langle \rangle) \delta_{ij} + \overline{\mathcal{C}}_j^{\dagger} [\mathcal{G}_j(\eta; \mathcal{R} - \langle \rangle) + \mathcal{F}_j(\eta; \mathcal{R} - \langle \rangle)] \right\} \end{aligned} \quad (\text{III.82})$$

is simplified by writing “pseudo-Wronskians” of the numerical solutions with the regular and the irregular Coulomb functions, that is, expressions giving Wronskians when h tends to 0. Writing the matrices :

$$\begin{aligned} A_i^k &= \frac{\xi_i^k(R+h) \mathcal{G}_j(\eta; \mathcal{R} - \langle \rangle) - \xi_j^{\parallel}(\mathcal{R} - \langle \rangle) \mathcal{G}_j(\eta; \mathcal{R} + \langle \rangle)}{\mathcal{F}_j(\eta; \mathcal{R} + \langle \rangle) \mathcal{G}_j(\eta; \mathcal{R} - \langle \rangle) - \mathcal{F}_j(\eta; \mathcal{R} - \langle \rangle) \mathcal{G}_j(\eta; \mathcal{R} + \langle \rangle)} \\ B_i^k &= \frac{\xi_i^k(R+h) \mathcal{F}_j(\eta; \mathcal{R} - \langle \rangle) - \xi_j^{\parallel}(\mathcal{R} - \langle \rangle) \mathcal{F}_j(\eta; \mathcal{R} + \langle \rangle)}{\mathcal{F}_j(\eta; \mathcal{R} + \langle \rangle) \mathcal{G}_j(\eta; \mathcal{R} - \langle \rangle) - \mathcal{F}_j(\eta; \mathcal{R} - \langle \rangle) \mathcal{G}_j(\eta; \mathcal{R} + \langle \rangle)} \end{aligned} \quad (\text{III.83})$$

the matching conditions become :

$$A_i^k = \sum_j \alpha_j^k \{ \delta_{ij} + i \overline{\mathcal{C}}_i^j \}, \quad B_i^k = - \sum_j \alpha_j^k \overline{\mathcal{C}}_i^j, \quad (\text{III.84})$$

and the solution is given by the linear system :

$$B_i^k = - \sum_j (A_j^k + i B_j^k) \overline{\mathcal{C}}_i^j. \quad (\text{III.85})$$

It should be noted that the solution of these equation for only one set of second members B_i^k for fixed i gives the $\overline{\mathcal{C}}_i^j$ for fixed i whereas the $\overline{\mathcal{C}}_i^j$ are needed for fixed j . Wronskian relations give :

$$\frac{k_i}{m_i} \overline{\mathcal{C}}_j^i = \frac{k_j}{m_j} \overline{\mathcal{C}}_i^j \quad (\text{III.86})$$

which can be used to obtain result from only one system of linear equations. This relation prove the symmetry of the matrix \mathcal{C} defined in Equ. (III.61) for $m_i = m_j$.

The error on the collision matrix is of the order of h^4 . They are equivalent to an error to the potential by a term :

$$\Delta V = \frac{h^4}{240} \frac{d^6}{dr^6}, \quad \Delta V = \frac{h^4}{240} \frac{d^6}{dr^6} - \frac{h^4}{144} V^3(r), \quad (\text{III.87})$$

for the Numerov method and for the Modified Numerov method respectively. **They have a coherent contribution.** The step size in kr can be as large as $\frac{\pi}{2}$ but smaller than the diffusenesses. A good value is $h = 0.5$.

C.2 Green function for Schrödinger equation

The differential equation written :

$$\left\{ -\frac{\hbar^2}{2m_i} \left[\frac{d^2}{dr^2} - \frac{2\eta_i k_i}{r} - \frac{l_i(l_i+1)}{r^2} + k_i^2 \right] - V_i^{opt}(r) + V_{ii}(r) \right\} f_i(r) = - \sum_j V_{ij}(r) f_j(r) \quad (\text{III.88})$$

and the boundary conditions for the solution :

$$f_i(r)_{r \rightarrow \infty} = \delta_{i,0} F_{l_i}(\eta_i; k_i r) + \overline{C}_i^0 [G_{l_i}(\eta_i; k_i r) + i F_{l_i}(\eta_i; k_i r)] \quad (\text{III.89})$$

where $\delta_{i,0}$ means only in the incident channel, can be replaced by an integral equation built with two kinds of solution of the single equations of the right member:

1) the **regular** or **optical** solution completely defined as vanishing at the origin and by its asymptotic value :

$$f_i^{reg}(r)_{r \rightarrow \infty} = F_{l_i}(\eta_i; k_i r) + C_i^{opt} [G_{l_i}(\eta_i; k_i r) + i F_{l_i}(\eta_i; k_i r)] \quad (\text{III.90})$$

obtained by numerical integration from the origin.

2) the **irregular** solution completely defined by its asymptotic value :

$$f_i^{irr}(r)_{r \rightarrow \infty} = G_{l_i}(\eta_i; k_i r) + i F_{l_i}(\eta_i; k_i r) \quad (\text{III.91})$$

obtained by numerical backward integration from the matching point.

These solutions are such that the Wronskian is :

$$f_i^{irr}(r) f_i'^{reg}(r) - f_i'^{irr}(r) f_i^{reg}(r) = k_i \quad (\text{III.92})$$

In practice, the irregular solution will be divided by k_i so that the Wronskian is unity. With these functions, the solution is :

$$\begin{aligned} f_i(r) = f_i^{reg}(r) \delta_{i,0} - \frac{2m_i}{\hbar^2 k_i} \left\{ f_i^{irr}(r) \int_0^r f_i^{reg}(r') \sum_j V_{ij}(r') f_j(r') dr' \right. \\ \left. + f_i^{reg}(r) \int_r^\infty f_i^{irr}(r') \sum_j V_{ij}(r') f_j(r') dr' \right\} \end{aligned} \quad (\text{III.93})$$

and the coefficient of the outgoing wave is :

$$\overline{C}_i^0 = C_i^{opt} \delta_{i,0} - \frac{2m_i}{\hbar^2 k_i} \int_0^\infty f_i^{reg}(r) \sum_j V_{ij}(r) f_j(r) dr \quad (\text{III.94})$$

The DWBA approximation is to replace $f_j(r)$ by $f_j^{reg}(r)$ in Equ. (III.94).

The derivative terms generated by the deformed spin orbit are in the second member of Equ. (III.88). In fact, it is $g(r) = r f'(r)$ which is needed because the form factor stored for the derivative term of Equ. (III.49) is $V_\lambda(r)/r^2$. It is obtained by numerical derivation related to the formulae giving the derivative y of the function y :

- for the 3 first points :

$$\begin{aligned} x_1 &= [147y_1 - 360y_2 + 450y_3 - 400y_4 + 225y_5 - 72y_6 + 10y_7]/(60h) \\ x_2 &= [10y_1 + 77y_2 - 150y_3 + 100y_4 - 50y_5 + 15y_6 - 2y_7]/(60h) \\ x_3 &= [-2y_1 + 24y_2 + 35y_3 - 80y_4 + 30y_5 - 8y_6 + y_7]/(60h) \end{aligned} \quad (\text{III.95})$$

- from the 4th point to $N - 3$ if there are N points :

$$x_i = [45(y_{i-1} - y_{i+1}) - 9(y_{i-2} - y_{i+2}) + y_{i-3} - y_{i+3}]/(60h) \quad (\text{III.96})$$

- for the 3 last points :

$$\begin{aligned}
x_{N-2} &= [y_{N-6} - 8y_{N-5} + 30y_{N-4} - 80y_{N-3} + 35y_{N-2} + 24y_{N-1} - 2y_N]/(60h) \\
x_{N-1} &= [-2y_{N-6} + 15y_{N-5} - 50y_{N-4} + 100y_{N-3} - 150y_{N-2} + 77y_{N-1} + 10y_N]/(60h) \\
x_N &= [10y_{N-6} - 72y_{N-5} + 225y_{N-4} - 400y_{N-3} + 450y_{N-2} - 360y_{N-1} + 147y_N]/(60h)
\end{aligned} \tag{III.97}$$

In fact, in this application, the formulae for the first three points are shifted, taking into account that the function to derive is zero at the origin.

Integrals with the Green function are not straightforward, due to the discontinuity of the integrand. With :

$$\int_0^h f(x)dx = \frac{h}{2}[f(0) + f(h)] + \frac{h^2}{12}[f'(0) - f'(h)] + \frac{h^5}{720}f^{(IV)}(r) \tag{III.98}$$

For Schrödinger equation, the integrand vanishes at the origin and at ∞ and its derivative has the discontinuity given by Equ. (III.92). Using :

$$F(nh) = \sum_{i=1}^n W(ih)f^{reg}(ih), \quad G(nh) = \sum_{i=n+1}^{\infty} W(ih)f^{irr}(ih) \tag{III.99}$$

where $W(ih)$ is the second member of Equ. III.85, the integral is :

$$f(nh) \propto f^{irr}(nh)F(nh) + f^{reg}(nh)G(nh) - \frac{h^2 k}{12}W(nh) \tag{III.100}$$

with an error of the order h^4 . In this equation k is the Wronskian given by Equ. (III.92) and reduces to unity in actual calculation with a redefinition of the irregular functions.

C.3 Green function for Dirac equation

The Dirac equation can be written as :

$$\begin{aligned}
\hbar\left[\frac{d}{dr} - A_i(r)\right]G_i(r) + E_i(r)F_i(r) &= S_i(r) \\
-\hbar\left[\frac{d}{dr} + A_i(r)\right]F_i(r) + D_i(r)G_i(r) &= T_i(r)
\end{aligned} \tag{III.101}$$

with :

$$A_i(r) = \frac{K_i}{r} + \frac{1}{2m} \frac{dV_T^0(r)}{dr}, \quad D_i(r) = E_i + m + V_S^0 - V_V^0, \quad E_i(r) = E_i - m - V_S^0 - V_V^0. \tag{III.102}$$

$S_i(r)$ and $T_i(r)$ are given by Equ. (III.60).

Two independent solutions of uncoupled equations are obtained by :

1. the two linear coupled equations are replaced by the equivalent Schrödinger equation,
2. its regular solution $f^{reg}(r)$ defined by Equ. (III.90) and its irregular solution $f^{irr}(r)$ defined by Equ. (III.91) are obtained,
3. they are multiplied by $D(r)^{\frac{1}{2}}$ to obtain $F^{reg}(r)$ and $F^{irr}(r)$ (there is no difficulty to obtain continuous square root of the complex function $D(r)$ as long as we assume that the real part is very large with respect to its imaginary part),
4. the small components $G^{reg}(r)$ and $G^{irr}(r)$ are obtained from the Dirac equation by numerical derivation using Equ. (III.95) to Equ. (III.97).

Due to the number of operations and their difficulties, all the reorientation terms are shifted in the second member of Equ. (III.58) (this is also done in the Schrödinger formalism if `LO(29)=TRUE.`) and these functions are kept if there are needed for the system with an higher total J .

With the relation (III.92) for $f_i^{reg}(r)$ and $f_i^{irr}(r)$ and the Dirac equation, one obtains the Wronskian :

$$F_i^{irr}(r) G_i^{reg}(r) - G_i^{irr}(r) F_i^{reg}(r) = \hbar k_i \quad (\text{III.103})$$

Here also, the irregular solution will be normalised such that the Wronskian is unity. With these functions, the system of equations (III.58) and its boundary conditions can be replaced by the integral equation :

$$\begin{aligned} F_i(r) &= F_i^{reg}(r) \delta_{i,0} - \frac{1}{\hbar^2 k_i} \left\{ F_i^{irr}(r) \int_0^r [F_i^{reg}(r') S_i(r') + G_i^{reg}(r') T_i(r')] dr' \right. \\ &\quad \left. + F_i^{reg}(r) \int_r^\infty [F_i^{irr}(r') S_i(r') + G_i^{irr}(r') T_i(r')] dr' \right\} \\ G_i(r) &= G_i^{reg}(r) \delta_{i,0} - \frac{1}{\hbar^2 k_i} \left\{ G_i^{irr}(r) \int_0^r [F_i^{reg}(r') S_i(r') + G_i^{reg}(r') T_i(r')] dr' \right. \\ &\quad \left. + G_i^{reg}(r) \int_r^\infty [F_i^{irr}(r') S_i(r') + G_i^{irr}(r') T_i(r')] dr' \right\} \end{aligned} \quad (\text{III.104})$$

and the coefficient of the outgoing wave is the coefficient of the regular functions at ∞ :

$$\overline{C}_i^0 = C_i^{opt} \delta_{i,0} - \frac{1}{\hbar^2 k_i} \int_0^\infty [F_i^{reg}(r') S_i(r') + G_i^{reg}(r') T_i(r')] dr' \quad (\text{III.105})$$

It is more difficult to obtain a precise value of this integral than in the Schrödinger formalism. The two terms of Equ. (III.98) introduce corrections and Equ. (III.100) does not hold.

To find the error coming from the first term of Equ. (III.98) let us add two elementary step sizes around $r = r'$:

$$\begin{aligned} \frac{h}{2} [F_i^{irr}(r) G_i^{reg}(r) T_i(r) + F_i^{reg}(r) G_i^{irr}(r) T_i(r)] &= h \left[F_i^{irr}(r) G_i^{reg}(r) - \frac{\hbar k_i}{2} \right] T_i(r) \\ \frac{h}{2} [G_i^{irr}(r) F_i^{reg}(r) S_i(r) + G_i^{reg}(r) F_i^{irr}(r) S_i(r)] &= h \left[G_i^{irr}(r) F_i^{reg}(r) + \frac{\hbar k_i}{2} \right] S_i(r) \end{aligned} \quad (\text{III.106})$$

which is an error of order h if this correction is neglected.

With the second term of Equ. (III.98), we obtain a correction :

$$\begin{aligned} & -\frac{\hbar^2 k_i}{12} \left[D_i(r) S_i(r) + \hbar \left\{ A_i(r) T_i(r) + \frac{d}{dr} T_i(r) \right\} \right], \\ & -\frac{\hbar^2 k_i}{12} \left[E_i(r) T_i(r) + \hbar \left\{ A_i(r) S_i(r) + \frac{d}{dr} S_i(r) \right\} \right], \end{aligned} \quad (\text{III.107})$$

for $F_i(r)$ and $G_i(r)$ respectively. All these correction give an error of the order h^4 . However, the derivatives of $S_i(r)$ and $T_i(r)$ are neglected in the code `ECIS` for the three first and the three last points.

C.4 Calculations

The actual calculations are quite different of what was presented here. In the Schrödinger formalism, the system of coupled equations (III.88) is replaced by :

$$\left[\frac{d^2}{dr^2} - \frac{2\eta_i k_i}{r} - \frac{l_i(l_i + 1)}{r^2} + k_i^2 + \frac{2m_i}{\hbar^2} \{ V_i^{opt}(r) - V_{ii}(r) \} \right] g_i(r) = \sum_j \mathcal{V}_{ij}(r) g_j(r) \quad (\text{III.108})$$

with :

$$f_i(r) = m_i^{\frac{1}{2}} g_i(r), \quad \mathcal{V}_{ij}(r) = \frac{2(m_i m_j)^{\frac{1}{2}}}{\hbar^2} V_{ij}(r) \quad (\text{III.109})$$

The result differs from the \overline{C} matrix defined by Equ. (III.61) : its matrix elements have to be multiplied by $\sqrt{k_f/k_i}$ to obtain the C_i^f .

Calculations are also simplified by using f_i^{irr} such that the Wronskian given by Equ. (III.92) and (III.103) is unity.

D Equations couplées en iterations séquentielles

The usual method of numerical integration of coupled differential equations can be applied only in the Schrödinger formalism without spin-orbit deformation. It needs a lot of computing time and involves all the solutions whereas we need only one or a few of them.

In the **ECIS method (Sequential Iteration for Coupled Equations)**, we suppose the equations ordered in decreasing order of coupling. The equations for the initial channel for which there is an incoming wave are in front.

The first one is labelled $f_0(r)$.

If there are more than one, the other solutions are obtained from a circular permutation of the equations.

With compound nucleus, all solutions are needed to compute the transmission coefficients and, if requested, to diagonalise the scattering matrix.

In its first presentation, there was the “differential” **ECIS** method which consist to solve single inhomogeneous second order differential equations and the “integral” **ECIS** method which has been generalised to Dirac formalism. They gave same results and they need same computational time and storage. The “differential” **ECIS** method have been suppressed because :

- there is no simple extension to the Dirac formalism,
- result is obtained as a difference of two numbers and can be quite wrong with high energy close channel.

D.1 0th order of iteration

We assume a multiplicative factor λ in front of the second members of the equations and do a power expansion with respect to λ and take the result for $\lambda = 1$.

For :

$$\left\{ -\frac{\hbar^2}{2m_i} \left[\frac{d^2}{dr^2} - \frac{2\eta_i k_i}{r} - \frac{l_i(l_i + 1)}{r^2} + k_i^2 \right] - V_i^{opt}(r) + V_{ii}(r) \right\} f_i(r) = \lambda W_i(r) \quad (\text{III.110})$$

the solution is :

$$f_i(r) = f_i^{reg}(r) \delta_{i,0} + \lambda \int_0^\infty \mathcal{G}_i(\nabla, \nabla') \mathcal{W}_i(\nabla') [\nabla' \quad (\text{III.111})$$

with :

$$\mathcal{G}_i(\nabla, \nabla') = \frac{\epsilon_{\nabla}^{\nabla}}{\epsilon_{\nabla}^{\nabla}} \{ \nabla \} \{ \nabla \} (\nabla <) \{ \nabla \} (\nabla >) \quad (\text{III.112})$$

the **0th order of iteration** is obtained with $\lambda = 0$. Therefore, it is :

$$f_0^{(0)}(r) = f_0^{reg}(r) \quad C_0^{(0)} = C_0^{opt} \quad f_i^{(0)}(r) = 0 \quad \overline{C}_i^{(0)} = 0 \quad (\text{III.113})$$

D.2 1st order iteration:**For $i = 1$**

$$\begin{aligned}
f_1^{(1)}(r) &= - \int_0^\infty \mathcal{G}_\infty(\nabla, \nabla') \mathcal{V}_{\infty'} \{^{(1)}(\nabla')\} [\nabla'] \\
\overline{C}_1^{(1)} &= - \frac{2m_i}{\hbar^2 k_i} \int_0^\infty f_1^{reg}(r') V_{10}(r') f_0^{(0)}(r') dr'
\end{aligned} \tag{III.114}$$

which gives the **DWBA** result.**For $i = 2$**

$$\begin{aligned}
f_2^{(1)}(r) &= - \int_0^\infty \mathcal{G}_\infty(\nabla, \nabla') \left[\mathcal{V}_{\infty'} \{^{(1)}(\nabla')\} + \mathcal{V}_{\infty} \{^{(\infty)}(\nabla')\} \right] [\nabla'] \\
\overline{C}_2^{(1)} &= - \frac{2m_i}{\hbar^2 k_i} \int_0^\infty f_2^{reg}(r') \left[V_{20}(r') f_0^{(0)}(r') + V_{21}(r') f_1^{(1)}(r') \right] dr'
\end{aligned} \tag{III.115}$$

which takes into account the result already obtained for $f_1^{(1)}(r)$.**and the same** up to the last equation.**After the last equation** go back to $i = 0$:

$$\begin{aligned}
f_0^{(1)}(r) &= f_0^{reg}(r) - \int_0^\infty \mathcal{G}_r(\nabla, \nabla') \left[\mathcal{V}_{r'} \{^{(1)}(\nabla')\} + \sum_{\downarrow} \mathcal{V}_{\eta}(\nabla') \{^{(\infty)}(\nabla')\} \right] [\nabla'] \\
C_0^{(1)} &= C_0^{opt} - \frac{2m_0}{\hbar^2 k_0} \int_0^\infty f_0^{reg}(r') \left[V_{00}(r') f_0^{(0)}(r') + \sum_i V_{0i}(r') f_i^{(1)}(r') \right] dr'
\end{aligned} \tag{III.116}$$

The result is obtained if $C_0^{(1)}$ and all the $\overline{C}_i^{(1)}$ are less than a given ϵ .**D.3 nth order iteration****For $1 \leq i \leq N$**

$$\begin{aligned}
f_i^{(n)}(r) &= - \int_0^\infty \mathcal{G}_i(\nabla, \nabla') \left[\mathcal{V}_{i'} \{^{(\infty)}(\nabla')\} + \sum_{|=\infty}^{\downarrow-\infty} \mathcal{V}_{|l}(\nabla') \{^{(\infty)}(\nabla')\} + \sum_{|=\downarrow}^{\mathcal{N}} \mathcal{V}_{|l}(\nabla') \{^{(\infty)}(\nabla')\} \right] [\nabla'] \\
\overline{C}_i^{(n)} &= - \frac{2m_i}{\hbar^2 k_i} \int_0^\infty f_i^{reg}(r') \left[V_{i0}(r') f_0^{(n-1)}(r') + \sum_{j=1}^{i-1} V_{ij}(r') f_j^{(n)}(r') \right. \\
&\quad \left. + \sum_{j=i}^N V_{ij}(r') f_j^{(n-1)}(r') \right] dr'
\end{aligned} \tag{III.117}$$

and for $i=0$

$$\begin{aligned}
f_0^{(n)}(r) &= f_0^{reg}(r) - \int_0^\infty \mathcal{G}_r(\nabla, \nabla') \left[\mathcal{V}_{r'} \{^{(\infty)}(\nabla')\} + \sum_{|=\infty}^{\downarrow-\infty} \mathcal{V}_{\eta} \{^{(\infty)}(\nabla')\} \right] [\nabla'] \\
C_0^{(n)} &= C_0^{opt} - \frac{2m_0}{\hbar^2 k_0} \int_0^\infty f_0^{reg}(r') \left[V_{00}(r') f_0^{(n-1)}(r') + \sum_{j=1}^N V_{0j}(r') f_j^{(n)}(r') \right] dr'
\end{aligned} \tag{III.118}$$

The result is obtained if $|C_0^{(n)} - C_0^{(n-1)}|$ and all the $|\overline{C}_i^{(n)} - \overline{C}_i^{(n-1)}|$ are less than a given ϵ .

D.4 Convergence and padé approximation

This method of iteration can give divergent results. The convergence can be accelerated and the possibility of divergence avoided by Padé approximants.

From the fourth iteration, if convergence is not obtained for the current equation but was obtained for all the previous equations in this iteration, the results $C_i^{(m)}$ for $1 \leq m \leq n$ are considered as :

$$C_i^{(m)} = \sum_{j=1}^m a_i^j \lambda^j \quad \text{for} \quad \lambda = 1 \quad (\text{III.119})$$

even if each iteration do not correspond to a definite power of the λ introduced in the equation.

The polynomial of coefficients a_i^j is replaced by the ratio of two polynomials $P(\lambda)/Q(\lambda)$ with the same number of coefficients. This ratio is obtained as a continued fraction in such a way that a new $C_i^{(n)}$ adds only a new coefficient. If the continued fraction evaluated with all the coefficients and the continued fraction evaluated without the last coefficient differ by less than ϵ , their value is assumed to be the result.

The smallest or the few smallest zeroes of the denominator polynomial $Q(\lambda)$ are the same for all the equations. There are complex values. If the coupling of the equations is multiplied by one of these zeroes, the coefficient of the outgoing wave of all the equations blows up. For this value, the set of coupled equations has a solution which is purely outgoing in all the channel: in an exact power expansion with respect to the coupling (that is without the sequential procedure) it is a **Weinberg state**. If λ_j are the Weinberg eigenvalues and x_i^j the amplitudes of the outgoing waves with a proper normalisation :

$$C_i(\lambda) = \sum_j x_i^j \frac{1}{\lambda_j - \lambda} \quad C_i^{(n)} = \sum_j \frac{x_i^j}{\lambda_j} \sum_{m=1}^n \frac{1}{\lambda_j^m} \quad (\text{III.120})$$

If λ_j is small, each iteration gives the result of last one multiplied by $1/\lambda_j$. Iterations are stopped if these results become too large, or by a maximum number of iterations.

The minimum value of $|\lambda_j|$ is the convergence radius of the Taylor expansion. **In practice, Padé approximants give good results up to 6 times the radius of convergence** : that is for 20 iterations, results as large as 10^{15} . Indications on the actions of the code were given in section (I.A.2.b).

Chapter IV

Long range interactions

The description of this topic in Ref. [7] applies only to **ECIS79**. From the code **ECIS88**, this method allows to use non relativistic Coulomb functions as already stated in Ref. [7] which needs no generalisation for that. The Coulomb excitation in the Dirac formalism needed a drastic change of formulae. The use of such methods for the Coulomb spin-orbit interaction of chargeless particles needs a completely different calculation, even in Schrödinger formalism. Note that, in the Schrödinger formalism, the integral related to Coulomb spin-orbit interaction cannot be computed by this method when the sum of the two angular momenta is equal to the angular momentum transfer; the code neglect this correction.

A Long range contributions

For large values of r with deformed Coulomb interaction, the coupling between equations is :

$$\frac{2m}{\hbar^2} V_{if}(r) = \sum_{\lambda} \alpha_{if}^{\lambda} r^{-\lambda-1} \quad (\text{IV.1})$$

where α_{if}^{λ} is the geometrical coefficient G_{if}^{λ} multiplied by some constant.

A.1 Iterations in the Schrödinger formalism

In Equ. (III.117) and (III.118) describing the n th iteration of **ECIS**, for sufficiently large values of r , we can assume that :

$$\begin{aligned} f_i^{reg}(r) &\rightarrow F_{l_i}(\eta_i; k_i r) + C_i^{opt} [G_{l_i}(\eta_i; k_i r) + iF_{l_i}(\eta_i; k_i r)] \\ f_i^{irr}(r) &\rightarrow G_{l_i}(\eta_i; k_i r) + iF_{l_i}(\eta_i; k_i r) \\ f_i^{(n)}(r) &\rightarrow F_{l_i}(\eta_i; k_i r) \delta_{i0} + C_i^{(n)} [G_{l_i}(\eta_i; k_i r) + iF_{l_i}(\eta_i; k_i r)] \end{aligned} \quad (\text{IV.2})$$

The long range contribution can be expressed with the integrals :

$$M(H, K, \bar{k}R)_{l_i l_f}^{-\lambda-1} = \int_{\bar{k}R}^{\infty} \frac{1}{\rho^{\lambda+1}} H_{l_i}(\eta_i; \epsilon \rho) K_{l_f}(\eta_f; \epsilon^{-1} \rho) d\rho \quad (\text{IV.3})$$

with $\bar{k} = \sqrt{k_i k_f}$, $\epsilon = \sqrt{k_i/k_f}$ and H and K are the regular or the irregular Coulomb function or any combination of them. We shall also use :

$$V(H, K)_{i,j} = \sum_{\lambda} \alpha_{i,j}^{\lambda} \frac{\bar{k}^{\lambda}}{k_i} M(H, K, \bar{k}R)_{l_i l_f}^{-\lambda-1} \quad (\text{IV.4})$$

The equations (III.117) and (III.118) can be replaced by :

$$\begin{aligned}
f_i^{(n)}(r) &= f_i^{reg}(r)\delta_{i0} + \frac{1}{k_i} f_i^{reg}(r) \left\{ \int_r^R f_i^{irr}(r') \sum_j V_{ij}(r') f_j^{(m)}(r') dr' + A \right\} \\
&\quad + \frac{1}{k_i} f_i^{irr}(r) \left\{ \int_0^r f_i^{reg}(r') \sum_j V_{ij}(r') f_j^{(m)}(r') dr' \right\} \\
C_i^{(n)} &= C^{(0)}\delta_{i0} + \frac{1}{k_i} \left\{ \int_0^R f_i^{reg}(r') \sum_j V_{ij}(r') f_j^{(m)}(r') dr' + B \right\}
\end{aligned} \tag{IV.5}$$

where $m = n$ if $j < i$ and $m = n - 1$ if $j \geq i$, the equation 0 being the last one. The corrections A are :

$$A = k_i \sum_j V(H^{(+)}, F)_{i,j} \delta_{j0} + C_j^{(m)} V(H^{(+)}, H^{(+)})_{i,j} \tag{IV.6}$$

and the corrections B are :

$$\begin{aligned}
B &= k_i \sum_j [V(F, F)_{i,j} + C_i^{opt} V(H^{(+)}, F)_{i,j}] \delta_{j0} \\
&\quad + C_j^{(m)} [V(F, H^{(+)})_{i,j} + C_i^{opt} V(H^{(+)}, H^{(+)})]
\end{aligned} \tag{IV.7}$$

where $H_{l_i}^{(+)}(\eta_i; k_i r) = G_{l_i}(\eta_i; k_i r) + iF_{l_i}(\eta_i; k_i r)$ is the Coulomb outgoing function.

All the integrals involved are for natural parity, that is with $l_i + l_f + \lambda$ even.

A.2 Usual coupled equations

For n equations and with \mathcal{F} and \mathcal{G} replaced by corrected values, the system of Equ. (III.82) cannot be solved as a system of n linear equations similar to Equ. (III.83). The linear system of equation would be of dimension $2n$. However, one can obtain the results solving successively two systems of n linear equations.

The first step is to obtain a \overline{C} -matrix by matching with the “uncorrected” Coulomb functions. In this step the values of $\mathcal{F}_{l_j}(\eta_j; \mathcal{R} \pm \langle \rangle)$ and $\mathcal{G}_{l_j}(\eta_j; \mathcal{R} \pm \langle \rangle)$ are obtained taking into account the long range term (IV.1) in the three Numerov integration steps and in Equ. (III.81). The method to use is described by Equ. (III.83) to (III.85). This is equivalent to the matching with Coulomb function and its derivative at the point R .

In the second step, the Equ. (III.61) at ∞ is rewritten at the matching point R . The Coulomb functions at ∞ are replaced by the values which they get at the point R , taking into account the coupling given by Equ. (IV.1), that is for equation j :

$$\begin{aligned}
F_{l_i}(\eta_i; \infty) &\Rightarrow F_{l_j}(\eta_j; k_j R) \left\{ \delta_{ij} - M(H^{(+)}, F)_{l_j, l_i} F_{l_j}(\eta_j; k_j R) \right\} + V(F, F)_{l_j, l_i} H_{l_j}^{(+)}(\eta_j; k_j R) \\
H_{l_i}^{(+)}(\eta_i; \infty) &\Rightarrow H_{l_j}^{(+)}(\eta_j; k_j R) \left\{ \delta_{ij} + M(F, H^{(+)})_{l_j, l_i} H_{l_j}^{(+)}(\eta_j; k_j R) \right\} - V(H^{(+)}, H^{(+)})_{l_j, l_i} F_{l_j}(\eta_j; k_j R)
\end{aligned} \tag{IV.8}$$

In the result, replacing :

$$F_{l_i}(\eta_i; k_i R) \Rightarrow - \sum_j \overline{C}_j^i H^{(+)}(\eta_j; k_j R) \tag{IV.9}$$

one obtains the system of linear equations :

$$\sum_l \left(\delta_{j,l} + V(F, H^{(+)})_{j,l} + \sum_k \overline{C}_k^j V(H^{(+)}, H^{(+)})_{k,l} \right) C_i^l = -V(F, F)_{j,i} + \sum_k \overline{C}_k^j \left(\delta_{ki} - V(H^{(+)}, F)_{ki} \right) \tag{IV.10}$$

which is also of dimension n .

The integrals involved are the same as with iterations.

A.3 Coulomb spin-orbit interaction

For this interaction, instead of Equ. (IV.1), the coupling between two equations is :

$$\begin{aligned} \frac{2m}{\hbar^2} V_{if}(r) &= \sum_{\lambda} \alpha_{if}^{\lambda} [x_1 r^{-\lambda-3} + x_2 \frac{1}{r} \frac{d}{dr} (r^{-\lambda-1}) + x_3 r^{-\lambda-2} \frac{d}{dr}] \\ &= \sum_{\lambda} \alpha_{if}^{\lambda} [\{x_1 - (\lambda+1)x_2\} r^{-\lambda-3} + x_3 r^{-\lambda-2} \frac{d}{dr}] \end{aligned} \quad (\text{IV.11})$$

where x_1 , x_2 and x_3 are the three coefficients which appear in Equ. (III.49) and (III.53) in front of the three form factors. The coefficient x_3 implies a derivation of the wave function which is a pure Coulomb function. One can use one of the two formulae :

$$\left(1 + \frac{\eta^2}{l^2}\right)^{\frac{1}{2}} F_{l-1}(\eta; \rho) = \left(\frac{d}{d\rho} + \frac{l}{\rho} + \frac{\eta}{l}\right) F_l(\eta; \rho), \quad \left(1 + \frac{\eta^2}{l^2}\right)^{\frac{1}{2}} F_l(\eta; \rho) = \left(-\frac{d}{d\rho} + \frac{l}{\rho} + \frac{\eta}{l}\right) F_{l-1}(\eta; \rho), \quad (\text{IV.12})$$

to express the derivative with the Coulomb function itself and another Coulomb function with the angular momentum increased by one unit (it could be decreased). The result involves three terms :

1. with the same angular momentum and a factor r^{-1} which can be summed with the contributions of x_1 and x_2 ,
2. with the same angular momentum and without power of r ,
3. with angular momentum increased by one unit.

So, three integrals are needed for spin-orbit interaction, one with $r^{-\lambda-3}$ and two with $r^{\lambda-2}$.

The integrals involved are of natural parity and also of non natural parity, with $l_i + l_f + \lambda$ odd.

A.4 Dirac formalism

In this formalism, the multipoles of the scalar and the vector potentials act only between large components or small components whereas the multipoles of a tensor potential acts between large and small components.

For the Coulomb potential, which is a vector potential, the geometrical coefficient for large components is the same as the geometrical coefficient between small components. However, there are quite different for a magnetic multipole excitation (in particular, one or the other can vanish). The coefficients between large and small components of equations i and j are multiplied by the normalisation factors :

$$N_F = \frac{1}{2} \left[\left(1 + \frac{E_i}{m}\right) \left(1 + \frac{E_j}{m}\right) \right]^{\frac{1}{2}}, \quad N_G = \frac{1}{2} \left[\left(1 - \frac{E_i}{m}\right) \left(1 - \frac{E_j}{m}\right) \right]^{\frac{1}{2}}, \quad (\text{IV.13})$$

respectively. The two integrals are computed independently because they involve different quantum numbers.

The tensor potential involve no derivative of the wave function. The two integrals are computed independently, with normalisations similar to those of Equ. (IV.13).

All the integral involved are of natural parity except for relativistic corrections of uncoupled equations.

In the Dirac formalism and in the iteration method for the Schrödinger formalism, the integrals M of Equ. (IV.3) must be corrected for finite step errors. For a step h , taking into account the numerical integration described by Equ. (III.99) and (III.100), the result is too large : in a next step between R and $R + h$, described by Equ. (III.98), the contribution at R has been already taken into account. The value needed is :

$$\begin{aligned} \overline{M}(H, K, \overline{k}R)_{l_i l_f}^{-\lambda-1} &= \int_{\overline{k}R}^{\infty} \frac{1}{\rho^{\lambda+1}} H_{l_i}(\eta_i; \epsilon \rho) K_{l_f}(\eta_f; \epsilon^{-1} \rho) d\rho - \frac{h}{2} \frac{1}{R^{\lambda+1}} H_{l_i}(\eta_i; \epsilon R) K_{l_f}(\eta_f; \epsilon^{-1} R) \\ &\quad - \frac{h^2}{12} \frac{d}{dr} \left\{ \frac{1}{R^{\lambda+1}} H_{l_i}(\eta_i; \epsilon R) K_{l_f}(\eta_f; \epsilon^{-1} R) \right\} \end{aligned} \quad (\text{IV.14})$$

On the contrary, with usual coupled equations, no finite step correction is needed.

A.5 Asymptotic J -region

Using the integrals defined by Equ. (IV.3), we divided the r -space into :

- $r < R$, the coupled channels region in which a numerical integration of the coupled equations is needed,
- $r > R$, a region in which only **DWBA** results are used.

When the total angular momentum J of the system increases, the result of the numerical integrals become smaller and smaller. For some critical value J_m , it does no more contribute, but the integrals of Equ. (IV.3) are still important because their mean contribution comes from $r \gg R$; the computation must be continued up to some $J = J_M$. So, in the J -space, there are also two regions :

- $J \leq J_m$, the coupled channels region in which a numerical integration of the coupled equations is needed,
- $J_m < J < J_M$, an asymptotic region in which the optical solutions of the diagonal equations are pure regular Coulomb functions.

In the asymptotic region, the integrals can be extended to the origin, keeping the expression given in Equ. (IV.1) because the wave functions are very small in the interior region. The result is :

$$C_i^{(1)} = \sum_{\lambda} \alpha_{i0}^{\lambda} M(F, F, 0)_{i,0}^{-\lambda-1} \quad (\text{IV.15})$$

These integrals are real. This first order does not respect unitarity. However, the coupled channel problem can be formulated differently, in terms of the reactance matrix K , which is the coefficient of the irregular function (instead of the outgoing function) in the solution, with the normalised regular function in one channel only. The matrix C is obtained, solving a linear system of equation :

$$K_{if} = \sum_{\lambda} \alpha_{if}^{\lambda} M(F, F, 0)_{i,f}^{-\lambda-1}, \quad C = \frac{K}{1 - iK} \quad (\text{IV.16})$$

The value of J_M can be very large. Valuable results can be obtained with a smaller value of J_M , using factorisation of $(1 - \cos \theta)$ in the amplitudes. Details will be given in section (VII.C.3).

B Recurrence relations

In the Schrödinger formalism, the product of the wave number k_i and the Coulomb parameter η_i of a channel i is a constant $C_i = k_i \eta_i$ independent of the channel i . All the considerations of Ref. [7] are based on this properties; the same restrictions are present also in the previous studies of the integrals from 0 to infinity in : [64] BIEDENHARN, L. C., McHALE, J., L. and THALER, R., M., “*Quantum calculation of Coulomb excitation. I*”, Phys. Rev. **100** (1955) page 376 as well in : [65] ALDER, K., BOHR, A., HUUS, T., MOTTELSON, B. and WINTHER, A., “*Study of Nuclear Structure by Electromagnetic Excitation with Accelerated Ions*”, Rev. Mod. Phys. **28** (1956) page 432. In the Dirac formalism, the products $\eta_i k_i$ are proportional to the total energy and all the considerations of these references have to be extended. However, the approach is quite similar.

B.1 General recurrence

In the Dirac phenomenology, the product $\eta_i k_i$ is proportional to the total energy of the incident particle, which varies from level to level. Therefore, the formula (36) of Ref. [7], which was a generalisation to

integrals from R to ∞ with regular or irregular Coulomb functions of formula (2.B.64) of Ref. [65] (which was written for integrals from 0 to ∞ involving only regular Coulomb functions) has to be replaced by :

$$\begin{aligned}
& x_1 \epsilon \frac{|l_f + 1 + i\eta_f|}{l_f + 1} M(H, K, \bar{k}R)_{l_i, l_f + 1}^{-\lambda-1} + x_2 \epsilon^{-1} \frac{|l_i + i\eta_i|}{l_i} M(H, K, \bar{k}R)_{l_i - 1, l_f}^{-\lambda-1} \\
& - x_3 \epsilon^{-1} \frac{|l_i + 1 + i\eta_i|}{l_i + 1} M(H, K, \bar{k}R)_{l_i + 1, l_f}^{-\lambda-1} - x_4 \epsilon \frac{|l_f + i\eta_f|}{l_f} M(H, K, \bar{k}R)_{l_i, l_f - 1}^{-\lambda-1} \\
& - \left(\epsilon \frac{x_1 \eta_f}{l_f + 1} + \epsilon^{-1} \frac{x_2 \eta_i}{l_i} - \epsilon^{-1} \frac{x_3 \eta_i}{l_i + 1} - \epsilon \frac{x_4 \eta_f}{l_f} \right) M(H, K, \bar{k}R)_{l_i, l_f}^{-\lambda-1} \\
& - [x_1(l_f - 1) + x_2 l_i - x_3(l_i + 1) - x_4(l_f + \lambda + 1)] M(H, K, \bar{k}R)_{l_i, l_f}^{-\lambda-2} \\
& = (x_1 + x_4) \frac{1}{(\bar{k}R)^{\lambda+1}} H_{l_i}(\eta_i, \epsilon \bar{k}R) K_{l_f}(\eta_f, \epsilon^{-1} \bar{k}R) \\
& + (x_1 + x_2 + x_3 + x_4) \int_{\bar{k}R}^{\infty} \frac{1}{\rho^{\lambda+1}} \left[\frac{d}{d\rho} H_{l_i}(\eta_i, \epsilon \rho) \right] K_{l_f}(\eta_f, \epsilon^{-1} \rho) d\rho
\end{aligned} \tag{IV.17}$$

where ϵ has been given below Equ. (IV.3) and x_1, x_2, x_3 and x_4 are independent parameters. Independent recurrence relations can be obtained by using various values of the x 's such that :

$$x_1 + x_2 + x_3 + x_4 = 0 \tag{IV.18}$$

This formula can be obtained from the Equ. (IV.12).

Other recurrence relations can be obtained by combining those obtained from Equ. (IV.13) with peculiar choices of the x 's.

Of particular interest are the recurrence relations between integrals with the same values for λ and the difference $q = l_f - l_i$. There are :

- three terms relations for $\lambda = q = 0$
- four terms relations for $\lambda = |q| = 0$
- five terms relations for $\lambda \neq q$.

In the non relativistic case, for which $\eta_i k_i = \eta_f k_f$, with $\lambda = 1$ and $l_f = l_i + 1$, $x_1 = x_2 = 0$, the coefficient of the fifth term $M(H, K, \bar{k}R)_{L_i, l_f}^{-\lambda-1}$ vanishes, allowing to express the dipole integral in terms of two monopole integrals. As the monopole integrals fulfill a three terms recurrence relation in all cases, the dipole integrals with $l_f = l_i + 1$ fulfill also a three terms recurrence relation, but only in the non-relativistic case.

B.2 Stored recurrence

In practice, we are interested in integrals with $\lambda \leq 4$. The structure of recurrence relation is such that any of them can be expressed in terms of four other ones. The simplest approach is to express all of them through local recurrences in terms of integrals with some fixed value of λ and q . **Integrals with the chosen values for λ and q will be evaluated beforehand, also by recurrence, with the best precision possible.**

One cannot choose the monopole integrals $\lambda = q = 0$, nor the integrals for which $\lambda = |q|$, because their recurrence involves less than five terms. Therefore, all the four integrals needed cannot be of one of these kinds.

For the stored integrals, we choose those with $q = 0$ to obtain symmetric expressions for the local recurrences with respect to the exchange of l_i and l_f and $\lambda = 1$ for simplicity of the expressions. Furthermore, the stored integrals used for diagonal terms are the corrections from non relativistic to relativistic Coulomb functions (for that, they can be obtained with equations much simpler than those needed in the general case and described below).

The five term recurrence relation for $\lambda = 1, q = 0$ is not easy to handle directly. It can be replaced by an inhomogeneous three terms recurrence relation involving the monopole integrals with $\lambda = 0, q = 0$ as inhomogeneous terms. This is similar to what was obtained for non relativistic corrections in Equ. (37a) and (37b) in Ref. [7], but the inhomogeneous terms $\lambda = 1, q = 1$ used in the non relativistic case have to be replaced by the monopole integrals because they do not have a three terms recurrence relation in the relativistic case.

The monopole integrals are given by :

$$\begin{aligned}
& \frac{2l+2}{2l+1} |l + i\eta_i| |l + i\eta_f| M(H, K, \bar{k}R)_{l-1, l-1}^{-1} - \left\{ 2\eta_i\eta_f + l(l+1)(\epsilon^2 + \epsilon^{-2}) \right\} M(H, K, \bar{k}R)_{l, l}^{-1} \\
& + \frac{2l}{2l+1} |l+1 + i\eta_i| |l+1 + i\eta_f| M(H, K, \bar{k}R)_{l+1, l+1}^{-1} \\
& = -\frac{l+1}{\bar{k}R} \left\{ \frac{1}{k_i} |l + i\eta_f| H_l(\eta_i; k_i R) K_{l-1}(\eta_f; k_f R) + \frac{1}{k_f} |l + i\eta_i| H_{l-1}(\eta_i; k_i R) K_l(\eta_f; k_f R) \right. \\
& \left. - \frac{1}{(\bar{k}R)^2} [l(2l+1) + \eta_i k_f + \eta_f k_i] H_l(\eta_i; k_i R) K_l(\eta_f; k_f R) \right\} \quad (IV.19)
\end{aligned}$$

The recurrence which replace formula (37a) of Ref. [7] is :

$$\begin{aligned}
& \frac{2l-1}{2l+1} |l + i\eta_i| |l + i\eta_f| [(l+1)^2(k_i^2 - k_f^2) + (k_i^2\eta_i^2 - k_f^2\eta_f^2)] M(H, K, \bar{k}R)_{l-1, l-1}^{-2} \\
& - \left[k_i^4\eta_i^4 - k_f^4\eta_f^4 + \frac{1}{k_i k_f} \{ (k_i^4\eta_i^2 - k_f^4\eta_f^2)(2l^2 + 2l + 1) + (k_i^4 - k_f^4)(l+1)^2 l^2 \} \right] M(H, K, \bar{k}R)_{l, l}^{-2} \\
& + \frac{2l+3}{2l+1} |l+1 + i\eta_i| |l+1 + i\eta_f| [l^2(k_i^2 - k_f^2) + (k_i^2\eta_i^2 - k_f^2\eta_f^2)] M(H, K, \bar{k}R)_{l+1, l+1}^{-2} \\
& = -\frac{1}{l} (k_i^2 - k_f^2)(\eta_i k_f + \eta_f k_i) |l + i\eta_i| |l + i\eta_f| M(H, K, \bar{k}R)_{l-1, l-1}^{-1} - \frac{1}{2l} (k_i^2 - k_f^2) \\
& \left[\{ l^2 + \eta_f^2 \} \{ l\eta_f k_f - (l+2)\eta_i k_i \} \frac{k_f}{k_i} + \{ l^2 + \eta_i^2 \} \{ l\eta_i k_i - (l+2)\eta_f k_f \} \frac{k_i}{k_f} \right] M(H, K, \bar{k}R)_{l, l}^{-1} \\
& + \frac{1}{l^2 \bar{k}R} (k_i^2\eta_i^2 - k_f^2\eta_f^2) |l + i\eta_i| |l + i\eta_f| H_{l-1}(\eta_i; k_i R) K_{l-1}(\eta_f; k_f R) - \frac{k_i^2(l^2 + \eta_i^2) - k_f^2(l^2 + \eta_f^2)}{2l^2 (\bar{k}R)^3} \\
& \left\{ \eta_f |l + i\eta_f| [2l(l+1)^2 + (l+2)\eta_i k_i - l\eta_f k_f] H_l(\eta_i; k_i R) K_{l-1}(\eta_f; k_f R) \right. \\
& \left. + \eta_i |l + i\eta_i| [2l(l+1)^2 + (l+2)\eta_f k_f - l\eta_i k_i] H_{l-1}(\eta_i; k_i R) K_l(\eta_f; k_f R) \right\} \\
& + \left\{ \frac{k_i^2(l^2 + \eta_i^2) + k_f^2(l^2 + \eta_f^2)}{2l^2 (\bar{k}R)^3} [2l^2(l+1)^2(2l+1) + l(l+2)(2l+1)(\eta_i k_i + \eta_f k_f) \right. \\
& \left. + 2(l+1)(\bar{k}R)^2 \eta_i \eta_f] - \frac{1}{2l \bar{k}R} [(k_i^4\eta_i^4 - k_f^4\eta_f^4) + l(l+2)(k_i^2\eta_i^2 + k_f^2\eta_f^2)(k_i^2 - k_f^2) \right. \\
& \left. - 2 \left(\frac{k_i^4\eta_i^2 - k_f^4\eta_f^2}{k_i k_f} \right) l] \right\} H_l(\eta_i; k_i R) K_l(\eta_f; k_f R) \quad (IV.20)
\end{aligned}$$

In practice, we introduced :

$$C_X = k_i^2\eta_i^2 + k_f^2\eta_f^2 \quad D_X = k_i^2\eta_i^2 - k_f^2\eta_f^2 \quad D_Y = \frac{k_i^4\eta_i^2 - k_f^4\eta_f^2}{k_i k_f} \quad D_T = k_i^2 - k_f^2 \quad (IV.21)$$

The Equ. (IV.20) is homogeneous in D_T, D_X and D_Y , which can be multiplied by a common factor :

- if $D_X = 0, DT = 1$ and $DY = \eta_i\eta_f$,
- if $D_X \neq 0$ and if $DT \neq 0$, D_X and D_Y are divided by DT and DT is replaced by 1.

With these notations, Equ. (IV.20) is :

$$\begin{aligned}
& \frac{2l-1}{2l+1} |l + i\eta_i| |l + i\eta_f| [(l+1)^2 D_T + D_X] M(H, K, \bar{k}R)_{l-1, l-1}^{-2} \\
& - \left[D_X C_X + D_Y (2l^2 + 2l + 1) + D_T (l+1)^2 l^2 (\epsilon^2 + \epsilon^{-2}) \right] M(H, K, \bar{k}R)_{l, l}^{-2} \\
& + \frac{2l+3}{2l+1} |l+1 + i\eta_i| |l+1 + i\eta_f| [l^2 D_T + D_X] M(H, K, \bar{k}R)_{l+1, l+1}^{-2} \\
& = -\frac{1}{l} D_T (\eta_i k_f + \eta_f k_i) |l + i\eta_i| |l + i\eta_f| M(H, K, \bar{k}R)_{l-1, l-1}^{-1} \\
& - \frac{1}{2l} D_T \left[\{l^2 + \eta_f^2\} \{l\eta_f k_f - (l+2)\eta_i k_i\} \epsilon^{-2} + \{l^2 + \eta_i^2\} \{l\eta_i k_i - (l+2)\eta_f k_f\} \epsilon^2 \right] M(H, K, \bar{k}R)_{l, l}^{-1} \\
& + \frac{1}{l^2 \bar{k}R} D_X |l + i\eta_i| |l + i\eta_f| H_{l-1}(\eta_i; k_i R) K_{l-1}(\eta_f; k_f R) \\
& - \frac{1}{2l^2 (\bar{k}R)^3} (l^2 D_T + D_X) \left\{ \eta_f |l + i\eta_f| [2l(l+1)^2 + (l+2)\eta_i k_i - l\eta_f k_f] H_l(\eta_i; k_i R) K_{l-1}(\eta_f; k_f R) \right. \\
& \left. + \eta_i |l + i\eta_i| [2l(l+1)^2 + (l+2)\eta_f k_f - l\eta_i k_i] H_{l-1}(\eta_i; k_i R) K_l(\eta_f; k_f R) \right\} \\
& + \frac{1}{2l^2 (\bar{k}R)^3} \left\{ [l^2 D_T + D_X] [2l^2(l+1)^2(2l+1) + l(l+2)(2l+1)(\eta_i k_i + \eta_f k_f) + 2(l+1)(\bar{k}R)^2 \eta_i \eta_f] \right. \\
& \left. - l(\bar{k}R)^2 [C_X D_X + l(l+2)C_X D_T - 2lD_Y] \right\} H_l(\eta_i; k_i R) K_l(\eta_f; k_f R) \quad (IV.22)
\end{aligned}$$

The four integrals $M(H, K, \bar{k}R)_{0,0}^{-1}$, $M(H, K, \bar{k}R)_{1,1}^{-1}$, $M(H, K, \bar{k}R)_{0,0}^{-2}$ and $M(H, K, \bar{k}R)_{1,1}^{-2}$ are needed to start the recurrence. As in : [66] RAWITSCHER, G. H. and RASMUSSEN, C. H., “*Error analysis of code AROSA for quantal Coulomb excitation calculations*”, Comput. Phys. Commun. **11** (1976) page 183, they can be obtained by integration of the product of the asymptotic expansions from the matching point to ∞ . For the asymptotic expansion, see for example in : [67] ABRAMOVITH, M. and STEGUN, I. A., “*Handbook of Mathematical Functions*” Dover, New-York, (1972) page 541. For $l = 0$ and $\eta < 30$, this asymptotic expansion can be used only (see Ref. [19]) for $kR > 5\eta/3 + 7.5$; for larger values of η , a look on the asymptotic formula shows a quadratic behaviour of the lower radius; $kR > 0.06\eta^2$ seems to be a safe limit. To use a smaller matching point than the larger value R_m of $k_i R_m = 5\eta_i/3 + 7.5$, $k_f R_m = 5\eta_f/3 + 7.5$ and $\bar{k}R_m = .06\eta_i\eta_f$, the integral between R and R_m is computed by 40 points Gauss integrations, each Gauss integration being for $\Delta\bar{k}R < 20$. This allows results for any R whereas the method of Ref. [66] gives results for a R which increases with angular momentum.

This procedure is convenient for $M(G, G, \bar{k}R)_{l,l}^{-2}$, which increases quickly with l . For $M(G, F, \bar{k}R)_{l,l}^{-2}$ and $M(F, G, \bar{k}R)_{l,l}^{-2}$, which remain of the same order of magnitude when η_i and η_f are not very different. On the contrary, $M(F, F, \bar{k}R)_{l,l}^{-2}$, which decreases when l increases, is not given accurately. The downward recurrence, starting from zero values for the integrals and using the usual procedure of downward recurrence for the regular Coulomb function of the inhomogeneous terms, give a very accurate value of the integral from the origin to the matching point; then $M(F, F, \bar{k}R)_{l,l}^{-2}$ is obtained by taking the difference with $M(F, F, 0)_{l,l}^{-2}$, which is needed anyway in the asymptotic region.

B.3 Local recurrence

Due to the complexity of the expressions in the general case, when $\eta_i k_j \neq \eta_j k_i$, the expression of the integrals had to be changed with respect to what was described in Ref. [7]. Any integral is expressed as :

$$\begin{aligned}
M(H, K, \bar{k}R)_{l,l}^{-\lambda-1} &= \alpha_1 M(H, K, \bar{k}R)_{n-1, n-1}^{-2} + \alpha_2 M(H, K, \bar{k}R)_{n, n}^{-2} \\
&+ \alpha_3 M(H, K, \bar{k}R)_{n+1, n+1}^{-2} + \alpha_4 M(H, K, \bar{k}R)_{n+2, n+2}^{-2}
\end{aligned}$$

$$\begin{aligned}
& + \beta_1 H_n(\eta_i; k_i R) K_n(\eta_f; k_f R) + \beta_2 H_n(\eta_i; k_i R) K_{n+1}(\eta_f; k_f R) \\
& + \beta_3 H_{n+1}(\eta_i; k_i R) K_n(\eta_f; k_f R) + \beta_4 H_{n+1}(\eta_i; k_i R) K_{n+1}(\eta_f; k_f R) \quad (\text{IV.23})
\end{aligned}$$

with :

$$n = \text{Integer part of } \frac{1}{2}(l_i + l_f - \lambda + 3) \quad (\text{IV.24})$$

If l_i and l_j are decreased by the same amount down to the first values for which the integral from 0 to ∞ converges, the value of n is 1. The four coefficients α depend upon λ , $q = l_f - l_i$, n , the Coulomb parameters and the wave numbers; the coefficients β depend upon the same parameters and also upon the radius R .

Using $\bar{\eta} = \sqrt{\eta_i \eta_f}$, $\epsilon' = \sqrt{\eta_f / \eta_i}$ ($\epsilon = \epsilon'$ in the non relativistic case) Equ. (IV.23) can be written :

$$\begin{aligned}
& \int_R^\infty \frac{1}{r^\lambda} H_{l_i}(\epsilon'^{-1} \bar{\eta}; \epsilon r) K_{l_f}(\epsilon' \bar{\eta}; \epsilon^{-1} r) dr \\
& = \sum_{i=1}^4 \alpha_i \int_R^\infty \frac{1}{r^2} H_{n+i-2}(\epsilon'^{-1} \bar{\eta}; \epsilon r) K_{n+i-2}(\epsilon' \bar{\eta}; \epsilon^{-1} \epsilon^{-1} r) dr + \sum_{i=1}^4 \beta_i(R) P_i(R) \quad (\text{IV.25})
\end{aligned}$$

with :

$$\begin{aligned}
P_1(R) &= H_n(\epsilon'^{-1} \bar{\eta}; \epsilon r) K_n(\epsilon^{-1} \bar{\eta}; \epsilon' r) \\
P_2(R) &= H_n(\epsilon'^{-1} \bar{\eta}; \epsilon r) K_{n+1}(\epsilon^{-1} \bar{\eta}; \epsilon' r) \\
P_3(R) &= H_{n+1}(\epsilon'^{-1} \bar{\eta}; \epsilon r) K_n(\epsilon^{-1} \bar{\eta}; \epsilon' r) \\
P_4(R) &= H_{n+1}(\epsilon'^{-1} \bar{\eta}; \epsilon r) K_{n+1}(\epsilon^{-1} \bar{\eta}; \epsilon' r) \quad (\text{IV.26})
\end{aligned}$$

Using the recurrence relation of the Coulomb functions :

$$(2l+1) \left[\eta + \frac{l(l+1)}{\rho} \right] F_l(\eta; \rho) = l|l+1+i\eta| F_{l+1}(\eta; \rho) + (l+1)|l+i\eta| F_{l-1}(\eta; \rho) \quad (\text{IV.27})$$

to express all the H with H_n and H_{n+1} and all the K with K_n and K_{n+1} (the power of ρ^{-1} always increases), we obtain :

$$\int_R^\infty \sum_{i=1}^4 C_i(r) P_i(r) dr = \sum_{i=1}^4 \beta_i(R) P_i(R) \quad (\text{IV.28})$$

where the $C_i(r)$ are known polynomials of r^{-1} and the $\beta_i(R)$ unknown polynomials in R^{-1} , that is :

$$C_i(r) = \sum_{n=2}^N C_i^{(n)} r^{-n}, \quad \beta_i(R) = \sum_{n=1}^{N-1} \beta_i^{(n)} R^{-n} \quad (\text{IV.29})$$

Differentiating with respect to R , we get :

$$\sum_{n=2}^N C_i^{(n)} R^{-n} = \sum_{n=1}^{N-1} \left[P_i(n) \beta_i^{(n)} R^{-n-1} + \sum_{j=1}^4 Q_{i,j} \beta_j^{(n)} R^{-n} \right] \quad (\text{IV.30})$$

with the diagonal matrix :

$$P(n) = \begin{vmatrix} n-2l-2 & 0 & 0 & 0 \\ 0 & n & 0 & 0 \\ 0 & 0 & n & 0 \\ 0 & 0 & 0 & n+2l+2 \end{vmatrix}, \quad (\text{IV.31})$$

and with the diagonal matrix :

$$Q = \frac{1}{2l+1} \begin{vmatrix} -\eta_i \epsilon - \eta_f \epsilon^{-1} & -|l+1+i\eta_f|\epsilon^{-1} & -|l+1+i\eta_i|\epsilon & 0 \\ |l+1+\eta_f|\epsilon^{-1} & \eta_f \epsilon^{-1} - \eta_i \epsilon & 0 & -|l+1+i\eta_i|\epsilon \\ |l+1+i\eta_i|\epsilon & 0 & \eta_i \epsilon - \eta_f \epsilon^{-1} & -|l+1+\eta_f|\epsilon^{-1} \\ 0 & |l+1+i\eta_i|\epsilon & |l+1+\eta_f|\epsilon^{-1} & \eta_i \epsilon + \eta_f \epsilon^{-1} \end{vmatrix}, \quad (\text{IV.32})$$

If the highest power of r^{-1} in the $C_i(r)$ is m , Equ. (IV.30) allows to obtain first the coefficients of R^{-m+1} in the β_i and, then, those of all the lower degree. As :

C_1 includes $[(\dots/r^2 + \dots/r + \dots)\alpha_1 + \alpha_2 + \dots\alpha_4]/r^2$

C_2 includes $[(\dots/r + \dots)\alpha_1 + (\dots/r + \dots)\alpha_4]/r^2$

C_3 includes $[(\dots/r + \dots)\alpha_1 + (\dots/r + \dots)\alpha_4]/r^2$

C_4 includes $[\dots\alpha_1 + \alpha_3 + (\dots/r^2 + \dots/r + \dots)\alpha_4]/r^2$

the coefficients $\beta_i^{(n)}$ for $n > 3$ do not depends on the α 's. The Equ. (IV.30) written for $n = 2$ and zero values for the $\beta_i^{(1)}$ is a linear system of equations giving the α 's. The coefficients $\beta_i^{(2)}$, which are in this equation, involve a specific dependence of l_i, l_j and λ and a very simple dependence upon α_1 and α_4 .

The dipoles integral are slightly different : in particular, when $\eta_i = \eta_j$ and $k_i = k_j$, a different formula has to be used to avoid a division by zero.

All the expressions are simpler in the non-relativistic case, when $X = \eta_i k_i - \eta_j k_j$ vanishes. They have been separated into a “non-relativistic” part, which is their value for $X = 0$, and a “relativistic” part, in which X can be factorised.

The linear system of equations expressing the α as functions of the $\beta_i^{(2)}$ can be inverted. The coefficients of the inverse system are quite complex when X do not vanish. The coefficients α_1 and α_4 depends only on $\beta_1^{(2)}$ and $\beta_4^{(2)}$ respectively, the coefficients α_2 and α_3 depends on all the $\beta^{(2)}$. A close expression of the coefficients α can be obtained only when some β vanish, that is for $l_j = l_i + 1$, $\lambda = 1$ and $l_i = l_j$, $\lambda = 2$. The subroutine **CORA** stores the values of the $\beta^{(2)}$ for the cases of interest and compute the coefficients α . It do not seem that this increase of intermediate computations decrease too much the precision of the results.

All these coefficients have been obtained up to $\lambda = 5$ for even parity integrals ($l_i + l_j + \lambda$ even) as well as odd parity integrals ($l_i + l_j + \lambda$ odd). For $\lambda = 6$, only the even parity integrals with $X = 0$ have been obtained except for $|l_i - l_j| = 6$. These last results allows the computation of $\lambda = 4$ for the Coulomb spin-orbit in the Schrödinger formalism. However, these results involve around 1100 cards and their use in **ECIS** is not absolutely needed. In **ECIS88** there are only the complete expressions for natural and unnatural parity up to $\lambda = 3$ and the expressions of natural parity integrals for $\lambda = 4$ with $X = 0$; this insure the use of Coulomb corrections up to $\lambda = 4$ for the central potential in any case, up to $\lambda = 2$ for the Coulomb spin-orbit in Schrödinger phenomenology and $\lambda = 3$ in Dirac phenomenology. In **ECIS94**, all the results obtained have been introduced, but the difference with **ECIS88** is on **COMMENT** cards which can easily been activated. In this domain, it is not possible to derive the expressions “by hand”. They were obtained using symbolic calculations on computer with **AMP** (“Algebraic Manipulation Program”) as described in : [68] DROUFFE, J.-M., “*AMP Language Reference Manual, Version 6*”, Note CEA-n-2297, (1982).

If $k_i = k_j$ and $\eta_i = \eta_j = \eta$, the matrix Q given by Equ. (IV.32) is singular of rank 2. If Equ. (IV.30) is used to compute some $\beta^{(1)}$, only two of them are independent because the coefficients of r^{-1} are :

$$\begin{aligned} 2\eta\beta_1^{(1)} + |l+1+i\eta|[\beta_2^{(1)} + \beta_3^{(1)}] &= 0 \\ \beta_1^{(1)} - \beta_4^{(1)} &= 0 \\ \beta_1^{(1)} - \beta_4^{(1)} &= 0 \\ 2\eta\beta_4^{(1)} + |l+1+i\eta|[\beta_2^{(1)} + \beta_3^{(1)}] &= 0 \end{aligned} \quad (\text{IV.33})$$

that is : $\beta_1^{(1)} = \beta_4^{(1)}$ and $\beta = \beta_2^{(1)} + \beta_3^{(1)}$ simple function of $\beta_1^{(1)}$. Consequently, Equ. (IV.23) holds with $\alpha_1 = \alpha_4 = 0$ and the first term of the polynomial expansion of the $\beta_i(r)$ is r^{-1} . In fact, the five terms recurrences reduce to three terms recurrence. This behaviour was used in ECIS79 but has been dropped since ECIS88.

B.4 Integrals from 0 to ∞

The integrals from 0 to ∞ are quite similar to those published in Ref. [65]: in general, the variables of the generalised hypergeometric functions of two variables depend upon the wave numbers and their parameters depend upon the Coulomb parameters. The formula (IIB.62), of Ref. [65], valid if $|l_i - l_f|$ is not equal to λ , becomes when $k_i > k_f$:

$$\begin{aligned}
M(F, F, 0)_{l_i, l_f}^{-\lambda-1} &= \pi \frac{k_f^{l_f}}{k_i^{l_i}} (2k_i)^{\lambda-2} \frac{\exp \pi \frac{\xi}{2}}{\sinh \pi \xi} (iX)^{\lambda+l_i-l_f-1} \frac{|\Gamma(l_f+1+i\eta_f)|}{|\Gamma(l_i+1+i\eta_i)|} \frac{\Gamma(2l_i+1)}{\Gamma(2l_f+1)\Gamma(l_i-l_f+1)} \\
&\times F_2(l_f-l_i+\lambda-1, l_f+1+i\eta_f, -l_i-i\eta_i, 2l_f+2, -2l_i; \frac{k_i}{k_f}X, -X) \\
&- \pi \frac{k_i^{l_i}}{k_f^{l_f}} (2k_f)^{\lambda-2} \frac{\exp -\pi \frac{\xi}{2}}{\sinh \pi \xi} (i\frac{k_i}{k_f}X)^{\lambda+l_f-l_i-1} \frac{|\Gamma(l_i+1+i\eta_i)|}{|\Gamma(l_f+1+i\eta_f)|} \frac{\Gamma(2l_f+1)}{\Gamma(2l_i+1)\Gamma(l_f-l_i+1)} \\
&\times F_2(l_i-l_f+\lambda-1, l_i+1+i\eta_i, -l_f-i\eta_f, 2l_i+2, -2l_f; -X, \frac{k_i}{k_f}X) \\
&+ \pi \frac{(k_i-k_f)^\lambda}{2k_i k_f} \frac{\exp -\pi \frac{\xi}{2}}{\sinh \pi \xi} \frac{|\Gamma(l_i+1+i\eta_i)|}{|\Gamma(l_f+1+i\eta_f)|} \\
&\times \Re \left\{ \frac{\Gamma(l_f+1+i\eta_f)}{\Gamma(\lambda+1+i\xi)\Gamma(l_i+1+i\eta_i)} i^{l_f-l_i-\lambda-1} (-X)^{i\eta_i} \left(-\frac{k_i}{k_f}X\right)^{-i\eta_f} \right. \\
&\left. \times F_3(-l_i+i\eta_i, -l_f-i\eta_f, l_i+1+i\eta_i, l_f+1-i\eta_f, \lambda+1+i\xi; X, \frac{k_i}{k_f}X) \right\} \quad (IV.34)
\end{aligned}$$

where :

$$\xi = \eta_i - \eta_f \quad X = \frac{k_i - k_f}{2k_f} \quad (IV.35)$$

Similarly, the formula (IIB.58) of Ref. [65] valid for $l_f = l_i + \lambda$, becomes :

$$\begin{aligned}
M(F, F, 0)_{l_i, l_i+\lambda}^{-\lambda-2} &= \exp \pi \frac{\xi}{2} \frac{|\Gamma(l_i+1+i\eta_i)|}{|\Gamma(l_f+1+i\eta_f)|} \left(\frac{k_i}{k_f}\right)^{l_i} (2k_f)^{\lambda-2} \left\{ \frac{|\Gamma(\lambda+i\xi)|^2}{\Gamma(2\lambda)} \right. \\
&\times F_2(-2\lambda+1, l_i+1-i\eta_i, l_i+1+i\eta_i, -\lambda+1+i\xi, -\lambda+1-i\xi; X, X) \\
&+ \Re \left[2 S X^{\lambda+i\xi} \frac{\Gamma(l_f+1-i\eta_f)\Gamma(-\lambda-i\xi)}{\Gamma(l_i+1+i\eta_i)} \right. \\
&\left. \left. \times F_2(-\lambda+1+i\xi, l_f+1-i\eta_f, l_i+1-i\eta_i, \lambda+1+i\xi, -\lambda+1-i\xi; X, X) \right] \right\} \quad (IV.36)
\end{aligned}$$

where :

$$\begin{aligned}
S &= -1, & \text{if } (\eta_i - \eta_f)(k_f - k_i) < 0 \quad \text{and } \lambda \text{ is odd;} \\
S &= 1, & \text{otherwise}
\end{aligned} \quad (IV.37)$$

These formulae holds even if $k_i = k_f$, that is $X = 0$: in the first formula, Equ. (IV.34), only the highest degree terms of the two generalised hypergeometric functions of two variables F_2 remain, all the others vanish; the generalised hypergeometric function F_3 of the first formula and one of the two F_2 of the Equ. (IV.25) reduce to unity; the second F_2 of the Equ. (IV.36) disappears. If $\eta_i = \eta_f$, careful limits of the expressions have to be taken, with derivative of gamma function inserted in each term of the generalised hypergeometric series. The special case $\eta_i = \eta_f = 0$ will be discussed in next section.

The monopole integrals are given by a different formula, corresponding to formula (IIB.56) of Ref. [65] :

$$M(F, F, 0)_{l,l}^{-1} = \left[\frac{4k_i k_f}{k_i + k_f} \right]^{l+1} \frac{1}{\sinh \pi \xi |\Gamma(l+1+i\eta_i)| |\Gamma(l+1+i\eta_f)|}$$

$$\Re \left\{ \left(\frac{k_f - k_i}{k_i + k_f} \right)^{i\xi} \frac{\Gamma(l+1+i\eta_i) \Gamma(l+1-i\eta_f)}{\Gamma(1+i\xi)} {}_2F_1(l+1-i\eta_f, l+1+i\eta_i, 1+i\xi; \left[\frac{k_f - k_i}{k_i + k_f} \right]^2) \right\}$$
(IV.38)

This formula do not hold if $k_i = k_f$ and has to be modified if $\eta_i = \eta_f$ but not zero; for $\eta_i = \eta_f = 0$, see below. All these formulae and their limits have been used to check the Coulomb integrals.

Equ. (IV.38) and (IV.34) are used to obtain the integrals for $l = 0$, $l = 1$ and $\lambda = 0$, $\lambda = 1$. All the other are obtained with the recurrence equations Equ. (IV.19) and (IV.20) (without the products of Coulomb functions). In fact, as downward recurrence is used to obtain stored integrals, only the integrals for $l = 0$ are necessary. Any integral is obtain by Equ. (IV.23), without the β 's.

If $\eta_i = \eta_f$ and $k_i = k_f$, the stored integrals are :

$$M(F, F, 0)_{l,l}^{-2} = \frac{1}{4l+2} \left\{ \pi - \pi \frac{\cosh \pi \eta}{\sinh \pi \eta} - \frac{1}{\eta} + \sum_{n=0}^l \frac{2\eta}{n^2 + \eta^2} \right\}$$
(IV.39)

which is used instead of the recurrence relation.

B.5 Coulomb corrections for chargeless particles

If there is no charge, the recurrence relations do not relate natural parity and unnatural parity integral, that is those for which $l_i + l_j + \lambda$ is even with those for which $l_i + l_j + \lambda$ is odd. The Equ. (IV.17) becomes :

$$x_1 \epsilon M(H, K, \bar{k}R)_{l_i, l_f+1}^{-\lambda-1} + x_2 \epsilon^{-1} M(H, K, \bar{k}R)_{l_i-1, l_f}^{-\lambda-1} - x_3 \epsilon^{-1} M(H, K, \bar{k}R)_{l_i+1, l_f}^{-\lambda-1}$$

$$- x_4 \epsilon M(H, K, \bar{k}R)_{l_i, l_f-1}^{-\lambda-1} - [x_1(l_f - 1) + x_2 l_i - x_3(l_i + 1) - x_4(l_f + \lambda + 1)] M(H, K, \bar{k}R)_{l_i, l_f}^{-\lambda-2}$$

$$= (x_1 + x_4) \frac{1}{(\bar{k}R)^{\lambda+1}} H_{l_i}(0, \epsilon \bar{k}R) K_{l_f}(0, \epsilon^{-1} \bar{k}R)$$

$$+ (x_1 + x_2 + x_3 + x_4) \int_{\bar{k}R}^{\infty} \frac{1}{\rho^{\lambda+1}} \left[\frac{d}{d\rho} H_{l_i}(0, \epsilon \rho) \right] K_{l_f}(0, \epsilon^{-1} \rho) d\rho$$
(IV.40)

The integrals needed are all of natural parity. Therefore, it is no more possible to compute them from stored integrals with $l_i = l_f$ and $\lambda = 1$. They have to be computed from stored integrals with $l_i = l_f$ and $\lambda = 2$. The recurrence relation is then :

$$2(l-1)M(H, K, \bar{k}R)_{l-1, l-1}^{-3} - (2l+1)(\epsilon^2 + \epsilon^{-2})M(H, K, \bar{k}R)_{l, l}^{-3} + 2(l+2)M(H, K, \bar{k}R)_{l+1, l+1}^{-3}$$

$$= \frac{1}{(\bar{k}R)^2} \left[H_{l-1}(k_i R) K_{l-1}(k_f R) - H_{l+1}(k_i R) K_{l-1}(k_f R) \right]$$
(IV.41)

The integral from 0 to ∞ , when $k_i > k_f$, is given by :

$$M(H, K, \bar{k}R)_{l_i, l_f}^{-\lambda-1} = \frac{\pi}{2^{\lambda+1}} \frac{k_f^{l_f}}{k_i^{l_f-\lambda+2}} \frac{\Gamma(\frac{1}{2}[l_i + l_f - \lambda + 2]) \Gamma(\lambda)}{\Gamma(\frac{1}{2}[l_i + l_f - \lambda + 1]) \Gamma(\frac{1}{2}[l_i - l_f + \lambda + 1]) \Gamma(\frac{1}{2}[l_i + l_f + \lambda + 1])}$$

$$\times {}_2F_1\left[\frac{1}{2}(l_i + l_f - \lambda + 2), \frac{1}{2}(l_i - l_f - \lambda + 1); 1 - \lambda; 1 - \left(\frac{k_f}{k_i}\right)^2\right]$$
(IV.42)

This expression has to be used with $\lambda = 2$ and $l_i = l_f = 1$ to start the recurrence; a second value is not needed because the recurrence can be started by assuming the product of the angular momentum with the integral to be $\frac{1}{3}$ for $l_i = l_f = 0$. When $k_i = k_f$, the integrals are simply :

$$M(H, K, \bar{k}R) = \frac{1}{2l_i(l_i + 1)}.$$
(IV.43)

To express any integral with the stored integrals, the coefficients described above in the general case are such that $\alpha_1 = \alpha_4 = 0$. The dipole integral cannot be expressed if $k_i = k_f$, but these integrals are used only for the Coulomb spin-orbit and it is the quadrupole integral which is needed for a dipole excitation. All the coefficients up to $\lambda = 6$ have been obtained, but only those up to $\lambda = 4$ are in **ECIS88** and all are on **COMMENT** cards in **ECIS94**.

C Coulomb excitations

The Coulomb potential as given by Equ. (III.21) is proportional to the product of charges of the particle and the target. In fact, it is the monopole folding of a charge distribution with the interaction $1/|r_1 - r_2|$. This charge distribution can be :

- a sphere with constant density,
- a plain Woods-Saxon distribution,
- a Woods-Saxon distribution multiplies by $1 + cr^2$,

of which the volume integral is the product of charges.

C.1 Coulomb deformation and heavy-ion option

This excitation is defined by the reduced electric transition probability $B(E\lambda)$ which is related to the Coulomb deformation β_c by :

$$B(E\lambda) = \left(\frac{3}{4\pi} Z e \beta_c R_c'^{\lambda+1} \right)^2 \quad (\text{IV.44})$$

where Z is the charge of the excited nucleus (see for example in : [69] HODGSON, P., E., “*Nucleon Heavy-Ion Reactions*”, Clarendon Press Oxford (1978) page 268 or in : [70] SATCHLER, G., R., “*Direct Nuclear Reactions*”, “*International series of monographs on physics 68*”, Clarendon Press Oxford, Oxford University Press New-York (1983) page 629. This relation do not give the sign of the deformation. A transition distribution is obtained with the deformation length $\beta_c R_c'$ and the transition form factor is the result of multipolarity λ obtained in the folding with $1/|r_1 - r_2|$.

In the usual notations, $R_c' = R_c A^{1/3}$ where R_c is the “reduced Coulomb radius” and A is the mass of the target. In the “heavy-ion option”, for a target of mass A and a particle of mass a , the radius R' is given in terms of the reduced radius R by :

$$R' = R(A^{1/3} + a^{1/3}) \quad (\text{IV.45})$$

and, as it is the deformation length which matters, the deformations are replaced by :

$$\beta' = \frac{A^{1/3}}{A^{1/3} + a^{1/3}} \beta, \quad (\text{IV.46})$$

except for the Coulomb deformations (central and spin-orbit) which are replaced by :

$$\beta_c' = \left(\frac{A^{1/3}}{A^{1/3} + a^{1/3}} \right)^\lambda \beta_c, \quad (\text{IV.47})$$

for a multipole λ . However, if **LO(6)=.TRUE.**, which means that deformations lengths are read instead of deformations, only Coulomb deformations are changed with a power $\lambda - 1$ instead of λ in Equ. (IV.47).

The introduction of magnetic multipole Coulomb excitation followed a question of Mme I. LINK from Strasbourg and was subject of discussions with Dr M. MERMAZ at Saclay and J. KIENER at Orsay. For relations between electric and magnetic excitation, we follow Ref. [65].

C.2 Electric multipole excitation

An electric multipole transition $B(E\lambda)$ is defined as an angular momentum transfer $\Delta L = \lambda$, a spin transfer $2\Delta S = 0$ and J transfer $2\Delta J = 2\lambda$ but can as well being defined with the same angular transfer, a spin transfer $2\Delta S = 2\lambda$ and a J transfer $2\Delta J = 0$. In Ref. [65], formulae (II B.32) and (II B.34) gives for an electric excitation :

$$d\sigma_{E\lambda} = \left(\frac{Z_1 e}{\hbar v_i} \right)^2 \frac{4k_i k_f}{(2\lambda + 1)^3} B(E\lambda) \sum_{\mu} | \langle \vec{k}_f | r_p^{-\lambda-1} Y_{\lambda\mu}(\theta_p, \phi_p) | \vec{k}_i \rangle |^2 d\Omega, \quad (\text{IV.48})$$

in the exterior region. This interaction needs geometrical factors given by Equ. (III.45) and (III.44).

C.3 Magnetic multipole excitation

A magnetic multipole transition $B(M\lambda)$ is defined as an angular momentum transfer $\Delta L = \lambda + 1$, a spin transfer $2\Delta S = 0$ and J transfer $2\Delta J = 2\lambda$. It is defined in the code by $2\Delta S = -2$ but $2\Delta S = -1$ in the input. The use of $\Delta L = \lambda + 1$ is necessary to go through the parity verifications.

In Ref. [65], formulae (II B.33) and (II B.35) gives for a magnetic excitation :

$$d\sigma_{M\lambda} = \left(\frac{Z_1 e}{\hbar c} \right)^2 \frac{v_f}{v_i} \frac{4}{\lambda^2 (2\lambda + 1)^3} B(M\lambda) \sum_{\mu} | \langle \vec{k}_f | \vec{l}_p \cdot \vec{\nabla}_p [r_p^{-\lambda-1} Y_{\lambda\mu}(\theta_p, \phi_p)] | \vec{k}_i \rangle |^2 d\Omega. \quad (\text{IV.49})$$

in the exterior region. The ratio of a magnetic excitation $\lambda - 1$ to an electric excitation λ given by Equ. (IV.47) is :

$$\frac{d\sigma_{M\lambda-1}}{d\sigma_{E\lambda}} = \frac{v_i v_f}{k_i k_f c^2} \frac{(2\lambda + 1)^3}{(\lambda - 1)^2 (2\lambda - 1)^3} \frac{B(M\lambda - 1)}{B(E\lambda)} \frac{\sum_{\mu} | \langle \vec{k}_f | \vec{l}_p \cdot \vec{\nabla}_p (r_p^{-\lambda} Y_{\lambda-1\mu}(\theta_p, \phi_p)) | \vec{k}_i \rangle |^2 d\Omega}{\sum_{\mu} | \langle \vec{k}_f | r_p^{-\lambda-1} Y_{\lambda\mu}(\theta_p, \phi_p) | \vec{k}_i \rangle |^2 d\Omega}, \quad (\text{IV.50})$$

that is the product of a coefficient of the potentials R_p^2 such that :

$$R_p = \frac{\hbar}{mc} \left(\frac{(2\lambda + 1)^3}{(\lambda - 1)^2 (2\lambda - 1)^3} \right)^{\frac{1}{2}} \left(\frac{B(M\lambda - 1)}{B(E\lambda)} \right)^{\frac{1}{2}} \quad (\text{IV.51})$$

obtained using :

$$\frac{v_i}{k_i} = \frac{v_f}{k_f} = \frac{\hbar}{m} \quad (\text{IV.52})$$

and a ratio R_g^2 such that :

$$R_g^2 = \frac{\sum_{\mu} | \langle \vec{k}_f | \vec{l}_p \cdot \vec{\nabla}_p (r_p^{-\lambda} Y_{\lambda-1\mu}(\theta_p, \phi_p)) | \vec{k}_p \rangle |^2 d\Omega}{\sum_{\mu} | \langle \vec{k}_f | r_p^{-\lambda-1} Y_{\lambda\mu}(\theta_p, \phi_p) | \vec{k}_p \rangle |^2 d\Omega} \quad (\text{IV.53})$$

The coefficient R_p is the product of three terms :

$$R_p = R_{p_1} R_{p_2} R_{p_3}, \quad R_{p_1} = \left(\frac{B(M\lambda - 1)}{B(E\lambda)} \right)^{\frac{1}{2}}, \quad R_{p_2} = \frac{\hbar}{mc}, \quad R_{p_3} = \left(\frac{(2\lambda + 1)^3}{(\lambda - 1)^2 (2\lambda - 1)^3} \right)^{\frac{1}{2}} \quad (\text{IV.54})$$

In practice :

- the factor R_{p_1} will be taken into account in the reduced nuclear matrix element multiplying by it the reduced nuclear matrix element of the electric transition to obtain the reduced nuclear matrix element of the magnetic transition,
- the factor R_{p_2} will be taken into account by multiplying by it a normalised Coulomb transition form factor of an electric excitation to obtain the magnetic transition form factor,

- the factor R_{p_3} will be factorised with R_g for convenience.

For the coefficient R_g , the gradient acting on $r^{-\lambda}Y_{\lambda\mu}(\hat{r})$ generates only terms in $r^{-\lambda-1}Y_{\lambda+1,\mu}(\hat{r})$. So, in the exterior region, the form factors of a magnetic excitation $B(M\lambda - 1)$ is the same as the form factor of an electric excitation $B(E\lambda)$. As the interior region has no importance in this case, we can use for the magnetic transition form factor a Coulomb electric transition form factor multiplied by the ratio R_{p_2} given in Equ. (IV.54). Between partial waves $|l_i >$ and $|l_f >$, the ratio of a magnetic multipole excitation $B(M\lambda - 1)$ to an electric multipole excitation $B(E\lambda)$ is :

$$\begin{aligned}
& \frac{\langle i^{l_f} Y_{l_f m_f}(\theta, \phi) | i^{\lambda} \vec{l} \cdot \vec{\nabla} (r^{-\lambda} Y_{\lambda-1\mu}(\theta, \phi)) | i^{l_i} Y_{l_i m_i}(\theta, \phi) \rangle}{\langle i^{l_f} Y_{l_f m_f}(\theta, \phi) | i^{\lambda} r^{-\lambda-1} Y_{\lambda\mu}(\theta, \phi) | i^{l_i} Y_{l_i m_i}(\theta, \phi) \rangle} \\
&= -(2\lambda - 1) [\lambda l_i (2l_i + 1)(l_i + 1)]^{\frac{1}{2}} \sum_q (-)^{\lambda+l_i-\mu-m_i+q} \begin{pmatrix} \lambda & 1 & \lambda-1 \\ -\mu-q & q & \mu \end{pmatrix} \\
&\quad \begin{pmatrix} l_i & 1 & l_i \\ -m_i+q & -q & m_i \end{pmatrix} \begin{pmatrix} l_i & \lambda & l_f \\ m_i-q & \mu+q & -m_f \end{pmatrix} \begin{pmatrix} l_i & \lambda & l_f \\ m_i & \mu & -m_f \end{pmatrix}^{-1} \\
&= (2\lambda - 1) [\lambda l_i (2l_i + 1)(l_i + 1)]^{\frac{1}{2}} \left\{ \begin{matrix} l_f & l_i & \lambda-1 \\ 1 & \lambda & l_i \end{matrix} \right\} \begin{pmatrix} l_i & \lambda-1 & l_f \\ m_i & \mu & -m_f \end{pmatrix} \begin{pmatrix} l_i & \lambda & l_f \\ m_i & \mu & -m_f \end{pmatrix}^{-1} \\
&= \frac{1}{2} \left[\frac{2\lambda-1}{2\lambda+1} \right]^{\frac{1}{2}} [(l_f + l_i + \lambda + 1)(l_f + l_i - \lambda + 1)(l_f - l_i + \lambda)(l_i - l_f + \lambda)]^{\frac{1}{2}} \\
&\quad \begin{pmatrix} l_i & \lambda-1 & l_f \\ m_i & \mu & -m_f \end{pmatrix} \begin{pmatrix} l_i & \lambda & l_f \\ m_i & \mu & -m_f \end{pmatrix}^{-1} \tag{IV.55}
\end{aligned}$$

using :

$$\left\{ \begin{matrix} l_f & l_i & \lambda-1 \\ 1 & \lambda & l_i \end{matrix} \right\} = (-)^{l_i+l_f+\lambda} \frac{1}{2} \left[\frac{(l_f + l_i + \lambda + 1)(l_f + l_i - \lambda + 1)(l_f - l_i + \lambda)(l_i - l_f + \lambda)}{l_i(2l_i + 1)(l_i + 1)(2\lambda - 1)\lambda(2\lambda + 1)} \right]^{\frac{1}{2}} \tag{IV.56}$$

and taking into account that $l_i + l_f + \lambda$ is even. This formula agrees with (II B.48) of Ref. [65] if $2l_i$ is replaced by $(2l_i + 1)$ in it.

The ratio of the couplings between a state $|(l_i s) j_i I_i J >$ and a state $|(l_f s) j_f I_f J >$ is :

$$\begin{aligned}
& -\frac{1}{2} \left[\frac{2\lambda-1}{2\lambda+1} \right]^{\frac{1}{2}} [(l_f + l_i + \lambda + 1)(l_f + l_i - \lambda + 1)(l_f - l_i + \lambda)(l_i - l_f + \lambda)]^{\frac{1}{2}} \\
& \left\{ \begin{matrix} j_f & \lambda-1 & j_i \\ I_i & J & I_f \end{matrix} \right\} \left\{ \begin{matrix} j_f & \lambda-1 & j_i \\ l_i & s & l_f \end{matrix} \right\} \left\{ \begin{matrix} j_f & \lambda & j_i \\ I_i & J & I_f \end{matrix} \right\}^{-1} \left\{ \begin{matrix} j_f & \lambda & j_i \\ l_i & s & l_f \end{matrix} \right\}^{-1} \tag{IV.57}
\end{aligned}$$

that is Equ. (III.44) written with $S = 0$, $I^t = L = \lambda$ but with $L = \lambda - 1$ instead of $L = \lambda$ in the $6j$ and the $9j$ coefficients, multiplied by the factor written in the first line of Equ. (IV.57).

For an excitation $1^+ \rightarrow 2^+$, the relative values of the coefficient el of the electric transition and ma of the magnetic one are :

$$\begin{aligned}
J = l_i = l_f & & el = 1/J & & ma = J \\
J = l_i = l_f \pm 1 & & el = 1 & & ma = 0 \\
J = l_i \pm 1, l_f = l_i & & el = 1 & & ma = J \\
J = l_i \pm 1, l_f = l_i \mp 2 & & el = 0 & & ma = 0
\end{aligned} \tag{IV.58}$$

C.4 Magnetic reduced matrix elements

The reduced matrix element of the magnetic excitation $B(M\lambda - 1)$ is the one of the electric excitation $B(E\lambda)$ multiplied by the ratio R_{p_1} reduced transition probabilities given in Equ. (IV.54). To do so, they should be expressed in compatible units (see [71] SKORKA, S., J₆, HERTEL, J. and RETZ-SCHMIDT, T., W., “*Compilation of electromagnetic transition rates in light nuclei ($A \leq 40$)*”, Nuclear Data Sheet **A2** (1967) page 347). Usually, the reduced electric probabilities $B(E\lambda)$ are given in units $e^2 \text{fm}^{2\lambda}$ and the reduced magnetic probabilities $B(M\lambda)$ in units $\mu_0^2 \text{fm}^{2\lambda-2}$: the value of $B(M\lambda)$ must be multiplied by .01106 to obtain it in units $e^2 \text{fm}^{2\lambda}$.

Chapter V

Excitation of particle and target

We consider the scattering of ^8Be on ^{64}Ni to have integer values of the masses at the power one third. We use the rotational model with a positive deformation for ^{64}Ni and a negative deformation for ^8Be . The following test case :

```
|BE8 + NI64   EXCITATION OF NI64   - 1 - (ECIS94)
|TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
|TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
|
|  2    3
|.25      10.
|
|      10.      8.      64.      0.
|2.      1.
|
|  2    4
|.45
|20.      1.      .5
|10.      1.      .5
|
|      1.      .5
|
|      1.      .5
|
|      1.      .5
|
|      1.      .5
|1.
|
|      1.      .5
|2.      2.      178.
|FIN
```

describes the excitation of a 2^+ state in ^{64}Ni at 1 MeV. The computation is limited to 3 J -values because the output is very detailed: the coefficients of each form-factor for all sets of equations are printed (LO(58).TRUE.) to compare one test with the others. The coulomb potential has been suppressed to get significant results. The usual coupled equations (LO(21)=.TRUE.) with Schmidt's orthogonalisation (LO(42)=.TRUE.) are used for more precision.

This Chapter is the transcription of the similar study in “Notes on ECIS79” which was starting with :

```
|BE8 + NI64   EXCITATION OF NI64   - 1 - (ECIS79)
|TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
|TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
|
|  2
|.25      10.      2.
|
|      10.      8.      64.      0.
|2.      1.
|
|  2    4      .45
```



```

|          10.          8.          64.          0.
|2.          1.
|2.          2.          178.
|  1    1    0
|  1    2    1
|  1    2    0    0    -1.00000000000000
|  2    2    2
|  1    2    0    0    1.195228338242
|  2    4    0    0    1.195228338242
|  1    1          1    2    -5
|20.          6.          .5
|          .3
|  1    1          2    2    -5
|10.          6.          .5
|          .3
|  1    1          3          -1
|
|  1    1          4          -1
|
|  1    1          7          -1
|0.          6.
|  2    2    1    1    2    -5
|20.          6.          .5
|          .3
|  2    2    1    2    2    -5
|10.          6.          .5
|          .3
|  2    2    1    3          -2
|
|  2    2    1    4          -2
|
|  2    2    2    1    2    -5
|20.          6.          .5
|          .3
|  2    2    2    2    2    -5
|10.          6.          .5
|          .3
|  2    2    2    3          -2
|
|  2    2    2    4          -2
|
|FIN

```

and the test case “- 2 -” can be done by changing only the energy in the laboratory system and the sign of the deformation and exchanging the masses :

```

|BE8 + NI64  EXCITATION OF BE8. EXTERNAL POTENTIALS  - 4 -
|TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
|TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
|  2    3
|.25          10.
|          10.          8.          64.          0.
|2.          1.
|2.          2.          178.
|  1    1    0
|  1    2    1

```

	1	2	0	0	-1.00000000000000
	2	2	2		
	1	2	0	0	1.195228338242
	2	4	0	0	1.195228338242
	1	1		1	2
20.		6.		.5	-5
		-.3			
	1	1		2	2
10.		6.		.5	-5
		-.3			
	1	1		3	-1
	1	1		4	-1
	1	1		7	-1
0.		6.			
	2	2	1	1	2
20.		6.		.5	-5
		-.3			
	2	2	1	2	2
10.		6.		.5	-5
		-.3			
	2	2	1	3	-2
	2	2	1	4	-2
	2	2	2	1	2
20.		6.		.5	-5
		-.3			
	2	2	2	2	2
10.		6.		.5	-5
		-.3			
	2	2	2	3	-2
	2	2	2	4	-2
FIN					

A.3 Change of spin for the particle

The excited state of the particle can be described as the scattering of a spin 2 particle on a target of spin 0 :

```

| BE8 + NI64 EXCITATION OF PARTICLE.EXTERNAL POTENTIALS - 5 -
| TFFFFFFTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
| TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
|      2      3
| .25      10.
|      10.      8.      64.      0.
|      1.      2.      8.      64.      0.
| 2.      2.      178.
|      1      1      0
|      1      2      1
|      1      2      4      0      -1.000000000000      (1)
|      2      2      2

```

```

|      1      2      4      0      1.195228338242      (1)
|      2      4      8      0      1.195228338242      (1)
|      1      1          1          2          -5
|20.          6.          .5
|          -.3
|      1      1          2          2          -5
|10.          6.          .5
|          -.3
|      1      1          3          -1
|
|      1      1          4          -1
|
|      1      1          5          -1
|
|      1      1          6          -1
|
|      1      1          7          -1
|0.          6.
|      1      1          8          -1
|
|      2      2      1      1          2          -5
|20.          6.          .5
|          -.3
|      2      2      1      2          2          -5
|10.          6.          .5
|          -.3
|      2      2      1      3          -2
|
|      2      2      1      4          -2
|
|      2      2      2      1          2          -5
|20.          6.          .5
|          -.3
|      2      2      2      2          2          -5
|10.          6.          .5
|          -.3
|      2      2      2      3          -2
|
|      2      2      2      4          -2
|
|FIN

```

The interaction which was $L = 2, S = 0, J = 2$ in the previous test cases is now $L = 2, S = 2, J = 0$ and the one which was $L = 4, S = 0, J = 4$ is now $L = 4, S = 4, J = 0$. The nuclear reduced matrix elements (1) are the same as the previous one, but were multiplied by the reduced “spin” matrix element which is $\sqrt{2S+1}$ in ECIS79 and ECIS88.

As the particle of one of the levels does not have a spin zero, a complex spin-orbit potential and a coulomb spin orbit potential must be read. As this last one is 0, use L0(46)=.TRUE. to avoid long range “Coulomb corrections” for this calculation.

B Excitation of the target and of the particle

We assume that there is no interaction between the level for which the target is excited and the one for which the particle is excited. In the rotational model, such an interaction could be only of second order

and it is smaller than the reorientation terms. In the second order vibrational model, such an interaction exists and is identical to the reorientation terms.

This is obtained by adding the test case “- 3 -” to the test case “- 4 -” or the test case “- 5 -”. However, the deformation of the optical potential of test case “- 3 -” is positive and the one of the other test cases is negative. We shall use an optical potential without deformation.

B.1 Excitation of the two nuclei

The three levels involved are the scattering of a spin zero particle on a spin zero target, of the spin zero particle on a 2^+ state as in test case “- 3 -”, and of a spin two particle on a spin zero target as in test case “- 5 -” :

```

| BE8 + NI64   EXCITATION OF PARTICLE AND TARGET  - 6 -
| TTTTTTTTFFFTFFFTFFFTFFFTFFFTFFFTFFFTFFFTFFFTFFFTFF
| TTTTTTTTFFFTFFFTFFFTFFFTFFFTFFFTFFFTFFFTFFFTFFFTFF
|
|   3   3
| .25   10.
|       10.           8.           64.           0.
| 2.     1.
|       1.           2.           8.           64.           0.
| 2.     2.          178.
|
|  1   1   0
|  1   2   1
|  1   2   0   0   -1.000000000000
|  1   3   1
|  2   2   4   0   -1.000000000000   (1)
|  2   2   2
|  1   2   0   0   1.195228338242
|  3   4   0   0   1.195228338242
|  2   3   0
|  3   3   2
|  2   2   4   0   1.195228338242   (1)
|  4   4   8   0   1.195228338242   (1)
|  1   1   1   2   -1
| 20.   6.   .5
|  1   1   2   2   -1
| 10.   6.   .5
|  1   1   3   -1
|
|  1   1   4   -1
|
|  1   1   5   -1
|
|  1   1   6   -1
|
|  1   1   7   -1
| 0.     6.
|  1   1   8   -1
|
|  2   2   1   1   2   -5
| 20.   6.   .5
|       .3
|  2   2   1   2   2   -5
| 10.   6.   .5

```

The reduced nuclear matrix elements are the ones of test cases “- 2 -” and “- 5 -”. The form factors “1” and “3” are the ones of test case “- 2 -” (positive deformation) and the form factors “2” and “4” are the ones of the test case “- 5 -” (negative deformation). The nuclear matrix elements (1) have no more to be multiplied by $\sqrt{2S+1}$ in ECIS94.

In this test case, the three levels are the scattering of a spin zero particle on a spin zero target and two 2^+ states of the target :

```

| BE8 + NI64      TWO EXCITATIONS OF TARGET  - 7 -
| TFFFFFFFFFFFFFTFFFFFFFTFFFFFFFFFFFFFFFFFFFFFFFFFTFTFF
| TTTTTTTTTFFFFFFFTFFFTFFFFFFFFFFFFFFFFFFFFFFFFFFFFF
|      3      3
| .25      10.
|           10.           8.           64.           0.
| 2.      1.
| 2.      1.

```

```

| 2.      2.      178.
| 1      1      0
| 1      2      1
| 1      2      0      0      -1.0000000000000
| 1      3      1
| 2      2      0      0      -1.0000000000000
| 2      2      2
| 1      2      0      0      1.195228338242
| 3      4      0      0      1.195228338242
| 2      3      0
| 3      3      2
| 2      2      0      0      1.195228338242
| 4      4      0      0      1.195228338242
| 1      1      1      2      -1
| 20.     6.      .5
| 1      1      2      2      -1
| 10.     6.      .5
| 1      1      3      -1
|
| 1      1      4      -1
|
| 1      1      7      -1
| 0.      6.
| 2      2      1      1      2      -5
| 20.     6.      .5
|
| .3
| ..... SAME CARDS AS IN - 6 - .....

```

The reduced nuclear matrix elements are the ones of the pure rotational model and there is no complex spin-orbit to be read with the optical potential because the particle has spin zero for all the levels.

C Double excitation

The double excitation can be described as the scattering of a particle with spin 2 on a target with spin 2, or as the sum of levels with spin 0 to 4 which are the channel spins.

C.1 One level description

The four channels are the scattering of a particle of spin zero and of spin two on a target with a spin zero and two. The coupling between the levels where the particle has a fixed spin and the target the spins zero and two is the same as in the rotational model for the excitation of a 2^+ state by a particle of spin zero or two. The coupling between the channels where the target has a given spin is the same as in test case “- 5 -”. The reorientation terms of the double excitation level are the sum of the reorientation terms of the 2^+ of ^8Be and the 2^+ state state of ^{64}Ni :

```

| BE8 + NI64  DOUBLE EXCITATION DESCRIPTION ONE LEVEL  - 8 -
| TFFFFFFFFFFFFFTFFFFFFFTFFFFFFFFFFFFFFFFFFFFFFFFFTFTFF
| TTTTTTTTFFFFFFFFFTFTFTFTFTFTFTFTFTFTFTFTFTFTFTFTFTFF
| 4      3
| .25     10.
|         10.      8.      64.      0.
| 2.      1.
|         2.      2.      8.      64.      0.

```



```

| 2.      3.
| 2.      2.      178.
| 1      1      0
| 1      2      1
| 1      2      0      0      -1.000000000000
| 1      3      1
| 2      2      4      0      -1.000000000000      (1)
| 1      4      0
| 2      2      2
| 1      2      0      0      1.195228338242
| 3      4      0      0      1.195228338242
| 2      3      0
| 2      4      1
| 2      2      4      0      -2.2360680      (2)
| 3      3      2
| 2      2      4      0      1.195228338242      (1)
| 4      4      8      0      1.195228338242      (1)
| 3      4      1
| 1      2      0      0      -1.000000000000
| 4      4      4
| 1      2      0      0      1.195228338242
| 3      4      0      0      1.195228338242
| 2      2      4      0      2.672612419      (2)
| 4      4      8      0      2.672612419      (2)
| ..... SAME CARDS AS IN - 6 - .....

```

The reduced nuclear matrix (2) are the usual nuclear matrix elements multiplied by the reduced matrix element in the spin space of the target $\sqrt{2I+1}$. This remains in ECIS94, whereas the reduced matrix elements (1) and (2) had to be multiplied by $\sqrt{2S+1}$ in ECIS79 and ECIS88.

C.2 Many levels description

The level where the particle and the target are in a 2^+ state can be replaced by five levels with spin 0^+ , 1^+ , 2^+ , 3^+ and 4^+ , resulting from the coupling of the spins of the particle and of the target (channel spin). The cross-section will be the sum of the five cross-sections. The data set is :

```

| BE8 + NI64 DOUBLE EXCITATION DESCRIPTION FIVE LEVELS - 9 -
| TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
| TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
| 8      3
| .25      10.
|          10.      8.      64.      0.
| 2.      1.
|          2.      2.      8.      64.      0.
|          3.      8.      64.      0.
| 1.      3.
| 2.      3.
| 3.      3.
| 4.      3.
| 2.      2.      178.
| 1      1      0
| 1      2      1
| 1      2      0      0      -1.000000000000
| 1      3      1
| 2      2      4      0      -1.000000000000      (1)

```

	1	4	0			
	1	5	0			
	1	6	0			
	1	7	0			
	1	8	0			
	2	2	2			
	1	2	0	0	1.195228338242	
	3	4	0	0	1.195228338242	
	2	3	0			
	2	4	1			
	2	2	0	0	-.4472135955	(3)
	2	5	1			
	2	2	0	0	.7745966692	(3)
	2	6	1			
	2	2	0	0	-1.	(3)
	2	7	1			
	2	2	0	0	1.183215957	(3)
	2	8	1			
	2	2	0	0	-1.341640786	(3)
	3	3	2			
	2	2	4	0	1.195228338242	(1)
	4	4	8	0	1.195228338242	(1)
	3	4	1			
	1	2	4	0	-0.447213596000	(4)
	3	5	1			
	1	2	4	2	0.774596669	(4)
	3	6	1			
	1	2	4	4	-1.0000000000	(4)
	3	7	1			
	1	2	4	6	1.183215957	(4)
	3	8	1			
	1	2	4	8	-1.341640786000	(4)
	4	4	0			
	4	5	0			
	4	6	2			
	1	2	0	0	.534522483825	(5)
	2	2	0	0	.534522483825	(5)
	4	7	0			
	4	8	2			
	3	4	0	0	.534522483825	(5)
	4	4	0	0	.534522483825	(5)
	5	5	2			
	1	2	0	0	-.547722555751	(5)
	2	2	0	0	-.547722555751	(5)
	5	6	2			
	1	2	0	0	-.462910049886	(5)
	2	2	0	0	.462910049886	(5)
	5	7	4			
	1	2	0	0	.585540043769	(5)
	2	2	0	0	.585540043769	(5)
	3	4	0	0	-.377964473009	(5)
	4	4	0	0	-.377964473009	(5)
	5	8	2			
	3	4	0	0	-.845154254729	(5)
	4	4	0	0	.845154254729	(5)

	6	6	4			
	1	2	0	0	-.256120416286	(5)
	2	2	0	0	-.256120416286	(5)
	3	4	0	0	.341493888381	(5)
	4	4	0	0	.341493888381	(5)
	6	7	4			
	1	2	0	0	-.808122035642	(5)
	2	2	0	0	.808122035642	(5)
	3	4	0	0	.714285714286	(5)
	4	4	0	0	-.714285714286	(5)
	6	8	4			
	1	2	0	0	.458162128993	(5)
	2	2	0	0	.458162128993	(5)
	3	4	0	0	.895404529326	(5)
	4	4	0	0	.895404529326	(5)
	7	7	4			
	1	2	0	0	.292770021885	(5)
	2	2	0	0	.292770021885	(5)
	3	4	0	0	-.886405260428	(5)
	4	4	0	0	-.886405260428	(5)
	7	8	4			
	1	2	0	0	-.958314847500	(5)
	2	2	0	0	.958314847500	(5)
	3	4	0	0	-.749149177244	(5)
	4	4	0	0	.749149177244	(5)
	8	8	4			
	1	2	0	0	1.20131123696	(5)
	2	2	0	0	1.20131123696	(5)
	3	4	0	0	.456568516759	(5)
	4	4	0	0	.456568516759	(5)
 SAME CARDS AS IN - 6 -					

The reduced nuclear matrix element in the spins spaces between a target of spin I_1 and a particle of spin s_1 , coupled to J_1 and a target of spin I_2 and a particle of spin s_2 , coupled to J_2 is :

$$\sqrt{(2I+1)(2S+1)(2L+1)(2J_1+1)(2J_2+1)} \left\{ \begin{array}{ccc} I_1 & I_2 & I \\ s_1 & s_2 & S \\ J_1 & J_2 & L \end{array} \right\} \quad (\text{V.1})$$

where $S = 0$ for the form factors 1 and 3, $I = 0$ for the form factors 2 and 4.

- For the reduced nuclear matrix elements (3) above, $s_1 = I = 0$ and the $9j$ coefficient reduces to $0.04(-)^{J_2}$.
- For the reduced nuclear matrix elements (4) above, $I_1 = S = 0$ and the $9j$ coefficient reduces to 0.04. They are multiplied by a reduced matrix element in the spin space which is $\sqrt{5}$. However, they must be multiplied by $(-)^{J_2} (?)$.
- For the reduced nuclear matrix elements (5) above, $I_1 = I_2 = s_1 = s_2 = 2$ and $I = 0$ or $S = 0$. In this reduced nuclear matrix element, the $9j$ coefficient becomes a $6-j$ coefficient and the result is :

$$\sqrt{(2L+1)(2J_1+1)(2J_2+1)/5} \left\{ \begin{array}{ccc} J_1 & J_2 & L \\ 2 & 2 & 2 \end{array} \right\} \quad (\text{V.2})$$

with a phase $(-)^{J_2}$ if $S = 0$ and $(-)^{J_1}$ if $I = 0$.

In fact, the explanation given above is what was given in the “Notes on ECIS79” and do not seem accurate. Perhaps, the result is due to the occurrence of 0 for S or I . Nevertheless, using the vibrational model with two different phonons 2^+ and with $0^+, 1^+, 2^+, 3^+, 4^+$ 2-phonons states :

	3	5	1			
	1	2	0	0	-0.774596669	(3)
	3	6	1			
	1	2	0	0	-1.0000000000	(3)
	3	7	1			
	1	2	0	0	-1.183215957	(3)
	3	8	1			
	1	2	0	0	-1.341640786	(3)
	4	4	0			
	4	5	0			
	4	6	2			
	1	2	0	0	.534522483825	(5)
	2	2	0	0	.534522483825	(5)
	4	7	0			
	4	8	2			
	3	4	0	0	.534522483825	(5)
	4	4	0	0	.534522483825	(5)
	5	5	2			
	1	2	0	0	-.547722555751	(5)
	2	2	0	0	-.547722555751	(5)
	5	6	2			
	1	2	0	0	-.462910049886	(5)
	2	2	0	0	.462910049886	(5)
	5	7	4			
	1	2	0	0	.585540043769	(5)
	2	2	0	0	.585540043769	(5)
	3	4	0	0	-.377964473009	(5)
	4	4	0	0	-.377964473009	(5)
	5	8	2			
	3	4	0	0	-.845154254729	(5)
	4	4	0	0	.845154254729	(5)
	6	6	4			
	1	2	0	0	-.256120416286	(5)
	2	2	0	0	-.256120416286	(5)
	3	4	0	0	.341493888381	(5)
	4	4	0	0	.341493888381	(5)
	6	7	4			
	1	2	0	0	-.808122035642	(5)
	2	2	0	0	.808122035642	(5)
	3	4	0	0	.714285714286	(5)
	4	4	0	0	-.714285714286	(5)
	6	8	4			
	1	2	0	0	.458162128993	(5)
	2	2	0	0	.458162128993	(5)
	3	4	0	0	.895404529326	(5)
	4	4	0	0	.895404529326	(5)
	7	7	4			
	1	2	0	0	.292770021885	(5)
	2	2	0	0	.292770021885	(5)
	3	4	0	0	-.886405260428	(5)
	4	4	0	0	-.886405260428	(5)
	7	8	4			
	1	2	0	0	-.958314847500	(5)
	2	2	0	0	.958314847500	(5)
	3	4	0	0	-.749149177244	(5)

D.1 Excitation of the particle

The same test case can be run with matrix elements and potentials read from cards with `LO(7)=.TRUE.` :

```

| BE8 + BE8   EXCITATION OF BE8 . EXTERNAL POTENTIALS   - 12 -
| TFFFFFFFFFTFFFTFFFTFFFFFFFFFFFFFFFFFFFFFFFFFFFFFTFF
| TTTTTTTTFFFFFFFFFTFFFTFFFFFFFFFFFFFFFFFFFFFFFFFFFF
|
|   2   5
| .25   10.
|       10.           8.           8.           0.
| 2.     1.
| 2.     2.     90.
|
| 1 1 0
| 1 2 1
| 1 2 0 0 -1.000000000000
| 2 2 2
| 1 2 0 0 1.195228338242
| 2 4 0 0 1.195228338242
| 1 1 1 2 -5
| 20. 6. .5
|     -.3
| 1 1 2 2 -5
| 10. 6. .5
|     -.3
| 1 1 3 -1
|
| 1 1 4 -1
|
| 1 1 7 -1
| 0. 6.
| 2 2 1 1 2 -5
| 20. 6. .5
|     -.3
| 2 2 1 2 2 -5
| 10. 6. .5
|     -.3
| 2 2 1 3 -2
|
| 2 2 1 4 -2
|
| 2 2 2 1 2 -5
| 20. 6. .5
|     -.3
| 2 2 2 2 2 -5
| 10. 6. .5
|     -.3
| 2 2 2 3 -2
|
| 2 2 2 4 -2
|
| FIN

```

The same result is obtained if the particle has a spin 2 in the final state and the target a spin 0 as in the test case “- 5 -” above. The interaction $L = 2, S = 0, J = 2$ between the ground state and the excited state is changed into an interaction $L = 2, S = 2, J = 0$. It is the same for the reorientation interaction $L = 2$. The reorientation interaction $L = 4, S = 0, J = 4$ becomes an interaction $L = 4, S = 4, J = 0$.

The nuclear nuclear matrix were multiplied by $\sqrt{2S+1}$ for ECIS79 and ECIS88 but not for ECIS94. As the excited state has a non-zero spin, a complex spin-orbit potential and a coulomb spin-orbit potential must be given and LO(46) set .TRUE..

```

| BE8 + BE8 EXCITATION OF BE8.EXCHANGE PARTICLE-TARGET - 13 -
| TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
| TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
|
| 2 5
| .25 10.
| 10. 8. 8. 0.
| 1. 2. 8. 8. 0.
| 2. 2. 90.
| 1 1 0
| 1 2 1
| 1 2 4 0 -1.000000000000 (1)
| 2 2 2
| 1 2 4 0 1.195228338242 (1)
| 2 4 8 0 1.195228338242 (1)
| 1 1 1 2 -5
| 20. 6. .5
| -.3
| 1 1 2 2 -5
| 10. 6. .5
| -.3
| 1 1 3 -1
| 1 1 4 -1
| 1 1 5 -1
| 1 1 6 -1
| 1 1 7 -1
| 0. 6.
| 1 1 8 -1
| 2 2 1 1 2 -5
| 20. 6. .5
| -.3
| 2 2 1 2 2 -5
| 10. 6. .5
| -.3
| 2 2 1 3 -2
| 2 2 1 4 -2
| 2 2 2 1 2 -5
| 20. 6. .5
| -.3
| 2 2 2 2 2 -5
| 10. 6. .5
| -.3
| 2 2 2 3 -2
| 2 2 2 4 -2

```


FIN

D.2 Excitation of the particle and the target

The two channels of excitation of the particle and of the target can be taken into account with the following data :

BE8 + BE8					
EXCITATION OF PARTICLE AND TARGET - 14 -					
TTTTTTTFFFFFTFFFTFFF					
TTTTTTTTTTTTTTTTTTTT					
.25	3	5			
		10.			
		10.	8.	8.	0.
2.		1.			
		1.	2.	8.	0.
2.	2.	90.			
1	1	0			
1	2	1			
1	2	0	0	-1.000000000000	
1	3	1			
1	2	4	0	-1.000000000000	(1)
2	2	2			
1	2	0	0	1.195228338242	
2	4	0	0	1.195228338242	
2	3	0			
3	3	2			
1	2	4	0	1.195228338242	(1)
2	4	8	0	1.195228338242	(1)
..... SAME CARDS AS IN - 13 -					

This data set is similar to the test case “- 6 -” when particle and target are not identical, but there is only one kind of deformation and the central potential is kept deformed. The interaction between the excited state of the target and the excited state of the particle have been neglected. The result of this test case is not the same as the one of test cases “- 11 -” to “- 12 -”.

As above, in test case “- 7 -”, the two excited states can be attributed to the target. As they are degenerated, the waves functions for the two excited states are identical and can be replaced by only one wave function in the equations after some modifications of the nuclear matrix elements. Data are the following :

```

BE8 + BE8      SIMPLIFIED EXCIT. OF PARTICLE AND TARGET  - 15 -
|TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
|TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
|      2      5
|.25      10.
|      10.      8.      8.      0.
|2.      1.
|2.      2.      90.
|      1      1      0
|      1      2      1
|      1      2      0      0      -1.414213562      (6)
|      2      2      2
|      1      2      0      0      1.195228338242
|      2      4      0      0      1.195228338242
|..... SAME CARDS AS IN - 12 - .....

```

The interaction between the ground state and the excited state has been multiplied by $\sqrt{2}$ and the reorientation terms are the sum of the reorientation terms and of the coupling terms between the two excited states, but we did not consider the interaction between the two excited states in this example.

As there is only one kind of deformations involved, this calculation can be done in the framework of standard rotational model by giving the nuclear matrix elements on card (LO(15)=.TRUE., LO(7)=.FALSE.). The data set is :

```
|BE8 + BE8   SIMPLIFIED EXCIT. OF PARTICLE AND TARGET  - 16 -
|TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
|TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
|      2      5
|.25      10.
|      10.      8.      8.      0.
|2.      1.
|      2      4
|-.6
|20.      1.5      .5
|10.      1.5      .5
|      1.5      .5
|      1.5      .5
|      1.5      .5
|      1.5      .5
|1.5
|      1.5      .5
|2.      2.      90.
|      1      1      0
|      1      2      1
| 3000      2      0      0      -1.414213562      (6)
|      2      2      2
| 3000      2      0      0      1.195228338242
| 5000      4      0      0      1.195228338242
|FIN
```

which almost as simple as the data set “- 11 -”.

D.3 Double excitation

The data set for the excitation of the target, the excitation of the particle and the simultaneous excitation of the target and the particle is obtained from the cases “- 12 -” and “- 13 -” when there is no interaction between the excited states. The nuclear matrix elements for the excitation of the target are independent of the excitation of particle and the one of the particle are independent of the the excitation of the target :

```
|BE8 + BE8   DOUBLE EXCITATION DESCRIPTION ONE LEVEL  - 17 -
|TTTTTTTFTFTFTFTFTFTFTFTFTFTFTFTFTFTFTFTFTFTFTFTFTFTFTFT
|TTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
|      4      5
|.25      10.
|      10.      8.      8.      0.
|2.      1.
|      1.      2.      8.      8.      0.
|2.      2.
|2.      2.      90.
|      1      1      0
|      1      2      1
```

```

| 1 2 0 0 -1.000000000000
| 1 3 1
| 1 2 4 0 -1.000000000000 (1)
| 1 4 0
| 2 2 2
| 1 2 0 0 1.195228338242
| 2 4 0 0 1.195228338242
| 2 3 0
| 2 4 1
| 1 2 4 0 -2.236067977 (2)
| 3 3 2
| 1 2 4 0 1.195228338242 (1)
| 2 4 8 0 1.195228338242 (1)
| 3 4 1
| 1 2 0 0 -1.000000000000
| 4 4 4
| 1 2 0 0 1.195228338242
| 2 4 0 0 1.195228338242
| 1 2 4 0 2.672612419 (2)
| 2 4 8 0 2.672612419 (2)
| ..... SAME CARDS AS IN - 13 - .....

```

where the interaction of the two single excitation channels with the double excitation were not the same with ECIS79 and ECIS88 but are the same for ECIS94.

One can use the channel spin description of the double excitation as for the test case “- 9 -” or “- 10 -”. One can see on the test case “- 10 -” that the 1^+ and the 3^+ states are not coupled to the other states, so they must not be taken into account. The equivalent to the test case “- 9 -” is :

```

| BE8 + BE8 DOUBLE EXCIT. DESCRIPTION THREE LEVELS - 18 -
| TTTTTTTTTTTTTFTFTFTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
| TTTTTTTTTTTTTTTTTFTFTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTT
| 6 5
| .25 10.
| 10. 8. 8. 0.
| 2. 1.
| 1. 2. 8. 8. 0.
| 2. 8. 8. 0.
| 2. 2.
| 4. 2.
| 2. 2. 90.
| 1 1 0
| 1 2 1
| 1 2 0 0 -1.000000000000
| 1 3 1
| 1 2 4 0 -1.000000000000 (1)
| 1 4 0
| 1 5 0
| 1 6 0
| 2 2 2
| 1 2 0 0 1.195228338242
| 2 4 0 0 1.195228338242
| 2 3 0
| 2 4 1
| 1 2 0 0 -.4472135955 (3)
| 2 5 1

```


2	5	1				
1	2	0	0	-1.	(3)	
2	6	1				
1	2	0	0	-1.341640786	(3)	
3	3	2				
1	2	0	0	1.195228338242		
2	4	0	0	1.195228338242		
3	4	1				
1	2	0	0	-0.447213596000	(3)	
3	5	1				
1	2	0	0	-1.0000000000	(3)	
3	6	1				
1	2	0	0	-1.341640786000	(3)	
4	4	0				
4	5	1				
1	2	0	0	1.06904496765	(5)	
4	6	1				
2	4	0	0	1.06904496765	(5)	
5	5	2				
1	2	0	0	-.512240832572	(5)	
2	4	0	0	.682987776762	(5)	
5	6	2				
1	2	0	0	.916324257986	(5)	
2	4	0	0	1.79080905865	(5)	
6	6	2				
1	2	0	0	2.40262247392	(5)	
2	4	0	0	.913137033518	(5)	
..... SAME CARDS AS IN - 12 -						

The absence of the double excitation channels 1^+ and 3^+ reduces the number of equations from 20 to 17. Furthermore, as the two single excitation levels are degenerated and have the same coupling to the other states, they can be replaced by a single level with reduced matrix elements multiplied by $\sqrt{2}$ as for the test case “- 16 -”. The number of coupled equations decreases to 13. The data are :

BE8 + BE8 DOUBLE EXCIT. DESCRIPTION THREE LEVELS - 20 -					
TFFFFFFTTTTFTTFTFFFFFFFFFFFFFFFFFFFFFTTFF					
TTTTTTTFFFFFFFFFTTFTFFFFFFFFFFFFFFFFFFFFF					
	5	5			
	.25	10.			
		10.	8.	8.	0.
	2.	1.			
		2.			
	2.	2.			
	4.	2.			
	2.	2.	90.		
	1	1	0		
	1	2	1		
	1	2	0	-1.414213562	(6)
	1	3	0		
	1	4	0		
	1	5	0		
	2	2	2		
	1	2	0	1.195228338242	
	2	4	0	1.195228338242	
	2	3	1		
	1	2	0	-.632455532034	(6)

```

| 2 4 1
| 1 2 0 0 -1.414213562 (6)
| 2 5 1
| 1 2 0 0 -1.8973665961 (6)
| 3 3 0
| 3 4 1
| 1 2 0 0 1.06904496765 (5)
| 3 5 1
| 2 4 0 0 1.06904496765 (5)
| 4 4 2
| 1 2 0 0 -.512240832572 (5)
| 2 4 0 0 .682987776762 (5)
| 4 5 2
| 1 2 0 0 .916324257986 (5)
| 2 4 0 0 1.79080905865 (5)
| 5 5 2
| 1 2 0 0 2.40262247392 (5)
| 2 4 0 0 .913137033518 (5)
| ..... SAME CARDS AS IN - 12 - .....

```

This calculation can be done with the usual rotational model by giving the nuclear reduced matrix elements on cards (LO(15)=.TRUE.) as in the test case “- 15 - ” with respect to the “- 14 - ”. This was already possible for the test case test case “- 19 -”, but not for the test case “- 18 -” which involves transfers of spin. The data set is :

```

| BE8 + BE8 DOUBLE EXCIT. DESCRIPTION THREE LEVELS - 21 -
| TTTTTTTTTFTTFTFTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTFF
| TTTTTTTTTTTTTTTTTFTFTTTTTTTTTTTTTTTTTTTTTTTTTTTTTTFF
| 5 5
| .25 10.
| 10. 8. 8. 0.
| 2. 1.
| 2.
| 2. 2.
| 4. 2.
| 2 4
| -.6
| 20. 1.5 .5
| 10. 1.5 .5
| 1.5 .5
| 1.5 .5
| 1.5 .5
| 1.5 .5
| 1.5
| 1.5 .5
| 2. 2. 90.
| 1 1 0
| 1 2 1
| 3000 2 0 0 -1.414213562 (6)
| 1 3 0
| 1 4 0
| 1 5 0
| 2 2 2
| 3000 2 0 0 1.195228338242
| 5000 4 0 0 1.195228338242
| 2 3 1

```


to the reorientation terms. By adding the nuclear matrix elements related to the same form factor as printed in the output of test case “- 23 -”, one get the test case :

BE8 + BE8						VIBRATIONAL SYMMETRISED LEVELS	- 24 -
FTFFFFFFFFFTFTTTTFTFTFTFFFFFFFFFFFFFFFFFFFFFFFFFFFFF							
TTTTTTTTTFTFTFFFFFFFFFTFTFTFFFFFFFFFFFFFFFFFFFFFFFFFFFFF							
	5	5					
	.25	10.					
		10.	8.	8.	0.		
	2.	1.5					
	1	1					
		3.					
	2	1	1				
	2.	3.					
	2	1	1				
	4.	3.					
	2	1	1				
	2	.6					
	20.	1.5	.5				
	10.	1.5	.5				
	1.5						
	2.	2.	90.				
	1	1	1				
	3	0	0	0	2.000000000000	(7)	
	1	2	1				
	1	2	0	0	-1.414213562373	(8)	
	1	3	1				
	3	0	0	0	0.894427191000		
	1	4	1				
	3	2	0	0	1.069044967650		
	1	5	1				
	3	4	0	0	1.434274331201		
	2	2	3				
	3	0	0	0	6.260990336000	(7)	
	3	2	0	0	2.138089935300	(7)	
	3	4	0	0	2.868548662402	(7)	
	2	3	1				
	1	2	0	0	-0.632455532034	(8)	
	2	4	1				
	1	2	0	0	-1.414213562373	(8)	
	2	5	1				
	1	2	0	0	-1.897366596101	(8)	
	3	3	1				
	3	0	0	0	2.800000000000	(7)	
	3	4	1				
	3	2	0	0	0.956182887468	(7)	
	3	5	1				
	3	4	0	0	1.282853961180	(7)	
	4	4	3				
	3	0	0	0	6.260990337000	(7)	

are more than one function y_2 , y_3 or y_4 , but the result is the same. The system of equations is :

$$\begin{aligned}
0 &= \frac{d^2}{dr^2} y_1 + V_1 y_1 + V_{11} y_1 + V_{12} y_2 + V_{12} y_3 + V_{14} y_4 \\
0 &= \frac{d^2}{dr^2} y_2 + V_2 y_2 + V_{12} y_1 + V_{22} y_2 + W_{22} y_3 + V_{24} y_4 \\
0 &= \frac{d^2}{dr^2} y_3 + V_2 y_3 + V_{12} y_1 + W_{22} y_2 + V_{22} y_3 + V_{34} y_4 \\
0 &= \frac{d^2}{dr^2} y_4 + V_4 y_4 + V_{14} y_1 + V_{24} y_2 + V_{24} y_3 + V_{44} y_4
\end{aligned} \tag{V.3}$$

where V_1, V_2, V_4 are the potentials, V_{11}, V_{22}, V_{44} the reorientation terms, $V_{12}, V_{14}, W_{22}, V_{24}$ the couplings between different channels. Using :

$$y_2 = \frac{1}{\sqrt{2}}(u_2 + u_3), \quad y_3 = \frac{1}{\sqrt{2}}(u_2 - u_3), \tag{V.4}$$

this system of equations can be written :

$$\begin{aligned}
0 &= \frac{d^2}{dr^2} y_1 + V_1 y_1 + V_{11} y_1 + \sqrt{2} V_{12} u_2 + V_{14} y_4 \\
0 &= \frac{d^2}{dr^2} u_2 + V_2 u_2 + \sqrt{2} V_{12} y_1 + (V_{22} + W_{22}) u_2 + \sqrt{2} V_{24} y_4 \\
0 &= \frac{d^2}{dr^2} y_4 + V_4 y_4 + V_{14} y_1 + \sqrt{2} V_{24} u_2 + V_{44} y_4 \\
0 &= \frac{d^2}{dr^2} u_3 + V_2 u_3 + (V_{22} - W_{22}) u_3
\end{aligned} \tag{V.5}$$

The last equation is not coupled to the others. As it is not the ground state, its solution is zero and $y_2 = y_3 = u_2/\sqrt{2}$. The C-matrix elements of y_2 or y_3 are those of u_2 divided by $\sqrt{2}$, so the cross-sections related to y_2 and y_3 are the half of the one computed with u_2 . The cross-section obtained with u_2 is the sum of the excitation of the target and of the particle. The symmetrisation, skipping odd J -values and doubling even J -values is independent of the problem of excitation of the target and of the particle.

In the three first equations above, the coupling between u_2 and y_1 or y_4 is the coupling of y_2 or y_3 with y_1 or y_4 multiplied by $\sqrt{2}$. The reorientation terms for u_2 is $V_{22} + W_{22}$, that is the sum of the reorientation terms of y_2 or y_3 and the coupling terms between y_2 and y_3 .

In the rotational model, the coupling between the two excited states must be smaller and of the order of the reorientation terms in second order vibrational model. They should be obtained by an angular integration on a multipole. We have neglected them here. The use of the multipole of order zero for the central potential can also be discussed to know if the deformation should be increased (multiplied by $\sqrt{2}$). With such an increase, it is not necessary to give the nuclear reduced matrix elements on cards as far as double excitation is not included. In the rotational model, it is not equivalent to increase the deformation or the nuclear matrix element.

In the vibrational model, after the correction introduced in January 1981 to the second order coupling between two one boson states, which makes it equal to the reorientation terms, the particle and target excitation can be taken into account by multiplying the deformation by $\sqrt{2}$ as well in the second order as in the first order if double excitation is not taken into account.

E.3 Summary of test cases

The output of these test cases varies from one to another. The total output is very large because the main value is 1200 lines by test.

The following test cases must give identical results:

1. “- 1 -” and “- 3 -” excitation of ^{64}Ni .

2. “- 2 -” , “- 4 -” and “- 5 -” excitation of 8Be .
3. “- 6 -” and “- 7 -” excitation of ${}^{64}Ni$ and 8Be .
4. “- 8 -” , “- 9 -” and “- 10 -” in summing the last five levels of the two last tests to obtain the last cross-section of the first one.
5. “- 11 -” , “- 12 -” and “- 13 -” excitation of one 8Be .
6. “- 14 -” , “- 15 -” and “- 16 -” excitation of the two 8Be .
7. “- 17 -” , “- 18 -” , “- 19 -” , “- 20 -” and “- 21 -” double excitation, but, in some cases, some cross-sections are the sum of cross-sections obtained in another tests.
8. “- 22 -” second order vibrational model for two 2^+ in 8Be .
9. “- 23 -” and “- 24 -” double excitation of 8Be in second order vibrational model (only these two examples deal with mutual excitation).
10. “- 25 -” first order vibrational model for double excitation in 2^+ of 8Be .

The nuclear reduced matrix elements are -1. between the ground state and the excited state and the two reorientation ones are $\sqrt{10/7}$ in the rotational model. The others couplings in the vibrational model are too various to be listed here. The signification of the indications between parenthesis is :

- (1) were multiplied by $\sqrt{2S+1}$ with $2S$ in column 15 in **ECIS79** and **ECIS88**, without change for **ECIS94**.
- (2) multiplied by $\sqrt{(2S+1)(2I+1)}$ where I is the spin of the target in **ECIS79** and **ECIS88**, by $\sqrt{2I+1}$ in **ECIS94**.
- (3) multiplied by a $9j$ coefficient as discussed after test case “- 9 -”.
- (4) multiplied by a $9j$ coefficient as discussed after test case “- 9 -”.
- (5) multiplied by a $9j$ coefficient as discussed after test case “- 9 -”.
- (6) multiplied by $\sqrt{2}$ to eliminate one level.
- (7) summed on equivalent form factors.
- (8) summation of coupling and reorientation terms.

Chapter VI

Transfer reactions

We consider the most general case of transfer reaction. We define the approximations which allows to use the codes **ECIS** and indicate the input of these codes. We shall follow : [72] OHMURA, T., IMANISHI, B., ICHIMURA, M., KAWAI, M., “*Study of Deuteron Stripping Reaction by Coupled Channel Theory. I – Variational Formulation and Discussion on Basic Equation –*”, Progress of Theor. Phys., **41** (1969) 391; ...*II – Properties of Interaction Kernel and Method of Numerical Solution –*”, Progress of Theor. Phys., **43** (1970) 347; “...*III – Numerical Results without Non Orthogonality –*”, Progress of Theor. Phys., **44** (1970) 1242 with slightly generalised notations.

Results of **ECIS79** have been compared with results of **CHUCK** and results of **ECIS94** with **DWUCK**.

A Notations

Let us consider a target of mass M_a at the point \vec{R}_a and a particle of mass M_d which is a bound state of a “neutron” of mass M_n at the point \vec{R}_n and of a “proton” of mass M_p at the point \vec{R}_p . In an other channel, the “neutron” can be bound to the nucleus of mass M_a : the target is then a nucleus of mass $M_b = M_a + M_n$ at the point \vec{R}_b and the particle is the proton.

A.1 Systems of coordinates

There are two systems of Jacobi coordinates after elimination of the centre of mass. The first, suitable for the first level is :

$$\begin{aligned}\vec{R}_{ad} &= \frac{M_n \vec{R}_n + M_p \vec{R}_p}{M_n + M_p} - \vec{R}_a = \vec{R}, \\ \vec{R}_{np} &= \vec{R}_p - \vec{R}_n,\end{aligned}\tag{VI.1}$$

The second one is more suitable for the second level :

$$\begin{aligned}\vec{R}_{bp} &= \vec{R}_p - \frac{M_n \vec{R}_n + M_a \vec{R}_a}{M_a + M_n} = \vec{r}, \\ \vec{R}_{an} &= \vec{R}_n - \vec{R}_a.\end{aligned}\tag{VI.2}$$

Any of these vectors can be expressed in terms of two others. These relations will be useful. They are :

$$\vec{R}_{ad} = -\frac{M_n(M_a + M_n + M_p)}{M_a(M_n + M_p)}\vec{R}_{np} + \frac{M_a + M_n}{M_a}\vec{R}_{bp}$$

$$\begin{aligned}
&= \frac{M_p}{M_n + M_p} \vec{R}_{bp} + \frac{M_n(M_a + M_n + M_p)}{(M_a + M_n)(M_n + M_p)} \vec{R}_{an} \\
&= \vec{R}_{an} + \frac{M_p}{M_n + M_p} \vec{R}_{np}, \\
\vec{R}_{np} &= -\frac{M_a(M_n + M_p)}{M_n(M_a + M_n + M_p)} \vec{R}_{ad} + \frac{(M_a + M_n)(M_n + M_p)}{M_n(M_a + M_n + M_p)} \vec{R}_{bp} \\
&= \vec{R}_{bp} - \frac{M_a}{M_n + M_p} \vec{R}_{an} \\
&= \frac{M_n + M_p}{M_p} (\vec{R}_{ad} - \vec{R}_{an}), \tag{VI.3}
\end{aligned}$$

and :

$$\begin{aligned}
\vec{R}_{bp} &= \frac{M_a}{M_a + M_n} \vec{R}_{ad} + \frac{M_n(M_a + M_n + M_p)}{(M_a + M_n)(M_n + M_p)} \vec{R}_{np} \\
&= \vec{R}_{np} + \frac{M_a}{M_a + M_n} \vec{R}_{an} \\
&= \frac{M_n + M_p}{M_p} \vec{R}_{ad} - \frac{M_n(M_a + M_n + M_p)}{M_p(M_a + M_n)} \vec{R}_{ap}, \\
\vec{R}_{an} &= \vec{R}_{ad} - \frac{M_p}{M_n + M_p} \vec{R}_{np} \\
&= \frac{M_a + M_n}{M_a} (\vec{R}_{bp} - \vec{R}_{np}) \\
&= \frac{(M_a + M_n)(M_n + M_p)}{M_n(M_a + M_n + M_p)} \vec{R}_{ad} - \frac{M_p(M_a + M_n)}{M_n(M_a + M_n + M_p)} \vec{R}_{bp}. \tag{VI.4}
\end{aligned}$$

A.2 Hamiltonian and wave function

The total Hamiltonian can be written in two alternative forms. For the deuteron channel :

$$H = H_a - a_{np} \Delta_{R_{np}} |R + V_{np}(\vec{R}_{np}) - a_d \Delta_R |R_{np} + V_{an}(\vec{R}_{an}) + V_{ap}(\vec{R}_{ap}) \tag{VI.5}$$

For the proton channel :

$$H = H_b - a_p \Delta_r |R_{np} + V_{ap}(\vec{R}_{ap}) + V_{np}(\vec{R}_{np}) \tag{VI.6}$$

with :

$$H_b = H_a - a_n \Delta_{R_{np}} |r + V_{an}(\vec{R}_{an}) \tag{VI.7}$$

Here the a 's are related to the reduced masses (or the reduced energies in the relativistic option). They are:

$$\begin{aligned}
a_d &= \frac{M_a + M_n + M_p}{2M_a(M_n + M_p)} \hbar^2, & a_{np} &= \frac{M_n + M_p}{2M_n M_p} \hbar^2, \\
a_p &= \frac{M_a + M_n + M_p}{2M_p(M_a + M_n)} \hbar^2, & a_n &= \frac{M_a + M_n}{2M_a M_n} \hbar^2. \tag{VI.8}
\end{aligned}$$

The intrinsic wave functions of the target A, the target B and the particle in the initial channel are respectively $F_a(\vec{x})$, $F_b(\vec{x}, \vec{R}_{an})$, $F_d(\vec{R}_{np})$ where \vec{x} are internal coordinates. They satisfy the equations :

$$\begin{aligned}
H_a F_a(\vec{x}) &= E_a F_a(\vec{x}), \\
H_b F_b(\vec{x}, \vec{R}_{an}) &= E_b F_b(\vec{x}, \vec{R}_{an}), \\
\left[-a_{np} \Delta_{R_{np}} + V_{np}(\vec{R}_{np}) \right] F_d(R_{np}) &= E_d F_d(R_{np}). \tag{VI.9}
\end{aligned}$$

Furthermore, we assume that :

$$\begin{aligned}
F_a(\vec{x}, \vec{R}_{an}) &= F_a(\vec{x}) F_n(\vec{R}_{an}), \\
\left[-a_n \Delta_{R_{an}} + V_{an}(\vec{R}_{an}) \right] F_n(\vec{R}_{an}) &= E_n F_n(\vec{R}_{an}). \tag{VI.10}
\end{aligned}$$

The total wave function is :

$$F(\vec{x}, \vec{R}_{an}, \vec{R}_{bp}) = F_a(\vec{x})F_d(\vec{R}_{np})F_1(\vec{R}) + F_b(\vec{x}, \vec{R}_{an})F_2(\vec{r}) \quad (\text{VI.11})$$

B System of coupled equations

The system of coupled equations is obtained from :

$$[H - E] F(\vec{x}, \vec{R}_{an}, \vec{R}_{bp}) = 0 \quad (\text{VI.12})$$

by projection on $F_a(\vec{x})F_d(\vec{R}_{np})$ and on $F_b(\vec{x}, \vec{R}_{an})$. In the first case, the total expression is integrated on the variables \vec{x} and \vec{R}_{np} and the result is a function of $\vec{R}_{ad} = \vec{R}$. In the second case, the total expression is integrated on \vec{x} and \vec{R}_{an} and the result is a function of $\vec{R}_{bp} = \vec{r}$.

B.1 Tri-dimensional equations

The first equation obtained is :

$$\begin{aligned} & -a_d(\Delta_R + k_d^2)F_1(\vec{R}) + U_{ad}(\vec{R})F_1(\vec{R}) \\ & + \int F_d(\vec{R}_{np}) \left[-a_p(\Delta_r|_{R_{an}} + k_p^2) + V_{ap}(\vec{R}_{ap}) + V_{np}(\vec{R}_{np}) \right] F_n(\vec{R}_{an})F_2(\vec{r})d\vec{R}_{np} = 0 \end{aligned} \quad (\text{VI.13})$$

with :

$$U_{ad}(\vec{R}) = \int F_d(\vec{R}_{np}) \left[V_{an}(\vec{R}_{an}) + V_{ap}(\vec{R}_{ap}) \right] F_d(\vec{R}_{np})d\vec{R}_{np} \quad (\text{VI.14})$$

The second equation is :

$$\begin{aligned} & -a_p(\Delta_r + k_p^2)F_2(\vec{r}) + U_{bp}(\vec{r})F_2(\vec{r}) \\ & + \int F_n(\vec{R}_{an}) \left[-a_p(\Delta_r|_{R_{an}} + k_p^2) + V_{ap}(\vec{R}_{ap}) + V_{np}(\vec{R}_{np}) \right] F_d(\vec{R}_{np})F_1(\vec{R})d\vec{R}_{an} = 0 \end{aligned} \quad (\text{VI.15})$$

with :

$$U_{bp}(\vec{r}) = \int F_n(\vec{R}_{an}) \left[V_{np}(\vec{R}_{np}) + V_{ap}(\vec{R}_{ap}) \right] F_n(\vec{R}_{an})d\vec{R}_{an} \quad (\text{VI.16})$$

In these integro-differential equations, all the variables can be expressed in terms of \vec{R} and \vec{r} : this introduces the Jacobian of the transformation from $\vec{R}_{ad}, \vec{R}_{np}$ or $\vec{R}_{bp}, \vec{R}_{an}$ to $\vec{R}_{ad}, \vec{R}_{bp}$ in front of these integrals. These two Jacobians are :

$$J = \left[\frac{(M_n + M_p)(M_a + M_n)}{M_n(M_a + M_n + M_p)} \right]^3 \quad (\text{VI.17})$$

We obtain:

$$\begin{aligned} & -a_d(\Delta_R + k_d^2)F_1(\vec{R}) + U_{ad}(\vec{R})F_1(\vec{R}) + \int K(\vec{R}, \vec{r})F_2(\vec{r})d\vec{r} = 0, \\ & -a_p(\Delta_r + k_p^2)F_2(\vec{r}) + U_{bp}(\vec{r})F_2(\vec{r}) + \int K(\vec{R}, \vec{r})F_1(\vec{R})d\vec{R} = 0 \end{aligned} \quad (\text{VI.18})$$

with :

$$K(\vec{R}, \vec{r}) = J F_d(\vec{R}_{np}) \left[-a_p(\Delta_r|_{R_{an}} + k_p^2) + V_{ap}(\vec{R}_{ap}) + V_{np}(\vec{R}_{np}) \right] F_n(\vec{R}_{an}) \quad (\text{VI.19})$$

The derivative term can be replaced by derivation of the bound functions $F_d(\vec{R}_{np})$ and $F_n(\vec{R}_{an})$ (see Ref. [72]). The result is :

$$K(\vec{R}, \vec{r}) = J \left[-a_p(\Delta_r|_R + k_p^2) + V_{ap}(\vec{R}_{ap}) + V_{np}(\vec{R}_{np}) \right] F_n(\vec{R}_{an})F_d(\vec{R}_{np}) \quad (\text{VI.20})$$

and can be divided into a non-orthogonality term :

$$K_N(\vec{R}, \vec{r}) = J \left[-a_p(\Delta_r |R + k_p^2) + U_{bp}(\vec{R}_{bp}) \right] F_n(\vec{R}_{an}) F_d(\vec{R}_{np}) \quad (\text{VI.21})$$

which do not exist in inelastic scattering and :

$$K_V(\vec{R}, \vec{r}) = J \left[V_{ap}(\vec{R}_{ap}) + V_{np}(\vec{R}_{np}) - U_{bp}(\vec{R}_{bp}) \right] F_n(\vec{R}_{an}) F_d(\vec{R}_{np}) \quad (\text{VI.22})$$

which is the only term which remains in **DWBA** because $K_N(\vec{R}, \vec{r})$ is identical to the homogeneous part of the equation for $F_2(\vec{r})$, multiplied by some function.

B.2 Radial equations

If the waves functions $F_1(\vec{R})$ and $F_2(\vec{r})$ are expanded into partial waves, a set of radial equations are obtained. The transition potential $K(\vec{R}, \vec{r})$ is characterised by :

- the transfer of an angular momentum \vec{L} which is the angular momentum \vec{L}_n of the “neutron” in $F_n(\vec{R}_{an})$,
- the transfer of a spin to the particle, which is the one of the “neutron”,
- the transfer of a total spin J to the nucleus, which is the total momentum of the “neutron” wave function $F_n(\vec{R}_{an})$.

The geometrical coefficients are the one implied by these three transfers. Note, in **ECIS**, the phase i^{L_n} and the extra phase i for odd values of L_n .

Without taking into account the spins, the wave functions F_1 and F_2 can be written :

$$\begin{aligned} F_1(\vec{R}) &= \sum_L i^L Y_L^M(\hat{R}) \frac{1}{R} F_1(L, R), \\ F_2(\vec{r}) &= \sum_l i^l Y_l^m(\hat{r}) \frac{1}{r} F_1(l, r). \end{aligned} \quad (\text{VI.23})$$

The system of coupled equations is :

$$\begin{aligned} &-a_d \left[\frac{d^2}{dR^2} + k_d^2 - \frac{L_i(L_i + 1)}{R^2} \right] F_1(L_i, R) + \sum_{L_j} U_{ad}(L_i, L_j, R) F_1(L_j, R) \\ &+ \sum_{l_j} \int G_{L_i, l_j} R r K(R, r) F_2(l_j, r) dr = 0 \\ &-a_p \left[\frac{d^2}{dr^2} + k_p^2 - \frac{l_i(l_i + 1)}{r^2} \right] F_2(l_i, r) + \sum_{l_j} U_{bp}(l_i, l_j, r) F_2(l_j, r) \\ &+ \sum_{L_j} \int G_{l_i, L_j} R r K(R, r) F_1(L_j, R) dR = 0 \end{aligned} \quad (\text{VI.24})$$

where $G_{L, l}$ is some geometrical coefficient.

C Zero-range approximation

We neglect the non-orthogonality term $K_N(\vec{R}, \vec{r})$ and we assume $V_{ap}(\vec{R}_{ap}) = U_{bp}(\vec{R}_{bp})$ in $K_V(\vec{R}, \vec{r})$. The zero-range approximation is obtained if we replace the product $V_{np}(\vec{R}_{np}) F_d(\vec{R}_{np})$ by a δ -function of \vec{R}_{np} .

C.1 Coupled differential equations

In the interaction :

$$K(\vec{R}, \vec{r}) = J V_{np}(\vec{R}_{np}) F_d(\vec{R}_{np}) F_n(\vec{R}_{an}) \quad (\text{VI.25})$$

the product $V_{np}(\vec{R}_{np}) F_d(\vec{R}_{np})$ is replaced by a δ -function of :

$$\vec{R}_{np} = -\frac{M_a(M_n + M_p)}{M_n(M_a + M_n + M_p)} \vec{R}_{ad} + \frac{(M_a + M_n)(M_n + M_p)}{M_n(M_a + M_n + M_p)} \vec{R}_{bp} \quad (\text{VI.26})$$

multiplied by some strength C_0 which is usually -122.5 meV (124.7 MeV in Ref. [72]). The δ -function can be replaced by a δ -function on the lengths divided by :

$$\frac{M_a(M_n + M_a)(M_n + M_p)^2 Rr}{[M_n(M_a + M_n + M_p)]^2} \quad (\text{VI.27})$$

with a δ -function on directions of R and r . So :

$$K(\vec{R}, \vec{r}) = C_0 \frac{(M_n + M_p)(M_a + M_n)^2}{M_a M_n (M_a + M_n + M_p)} F_n(\vec{R}_{an}) \frac{\delta(R_{np})}{Rr} \quad (\text{VI.28})$$

and after integration the system of coupled equations becomes :

$$\begin{aligned} & -a_d \left[\frac{d^2}{dR^2} + k_d^2 - \frac{L_i(L_i + 1)}{R^2} \right] F_1(L_i, R) + \sum_{L_j} U_{ad}(L_i, L_j, R) F_1(L_j, R) \\ & + \sum_{l_j} C_0 G_{L_i, l_j} \frac{M_a + M_n}{M_a} F_n(R) F_2(l_j, \frac{M_a}{M_a + M_n} R) = 0, \\ & -a_p \left[\frac{d^2}{dr^2} + k_p^2 - \frac{l_i(l_i + 1)}{r^2} \right] F_2(l_i, r) + \sum_{l_j} U_{bp}(l_i, l_j, r) F_2(l_j, r) \\ & + \sum_{L_j} C_0 G_{l_i, L_j} \left(\frac{M_a + M_n}{M_a} \right)^2 F_n\left(\frac{M_a + M_n}{M_a} r \right) F_1(L_j, \frac{M_a + M_n}{M_a} r) = 0. \end{aligned} \quad (\text{VI.29})$$

These equations are not symmetric. Note that M_a and $M_a + M_p$ are the masses of the target for the two levels. From now on, we shall use:

$$M_1 = M_a \quad M_2 = M_a + M_n \quad (\text{VI.30})$$

C.2 Symmetric equations

To obtain more symmetric equations, we shall use different units of lengths in each channel: in channel i , we use $s = M_i R_i / M_0$, that is the step size $h_i = h M_0 / M_i$, where M_0 is the mass of the target in some first channel labelled by 0. The equations are :

$$\begin{aligned} & -a_d \left(\frac{M_1}{M_0} \right)^2 \left[\frac{d^2}{ds^2} + \left(\frac{M_0}{M_1} \right)^2 k_d^2 - \frac{L_i(L_i + 1)}{s^2} \right] F_1(L_i, s) + \sum_{L_j} U_{ad}(L_i, L_j, s) F_1(L_j, s) \\ & + \sum_{l_j} C_0 G_{L_i, l_j} \frac{M_2}{M_1} F_n(s) F_2(l_j, s) = 0, \\ & -a_p \left(\frac{M_2}{M_0} \right)^2 \left[\frac{d^2}{ds^2} + \left(\frac{M_0}{M_2} \right)^2 k_p^2 - \frac{l_i(l_i + 1)}{s^2} \right] F_2(l_i, s) + \sum_{l_j} U_{bp}(l_i, l_j, s) F_2(l_j, s) \\ & + \sum_{L_j} C_0 G_{l_i, L_j} \left(\frac{M_2}{M_1} \right)^2 F_n(s) F_1(L_j, s) = 0. \end{aligned} \quad (\text{VI.31})$$

With the following notations :

$$V_{11} = U_{ad}(s), \quad V_{12} = GC_0 \left[\frac{M_2}{M_1} \right]^{\frac{3}{2}} F_n(s), \quad V_{22} = U_{bp}(s), \quad (\text{VI.32})$$

and using the coefficients :

$$A_1 = \left[\frac{M_0}{M_1 a_d} \right]^{\frac{1}{2}} \quad A_2 = \left[\frac{M_0}{M_2 a_p} \right]^{\frac{1}{2}} \quad B_1 = \frac{M_0}{M_1 a_d^{\frac{1}{2}}} \quad B_2 = \frac{M_0}{M_1 a_p^{\frac{1}{2}}} \quad (\text{VI.33})$$

wave functions and potentials can be redefined with :

$$F_i(s) = A_i G_i(s) \quad U_{ij}(s) = B_i B_j V_{ij}(s) \quad (\text{VI.34})$$

to get the symmetric set of equations :

$$\begin{aligned} -\left[\frac{d^2}{ds^2} + \left(\frac{M_0}{M_1} \right)^2 k_d^2 - \frac{L(L+1)}{s^2} \right] G_1(s) + U_{11}(s)G_1(s) + U_{12}(s)G_2(s) &= 0 \\ -\left[\frac{d^2}{ds^2} + \left(\frac{M_0}{M_2} \right)^2 k_p^2 - \frac{l(l+1)}{s^2} \right] G_2(s) + U_{22}(s)G_2(s) + U_{12}(s)G_1(s) &= 0 \end{aligned} \quad (\text{VI.35})$$

in which $[M_0/M_1]k_d$ and $[M_0/M_2]k_p$ appear instead of the wave numbers. In **ECIS**, the mass of the target in the first channel is chosen as M_0 . In the subroutine **COLF** the wave number is multiplied by M_0/M_I , because this product plays the same role in all parts of the code.

D Use of ECIS94

D.1 Possibilities of the code

The code **ECIS94** as well as **ECIS79** allows zero-range calculations without the non-orthogonality term using external nuclear matrix elements and form factors (**LO(7)=.TRUE.**). The step sizes for each levels are related with respect to the difference of masses, unless **tt LO(93)=.TRUE.** (**LO(48)=.TRUE.** in **ECIS79**): this control can be used to interchange particle and target in some level, in which case there is no recoil correction. For the transfer of a neutron with angular momentum L_n and total spin J_n , the data needed for the reduced matrix elements are :

- **columns** 1- 5 Sequence number or blank.
- **columns** 6-10 angular momentum L_n .
- **columns** 11-15 1 (twice the spin of the neutron).
- **columns** 16-20 $2J_n$ twice the total spin.
- **columns** 21-40 the nuclear matrix element which depends upon the form factor, but includes :
 - a reduced nuclear matrix element in the spin space which is not the same in **ECIS79** and **ECIS94**,
 - a reduced nuclear matrix element in the target space,
 - a recoil factor $[(a+1)/a]^{3/2}$ if **LO(48)=.FALSE.**

The form factor can be given by points. However, they are two “standard” possibilities:

- 7 Laguerre polynomial,
- 8 solution in a real Woods-Saxon potential of which the depth is searched for a given binding energy.

In these two cases, there is a discrepancy in the code between these form factors and the macroscopic ones such that the reduced nuclear matrix element has to be divided by $\sqrt{4\pi}$.

D.2 Example

The following example has been chosen to compare **ECIS79** with the code **CHUCK**. The original test case was intended for *Zr* but the mass of the target has been decreased to enforce the recoil effects.

For **ECIS94**, this test case is :

```

|COMPARAISON AVEC CHUCK POUR BE8
|FFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFF
|TTTTTFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFFF
|
|      2      1      2      3      3
|.12      18.
|      15.      1.      2.      8.      4.
|3.5      2 -5.      .5      1.      9.      4.
|0.      2.5      137.5
|
|      1      1
|      1      2      1
|      1      4      1      7      142.84024
|      2      2
|      1      1      0      1      0      0      0      -1
|104.3      2.14      .78
|      1      1      0      2      0      0      0      -1
|
|      1      1      0      3      0      0      0      -1
|
|      1      1      0      4      0      0      0      -1
|13.9      2.52      .8
|      1      1      0      5      0      0      0      -1
|3.5      2.14      .78
|      1      1      0      6      0      0      0      -1
|
|      1      1      0      7      0      0      0      -1
|4.      2.5
|      1      1      0      8      0      0      0      -1
|
|      2      2      0      1      0      0      0      -1
|54.9      2.3713      .7
|      2      2      0      2      0      0      0      -1
|.41      2.7249      .63
|      2      2      0      3      0      0      0      -1
|
|      2      2      0      4      0      0      0      -1
|8.7875      2.7249      .63
|      2      2      0      5      0      0      0      -1
|8.3825      2.609      .53
|      2      2      0      6      0      0      0      -1
|
|      2      2      0      7      0      0      0      -1
|4.      2.6001
|      2      2      0      8      0      0      0      -1
|
|      1      2      1      1      0      0      0      -8
|7.2      9.      1.      0.      0.      1.25      .65
|0.      1.      1.      1.
|FIN

```

There is no spin-orbit for the bound wave-functions because **CHUCK** takes is proportional to the central

potential and **ECIS** keeps it constant. As the first channel is the deuteron channel, there is no need to indicate to the code to use the deuteron step size : for the inverse reaction, a “2” should be in column 45 of the third card before **FIN**. The “deformation” in the test with the code **CHUCK** was 122.5 . The value (1) used here is the same

- divided by $\sqrt{4\pi}$, due to the normalisation of bound-states in **ECIS**,
- multiplied by the reduced matrix element in the spin space which is $\sqrt{3}/2$ in **ECIS94** but was $\sqrt{3}$ for **ECIS79**,
- multiplied by the reduced matrix element in the target space, which is $\sqrt{8}$,
- multiplied by the recoil factor which is $(9./8.)^{3/2}$.

Chapter VII

Miscellaneous

A Relativistic cinematics

This section is to explain some expressions which are used in the programme when relativistic cinematics are used (`LO(8)=.TRUE.`) and compare then to the non-relativistic ones (`LO(8)=.FALSE.`). The code includes in subroutine `CALC` the following values :

	<code>CM=931.5017646D0</code>	<code>CALC-085</code>
	<code>CHB=197.328604D0</code>	<code>CALC-086</code>
	<code>CZ=137.0360411D0</code>	<code>CALC-087</code>

in the `IBM` version (single precision in the `CDC` version). They are the values of the Atomic Mass Unit M_u in MeV fm/c² and of $\hbar c$ in MeV fm respectively where c is the speed of light in vacuum. In these explanations, we supposes the masses multiplied by `CM` and the wave numbers by `CHB`.

A.1 Notations

Let us consider a particle of mass m_i and energy E_{lab} in the Laboratory system and a target of mass M_i ; in the final state with an excitation energy Q , an outgoing particle of mass m_f and a residual nucleus of mass M_f :

- in non-relativistic cinematics, the mass M_f is unaltered;
- in relativistic cinematics :

$$M_f = m_i + M_i - m_f + Q/M_u \quad (\text{VII.1})$$

in such a way that the total energy in the center of mass system is a constant. However, this correction is dropped in the Schrödinger formalism if they are Coulomb corrections.

A.2 Energy in the center of mass system

The relativistic energy of the incoming particle in the Laboratory System is given by $e_i = m_i + E_{lab}$, as a pure extension of the non-relativistic case. Its momentum p_i is such that $e_i^2 = m_i^2 + p_i^2$. A change of frame along the quantification axis is obtained by :

$$P = p \cosh x + e \sinh x, \quad E = p \sinh x + e \cosh x \quad (\text{VII.2})$$

The change to the Center of mass system is obtained with :

$$P_{cm} = p_i \cosh x + e_i \sinh x + M_i \sinh x = 0, \quad E_{cm} = p_i \sinh x + e_i \cosh x + M_i \cosh x \quad (\text{VII.3})$$

from which $\sinh x$ and $\cosh x$ can be eliminated to obtain :

$$E_{cm_i} = E_{cm} - m_i - M_i = [(m_i + M_i)^2 + 2E_{lab}M_i]^{\frac{1}{2}}, \quad E_{cm_f} = E_{cm_i} - Q/M_u \quad (\text{VII.4})$$

for each channel.

A.3 Wave numbers - Momenta

From :

$$E_{cm} = (m_f^2 + p_f^2)^{\frac{1}{2}} + (M_f^2 + p_f^2)^{\frac{1}{2}} \quad (\text{VII.5})$$

we get :

$$\begin{aligned} p_f^2 &= \frac{[(E_{cm} - m_f - M_f)(E_{cm} + m_f + M_f)(E_{cm} + m_f - M_f)(E_{cm} - m_f + M_f)]}{4E_{cm}^2} \\ &= \frac{(E_{cm_f}(E_{cm_f} + 2m_f + 2M_f)(E_{cm_f} + 2m_f)(E_{cm_f} + 2M_f))}{4(E_{cm_i} + m_i + M_i)^2} \end{aligned} \quad (\text{VII.6})$$

With relativistic cinematics without Dirac formalism, the reduced mass μ_{nr} is replaced by the reduced energy μ_r :

$$\mu_{nr} = \frac{m_f M_f}{m_f + M_f}, \quad \mu_r = \frac{(m_f^2 + p_f^2)^{\frac{1}{2}}(M_f^2 + p_f^2)^{\frac{1}{2}}}{E_{cm}} = \frac{E_{cm}^4 - (m_f^2 - M_f^2)^2}{4E_{cm}^3}, \quad (\text{VII.7})$$

but, with Coulomb corrections, M_i is used instead of M_f in order to have the same reduced mass for all the levels : until their generalisation, the Coulomb corrections are valid only with a product of the wave number by the Coulomb parameter constant in all the channels.

Without relativistic cinematics, the residual target mass M_f is not corrected by the excitation energy. The center of mass energy of each levels are given by :

$$E_{cm_i} = \frac{E_{lab}M_i}{m_i + M_i}, \quad E_{cm_f} = E_{cm_i} - Q, \quad p_f^2 = \frac{2M_u}{\hbar^2} \frac{E_{cm_f}m_f M_f}{m_f + M_f}. \quad (\text{VII.8})$$

A.4 Observables in the laboratory system

We shall note the incoming particle, the initial target, the outgoing particle and the residual target as particles 1, 2, 3 and 4 respectively. In order to compute the rotation to be applied to the S matrix to obtain a description of polarisation phenomena in the Laboratory system, we have to consider the following frames :

1. the center of mass with Oz in incoming direction;
2. the center of mass with Oz in outgoing direction of particle 3;
3. the system in which particle 3 is at rest and Oz opposite to outgoing direction of particle 4; the description of polarisations is invariant in this change of frame;
4. the system in which particle 3 is at rest and Oz opposite to incoming particle 2; the description of polarisations has to be rotated for this change of frame;
5. the system where particle 2 is at rest and Oz along the outgoing particle 3; the polarisations are invariant;
6. the system where particle 2 is at rest and Oz along the incoming particle m_i ; the description in this frame can be obtained easily from the description in frame (1).

We note by p_n the momentum of the particle n in the center of mass frame; by e_n their relativistic energy and by m_n their mass. In fact, $p_1 = p_2$ and $p_3 = p_4$. We shall give z , x and time components of the quadri-vector (\vec{p}, e) for each particle in each frame.

A.4.a Center of mass system - Oz incoming direction

For an outgoing particle 3 at the angle α in the center of mass system, the 4 quadri-vectors are:

$$(1) \begin{vmatrix} p_1 \\ 0 \\ e_1 \end{vmatrix} \quad (2) \begin{vmatrix} -p_1 \\ 0 \\ e_2 \end{vmatrix} \quad (3) \begin{vmatrix} p_3 \cos \alpha \\ p_3 \sin \alpha \\ e_3 \end{vmatrix} \quad (4) \begin{vmatrix} -p_3 \cos \alpha \\ -p_3 \sin \alpha \\ e_4 \end{vmatrix} \quad (\text{VII.9})$$

A.4.b Center of mass system - Oz outgoing direction

This new frame is obtained by a rotation in the ordinary space :

$$(1) \begin{vmatrix} p_1 \cos \alpha \\ -p_1 \sin \alpha \\ e_1 \end{vmatrix} \quad (2) \begin{vmatrix} -p_1 \cos \alpha \\ p_1 \cos \alpha \\ e_2 \end{vmatrix} \quad (3) \begin{vmatrix} p_3 \\ 0 \\ e_3 \end{vmatrix} \quad (4) \begin{vmatrix} -p_3 \\ 0 \\ e_4 \end{vmatrix} \quad (\text{VII.10})$$

A.4.c Particle 3 rest system - Oz outgoing direction

This new frame is obtained by a special Lorentz transformation with $\cosh x = e_3/m_3$ and $\sinh x = -p_3/m_3$:

$$(1) \begin{vmatrix} (p_1 e_3 \cos \alpha - e_1 p_3)/m_3 \\ -p_1 \sin \alpha \\ (e_1 e_3 - p_1 p_3 \cos \alpha)/m_3 \end{vmatrix} \quad (2) \begin{vmatrix} -(p_1 e_3 \cos \alpha + e_2 p_3)/m_3 \\ p_1 \sin \alpha \\ (e_2 e_3 + p_1 p_3 \cos \alpha)/m_3 \end{vmatrix} \quad (3) \begin{vmatrix} 0 \\ 0 \\ m_3 \end{vmatrix} \quad (4) \begin{vmatrix} -p_3(e_3 + e_4)/m_3 \\ 0 \\ (e_3 e_4 + p_3^2)/m_3 \end{vmatrix} \quad (\text{VII.11})$$

A.4.d Particle 3 rest system - Oz opposite to particle 2

The axis Oz must be opposite to the momentum of particle 2. This rotation β is such that :

$$\begin{aligned} \sin \beta &= -\frac{\sin \alpha}{y} & \cos \beta &= \frac{e_3 p_1 \cos \alpha + e_2 p_3}{p_1 m_3 y} \\ y^2 &= \frac{m_3^2 p_1^2 + p_3^2 p_1^2 \cos^2 \alpha + 2 p_1 p_3 e_2 e_3 \cos \alpha + p_3^2 e_2^2}{p_1^2 m_3^2} \end{aligned} \quad (\text{VII.12})$$

The expressions obtained are :

$$(1) \begin{vmatrix} z_{14} \\ -(e_1 + e_2) p_1 \sin \alpha / (m_3 y) \\ (e_1 e_3 - p_1 p_3 \cos \alpha) / m_3 \end{vmatrix} \quad (2) \begin{vmatrix} -p_1 y \\ 0 \\ (e_2 e_3 + p_1 p_3 \cos \alpha) / m_3 \end{vmatrix} \quad (3) \begin{vmatrix} 0 \\ 0 \\ m_3 \end{vmatrix} \quad (4) \begin{vmatrix} z_{44} \\ -(e_3 + e_4) p_3 \sin \alpha / (m_3 y) \\ (e_3 e_4 + p_3^2) / m_3 \end{vmatrix} \quad (\text{VII.13})$$

where :

$$z_{14} = y p_1 - \frac{(e_1 + e_2) p_3 (e_3 p_1 \cos \alpha + e_2 p_3)}{m_3^2 p_1 y}, \quad z_{44} = -\frac{(e_3 + e_4) p_3 (e_3 p_1 \cos \alpha + e_2 p_3)}{m_3^2 p_1 y}. \quad (\text{VII.14})$$

The angle β is the angle of transformation of the scattering matrix in the helicity formalism. In the non-relativistic case, $e_3 = m_3$ and $e_2 = m_2$; the angle β is then the angle in the Laboratory system minus the angle in the Center of mass system. For the target, the quantification axis is opposite to the momentum in the Center of mass system: in the formulae above, p_3 must be changed into $-p_3$, e_3 and m_3 in e_4 and m_4 respectively.

A.5 Angles in the laboratory system

The transformations can be continued :

A.5.a Laboratory system - Oz outgoing direction

This frame is obtained by a special Lorentz transformation with :

$$\cosh x = \frac{e_2 e_3 + p_1 p_3 \cos \alpha}{m_2 m_3}, \quad \sinh x = \frac{y p_1}{m_2} \quad (\text{VII.15})$$

The expressions obtained are :

$$(1) \begin{vmatrix} z_{15} \\ -(e_1 + e_2)p_3 \sin \alpha / (m_3 y) \\ (e_1 e_2 + p_1^2) / m_2 \end{vmatrix} \quad (2) \begin{vmatrix} 0 \\ 0 \\ m_2 \end{vmatrix} \quad (3) \begin{vmatrix} m_3 p_1 y / m_2 \\ 0 \\ (e_2 e_3 + p_1 p_3 \cos \alpha) / m_2 \end{vmatrix} \quad (4) \begin{vmatrix} z_{45} \\ -(e_3 + e_4)p_3 \sin \alpha / (m_3 y) \\ (e_2 e_4 - p_1 p_3 \cos \alpha) / m_2 \end{vmatrix} \quad (\text{VII.16})$$

where :

$$z_{15} = \frac{(e_1 + e_2)(e_2 p_3 \cos \alpha + e_3 p_1)}{m_2 m_3 y}, \quad z_{45} = \frac{p_1 p_3 e_2 (e_4 - e_3) \cos \alpha + p_1^2 e_3 e_4 - p_3^2 e_2^2 + p_1^2 p_3^2 \sin^2 \alpha}{m_2 m_3 p_1 y}. \quad (\text{VII.17})$$

The scattering angle in the Laboratory system γ corresponding to the scattering angle α in the center of mass system is such that :

$$\sin \gamma = \frac{\sin \alpha}{z}, \quad \cos \gamma = \frac{e_2 p_3 \cos \alpha + e_3 p_1}{p_1 m_2 z} \\ z^2 = \frac{p_1^2 p_3^2 \cos^2 \alpha + 2 p_1 p_3 e_2 e_3 \cos \alpha + p_3^2 m_2^2 + e_3^2 p_1^2}{p_3^2 m_2^2} \quad (\text{VII.18})$$

A.5.b Laboratory system - Oz incoming direction

This frame is obtained by a rotation of angle γ in the tri-dimensional space or from the first frame by a special Lorentz transformation with $\cosh x = e_2 / m_2$ and $\sinh x = p_1 / m_2$:

$$(1) \begin{vmatrix} p_1(e_1 + e_2) / m_2 \\ 0 \\ (e_1 e_2 + p_1^2) / m_2 \end{vmatrix} \quad (2) \begin{vmatrix} 0 \\ 0 \\ m_2 \end{vmatrix} \quad (3) \begin{vmatrix} (p_3 e_2 \cos \alpha + p_1 e_3) / m_2 \\ p_3 \sin \alpha \\ (e_2 e_3 + p_1 p_3 \cos \alpha) / m_2 \end{vmatrix} \quad (4) \begin{vmatrix} (p_1 e_4 - p_3 e_2 \cos \alpha) / m_2 \\ -p_3 \sin \alpha \\ (e_2 e_4 - p_1 p_3 \cos \alpha) / m_2 \end{vmatrix} \quad (\text{VII.19})$$

The angle γ in the laboratory system is given in function of the angle α in the center of mass system by :

$$\sin \gamma = x \sin \alpha, \quad \cos \gamma = x \frac{e_2 p_3 \cos \alpha + e_3 p_1}{m_2 p_3} \quad (\text{VII.20})$$

Using :

$$A = \frac{e_2}{m_2}, \quad B = \frac{e_3 p_1}{m_2 p_3}, \quad (\text{VII.21})$$

we have :

$$\tan \gamma = \frac{\sin \alpha}{A \cos \alpha + B} \quad (\text{VII.22})$$

This relation can be inverted by :

$$\tan \alpha = \sin \gamma \frac{B \cos \gamma + AC}{C \cos \gamma - AB \sin^2 \gamma}, \quad C^2 = (A^2 - B^2) \sin^2 \gamma + \cos^2 \gamma \quad (\text{VII.23})$$

The factor multiplying the cross-section in the center of mass system to obtain the cross-section in the Laboratory system is:

$$\frac{d(\cos \alpha)}{d(\cos \gamma)} = \frac{[(A \cos \alpha + B)^2 + \sin^2 \alpha]^{\frac{3}{2}}}{A + B \cos \alpha} \quad (\text{VII.24})$$

In the non relativistic case, these formulae are simplified only by the fact that $A = 1$ because $e_2 = m = 2$.

A.6 Conclusions

If the observables are not defined in the center of mass system, the scattering matrix has to be rotated. For a scattering angle α in the center of mass system, the rotation needed to describe the polarisation of the outgoing particle is given by Equ. (VII.12), that is :

$$\tan \beta = -\frac{p_1 m_3 \sin \alpha}{e_3 p_1 \cos \alpha + e_2 p_3} \quad (\text{VII.25})$$

and the rotation needed to describe the polarisation of the target nucleus is :

$$\tan \beta = -\frac{p_1 m_4 \sin \alpha}{e_4 p_1 \cos \alpha - e_2 p_4} \quad (\text{VII.26})$$

with relativistic kinematics. In the non relativistic approximation, the energies e are replaced by the masses m in Equ. (VII.24) and (VII.26).

If the observables are defined with only one axis of quantification along the incoming direction, the collision matrix must be rotated with the angle $-\alpha$.

If an angular distribution is requested in the Laboratory system, the angle α in the center of mass system is given by Equ. (VII.23) for the angle γ in the Laboratory system. Cross-sections only must be multiplied by the factor given by Equ. (VII.24). Here also, the non relativistic approximation is obtained by replacing the energies e by the masses m .

B Coefficients 3j, 6j and 9j

Clebsch-Gordan coefficients are used only in the computation of reduced nuclear matrix elements, with the function **DJCG**. These reduced matrix elements need also some 6j computed by the function **DJ6J**. Clebsch-Gordan coefficients with magnetic quantum numbers zero or $(1/2, 0, -1/2)$ are given by the function **DCGS**. The subroutine **QUAN**, which computes the couplings of Equ. (III.44) needs the 9j coefficients given by the function **DJ9J** only for transitions with $\Delta S \neq 0$ and functions **DCGS** and **DJ6J** for all transitions.

On the contrary, the Clebsch-Gordan needed in subroutine **SCHE** to transform the results into the helicity formalism by Equ. (III.64) are obtained by recurrence.

An array of logarithms of factorials is used by the functions **DJCG**, **DJ6J** and **DCGS**.

B.1 The functions for geometric coefficients

All the arguments of these four functions are integer double values of the quantum numbers.

B.1.a Clebsch-Gordan coefficients

If one of the angular momenta is zero, the simple result is returned. If one of the magnetic quantum numbers is zero and the others 0 and $\pm \frac{1}{2}$, the simpler formula of function **DCGS** is used.

In the general case, the subroutine use the formula (21) of Appendix C in : [73] MESSIAH, A., "*Mécanique quantique*", Dunod, Paris (1960) page 910. The last term of the sum is computed. The sum is then obtained by the **DO LOOP** :

```
|C K2,K3,K4,N1,N2,N3 ARE THE ARGUMENTS OF THE FACTORIALS IN THE LAST TERMDJCG-061
|      DO 2 I=1,K                                  DJCG-062
|      A1=(K2-I)*(K3-I)*(K4-I)                     DJCG-063
|      A2=(N1+I)*(N2+I)*(N3+I)                     DJCG-064
|      2 DJCG=A4-DJCG*A2/A1                          DJCG-065
```

Exponential is used only for the first term.

B.1.b 6j coefficients

If one of the angular momenta is zero, a simple result is returned.

In the general case, the subroutine use the formula (36) of Appendix C in Ref. [73] page 915. As for Clebsch-Gordan coefficients, the last term of the sum is computed. The sum is then obtained by a similar `D0 LOOP`.

B.1.c Simple 3j coefficients

Due to the relation :

$$\begin{pmatrix} j' & J & j \\ -\frac{1}{2} & 0 & \frac{1}{2} \end{pmatrix} = -\sqrt{(2l+1)(2j+1)} \begin{pmatrix} l & l' & J \\ 0 & 0 & 0 \end{pmatrix} \begin{Bmatrix} l & l' & J \\ j' & j & \frac{1}{2} \end{Bmatrix}, \quad (\text{VII.27})$$

such a 3j coefficient can be used to compute Equ. (III.44) when $S = 0$ and the spins are $\frac{1}{2}$. This relation shows up in the helicity formalism for multipole expansion of an interaction (see Ref. [62]).

The function `DCGS` computes :

$$CGS = (-)^{j'+\frac{1}{2}} \sqrt{(2j+1)(2j'+1)} \begin{pmatrix} j & L & j' \\ -\frac{1}{2} & 0 & \frac{1}{2} \end{pmatrix} \quad (\text{VII.28})$$

if j and j' are half integers and :

$$CGS = (-)^{j'+1} \begin{pmatrix} j & L & j' \\ 0 & 0 & 0 \end{pmatrix} \quad (\text{VII.29})$$

if j and j' are integers.

Using :

$$g(n) = \frac{m!}{\sqrt{(n!)}} = \frac{1}{2^m} \begin{cases} \sqrt{\frac{2.4.6\dots n}{3.5.7\dots(n-1)}} & \text{if } n \text{ is even} \\ \sqrt{\frac{2.4.6\dots(n-1)}{3.5.7\dots n}} & \text{if } n \text{ is odd} \end{cases} \quad (\text{VII.30})$$

where m is the integer part of $n/2$, these coefficients are :

$$CGS = (-)^{\text{Int}\{(L+j-j'+2)/2\}} \frac{g(L+j+j'+1)}{g(L+j-j')g(L-j+j')g(j+j'-L)} \quad (\text{VII.31})$$

If j and j' are integer, this result must be :

- multiplied by 2 if $L+j+j'$ is even,
- 0 if $L+j+j'$ is odd.

These two conditions are $\text{Int}\{(L+j+j'+1)/2\} - \text{Int}\{(L+j-j')/2\} - \text{Int}\{(L-j+j')/2\} - \text{Int}\{(j+j'-L)/2\}$ positive or zero respectively. If this quantity is negative, j and j' are half integers.

B.1.d 9j coefficients

The subroutine use the formula (41) of Appendix C in Ref. [73] page 917, which is :

$$\begin{Bmatrix} j_1 & j_2 & j_3 \\ j_4 & j_5 & j_6 \\ j_7 & j_8 & j_9 \end{Bmatrix} = \sum_x (-)^{2x} \begin{Bmatrix} j_1 & j_2 & j_3 \\ j_6 & j_9 & x \end{Bmatrix} \begin{Bmatrix} j_4 & j_5 & j_6 \\ j_2 & x & j_8 \end{Bmatrix} \begin{Bmatrix} j_7 & j_8 & j_9 \\ x & j_1 & j_4 \end{Bmatrix} \quad (\text{VII.32})$$

A cyclic permutation of the lines is done to get the shortest sum on x in Equ. (VII.32). Then, the three series of $6j$ coefficients are computed by recurrence. For the first of them :

$$(x+1)A(x) \begin{Bmatrix} j_1 & j_2 & j_3 \\ j_6 & j_9 & x-1 \end{Bmatrix} + B(x) \begin{Bmatrix} j_1 & j_2 & j_3 \\ j_6 & j_9 & x \end{Bmatrix} + xA(x+1) \begin{Bmatrix} j_1 & j_2 & j_3 \\ j_6 & j_9 & x+1 \end{Bmatrix} = 0 \quad (\text{VII.33})$$

with :

$$\begin{aligned} A(x) &= \left\{ (j_1 - j_9 - x)(j_1 - j_9 + x)(j_1 + j_9 - x + 1)(j_1 + j_9 + x + 1)(j_2 - j_6 - x)(j_2 - j_6 + x) \right. \\ &\quad \left. (j_2 + j_6 - x + 1)(j_2 + j_6 + x + 1) \right\}^{\frac{1}{2}} \\ B(x) &= (2x+1) \left\{ (j_1 - j_9)(j_1 + j_9 + 1)(j_6 - j_2)(j_2 + j_6 + 1) + [j_1(j_1 + 1) + j_2(j_2 + 1) \right. \\ &\quad \left. + j_6(j_6 + 1) + j_9(j_9 + 1) - 2j_3(j_3 + 1) - x(x+1)]x(x+1) \right\} \end{aligned} \quad (\text{VII.34})$$

This is not a relation between “contiguous” $6j$ coefficients as defined in : [74] RAYNAL, J., *On the definition and properties of generalized 6-j symbols*, J. Math. Phys. **20** (1979) page 2398, but can be obtained by writing three such relations and eliminating two $6j$ coefficients between them. This relation is used to compute the $6j$ coefficients for $x_m \leq x \leq x_M$ with $x_m = \max\{|j_1 - j_2|, |j_6 - j_9|\}$ and $x_M = \min\{j_1 + j_2, j_6 + j_9\}$ starting from $x = x_M$ with the value 1. The first recurrence reduces to a two terms relations because the coefficient of the $6j$ coefficient for $x = x_M + 1$, which does not exist, vanishes. The sum N_1 of the values obtained multiplied by $(2x+1)$ is computed; if it is larger than 10^{15} , all the values obtained before are divided by 10^{15} and N_1 is divided by 10^{30} . This method needs a working array to store all these unnormalised $6j$ coefficients and no table of logarithm. As used in **ECIS**, the series of $6j$ coefficients are short because two of the three quantum numbers L , S or I^t are in each of them.

The result of Equ. (VII.32) must be normalised by dividing by the square root of the product $N_1 N_2 N_3$. As the sign of the $6j$ coefficient with $x = x_M$ is $(-)^{2x_M}$ as can be seen in formula (36) in Ref. [73] when the sum reduces to only one term and as all the three x_M and x are integer or half integer at the same time, the product of all these signs is plus. The result has the correct sign.

B.2 Recurrence for 3j coefficients

Two different recurrence relations are used in the code **ECIS94** for two very different purposes.

B.2.a Transformation to helicity formalism

To use conveniently Equ. (III.64), for a given total angular momentum J and parity π , one needs a tri-dimensional array of which the dimensions correspond to :

1. all sets of quantum numbers (l_i, s_i, j_i, I_i) or (l_f, s_f, f_f, I_f) involved for this value of J and π ,
2. all the $(2s+1)$ Clebsch-Gordan $\langle l, s, 0, \sigma | j, \sigma \rangle$ for these values of (l, s, j) ,
3. all the $(2I+1)$ Clebsch-Gordan $jI\sigma - \mu JM$ for these values of (j, σ, I) .

In practice, this array is built with the dimensions in the order (2),(3),(1).

The Clebsch-Gordan coefficients are ordered by increasing helicity. They can be obtained by elementary recurrence relations between “contiguous” coefficients as defined in : [75] RAYNAL, J., *On the definition and properties of generalized 3-j symbols*, J. Math. Phys. **19** (1978) page 467. More explicitly, a part of the Equ. (A7) in Ref. [74] is :

$$C^{te} = \left\{ -\frac{1}{2}[a(a+1) + b(b+1) - c(c+1)] - \alpha\beta - \frac{1}{3}(\alpha - \beta) \right\} \begin{pmatrix} a & b & c \\ \alpha & \beta & \gamma \end{pmatrix}$$

$$\begin{aligned}
& -\left\{(a+\alpha)(a-\alpha+1)(b-\beta)(b+\beta+1)\right\}^{\frac{1}{2}} \begin{pmatrix} a & b & c \\ \alpha-1 & \beta+1 & \gamma \end{pmatrix} \\
= & \left\{\frac{1}{2}[a(a+1)+b(b+1)-c(c+1)]+\alpha\beta-\frac{1}{3}(\alpha-\beta)\right\} \begin{pmatrix} a & b & c \\ \alpha & \beta & \gamma \end{pmatrix} \\
& +\left\{(a-\alpha)(a+\alpha+1)(b+\beta)(b-\beta+1)\right\}^{\frac{1}{2}} \begin{pmatrix} a & b & c \\ \alpha+1 & \beta-1 & \gamma \end{pmatrix} \tag{VII.35}
\end{aligned}$$

which can be simplified in this simple case into :

$$\begin{aligned}
0 = & \left\{(a+\alpha)(a-\alpha+1)(b-\beta)(b+\beta+1)\right\}^{\frac{1}{2}} \begin{pmatrix} a & b & c \\ \alpha-1 & \beta+1 & \gamma \end{pmatrix} \\
& +\left\{[a(a+1)+b(b+1)-c(c+1)]+2\alpha\beta\right\} \begin{pmatrix} a & b & c \\ \alpha & \beta & \gamma \end{pmatrix} \\
& +\left\{(a-\alpha)(a+\alpha+1)(b+\beta)(b-\beta+1)\right\}^{\frac{1}{2}} \begin{pmatrix} a & b & c \\ \alpha+1 & \beta-1 & \gamma \end{pmatrix} \tag{VII.36}
\end{aligned}$$

and give for the $\langle j, I, \sigma, -\mu | J, \sigma - \mu \rangle$ the following recurrence relation :

$$\begin{aligned}
0 = & \left\{(I-\mu)(I+\mu+1)(J-\mu+\sigma)(J+\mu-\sigma+1)\right\}^{\frac{1}{2}} \langle j, I, \sigma, -\mu-1 | J, \sigma - \mu - 1 \rangle \\
& -\left\{[I(I+1)+J(J+1)-j(j+1)]+2\mu(\sigma-\mu)\right\} \langle j, I, \sigma, -\mu | J, \sigma - \mu \rangle \\
& +\left\{(I+\mu)(I-\mu+1)(J+\mu-\sigma)(J-\mu+\sigma+1)\right\}^{\frac{1}{2}} \langle j, I, \sigma, -\mu+1 | J, \sigma - \mu + 1 \rangle \tag{VII.37}
\end{aligned}$$

and a simpler recurrence relation for $\langle l, s, 0, \sigma | j, \sigma \rangle$. The sign of $\langle l, s, 0, -s | j, -s \rangle$ is positive and if this coefficient does not exist, the sign of $\langle l, s, 0, -j | j, -j \rangle$ is $(-)^{s-j}$. Similarly, the sign of $\langle j, I, \sigma, -I | J, -I \rangle$ is also positive and if this coefficient does not exist the sign of $\langle j, I, \sigma, -J-\sigma | J, -J \rangle$ is $(-)^{I-J-\sigma}$. Consequently, starting the recurrences from the most negative values, the sign is given by the number of coefficients which do not verify the usual relations on quantum numbers. Normalised values are easily obtained, summing the squares of the unnormalised coefficients for each recurrence.

B.2.b Cross-section expansion in Legendre polynomials

Two similar recurrence computations of Clebsch-Gordan coefficients are needed to obtain the coefficients of Legendre polynomials describing cross-sections.

The first of them is to obtain the product of the two reduced rotation matrix elements in the helicity Coulomb amplitude as given by Equ. (III.65). As given by Equ. (69) of Appendix C in Ref. [73] page 920, their product is :

$$r_{\sigma_f, \sigma_i}^{(s)}(\theta) r_{-\mu_f, -\mu_i}^{(I)}(\theta) = \sum_{j=|I-s|}^{I+s} \langle s, I, \sigma_i, -\mu_i | j, \sigma_i - \mu_i \rangle \langle s, I, \sigma_f, -\mu_f | j, \sigma_f - \mu_f \rangle r_{\sigma_f - \mu_f, \sigma_i - \mu_i}^{(j)}(\theta) \tag{VII.38}$$

The recurrence relation needed here is not between “contiguous” coefficients, but similar to the one used for $6j$ coefficients when computing $9j$ coefficients as shown in section (B.1.d) by Equ. (VII.33) and (VII.34). The recurrence relation can be found of Appendix C in Ref. [73], formula (20) page 910, which is :

$$\begin{aligned}
& A(J) \langle j_1, j_2, m_1, m_2 | J-1, M \rangle + B(J) \langle j_1, j_2, m_1, m_2 | J, M \rangle \\
& + A(J+1) \langle j_1, j_2, m_1, m_2 | J+1, M \rangle = 0 \\
& A(J) = \frac{1}{J} \sqrt{\frac{[J^2 - M^2][(j_1 + j_2 + 1)^2 - J^2][J^2 - (j_1 - j_2)^2]}{(2J-1)(2J+1)}}, \\
& B(J) = m_1 - m_2 + M \frac{j_2(j_2+1) - j_1(j_1+1)}{J(J+1)} \tag{VII.39}
\end{aligned}$$

This recurrence relation is used for increasing values of j . The normalisation is obtained by dividing by the square root of the sum of unnormalised values. The sign is such that the product of Clebsch-Gordan coefficients is positive for the highest value of j in Equ. (VII.38).

The second occurrence is for the angular distribution of compound nucleus, which is given by :

$$\left(\frac{d\sigma(\theta)}{d\Omega}\right)_{c.n.} = \frac{1}{(2s_i + 1)(2I_i + 1)} \sum_{\sigma_i \sigma_f \mu_i \mu_f} \sum_J f_{\mu_f \sigma_f \mu_i \sigma_i}^{(J) \text{ hel c.n.}} |r_{\sigma_f - \mu_f, \sigma_i - \mu_i}^{(J)}(\theta)|^2, \quad (\text{VII.40})$$

which can be expressed in terms of Legendre polynomials with the relation :

$$|r_{m, m'}^{(J)}(\theta)|^2 = (-)^{m-m'} \sum_{\lambda=0}^{2J} \langle J, J, m, -m | \lambda, 0 \rangle \langle J, J, m', -m' | \lambda, 0 \rangle P_\lambda(\cos \theta) \quad (\text{VII.41})$$

The recurrence relation needed here is the same. After simplification for this application, Equ. (VII.39) is :

$$\begin{aligned} & A(J) \langle j, j, m, -m | J-1, 0 \rangle + B(J) \langle j, j, m, -m | J, 0 \rangle \\ & + A(J+1) \langle j, j, m, -m | J+1, 0 \rangle = 0 \\ & A(J) = J \sqrt{\frac{(2j+1)^2 - J^2}{(2J-1)(2J+1)}}, \quad B(J) = 2m \end{aligned} \quad (\text{VII.42})$$

Recurrences start from values 1 for $\lambda = 0$ because :

$$(-)^{m-m'} \langle J, J, m, -m | 0, 0 \rangle \langle J, J, m', -m' | 0, 0 \rangle = \frac{1}{2J+1} \quad (\text{VII.43})$$

Contributions of odd values of J disappear in the sum on helicities.

C Reduced rotation matrix elements

The subroutine **EMRO** computes an array of reduced rotation matrix elements $r_{m, m'}^{(J)}(\theta)$ for $J_m \leq J \leq J_M$ where $J_m = \max\{m, m'\}$ and J_M is the maximum value involved. But reduced matrix elements are also computed by recurrence to rotate the scattering matrix, to factorise $(1 - x \cos \theta)$ in the amplitudes and to express the cross-sections with Legendre polynomials. For reduced rotation matrix elements, see Ref. [73] and in : [76] VARSHALOVICH, D., A., MOSKALIEV, A., N. and KHERSONSKII, V., K., “*Kvantovaya Teoriya Uglovogo Momenta*”, (“*Quantum Theory of Angular Momenta*”), Nauka, Leningrad (1975) Chapter 4.

C.1 Recurrence with respect to J

This recurrence is a simplification of Equ. (1) of section (4.8) of Ref. [76], page 79. It can be rewritten as :

$$\begin{aligned} & (J+1) \sqrt{(J+m)(J-m)(J+m')(J-m')} r_{m, m'}^{(J-1)}(\theta) + (2J+1) [mm' - J(J+1) \cos \theta] r_{m, m'}^{(J)}(\theta) \\ & + J \sqrt{(J+1+m)(J+1-m)(J+1+m')(J+1-m')} r_{m, m'}^{(J+1)}(\theta) = 0 \end{aligned} \quad (\text{VII.44})$$

which allows the computation of $r_{m, m'}^{(J_m+1)}(\theta)$ for $J_m = \max\{m, m'\}$ if $r_{m, m'}^{(J_m)}(\theta)$ is known because the coefficient of $r_{m, m'}^{(J_m-1)}(\theta)$ vanishes. This starting value is given by Equ. (73) of Appendix C in Ref. [73] page 922 :

$$r_{m, j}^{(j)}(\theta) = (-)^{j-m} r_{j, m}^{(j)}(\theta) = \sqrt{\frac{(2j)!}{(j+m)!(j-m)!}} \xi^{j+m} \eta^{j-m}, \quad \xi = \cos \frac{1}{2} \theta, \quad \eta = \sin \frac{1}{2} \theta \quad (\text{VII.45})$$

which is obtained easily because the arguments of the subroutine **EMRO**, which computes the reduced rotation matrix elements include ξ , η , $\cos \theta$ and an array of square roots of integers. The recurrence given by Equ. (VII.44) used upwards seems to be stable without limitation.

If $m' = 0$, the recurrence of Equ. (VII.44) can be simplified to give :

$$\sqrt{(J+m)(J-m)} r_{m,0}^{(J-1)}(\theta) - (2J+1) \cos \theta r_{m,0}^{(J)}(\theta) + \sqrt{(J+1+m)(J+1-m)} r_{m,0}^{(J+1)}(\theta) = 0 \quad (\text{VII.46})$$

If $m = 0$, we obtain the same equation with m' instead of m . The starting value are :

$$r_{0,j}^{(j)}(\theta) = (-)^j r_{j,0}^{(j)}(\theta) = \sqrt{\frac{(2j)!}{j!j!}} (\xi\eta)^j \quad (\text{VII.47})$$

If $m = m' = 0$, the subroutine uses :

$$J r_{0,0}^{(J-1)}(\theta) - (2J+1) \cos \theta r_{0,0}^{(J)}(\theta) + (J+1) r_{0,0}^{(J+1)}(\theta) = 0 \quad (\text{VII.48})$$

starting with $r_{0,0}^{(0)}(\theta) = 1$.

For the angular distribution of a level, this calculation is done only once for all the amplitudes involving the same values of m and m' , but also for $-m$ and $-m'$, using the relation :

$$r_{m,m'}^{(J)}(\theta) = (-)^{m-m'} r_{-m,-m'}^{(J)}(\theta) \quad (\text{VII.49})$$

C.2 Recurrence with respect to m

It can be necessary to multiply the scattering matrix by the reduced rotation matrix $r^{(j)}(\alpha)$ where j is the spin of the outgoing particle or of the recoil nucleus and α the angle given by Equ. (VII.25) or (VII.26) respectively or $\alpha = -\theta$ for observables described with only one axis of quantification. The recurrence relation is Equ. (16) or (17) of section (4.8) in [76], page 81. It is :

$$\sqrt{(j+m')(j-m'+1)} r_{m,m'-1}^{(j)}(\alpha) + 2 \frac{m-m' \cos \alpha}{\sin \alpha} r_{m,m'}^{(j)}(\alpha) + \sqrt{(j-m')(j+m'+1)} r_{m,m'+1}^{(j)}(\alpha) = 0 \quad (\text{VII.50})$$

which can start with $r_{m,j}^{(j)}(\alpha)$ given by :

$$\sqrt{(j+m)(j-m+1)} r_{m-1,j}^{(j)}(\alpha) - 2 \frac{j-m \cos \alpha}{\sin \alpha} r_{m,j}^{(j)}(\alpha) + \sqrt{(j-m)(j+m+1)} r_{m+1,j}^{(j)}(\alpha) = 0 \quad (\text{VII.51})$$

This relation is used to compute $r_{m,j}^{(j)}(\alpha)$ for m varying from $-j$ to j , starting with :

$$r_{-j,j}^{(j)} = (-\sin \frac{1}{2} \alpha)^{2j} \quad (\text{VII.52})$$

and Equ. (VII.50) is used to compute $r_{m,m'}^{(j)}$ for m' varying from j to m . The other reduced rotation matrix elements are obtained from Equ. (VII.49).

C.3 Factorisation of $(1 - x \cos \theta)$

To accelerate convergence with respect to the total spin J , it is convenient to replace the amplitude f by an amplitude f' such that :

$$\sum_J f_{\mu_f \sigma_f, \mu_i \sigma_i}^{(J)} r_{m,m'}^{(J)}(\theta) = \frac{1}{1 - x \cos \theta} \sum_J f'_{\mu_f \sigma_f, \mu_i \sigma_i}^{(J)} r_{m,m'}^{(J)}(\theta) \quad (\text{VII.53})$$

with $m = \sigma_f - \mu_f$ and $m' = \sigma_i - \mu_i$. Multiplying the two sides by $(1 - x \cos \theta)$ and using the Equ. (VII.44) to express the product of a reduced rotation matrix element by $\cos \theta$, we can identify the coefficients of the same $r_{m,m'}^{(j)}(\theta)$ to get :

$$\begin{aligned} f_{\mu_f \sigma_f, \mu_i \sigma_i}^{(j)} &= -x \frac{\sqrt{(j-m)(j+m)(j-m')(j+m')}}{j(2j+1)} f_{\mu_f \sigma_f, \mu_i \sigma_i}^{(j-1)} + \left[1 - x \frac{mm'}{J(J+1)}\right] f_{\mu_f \sigma_f, \mu_i \sigma_i}^{(j)} \\ &\quad - x \frac{\sqrt{(j+1-m)(j+1+m)(j+1-m')(j+1+m')}}{(j+1)(2j+1)} f_{\mu_f \sigma_f, \mu_i \sigma_i}^{(j+1)} \end{aligned} \quad (\text{VII.54})$$

which allows to obtain $f^{(j)}$ up to $j = J_M - 1$ if the values of $f^{(j)}$ are known up to $j = J_m$. Writting this new amplitude $f'^{(j)} = f^{(j)} - x f''^{(j)}$, the value of x is obtained by :

$$D1 = \frac{1}{2} \sum_{j=J_M-5}^{j=J_M-1} (2j+1)^2 [f^{(j)} f''^{(j)*} + f^{(j)*} f''^{(j)}], \quad D2 = \sum_{j=J_M-5}^{j=J_M-1} (2j+1)^2 [f'^{(j)} f''^{(j)*}], \quad x = \frac{D1}{D2} \quad (\text{VII.55})$$

but with the restriction $|x| < 1$. The value of x is stored instead of the real part of $f^{(J_M)}$. This process can be done again several times.

C.4 Expression of cross-sections in Legendre polynomials

If the inelastic scattering amplitudes or the elastic scattering amplitude of a chargeless particle is known up to J_M , the cross-section can be expressed as a sum of Legendre polynomials $P_l(\cos \theta)$ with l from $l = 0$ to $l = 2J_M$. For the elastic cross-section of a charged particle, the Coulomb cross-section must be subtracted; the remaining cross-section cannot be expanded into Legendre polynomials because it involves an interference term with the Coulomb amplitude and the Legendre expansion of the Coulomb amplitude does not converge. So, the difference between elastic and pure Coulomb cross-section must be multiplied by some power of $(1 - \cos \theta)$ before being expanded in Legendre polynomials.

C.4.a Legendre expansion without Coulomb amplitude

The contribution to the cross-section of an helicity amplitude being :

$$\frac{d\sigma(\theta)}{d\Omega} = \frac{1}{(2s_i + 1)(2I_i + 1)} \left| \sum_J f_{\mu_f \sigma_f, \mu_i \sigma_i}^{(J) \text{ hel.}} r_{\sigma_f - \mu_f, \sigma_i - \mu_i}^{(J)}(\theta) \right|^2 \quad (\text{VII.56})$$

its expression as an expansion on Legendre polynomials and the coefficients are :

$$\frac{d\sigma(\theta)}{d\Omega} = \sum_L (2L+1) A_L P_L(\cos \theta), \quad A_L = \frac{1}{2} \int \frac{d\sigma(\theta)}{d\Omega} P_L(\cos \theta) d\cos \theta \quad (\text{VII.57})$$

The coefficient of $P_0(\cos \theta)$ is quite trivial :

$$A_0 = \frac{1}{(2s_i + 1)(2I_i + 1)} \sum_J \frac{1}{2J+1} \left| f_{\mu_f \sigma_f, \mu_i \sigma_i}^{(J) \text{ hel.}} \right|^2 \quad (\text{VII.58})$$

and can be generalised to any $P_L(\cos \theta)$ if one defines :

$$F_L = P_L(\cos \theta) f_{\mu_f \sigma_f, \mu_i \sigma_i}^{\text{hel.}} = \sum_J F_L^{(J)} r_{\sigma_f - \mu_f, \sigma_i - \mu_i}^{(J)}(\theta) \quad (\text{VII.59})$$

and use :

$$A_L = \frac{1}{(2s_i + 1)(2I_i + 1)} \sum_J \frac{1}{2J+1} f_{\mu_f \sigma_f, \mu_i \sigma_i}^{(J) \text{ hel.}} F_L^{(J)*}. \quad (\text{VII.60})$$

The F_L can be obtained with the same recurrence relations as the $P_L(\cos\theta)$. This give for their coefficients, using again $m = \sigma_f - \mu_f$ and $m' = \sigma_i - \mu_i$:

$$F_L^{(J)} = -\frac{L-1}{L}F_{L-2}^{(J)} - \frac{2L-1}{L}\left\{\frac{\sqrt{(J-m)(J+m)(J-m')(J+m')}}{J(2J-1)}F_{L-1}^{(J-1)} + \frac{mm'}{J(J+1)}F_{L-1}^{(J)} + \frac{\sqrt{(J+1-m)(J+1+m)(J+1-m')(J+1+m')}}{(J+1)(2J+3)}F_{L-1}^{(J+1)}\right\} \quad (\text{VII.61})$$

easy to use because the dependence on L is very simple and the dependence in J is the same for every L and can be stored. In this process, the number of J values increases with L .

C.4.b Legendre expansion of Coulomb amplitude

To generalise this calculation to the elastic scattering of charged particle, the helicity amplitude in Equ. (VII.59) does not include the coulomb amplitude. To obtain F_0 :

1. the helicity amplitude without Coulomb amplitude is multiplied by $(1 - \cos\theta)^2$, using recurrence relations on the reduced rotation matrix elements and also by $(1 + \cos\theta)$ for identical particles, giving $h(\theta)$,
2. the two reduced matrix elements in the spin space are multiplied as shown by Equ. (VII.38), giving some function $g(\theta)$,
3. the Legendre expansion of the Coulomb amplitude multiplied by $(1 - \cos\theta)^n$ or $(1 - \cos^2\theta)^n$ is generated as :

$$c(\theta) = \sum_L (2L+1)C_L^{(n)} P_L(\cos\theta) \quad (\text{VII.62})$$

4. $g(\theta)$ is multiplied by $c(\theta)$ and added to $h(\theta)$ to get the function F_0 used in Equ. (VII.60) to compute A_0 :

$$F_0 = h(\theta) + g(\theta) c(\theta) = h(\theta) + \sum_L (2L+1)C_L^{(n)} G_L \quad (\text{VII.63})$$

where the G_L are defined with respect to $g(\theta)$ as F_L was defined with the helicity amplitude in Equ. (VII.59).

Then, the computation of the coefficients A_L can be done as in the absence of Coulomb amplitude. This expansion must be done up to $J_M + L_M$ where L_M is the maximum L value for the A_L requested and J_M the maximum J value of the amplitudes.

The Legendre expansion of the coulomb amplitude multiplied by any power of $(1 - \cos\theta)$ is :

$$(1 - \cos\theta)^n f_c(\theta) = \sum_l (2l+1)C_l^{(n)} P_l(\cos\theta) \quad (\text{VII.64})$$

For a symmetrised amplitude, we can use the relation :

$$(1 - \cos^2\theta)^n = \sum_{m=0}^{m=n} \frac{n!}{m!(n-m)!} (-)^m 2^{n-m} (1 - \cos\theta)^{n+m} \quad (\text{VII.65})$$

The coefficients $C_l^{(n)}$ have quite simple recurrence relations. For a given value of n , all the coefficients needed are given by :

$$C_{l+1}^{(n)} = \frac{l+1-n+i\eta}{l+1+n-i\eta} C_l^{(n)} \quad (\text{VII.66})$$

starting with :

$$C_0^{(n)} = -2^n \frac{i\eta}{n-i\eta} C_0^{(0)} \quad (\text{VII.67})$$

With the convention $\sigma_0 = 0$ used in ECIS, $C_0^{(0)} = 1/(2ik)$.

Chapter VIII

Description of the subroutines

We shall describe the subroutines in the order of their appearance in the code **ECIS94**. The Table of Contents of these paragraph will reproduce a flow chart of the programme. *Italic characters shifted to the right side indicates comments related to a calling subroutine.*

All the steps of the calculation are managed by 240 logicals from **L0(1)** to **L0(250)**. They are :

- (a) from **L0(1)** to **L0(100)**, the values read in the second and the third cards of the input; the meaning of those which are used can be found in the “**DESCRIPTION OF INPUT**”.
- (b) from **L0(101)** to **L0(200)**, the opposite values of the preceding one.
- (c) from **L0(201)** to **L0(250)**, internal controls of the code; the meaning of those which are used is given as comment cards in subroutine **CALC** reproduced below.

C L0(201) IS TRUE IF THERE IS NO REAL SPIN-ORBIT POTENTIAL	CALC-038
C L0(202) IS TRUE IF THERE IS NO IMAGINARY SPIN-ORBIT POTENTIAL	CALC-039
C L0(203) IS TRUE IF THERE IS NO COULOMB SPIN-ORBIT POTENTIAL	CALC-040
C L0(204) IS TRUE IF CONVERGENCE IS OBTAINED IN THE ITERATION	CALC-041
C L0(205) IS TRUE IF CONVERGENCE IS OBTAINED FOR THIS EQUATION	CALC-042
C L0(206) IS TRUE WHEN THE ITERATION IS NOT THE LAST ONE PERMITTED	CALC-043
C L0(207) IS TRUE IF ALL THE COUPLINGS HAVE TO BE CALCULATED BEFOREHAND	CALC-044
C L0(208) IS TRUE IF THE DIAGONAL COULOMB CORRECTIONS ARE NEEDED	CALC-045
C L0(209) IS TRUE FOR DIRAC POTENTIALS	CALC-046
C L0(210) IS TRUE IF DERIVATIVES ARE NEEDED	CALC-047
C L0(211) IS TRUE IF DEFORMATIONS ARE CHANGED IN SEARCH	CALC-048
C L0(212) IS TRUE IF NUCLEAR PARAMETERS ARE CHANGED IN SEARCH	CALC-049
C L0(213) IS TRUE IF NUCLEAR MATRIX ELEMENTS ARE CHANGED IN SEARCH	CALC-050
C L0(214) IS TRUE IF SPIN-ORBIT PARAMETRISATION IS CHANGED IN SEARCH	CALC-051
C L0(215) IS TRUE FOR NO OUTPUT	CALC-052
C L0(216) IS TRUE FOR A STOP, FOR EXAMPLE THE WORKING FIELD IS TOO SMALL	CALC-053
C L0(217) IS TRUE FOR ALL THE CALCULATIONS EXCEPT THE FIRST	CALC-054
C L0(218) IS TRUE FOR LAST RESULTS	CALC-055
C L0(219) IS TRUE FOR RESULTS WITHOUT DOING THE CALCULATION AGAIN	CALC-056
C L0(220) IS TRUE FOR OUTPUT AND LAST CALCULATION IS THE BEST ONE	CALC-057
C L0(221) IS TRUE FOR OPTICAL MODEL WITHOUT COUPLING	CALC-058
C L0(222) IS TRUE IF IT IS THE FIRST COMPUTATION FOR THIS ENERGY	CALC-059
C L0(223) IS TRUE IF L0(18) IS .TRUE. AND NO SPIN IN THE INITIAL STATE	CALC-060
C L0(224) IS TRUE FOR COMPOUND NUCLEUS OR PUNCH OF TRANSMISSION COEFF.	CALC-061
C L0(225) IS TRUE IN CAL1 FOR A CALL TO USUAL COUPLED EQUATIONS SUBR.	CALC-062
C L0(226) IS TRUE IF THERE ARE OBSERVABLES IN THE LABORATORY SYSTEM	CALC-063
C L0(227) IS TRUE FOR COULOMB CORRECTIONS WITH PURE REGULAR FUNCTIONS	CALC-064
C L0(228) IS TRUE FOR NO COPY OF UNCOUPLED FUNCTIONS AND PHASE-SHIFT	CALC-065

A Subroutines ECIS/CALS, HORA, STIM and MEMO

For the **MAIN**, the subroutines **ECIS/CALS**, **HORA**, **STIM** and **MEMO**, see Chapter II, “*Use on various Computers*”.

1. The **MAIN** defines only the working array and calls **CALC**, directly or via the Assembler subroutine **ECIS** on **IBM**, or via **CALS** on **CDC 7600**.
2. The subroutine **HORA** is called from different places to give the elapsed **CPU** time for the **JOB**. It uses the subroutine **STIM** to get the allowed remaining **CPU** time.
3. The subroutine **STIM** should give the time remaining for the **JOB**. This subroutine is very machine dependent. It is called directly by **CALC** to obtain the time needed by a single calculation to stop an automatic search when **LO(34)=.TRUE.** and to save results or a search before being stopped by time limitation. **STIM** allows to the subroutine **HORA** to print the elapsed time.
4. The subroutine **MEMO** is called from anywhere to increase the size of the working array. On computers other than **CDC** or **UNIVAC** this subroutine stops the calculation if the request is absolute. This subroutine is also very machine dependent.

We shall not quote the calls to **HORA** or **MEMO** in the description of the subroutines.

B Subroutine CALC

The main subroutine is **CALC**, from which the calculation never returns. Its arguments are the working array and its length. The same working array is floating values and integers in the **CDC** version, also double precision in the **IBM** version. This subroutine calls the subroutines **CALX**, **COLF**, **REDM**, **EXTP**, **CAL1**, **VARI**, **FITE**, **REST**, **EVAL** which we describe below. Meaning of addresses stored in the common **/DECOU/** and defined in **CALC** after the return from **CALX** are given on **COMMENT** cards.

The computation is normally stopped by a control word read in subroutine **CALX**, so there is usually a warning at the compilation “subprogram exit cannot be reached from entry”.

The calculation never returns to the first instructions which are values of the Atomic Mass Unit M_u , $\hbar c$ and fine structure constant, as shown in section (VII.A) and the calculation of $M/(\hbar c)$, $2M/(\hbar c)^2$ and $\hbar c\alpha$. Then **CALC** calls **CALX**.

C Subroutine CALX

At the beginning of this subroutine is given on comment cards the meaning of the arrays **SP**, **IPI**, **WV** which store all the informations relative to the levels. They are followed by the meaning of data and addresses stored in the common **/DECOU/** and some informations on other commons. The subroutine calls **INPA**, **INPB**, **INPC**, **LECL**, **LECT**, **DEPH**, **LECD** and **OBSE**.

This subroutine reads all the input except for nuclear matrix elements and external form factors and new parameters if **LO(37)=.TRUE.**. This input starts by a title card :

- if this title is “**FIN** ”, the calculation is stopped,
- if this title is “**DESCRIPTION** ”, the description of the input is printed.

C.1 Subroutines INPA, INPB, INPC

These three subroutine are called one after the other if the title is “**DESCRIPTION**”. They include only **WRITE** statements and they have been generated from the text written on cards with a special programme. Output can be listed as **FORTRAN** or by “enscript -2Bhr -L58” after suppression of the first column. It is 27 pages long.

After the printing, a new title card is read in CALX. If the title card is neither “FIN” or “DESCRIPTION”, the subroutine CALX reads the first card of logical control (card 2) : if LO(36)=.TRUE. (restart of a computation from file 8), the subroutine returns to CALC which calls to the subroutine REST. If LO(36)=.FALSE., the second card of logical control (card 3) and card 4 (integers) are read. Some logical controls can be changed (for example, the real spin-orbit must be deformed if the imaginary spin-orbit is deformed) before being printed. Then the card 5 (floating values) is read and default values for card 4 and 5 are set and the values printed. The logicals from LO(211) to LO(220) are set .FALSE. and LO(222) is set .TRUE..

Then, informations needed when some logicals are .TRUE. are read :

- *limitation on angular momenta for coulomb corrections if LO(45)=.TRUE.,*
- *compound nucleus data or indication to calculate the Legendre Polynomial expression of cross-sections if LO(84), LO(85), LO(86) or LO(65) is .TRUE.,*
- *interpolation on total spin if LO(43)=.TRUE..*

Some default options are set. The subroutine computes storage requirements as soon as it can be done and calls LECL, LECT and DEPH with only some storage calculations between them.

C.2 Subroutine LECL

This subroutine reads energies, masses, spins and product of charges for all the levels. Informations relevant of the nuclear model are read for coupled states but not for uncoupled states introduced in a compound nucleus calculation. For the level **I**, the informations are stored in SP(3,I), IPI(3,I) and WV(11,I) in the following order :

- SP(1,I) s_I spin of incident particle,
- SP(2,I) I_I spin of the target,
- SP(3,I) Z_I product of the charges of the particle and the target,
- IPI(1,I) π_I parity (0 for “+” and 1 for “-”),
- IPI(2,I) $n_I = 2s_I + 1$ multiplicity of the incident particle,
- IPI(3,I) $N_I = 2I_I + 1$ multiplicity of the target,
- IPI(4,I) position of the potential used,
- IPI(5-6,I) beginning and end in the description of amplitudes (defined in DEPH),
- IPI(7-8,I) beginning and end in the description of observables (defined in DEPH),
- IPI(9,I) cross-reference to potentials (if there are more potentials than levels, this table is extended to the number of potentials),
- WV(1,I) m_I mass of the incident particle,
- WV(2,I) M_I mass of the target (for relativistic kinematics, LO(8)=.TRUE., the target mass is corrected by addition of Q/M_u but this correction is dropped if LO(44)=.TRUE. to compute the wave number in the Schrödinger formalism),

WV(3,I) E_I energy in the center of mass system; for the incident channel, this energy is given by Equ. (VII.8) if **LO(8)=.FALSE.** (non relativistic kinematics) or by Equ. (VII.4) if **LO(8)=.TRUE.** (relativistic kinematics); for the other channels $E_I = E_1 - Q$.

For the other entries of **WV**, see description of **COLF**.

The wave number of the ground state is computed only to obtain the default options of matching radius and step size in **LECT**. The same computation will be done in **COLF** for all the levels.

C.3 Subroutine LECT

This subroutine reads the description of the phonons if there are some in the nuclear model, the deformations for the rotational model and the optical potentials if **LO(7)=.FALSE.** (no external potential). The default value of the matching radius is so that the largest potential is less than the data **ACONV** of card 5 and the default value of the step size to the smaller of half diffuseness and $1/(2k)$ where k is the wave number of the ground state. The step size is modified to be the matching radius divided by an integer. The logical asking for deformation of the real spin-orbit, the imaginary spin-orbit, the coulomb or the spin-orbit coulomb potentials is set **.FALSE.** if these interactions vanish.

If **LO(7)=.TRUE.**, the default value of the matching radius is 20fm and the default value of the step size is given only by the wave number. The logicals for deformation of spin-orbit potentials can be changed if there is no spins in the levels. However, without spin-orbit coulomb potential and no coulomb corrections, one must use **LO(46)=.TRUE.**. This subroutine returns also in the **COMMON TITLE** the factor **XZ** equal to 10 divided by the multiplicity of the initial channel (**IPI(2,1)** and **IPI(3,1)** defined in **LECL**), used by the subroutine **SCAT** to express the cross-sections in millibarns.

Then the subroutine reads :

- angles for equidistant angular distributions if **LO(66)=.FALSE.**,
- spin-orbit parametrisation if **LO(4)=.TRUE.**,
- Hauser-Feshbach parameters if **LO(82)=.TRUE.**,
- Fission penetrabilities if **LO(85)=.TRUE.**,
- Gamma penetrabilities if **LO(86)=.TRUE.**.

C.4 Subroutine DEPH

There are two parts with comment cards before each to explain how are stored informations on each amplitude. The levels taken into account are those which are coupled plus those which are not coupled but of which an angular distribution is wanted.

In the first one, the amplitudes are counted and arrays of quantum number stored. Indications to compute only once the reduced rotation matrix elements when they can be used for more than one amplitude to a sign are also stored. The number of solutions and of equation for each parity is obtained in **NCT**. For each level, beginning and end addresses are stored in **IPI(5,I)** and **IPI(6,I)** respectively.

In the second part, the default output is generated unless **LO(94)=.TRUE.**, in which case the instructions are read. For each level, beginning and end addresses of instructions are stored in **IPI(7,I)** and **IPI(8,I)** respectively. This subroutine reads also instructions for restricted coulomb corrections if **LO(44)** and **LO(46)=.TRUE.**. Addresses of coulomb corrections are stored after the array of beginning and end of coupling **NIV** (see **REDM**) in **NIV(I,J,3)**.

*If **LO(31)=.TRUE.**, the subroutine **CALX** reads number of angular distributions and of parameters in search.*

C.5 Subroutine LECD

This subroutine reads the angular distributions.

The subroutine CALX reads identifications and step sizes of the parameters in search if LO(32)=.TRUE. and if their number and the number of angular distributions are non zero.

C.6 Subroutine OBSE

This subroutine computes for all the observables the indications (described on **COMMENT** cards) for the do-loops and the geometrical coefficients which will be needed in **SCAT** to obtain the observables requested. A first part reads indications for non standard observables and transform them into tensor notations. A second part computes geometrical coefficients and do-loop limits. This part is run twice, the first time to obtain storage requirements, the second one for effective computation.

After return from OBSE, the subroutine CALX computes some storage requirement and a table of logarithms of factorial as long as needed.

After return to CALC from CALX, the subroutine REST is called with KF=1 if LO(36)=.TRUE. (restart of a search). If LO(36)=.FALSE. the subroutine COLF is called.

D Subroutine COLF

This subroutine and the other subroutines called in it compute the matching values and the stored integrals for coulomb corrections. The array **WV(11,I)** is completed for all the levels, coupled or uncoupled. Using the recoil ratio R_{ec} to take into account recoil corrections for a zero-range interaction when the masses are not the same in all the channels ($R_{ec} = 1$ for the incident channel, but for the excited channels $R_{ec} = M_1/M_i$ if the mass of the particle is not the same and if **LO(93)=.FALSE.**), the content of this array is :

- WV(4,I)** k_I wave number; if **LO(8)=.FALSE.** : it is given by Equ. (VII.8) and if **LO(8)=.TRUE.** by Equ. (VII.6),
- WV(5,I)** η_I coulomb parameter,
- WV(6,I)** $\sqrt{R_{ec}k_I/k_1}$, square root of ratio of wave numbers,
- WV(7,I)** E_{Irel} relativistic energy in the center of mass system in Dirac formalism in Mev,
- WV(8,I)** square root of $2m/\hbar^2$,
- WV(9,I)** $k_i R_{ec}$, wave number multiplied by ratio of step sizes,
- WV(10,I)** $h_i^2 k_i^2$ where h_i is the last item,
- WV(11,I)** $h_i = h_1 R_{ec}$ step size of this channel.

The subroutine **COCL** is called for closed channels and the subroutine **FCOU** for open channels.

D.1 Subroutine COCL

This subroutine computes matching values for closed channels. For $L = 0$, it uses asymptotic expansion with/without backwards integration. For the other values of L , it uses upwards recurrence. This subroutine comes from the Buck and Hill's code **INCH**.

D.2 Subroutine FCOU

This subroutine and the subroutines called by it are described in Ref. [18] and Ref. [19]. The calculation of phase-shifts has been suppressed except for $L = 0$, the factorisation of some power of 10 has been changed from modulo 60 to modulo 15 for VAX computers to handle squares of Coulomb functions.

The subroutine **FCOU** calls **FCZO** to obtain coulomb functions and their derivatives for $L = 0$ and computes the other ones by recurrence. It uses upwards recurrence for irregular to the end and for regular functions as long as $\rho > \eta + [\eta^2 + L(L + 1)]^{\frac{1}{2}}$, downwards recurrence for the regular function starting from the maximum L requested plus $25 + 5|\eta|$ normalised with the last value obtained from upwards recurrence.

D.2.a Subroutines FCZO

The subroutine **FCZO** computes the coulomb functions for $L = 0$. It calls the function **SIGM** to obtain the phase-shift. For some values of the coulomb parameter and the radius, it calls the subroutines **YFRI** or **YFAS**. For some other values near the origin, it computes the regular function by expansion into Chebyshev polynomial (Clenshaw expansion) or series expansion and calls **YFIR** for the irregular one.

D.2.b Subroutine YFRI

For some values of the matching radius and of the coulomb parameter, it calls **YFCL**. For other values, it uses Riccati methods, “Riccati at the origin” or “asymptotic Riccati”.

D.2.c Subroutine YFAS

Computes coulomb functions with the asymptotic expansions.

D.2.d Subroutine YFIR

Computes irregular coulomb functions by Taylor expansion around the origin or around the point $R = 7.5 + 4/3\eta$ at which the functions and their derivatives are obtained with the subroutine **YFAS**. For the expansion around the origin, this subroutine calls the function **PSI**.

D.2.e Subroutine YFCL

This subroutine is called by **YFRI**. It uses an expansion on Chebyshev polynomials in the asymptotic region or a MacLaurin series expansion near the origin for which it needs the function **PSI**.

D.2.f Functions SIGM, PSI

These functions compute respectively the coulomb phase-shift for $L = 0$ and the real part of the logarithmic derivative of the gamma function for a complex argument.

*The subroutine **COLF** checks if some coulomb functions are too large. If the power of 10^{15} returned by **FCOU** is not zero, it limits the number of Coulomb functions to avoid them (with a ratio 10^{30} between regular and irregular functions, the phase-shift obtained by the code cannot be significative). If coulomb corrections have to be used, it calls subroutine **CORI** to compute the stored integrals with all the previous channel up this one, if the pair of channels are both open and the corrections requested.*

D.3 Subroutine CORI

If $\eta_i > 2.2\eta_f$, the code is unable to do the calculation and stops. If $\eta_i = 0$, η_f must be also zero : the subroutine **CORZ** is called and **CORI** returns.

In any other case, the subroutine **CORI** calls the subroutine **CORO** to obtain the integrals from R_m to ∞ of product of regular and irregular Coulomb functions with $L = 0$ and $L = 1$, divided by r and r^2 . Then, if $\eta_i = \eta_f$ and $k_i = k_f$, the integrals from 0 to infinity of the product of two regular Coulomb functions divided by r^2 are obtained by Equ IV.23. Details are given in section (IV.B).

D.3.a Subroutine CORO

This subroutine computes the integrals from R_m to ∞ of the product of two different regular or irregular coulomb functions with the same angular momentum $L = 0$ and $L = 1$ divided by r and r^2 (16 values). With $3R_s = \max[(5|\eta_i| + 22.5)\sqrt{k_f/k_i}, (5|\eta_f| + 22.5)\sqrt{k_i/k_f}, .18\sqrt{\eta = i\eta_f}]$, these integrals are computed from $\max(R_m, R_s)$ to ∞ by a generalisation of the method described in Ref. [66]. If $R_s > R_m$, the integral between R_m and R_s is computed by 40 points Gauss integrations, each Gauss integration being for $\Delta r < 20$.

D.3.b Subroutine CORZ

This subroutine computes integrals from R_m to ∞ of products of regular and irregular functions of the same L divided by r^3 and integrals from 0 to ∞ of regular functions of the same L divided by r^3 .

*Then the coulomb phase-shift minus the one for $L = 0$ are computed in COLF.
After return from COLF, the subroutine CALC calls REDM.*

E Subroutine REDM

This subroutine generates or reads the reduced nuclear matrix elements between the coupled levels. If **LO(7)=.FALSE.** (standard nuclear model) and **LO(15)=.FALSE.** (reduced nuclear matrix elements not read), **REDM** calls the subroutines **VIBM**, **ROTM** or **ROAM**. If **LO(7)=.TRUE.** or **LO(15)=.TRUE.**, this subroutine reads the reduced matrix elements. Then, the informations are stored in **IQ(4,I)** and **T(4,I)** which are in equivalence, for **I=1** to **IT**.

IQ(1,I) Form factor identification (see **DESCRIPTION OF INPUT**),

IQ(2,I) L transferred angular momentum,

IQ(3,I) $10000(2\Delta S) + 10(2\Delta J) + 10a + b$, where $a = b = 0$ usually, but $a = 1$ for existence of spin-orbit deformation and $b = 1$ for a magnetic Coulomb excitation with $\lambda = L - 1$,

T(4,I) Reduced nuclear matrix element multiplied by $(-1)^{\text{Int}(L/2)}$.

The address of the first and the last reduced nuclear matrix element between level **I** and level **J** are in **NIV(1,I,J)** and **NIV(2,I,J)** respectively. The table **NIV** is completed if there are uncoupled states for which an angular distribution is requested and the reduced nuclear matrix elements written on file 7 as they should be read if **LO(61)=.TRUE.** .

Then, the table of **IM** multipoles, **IVQ(3,I)** is generated :

IVQ(1,I) L transferred angular momentum,

IVQ(2,I) $2\Delta S$ twice the transferred spin to the particle,

IVQ(3,I) $2\Delta J$ twice the transferred spin to the target or 0 if $\Delta S = 0$.

followed by a table of form factors **IVZ(7,I)** including correction terms (with external potentials and monopole or dipole corrections to vibrations in the rotational model) :

IVZ(1,I) Form factor control number,

IVZ(2,I) Address in the table of multipoles,

IVZ(3,I) 0 or address of the first spin-orbit form factor,

IVZ(4,I) 0 or address of the temporary Coulomb form factor,

IVZ(5,I) 0 or address of the temporary Coulomb spin-orbit form factor,

IVZ(6,I) 0 or address of correction term (positive for corrected term, negative for correction term),

IVZ(7,I) L transferred angular momentum.

The array **IQ** has been changed to be :

IQ(1,I) Address in the table of form factors **IVZ**

IQ(2,I) Address in the table of multipoles **IVQ**

IQ(3,I) 0 or address of the spin-orbit form factor

but **T(4,I)** is unchanged. A table **ITX** of the 8 starting addresses of potentials and 8 starting addresses of transition form-factors is built with :

NV Number of real form factors (**INVT+2 INLS**),

MV Number of imaginary form factors,

INVT Number of transitions, without correction terms,

INLS Number of spin-orbit form factors not taking into account multiplication by 2,

INVC Number of coulomb transition form factors,

INVD Idem for coulomb spin-orbit,

INTC Same as **INVT** but including correction terms.

These numbers include dummy central form factors for magnetic coulomb excitations.

The reduced nuclear matrix elements are written and/or punch on file 7 on request.

In a search, this subroutine is called again if some nuclear parameter is changed (there is no nuclear parameter if **LO(7)=.TRUE.**). At the end of a search is called again if reduced nuclear matrix elements have to be printed or punched.

E.1 **Subroutine VIBM**

Computes the reduced nuclear matrix elements of the vibrational model, taking into account that the phonons amplitudes and a factor $(n!\sqrt{4\pi})^{-1}$ is shifted to the radial form factors. This subroutine is limited to two-phonons states and second order vibrations and uses the functions **DJCG** and **DJ6J**. There is a transposition relation :

E.2 Subroutine ROTM

Computes the reduced nuclear matrix elements of the symmetric rotational model. This subroutine calls the function **DJCG**.

E.3 Subroutine ROAM

Computes the reduced nuclear matrix elements of the asymmetric rotational model. This subroutine calls the function **DJCG**. If **LO(2)=.TRUE.**, the mixing parameter of the second level is used as the asymmetry gamma angle and the mixing parameter of the first levels are computed from it.

E.4 Function DJCG

Computes Clebsh-Gordan coefficients. The subroutine verifies triangular relations. For more details see section (VII.B.1.a).

E.5 Function DJ6J

Computes $6j$ coefficients. The function tests if some angular momentum is zero to use simpler expression. For more details see section (VII.B.1.b).

*After the call to **REDM**, at the first calculation (**LO(225)=.TRUE.**), if **LO(7)=.TRUE.** (external potentials), the subroutine **EXTP** is called. After return from it, informations from **REDM** are copied after those of **EXTP** and the whole is shifted down (to prevent against variation of the size of informations from **REDM** in a search).*

F Subroutine EXTP

The subroutine **EXTP** calls only the function **DCGS** to obtain the geometrical coefficient of a particle-hole excitation. The form factors are read in any order. In case of error, the subroutine prints the list of form factors not yet read under the form (**I,J**) where **I** varies from 1 to 8 and **J** is the potential if less or equal to the number n_p of potentials and **J**- n_p is the form factor in the order printed by **REDM** if $j > n_p$; the subroutine stops after this output.

*The subroutine **CALC** computes storage requirements for potentials, calls the subroutine **STIM** to obtain remaining time before calculation. Then, **CALC** calls **CAL1**.*

G Subroutine CAL1

This subroutine computes from form factors to final results. It calls the subroutines **POTE**, **CONU**, **QUAN**, **MTCH**, **INTI**, **INTR**, **INCH**, **SCAM**, **SCHE** and **RESU**. If the computation is the last one of a search :

- if no printing is requested, **LO(219)=.TRUE.**,
- if the printing related to **LO(53)** or **LO(55)** to **LO(58)** and the punching related to **LO(62)**, **LO(64)** or **LO(65)** (this last one if **LO(41)=.TRUE.**) are not requested, and if the last χ^2 is the best one, **LO(220)=.TRUE.**,
- if **LO(220)=.FALSE.**, all the calculation has to be done again.

If **LO(220)=.TRUE.** and the printing of potentials is requested, **CAL1** calls the subroutine **POTE** and the subroutine **SCHE** after return. If **LO(219)=.TRUE.**, **CAL1** calls the subroutine **SCHE**.

First, **CAL1** set to 0 all the memories needed for the computation of potentials and calls **POTE**.

G.1 Subroutine POTE

This subroutine computes the form factors.

It calls the subroutines **ROTP**, **FOLD** and **STDP**.

If **LO(220)=.TRUE.** and **LO(51)=.TRUE.**, this subroutine prints only the form factors.

For standard models, there is a do-loop on the potentials, of which only the first is deformed. For each potential, **POTE** calls the subroutine **ROTP** and, if **LO(17)=.TRUE.** (folding), the subroutine **FOLD**. Instead of that, if **LO(7)=.TRUE.** (external potentials), **POTE** calls the subroutine **STDP**. If **LO(61)=.TRUE.**, the potentials are punched on file 7 with the format needed by subroutine **EXTP**. In the Schrödinger formalism, volume and surface potentials are added together.

If there are some, corrections are done. Separately for central, spin-orbit, coulomb and spin-orbit coulomb transition form factor in the Schrödinger formalism, the scalar, vector, tensor, coulomb transition form factors in the Dirac formalism, the correction form factor is added to the form factor to be corrected in such a way that the integral with r^{L+2} vanishes. For spin-orbit transition form factors in the Schrödinger formalism, it is the form factor V_λ/r^2 which is used with a factor r^{L+4} .

From there, the subroutine proceeds quite differently for the two formalism :

- **Schrödinger formalism** including equivalent for a Dirac formalism.

The coulomb potentials are added to the central ones, keeping their strengths in **VCO** for the potentials (correction to r^{-1} behaviour) and in **VDO** for the transitions. The coulomb magnetic form factors are multiplied by $\hbar c/m_r M_u$. The potentials are printed on request.

- **Dirac formalism**

The Schrödinger potential for each level j is computed separately and stored in the array **V(*,I,J)** with **I=1** to **I=14**, real part for odd **I** and imaginary part for even **I** :

V(*, 1,J)	Central potential,
V(*, 3,J)	Spin-orbit potential,
V(*, 5,J)	$D = E + m - V_v + V_s$ divided by $\hbar c$,
V(*, 7,J)	$D' = E - m - V_v - V_s$ divided by $\hbar c$,
V(*, 9,J)	$(E + m - V_v + V_s)^{\frac{1}{2}}$ more precisely square root of V(*, 5,J) ,
V(*,11,J)	$(E + m - V_v + V_s)^{-1}$ multiplied by $\hbar c$,
V(*,13,J)	Tensor potential multiplied by $\hbar c$.

The transitions are stored after that beginning by the scalar and vector ones :

V(*, 1,J)	$V_v^\lambda + V_s^\lambda$ to use between large components,
V(*, 3,J)	$V_v^\lambda - V_s^\lambda$ to use between small components.

followed by the tensor ones :

V(*, 1,J)	derivative of the tensor potential $\frac{d}{dr}V_T^\lambda$,
V(*, 3,J)	tensor potential V_T^λ divided by r .

The strengths of the coulomb potentials are kept in **VCO** and those of coulomb transitions in **VDO**. The coulomb magnetic form factors are multiplied by $\hbar c/m_r M_u$. The potentials are printed on request.

G.1.a Subroutine ROTP

This subroutine computes the potentials and the transition form factors for all the standard models. It calls the subroutines **ROTD**, **ROTZ**, **COPO** and **DERI**.

If a diffuseness, except for the coulomb potentials, is less than .02 times the step size, it is changed into the maximum of its opposite and .02*step size; if a radius is less than the step size and symmetrised Woods-Saxon potentials are not used, it is changed into the maximum of its opposite and the step size (this can be useful in a search).

For the asymmetric rotational model, this subroutine uses a large array of weights and rotation matrix elements which are **DATA** in **ROTD**. The symmetric rotational model uses one half of a 20 points Gauss-Legendre integration of which abscissae and weights are **DATA** in **POTE**. A table of deformation is built for all the transition for factors and the subroutine **ROTD** is called to get the number of points of angular integration, starting values, increments and so on.

For each value of r and for each angle (see **ROTD**), up to the fourth derivative can be needed (spin-orbit transition for the third order of the anharmonic vibrational model). They are stored in **VR(6,10)** (10 for the 8 components of the potential plus 2 Schrödinger spin-orbit, 6 for up to the second derivative and third order vibration in the Dirac formalism). In the Dirac formalism using Schrödinger equivalent (**LO(99)=.TRUE.**), the subroutine **ROTZ** is called. In other formalisms, the coulomb potentials and transitions for a non diffuse charge are computed; the charge distributions of diffuse charges are obtained as the other components of the potential. The angular integrated values are stored differently in the two formalism.

If there is a charge diffuseness, **ROTP** calls **COPO** to obtain the coulomb potential or transition form factor from the charge distribution or transition. For spin-orbit coulomb potentials, **ROTP** has also to call the subroutine **DERI**.

G.1.a.i Subroutine ROTD

This subroutine returns the number of angular integration points, the weight, the radii and its two first derivatives at each point of each form factor.

- In the symmetric rotational model, 20 points Gauss-Legendre integration method is used, reduced to 10 points by symmetry. For a vibration, the angular dependence of the boson is included in the weight.
- In the asymmetric rotational model, 36 values of (θ, ϕ) have been chosen : the values of θ are $N\pi/14$ and for each value of N , there are $N + 1$ values of ϕ equidistant from 0 to $\pi/2$. The matrix of which the elements are the rotation matrix elements for these angles and $L \leq 14$ with L and K even has been inverted. The coefficients of the 15 lowest (L, K) are stored as integration weights.
- In the vibrational model, the angular integration reduces to one point.

G.1.a.ii Subroutine ROTZ

The Schrödinger equivalent equation to the Dirac equation is supposed to describe also the inelastic scattering. For the vibrational model, the result of elastic scattering should be the same than the pure Dirac formalism; inelastic scattering is described by derivatives of these potential. In the rotational model, result should be different because gradient and Laplacian are computed on potentials depending on θ instead of its monopole part.

G.1.a.iii Subroutine COPO

Computes the coulomb potential and the coulomb interactions from the density distribution of charge and the form factors of charge transition.

G.1.a.iv Subroutine DERI

This subroutine returns the derivative $x(r) = -\frac{d}{dr}y(r)$ of a function $y(r)$ known by n equidistant points by steps h or this result divided by r . The formulae used are given by Equ. (III.95) to Equ. (III.97). The subroutine stops if there are less than 7 points.

G.1.b Subroutine FOLD

When called directly from **PUTE**, this subroutine folds separately the real potential, the imaginary potential and the coulomb potential with a particle density which can be a Woods-Saxon distribution or a Gaussian one. This subroutine can be called from the subroutine **STDP** to perform the same operations on each components of the potentials and of the transition form factors independently. To obtain spin-orbit form factors, and derivatives in Dirac formalism, the subroutine **DERI** is used.

G.1.c Subroutine STDP

This subroutine computes the external form factors of which the parameters are read in **EXTP** and calls the subroutines **INTP**, **STBF**, **COPO**, **DERI** and **FOLD**. The form factors can be :

- Copy of another form factor,
- Interpolated from values read on cards by subroutine **INTP**, eventually derived with subroutine **DERI** for spin-orbit form factors or Dirac formalism,
- Woods-Saxon volume and surface form factors at some power and their derivatives which can be deformed with even or odd deformation,
- Single or product of two Laguerre polynomials or solutions in a Woods-Saxon potential computed by subroutine **STBF**,
- Sums of Bessel functions or their derivatives. For $L = 0$, the zeros are $z_n = n\pi$; for $L = 1$ to $L = 11$, the $L + 1$ first zeros are stored as data; all the others are given by Mc Mahon formula which can be written here :

$$z_n = (n + \frac{1}{2}L)\pi - \frac{L(L+1)}{(2n+L)\pi} \left[1 + \frac{7L(L+1)-6}{3(2n+L)^2\pi^2} + \frac{166L^2(L+1)^2 - 408L(L+1) + 360}{15(2n+L)^4\pi^4} + \frac{6949L^3(L+1)^3 - 3352L^2(L+1)^2 + 81180L(L+1) - 75600}{105(2n+L)^6\pi^6} \right] \quad (\text{VIII.1})$$

(see Ref. [67], page 371).

If folding is requested for some form-factors, **STDP** calls **FOLD**.

G.1.c.i Subroutine INTP

Interpolates the form factors between the points given multiplied by a normalisation factor g . It is :

$$V(r) = g \sum_{i=1}^{i=4} V(r_i) \prod_{j \neq i} \frac{r - r_j}{r_i - r_j} \quad (\text{VIII.2})$$

with, as long as possible, $x_1 < x_2 < x < x_3 < x_4$.

G.1.c.ii Subroutine STBF

This subroutine computes a Laguerre polynomials or search on the depth of a real Woods-Saxon potential for a bound state with a given binding energy. With negative binding energy, this subroutine returns a real scattering state normalised to $\text{sink}r + \delta$ at infinity.

For the last calculation, after return from POTE, CAL1 calls SCHE if LO(220)=.TRUE.. In all other cases, the memories needed for the S-matrix, the compound nucleus results and the integrated cross-sections are set to 0. If there is compound nucleus (LO(81)=.TRUE.), the subroutine CONU is called.

G.2 Subroutine CONU

This subroutine does the preliminary calculations for compound nucleus, as introduced in **ANLECIS** by [77] MOLDAUER, P., A., “*ANLECIS*”, presented at the Workshop on Nuclear Model Computer Codes, Trieste (1984). If there are uncoupled states, **CONU** computes their transmission coefficients, using the subroutines **MTCH** and **INSH** or **INRH** in the Schrödinger and the Dirac formalism respectively.

There is a pseudo-loop on parities (first, same parity for the levels and the angular momentum and, last, opposite parities) and inside it, another do-loop on the total angular momentum. Inside these two pseudo-loops, CAL1 calls the subroutine QUAN.

G.3 Subroutine QUAN

This subroutine find the quantum numbers of all the equations coupled for a given total angular momentum J and a given parity. If there are none for the first level, it returns to **CAL1**. The subroutine computes all the coupling between equations and scans if some coefficients are zeros or can be summed up. The coefficients between levels **I1** and **I2** are multiplied by **WV(8,I1)*WV(8,I2)**. Results less than 10^{-10} are eliminated. The address of the first coefficient is stored in **NVI(1,I1,I2)**, the address of the last one in **NVI(2,I1,I2)** and the address of the last derivative one in **NVI(3,I1,I2)**. The subroutine counts the number of couplings coupling between the equations and tests if there will be derivatives of the functions. This subroutine calls the functions **DCGS**, **DJ6J** and **DJ9J**.

G.3.a Function DCGS

This subroutine computes special Clebsch-Gordan coefficients for which the formula involves no summation. For more details, see section (VII.B.1.c).

G.3.b Function DJ9J

Computes 9-j coefficients. The subroutine verifies triangular relations. For more details see section (VII.B.1.d).

After the return from QUAN, if there is no equation for the first level, the pseudo-loop is ended; otherwise, CAL1 calls the subroutine MTCH. Up to the call to subroutine SCAM, all the levels are considered as incoming if compound nucleus or output of transmission coefficients are requested.

G.4 Subroutine MTCH

If `LO(227)=.FALSE.`, as in the beginning of the do-loop on angular momenta of the subroutine `CAL1`, this subroutine computes matching values for each equation and corrections to the S-matrix by integrals from the matching radius to infinity of regular and irregular Coulomb functions divided by some power of r if Coulomb corrections are requested. At the end of the do-loop, it can be called with `LO(227)=.TRUE.` and computes only integrals from 0 to infinity of products of regular Coulomb functions to obtain the reactance K-matrix. `MTCH` calls the subroutines `CORA` to get the coefficients α and β of Equ IV.23 and `LINS` to obtain the S-matrix from a K-matrix. For more details, see section (IV.A).

G.4.a Subroutine CORA

For given angular momenta, this subroutine returns the four coefficients α needed in the asymptotic region in Equ IV.23, if its last argument is `.TRUE.` (this argument is `LO(227)`). When this last argument is `.FALSE.`, it returns also the four other coefficients β needed for finite integrals. For limitations, see `INPUT DESCRIPTION`.

G.4.b Subroutine LINS

This subroutine solves a complex linear system of equations with real and imaginary parts in different array. This subroutine is also called by `INCH`.

After return from MTCH and after some storage evaluations :

- *If `LO(227)=.TRUE.`, the subroutine `CAL1` skips the numerical integration and calls `SCAM` at the end of the loop on total angular momentum,*
- *If `LO(100)=.TRUE.` (Dirac formalism), the subroutine `CAL1` calls first `INTR` and then `SCAM`,*
- *If `LO(21)=.TRUE.` or `LO(225)=.TRUE.` (Schrödinger formalism with usual coupled equations), the subroutine `CAL1` calls first `INCH` and then `SCAM`.*
- *In other cases, the subroutine `CAL1` calls `INTI` and then :*
 - *If convergence is not obtained (`LO(204)=.FALSE.`) and if `LO(23)=.TRUE.` to allow that, `LO(225)` is set `.TRUE.` and the subroutine `QUAN` is called again with the same quantum numbers.*
 - *In any other case, the subroutine `CAL1` calls `SCAM`.*

G.5 Subroutine INTI

This subroutine solves the set of coupled equations by iterations. First, it calls the subroutine `INSH` to solve the diagonal homogeneous equations, unless there is indications from `QUAN` that this solution has been conserved. If all the coupling have to be computed (`LO(207)=.TRUE.`), their are computed and the address of the real and the imaginary coupling between equation `I1` and `I2` are stored in `NVI(1,I1,I2)` and `NVI(2,I1,I2)`, those of the derivative coupling in `NVI(3,I1,I2)` and `NVI(4,I1,I2)`. The system is solved for each equations related to the first level :

- This uncoupled solution and its phase-shift is transferred as the corresponding function of the complete solution, eventually with a preliminary call to the subroutine `INSI` to derive the uncoupled function,
- The iteration procedure is started :
 - If there are Coulomb corrections, the matching conditions are computed, using the phase-shifts known at that stage,

- The subroutine **INSI** is called,
- The equations of which the coupling is too small are counted,
- The maximum phase-shift is computed to be compared to 10^{10} to decide that it is the last iteration.

G.5.a Subroutine INSH

This subroutine solves a single homogeneous equation. It computes the potential and solve the equation with the Numerov or the Modified Numerov method, with a correction term if **LO(26)=.TRUE.**. The matching is with two points. If there is only elastic scattering, computation of transmission coefficients for uncoupled state, the subroutine returns. Otherwise, the solution is normalised, the point from which the solution is larger than a minimum value obtained. For **DWBA (LO(92)=.TRUE.)**, the subroutine returns at this point; otherwise, the irregular solution computed by backward integration down to this point.

G.5.b Subroutine INSI

This subroutine solves the inhomogeneous equation. The second members are obtained by multiplication of the coupled solution with the potentials and the geometrical coefficient or with the coupling if **LO(207)=.TRUE.** (usually quicker). The contribution of the derivative of coupled solution is added. the subroutine compute from which point the inhomogeneous term is not too small. Unless there is no inhomogeneous term :

- the integral of the regular uncoupled function with the inhomogeneous term is computed,
- the phase-shift is obtained,
- if **LO(92)=.FALSE.** (not pure DWBA) the backward integral with the irregular uncoupled function and the solution taking into account correction for the Green's function are obtained in the same loop,
- the subroutine computes the point from which the solution is not too small.

If the difference with last iteration is too large, that the iteration is at least the fourth one, that convergence was obtained for all the other equations in this iteration, **INSI** calls the subroutine **PADE**. If needed, it computes the derivative multiplied by r with formulae analogous to those of **DERI**. This subroutine can be used to do only this last derivation.

G.5.c Subroutine PADE

This subroutine is called by **INSI** and **INRI**. It computes Pade approximants with the results of all the iterations. It evaluates them with all the partial series and returns the mean value of the two nearest successive results if convergence is obtained.

G.5.d Subroutine INCH

This subroutine is called by **CAL1** as an alternative to the call to subroutine **INTI** to use usual coupled equations instead of iterations. It solves the system of coupled equations with the Modified Numerov method. It computes the matrices of coupling in as many points as possible, leaving free the place of a first matrix. Then it computes $V'_{i,j} = V_{i,j} - h^2 \sum_k V_{i,k} V_{k,j} / 12$, shifting the result downwards. These values are used in the integration procedure. If **LO(42)=.TRUE.**, there is a Schmidt's orthogonalisation procedure between the solutions every **ITERM** points. The matching uses the values at two points. The C-matrix is obtained by a call to the subroutine **LINS**. Lines of the C-matrix are computed instead of columns. However, if coulomb corrections, Hauser-Feshbach corrections or punch of penetrabilities are requested, the complete matrix is computed. For Coulomb corrections, another system of linear equations is built and solved with the subroutine **LINS** (see Ref. [7], Equ (35) and explanations above it).

G.6 Subroutine INTR

This subroutine is similar to **INTI**. It calls the subroutines **INRH** and **INRI** instead of **INSH** and **INSI**. The main difference is the storage of tensor form factors.

G.6.a Subroutine INRH

This subroutine solves a single homogeneous equation with the Schrödinger equivalent potential obtained in **POTE**. It computes the potential and solve the equation with the Numerov or the Modified Numerov method, with a correction term if **LO(26)=.TRUE.**. The matching is with two points. If there is only elastic scattering, computation of transmission coefficients for uncoupled state, the subroutine returns. Otherwise, the large component is computed and the point from which it is larger than a minimum value obtained, the two solutions are derived and the small components obtained with the Dirac equation.

G.6.b Subroutine INRI

This subroutine solves the inhomogeneous equation in the Dirac formalism. With tensor interaction instead of deformed spin-orbit, the operations are very similar to those of subroutine **INSI**. The subroutine **INRI** calls the subroutine **PADE** in the same circumstances as subroutine **INSI**. Corrections for finite step size are given by Equ. (III.106) and (III.107).

G.7 Subroutine SCAM

This subroutine computes the transmission coefficients if they are requested. For compound nucleus, **SCAM** computes gamma, fission and find transmission coefficients of uncoupled states related to this total angular momentum. This subroutine stores the results of coupled equations in a two dimensional array in such a way that one index is the total angular momentum and, for a fixed values of the other, the quantum numbers are increased by the difference of total angular momentum. This disposition will make interpolation easier. If iterations are used, the crude results are multiplied by **WV(6,I)** where **I** is the level, to avoid any further factor in the computation of cross-sections. If usual coupled equations are used, they are divided by this quantities to take into account that a line of the matrix has been obtained instead of a column. If compound nucleus is used, results related to this total angular momentum are computed and stored. **SCAM** calls the subroutine **DIAG** to diagonalise the S-matrix if **LO(83)=.FALSE.**.

G.7.a Subroutine DIAG

This subroutine diagonalises a hermitian complex matrix with an extension of Jacobi's method. Real and imaginary values are in different array.

*In subroutine **CAL1**, there are the tests of convergence to end the pseudo-loop on total angular momentum or to shift from nuclear to asymptotic region when there are coulomb corrections.*

*If the pseudo-loops are ended, **CAL1** calls the subroutine **SCHE**, **RESU**, set **LO(217)=.TRUE.**, **LO(222)=.FALSE.** and returns.*

G.8 Subroutine SCHE

This subroutine computes an array of square roots of integer. If **LO(219)=.TRUE.**, the subroutine returns; if **LO(220)=.TRUE.**, it calls **LCSP** if requested and returns. In other cases, this subroutine replaces the values stored by the subroutine **SCAM** by the scattering coefficients in the helicity formalism. First, the subroutine computes a table of quantum numbers. Then, the interpolations are done if they are needed (on the complex values or on their polar expression, by the method used in subroutine **INTP**).

Then, for each total angular momentum and parity, the subroutine computes by recurrence the Clebsch-Gordan coefficients needed to go to the helicity formalism, transforms the scattering coefficients and computes the Hauser-Feshbach corrections if they are requested. The scattering coefficients are symmetrised for identical particle and target with non zero spins. Then, the reactions cross-sections are computed. If `LO(65)=.TRUE.` (Legendre expansions printed and punched), the subroutine `LCSP` is called. If `LO(220)=.FALSE.` and `LO(41)=.TRUE.` (factorisation of $1 - x \cos \theta$), new amplitudes f'^J are computed as described in section (VII.C.3).

G.8.a Subroutine LCSP

This subroutine computes the coefficients of the expansion of the cross-sections in terms of Legendre polynomials. For charges particles elastic scattering, the cross-section is multiplied by $(1 - \cos \theta)^n$ where n is given (default option $n = 2$) for different particle and target, by $(1 - \cos^2 \theta)^n$ for identical particle and target. Only even coefficients are given for compound nucleus contribution. See section (VII.C.4).

G.9 Subroutine RESU

This subroutine prints comparison of computed values and experimental values, computes experimental normalisations and χ^2 , prints results at equidistant angles. In a search, if the computation is ended, the subroutine copies the scattering matrices related to the best χ^2 . Then, for each angular distribution, `RESU` calls the subroutine `SCAT` for each data and computes the partial χ^2 . If the χ^2 is the best obtained up here, the scattering matrices are copied in another array. If complete output is requested, results are printed and a graph can be drawn by a call to the subroutine `GRAL`; cross-sections and polarisations are computed at equidistant angles by the subroutine `SCAT` and a graph can be drawn by the subroutine `GRAL`. If results are requested in the Laboratory system, the angles are transformed by Equ. (VII.23) before the call to subroutine `SCAT` and the cross-sections multiplied by the factor given in Equ. (VII.24).

G.9.a Subroutine SCAT

This subroutine computes the observables. In a first part it computes the helicity amplitudes with the reduced rotation matrix elements obtained by a call to subroutine `EMRO`. A simple loop gives the cross-section in all cases (they are the sum of the squares of the amplitudes multiplied by `XZ` computed in subroutine `LECT`) and the vector polarisation and analysing power for spins 1/2 and 1. The other observables involve do-loops and coefficients which were computed in subroutine `OBSE` and are used after transfer in some memories in equivalence. If observables are in the Laboratory system or with axis of quantification along the incoming direction, the collision matrix is rotated as indicated by Equ. (VII.25), (VII.26) or the angle $-\theta$.

G.9.b Subroutine EMRO

This subroutine computes a series of reduced rotation matrix elements with increasing angular momentum and the same magnetic quantum numbers. It uses upwards recurrence relations. Square roots are avoided by a table of square roots of integer as argument. In fact, there are three independent computations: one if none of the two magnetic quantum numbers are zeros, the second if one of them is zero and the third if both of them are zeros.

G.9.c Subroutine GRAL

This subroutine prints graphs of cross-sections or polarisations. For different values of its last argument, it can print or not experimental data on the same graph.

After the return from CAL1, in the subroutine CALC :

- if there is no automatic search of if it is to its end,
 - if `LO(37)=.TRUE.` the subroutine `EVAL` is called,
 - if `LO(37)=.FALSE.` the working array is given back and the subroutine `CAL1` is called; the computation can be stopped by a card “`FIN`”.
- if there is an automatic search which is not finished,
 - the subroutine `STIM` is called to get the time needed in `CAL1`,
 - the subroutine `VARI` is called with `KF=1`,
 - if `LO(35)=.TRUE.` and it is the last run allowed for the job, the subroutine `REST` is called with `KF=0`,
 - then the subroutine `FITE` is called,
 - the subroutine `VARI` is called with `KF=0`.

H Subroutine VARI

This subroutine is called before `FITE` with its first argument `KF=1` and after the return from `FITE` with `KF=.` Its second argument is the control `KE` of `FITE`. At the first call (`KF=1` and `KE=0`), the subroutine defines the variables in search. For all the calls with `KF=1`, the χ^2 value are only printed. When called after `FITE`, the parameters are set to their values for the next evaluation and if the calculation has to be stopped for any reason, the final parameters are printed.

I Subroutine REST

The first argument of this subroutine, `KF` can be 0 or 1. If `KF=0`, the commons and the beginning of the working array which is needed to restart the calculation are written on tape 8. If `KF=1`, these commons and the working array are read from the tape to restart the search.

J Subroutine FITE

This is the χ^2 minimising subroutine `FITEX` written by G. SCHWEIMER, slightly modified. It calls the subroutines `FIT1` and `FIT2`.

J.1 Subroutine FIT1

Minimisation of the χ^2 for one variable. It is used if there is only one variable in search but also if the subroutine `FITE` find a linear dependency of the χ^2 on the variables.

J.2 Subroutine FIT2

This subroutine replaces the subroutines `LILESQ` and `INVATA` of G. SCHWEIMER. If the last argument `LLO=.TRUE.`, it solves a linear least square problem $\|B - A * D\| = \text{Min}$. If `LLO=.FALSE.`, it inverses the product matrix $A^\dagger A$. The same Householder transformation is used in the two cases.

After return from the second call to `VARI`, `CALC` calls `REDM` if a nuclear parameter has been changed or at the end of a search if the nuclear matrix elements have to be printed or punched. Then, `CALC` calls again `CAL1`.

If `LO(36)=.TRUE.` after return from `CALX`, `CAL1` calls `REST` with `KF=1` to restart a search. Then `CAL1` calls `FITE` if the search is not ended or goes to the second call of `VARI` if the search was ended.

At the end of a search, `CALC` calls `REST` with `KF=0` if `LO(35)=.TRUE.` and calls `EVAL` if `LO(37)=.TRUE.` Otherwise, `CALC` calls again `CALX`.

K Subroutine EVAL

This subroutine is called if `LO(37)=.TRUE.` . All the parameters on which there can be automatic search can be changed for a new calculation. The total energy can also be changed in this subroutine, to simplify the input for the calculation of an excitation function.

After the call to EVAL, the subroutine CALC calls again the subroutine COLF if the total energy has been changed, the subroutine REDM if some nuclear parameters or nuclear matrix elements are modified, the subroutine CAL1 in the other cases.

L Summary

Without the calls to `MEMO`, `HORA` and `STIM`, the structure of the code is given by the following table :

	CALC	CALX	(INPA, INPB, INPC, LECL, LECT, DEPH, LECD, OBSE)
	''	COLF	FCOU FCZO.....
	''	''	COCL
	''	''	CORI CORO FCOU FCZO.....
	''	''	CORZ
	''	REDM	VIBM (DJCG, DJ6J)
	''	''	ROTM DJCG
	''	''	ROAM DJCG
	''	EXTP	DJCG
	''	CAL1	POTE ROTP (ROTD, ROTZ, COPO, DERI)
	''	''	STDP STBF FCOU FCZO.....
	''	''	'' FOLD DERI
	''	''	'' (INTP, COPO, DERI)
	''	''	FOLD DERI
	''	''	CONU MTCh (CORA, LINS)
	''	''	'' (INSH, INRH)
	''	''	QUAN (DCGS, FJ6J, DJ9J)
	''	''	MTCH (CORA, LINS)
	''	''	INTI INSH
	''	''	'' INSI PADE
	''	''	INCH (INSH, LINS)
	''	''	INTR INRH
	''	''	'' INRI PADE
	''	''	SCAM DIAG
	''	''	SCHE (DJ6J, LCSP)
	''	''	RESU SCAT EMRO
	''	''	'' GRAL
	''	VARI	
	''	REST	
	''	FITE	(FIT1, FIT2)
	''	EVAL	

where (PR01,PR02,PR03) means that these 3 subroutine are called by the subroutine written on the left and call nothing by themselves. For Coulomb functions, this Table must be complete by :

	FCOU	FCZO	YFAS
	''	''	YFIR (YFAS,PSI)
	''	''	YFRI YFCL PSI
	''	''	SIGM

which occurs in 3 places.

Chapter IX

Conclusion

A Test cases

Nine test cases have been chosen to help to use **ECIS94**. There are quite arbitrary. There are identified by **TSTx** in columns 73-80.

The first of them, with the titles :

ESSAI DU NOYAU COMPOSE - USUAL COUPLED EQUATIONS	TST1-000
ESSAI DU NOYAU COMPOSE - USUAL ITERATIONS	TST1-024
ESSAI DU NOYAU COMPOSE - ITERATIONS WITH DEFORMED SPIN-ORBIT POTENTIAL	TST1-048

is compound nucleus at low energy. The same calculation with 3 levels and 2 uncoupled levels is done :

- without spin-orbit deformation, without iterations and without angular distribution of uncoupled states,
- without spin-orbit deformation, with iterations and with angular distribution of the first uncoupled states,
- with spin-orbit deformation, with iterations and with angular distribution of the two uncoupled states,

Results show that the spin-orbit deformation does not matter in this case.

The second test case, with the title :

28 SI + 104 MEV ALPHAS, SYMMETRIC ROTATOR, NON STANDARD OBSERVABLES	TST2-000
---	----------

shows different ways to define observables.

The third test case, with the title :

CA40 A 497 MEV RELATIVISTE SAME AS ECIS88	TST3-000
---	----------

shows the use of Dirac formalism.

The fourth test case, with the title :

12N+208PB 780 MEV ELECTIC AND MAGNETIC COULOM INTERACTION	TST4-000
---	----------

use a magnetic $M1$ and electric $E2$ Coulomb interactions.

The fifth test case, with the title :

COMPARAISON AVEC CHUCK POUR BE8	DWBA	TST5-000
COMPARAISON AVEC CHUCK POUR BE8	COUPLED EQUATIONS	TST5-047

is the example of section (VI.D.2) and shows that the result with coupled equations is quite different of the result of DWBA.

The sixth test case, with the title :

TEST EXTERNAL POTENTIALS	INPUT AS ECIS88	TST6-000
TEST EXTERNAL POTENTIALS	NEW INPUT	TST6-106

compare an input with external potentials as needed by ECIS88 with one of the new possibilities of ECIS94.

The seventh test case, with the title :

ESSAI PARTICULES IDENTIQUES ET POLYNOMES DE LEGENDRE	TST7-000
--	----------

is a calculation with identical particle of spin one half. The expansion in Legendre polynomials of the cross-sections are written on file 7.

The eighth test case, with the title :

28SI + 160	56.MEV	G.S. , 1.78 MEV.	HEAVY IONS OPTION	TST8-000
28SI + 160	56.MEV	G.S. , 1.78 MEV.	RESTART	TST8-110

is an automatic search on 3 parameters which needs about 24 calculations. After 12 calculations, the job is saved on file 8 and restarted to the end.

The ninth test case, with the title :

NEUTRON DE 14. MEV SUR GADOLINIUM	COUL. SPIN-ORBITE DEFORME	TST9-000
NEUTRON DE 14. MEV SUR GADOLINIUM	COUL. SPIN-ORBITE NON DEFORME	TST9-019

is a calculation for neutron with spin-orbit Coulomb potential. The two calculations, with and without deformation of this interaction gives almost the same results.

B Missing topics

Many points quite important are not in this report :

- a detailed description of the reduced matrix elements used in the macroscopic models,
- relation between form factors of ECIS94 and of DWBA91
- more details on identical particle and target,
- more details on the use of form factors expanded with Bessel functions,
- details on the manipulation of symmetrised Woods-Saxon form factors,
- explanations on the interpolations,
- and so on.

The most important point missing here is **Compound Nucleus**.

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