Entrance-channel potentials in the synthesis of the heaviest nuclei

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Abstract. Entrance-channel potentials in nucleus-nucleus collisions, relevant for the synthesis of superheavy elements, are systematically studied within a semi-microscopic approach, where microscopic nuclear densities of the colliding spherical or deformed nuclei are used in semi-classical expressions of the energy-density functional. From experimental data on fusion windows evidence is found that the existence of pockets in the entrance-channel potentials is crucial for fusion. Criteria for the choice of best collision systems for the synthesis of superheavy elements are discussed.

PACS. 25.60.Pj Fusion reactions – 25.70.Jj Fusion and fusion-fission reactions

1 Introduction and summary

The formation of compound nuclei by fusion of very heavy nuclei is one of the outstanding problems of low-energy nuclear reactions. Such processes play a key role in the production of superheavy elements (SHEs) \cite{1-19} and of heavy nuclei far from the $\beta$-stability line, as for example $^{217}$U \cite{20}.

Recently, various properties of SHEs have been studied experimentally as well as theoretically \cite{1-8,20-29}. However, the dynamical processes leading to SHEs in heavy-ion fusion reactions are not yet well understood. Few attempts have been made to develop models, which yield cross-sections for SHE production \cite{11-13,14-16,17-18}. These models are mainly concerned with the evolution from the touching configuration of the colliding nuclei to the compound nucleus. The dynamics has been treated as diffusion in mass asymmetry at the touching point \cite{11-13}, as collective motion in 3-dimensions within Langevin equations \cite{14}, as a tunneling process of outer barrier \cite{15,16}, and as shape evolution and tunneling \cite{17,18}.

We suggest to consider the fusion process as part of the quantum-mechanical collision process \cite{30}, where the features of the entrance channel are essential for the scattering solution. In such a formulation the population of quasi-bound states near touching is regarded as the first decisive step for fusion and is similar to the doorway-state mechanism in resonance scattering \cite{31,32}. The existence of such capture states depends crucially on the detailed properties of the entrance-channel potential as a function of the distance $R$ between the nuclei, in particular on the existence of a pocket which serves as a source for the quasi-bound states. If the capture pocket is deep and wide, many quasi-bound states exist. Then the coupling to complex states is strong and the probability of compound-nucleus formation is much larger than for shallow pockets. Because of its crucial role on the initial stage of the fusion process, a precise and systematic knowledge of the interaction potentials between the colliding nuclei is needed.

For determining the interaction potentials between two nuclei various methods have been introduced. Early attempts are due to Bass \cite{33,34}, who parametrized a simple analytical expression for the potential, to Swiatecki and coworkers \cite{35}, who introduced the proximity interaction of leptodermous systems and to Krappe, Nix and Sierk, who used the folding procedure with a phenomenological Yukawa-plus-exponential interaction \cite{36}. Recently, a folding procedure using the Migdal force \cite{37} has been used \cite{11,12,13,14} together with phenomenological density distributions. Here like in \cite{12,13,14,38} a simple point-point interaction $Z_1 Z_2 e^2 / R$ (charges $Z_1$, $Z_2$) has been used for the Coulomb part at distances larger than the touching point. A quite large variance of the interaction potentials near touching are encountered, when these different methods are used. These uncertainties allow for quite different conclusions on the fusion mechanism in the synthesis of SHEs \cite{11,12,13,14}. Therefore, the precise knowledge of the nucleus-nucleus potential near the touching point may also improve our qualitative understanding of the fusion process.

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We evaluate the interaction potential by keeping the densities of the colliding nuclei fixed. Due to the short collision time this frozen-density approximation is appropriate for the definition of the entrance-channel potential (cf. section 2). The interaction energy between the nuclei is obtained from the energy density functional according to a definition which has been introduced by Brueckner et al. [39]. The Thomas-Fermi approximation, extended to all second-order gradient terms (ETF), is used for the evaluation of the kinetic-energy density functional [14,11], while the nuclear interaction energy is obtained from a Skyrme energy-density functional. The nucleon densities of the colliding spherical or deformed nuclei are calculated in the microscopic Hartree-Fock-Bogoliubov approximation. Our method is semi-microscopic, because we use the microscopic nucleon densities and the semiclassical ETF approach and energy densities. We refer to the interaction potential, which is obtained from this semi-microscopic approach, as the semi-microscopic potential (SMP). Since the microscopic HFB densities are used in the extended semi-classical energy-density-functional for the semi-classical wave function to more compact configurations. Therefore, we are interested to understand as a first step the basic properties of the entrance-channel wave function to more compact configurations. It is obvious that the energy dependence of the entrance-channel wave function strongly affects the crucial steps in such a description is the coupling of a large number of channels which include inelastic excitations and transfer, quasi-fission, compound-nucleus formation and fission. In order to cope with this complexity one has to formulate the stationary scattering problem in a suitable way, which is the subject of a forthcoming paper [30]. One of the crucial steps in such a description is the coupling of the entrance-channel wave function to more compact configurations. It is obvious that the energy dependence of the entrance-channel wave function strongly affects the cross-sections for all processes including those which involve compact-shape configurations. Therefore, we are interested to understand as a first step the basic properties of the entrance channel in fusion reactions. We evaluate targets on the waist-line (90° with respect to its symmetry axis). If this is assumed also for the experimentally studied fusion reactions with 48Ca on prolate uranium and transuranium nuclei we find the same location of the fusion window as in the lead-target systems, i.e. 5 to 10 MeV below the SMP barrier. 48Ca projectiles are less favorable, because they lead to higher excitation energies, which reduces the evaporation-residue cross-section. The systematic study of 238U and 252Cf target systems show deeper pockets for 238U and even more for 252Cf than for 208Pb target systems, however at higher compound-nucleus excitation energies. Only for very heavy systems like 82Se+238U=320126 or 68Ni+252Cf=320126 we encounter low excitation energies like in cold-fusion systems.

The warm-fusion systems, formed with the oblate 198Pt target, yield compound-nucleus excitation energies which lie in between those with 208Pb and with 238U or 252Cf. Pockets are obtained up to 106Mo+198Pt=298120, where the compound-nucleus excitation energy is comparable to cold-fusion systems (section 3.3).

In conclusion, there are indications that the existence of quasi-bound capture states in the entrance-channel potential pocket is crucial for the synthesis of SHEs. These capture states are only populated in an energy window with a mean value 5 to 10 MeV below our SMP barrier and a width roughly given by the depth of the pocket. Whenever deformed nuclei are involved, fusion is expected through the most compact capture states (pockets). Applying these considerations to the synthesis of SHE 118, we conclude that there are two favorable systems: the hot-fusion system (which actually is more like a warm-fusion system) 48Ca+252Cf at about 206 MeV center-of-mass energy and the warm-fusion system (which actually is an almost cold-fusion system) 96Zr+198Pt at about 330 MeV center-of-mass energy (section 4).

2 Definition of a semi-microscopic potential (SMP) in the entrance channel

The ultimate aim is the description of fusion as part of the quantum-mechanical scattering process between two nuclei. In such a formulation one has to consider all kinds of channels which are coupled (in general non-perturbatively) to the entrance channel of the colliding nuclei. Thus one faces the problem of describing the coupling of a large number of channels which include inelastic excitations and transfer, quasi-fission, compound-nucleus formation and fission. In order to cope with this complexity one has to formulate the stationary scattering problem in a suitable way, which is the subject of a forthcoming paper [30]. One of the crucial steps in such a description is the coupling of the entrance-channel wave function to more compact configurations. It is obvious that the energy dependence of the entrance-channel wave-function strongly affects the cross-sections for all processes including those which involve compact-shape configurations. Therefore, we are interested to understand as a first step the basic properties of the entrance channel in fusion reactions. We evaluate
the interaction potential of the entrance channel in a semi-
microscopic approach keeping the densities of the colliding
nuclei fixed. In the following, details of definitions and just-
ifications are given for these semi-microscopic potentials
(SMPs).

2.1 Entrance channel dynamics

Whereas the entrance channel (elastic channel) is uniquely
defined outside the range of nucleus-nucleus interactions, there
is considerable freedom inside. In principle, any def-
inition, which asymptotically describes two nuclei in their
ground states, is possible if all relevant inelastic channels
are included in the solution of the scattering problem.
However, from a physical point of view a reasonable def-
inition should account for the essential collective dynamics
in the region of overlap.

A crucial quantity, which characterizes the collective
dynamics in the entrance channel during the capture pro-
cess, is the nuclear interaction time \( \tau_{\text{coll}} \) (collision time)
as compared to the characteristic time \( \tau_{\text{relax}} \) for the relax-
ation of the intrinsic nuclear state due to nucleon-nucleon
interactions. The collision time \( \tau_{\text{coll}} \) may be estimated by

\[
\tau_{\text{coll}} \approx \frac{\pi}{\omega_{\text{pocket}}} = \pi \left[ \frac{mA_1 A_2}{(A_1 + A_2)V_n(R_{\text{pocket}})} \right]^{1/2},
\]

where \( \omega_{\text{pocket}} \) denotes the “oscillator frequency” of the
interaction potential pocket \( V(r) \), \( R_{\text{pocket}} \) is the center-to-
center distance at the pocket, while \( m, A_1 \\text{ and } A_2 \) are the
bare nucleon mass and the number of nucleons in the nu-
clei. A typical value for \( \omega_{\text{pocket}} \) in reactions used for SHE
production is given by \( \hbar \omega_{\text{pocket}} \approx 4 \text{ MeV} \). Therefore, typi-
cal collision times for such cases are \( \tau_{\text{coll}} \approx 5 \cdot 10^{-22} \text{ s} \). This
value has to be compared to typical times for the relaxation
of the intrinsic nuclear state due to nucleon-nucleon
interactions. This time is estimated as \([4]\]

\[
\tau_{\text{relax}} \approx \frac{\epsilon_F}{3.2 \sigma v_F \rho_0 \epsilon^*} \approx \frac{2 \cdot 10^{-22}}{\epsilon^*/\text{MeV}} \text{s},
\]

where \( \epsilon_F \) and \( v_F \) denote the Fermi energy and velocity,
respectively, \( \sigma \) the averaged nucleon-nucleon cross-section,
\( \rho_0 \) the normal density of nuclear matter and \( \epsilon^* \) the ex-
citation energy per nucleon. For reactions leading to SHEs
we have at touching \( \epsilon^* \lesssim (5 \text{ MeV})/250 \approx 0.02 \text{ MeV} \), and
hence \( \tau_{\text{relax}} \gtrsim 10^{-20} \text{ s} \), i.e. more than one order of magni-
tude larger than the collision time \( \tau_{\text{coll}} \). Thus we conclude
that the entrance channel in the region of nucleus-nucleus
overlap is well defined by the fixed configuration of the
colliding nuclei. With respect to the nucleus-nucleus poten-
tial in the entrance channel this means that the interact-
ion energy of two overlapping nuclei with frozen densities
is relevant.

The frozen-density potential should however not be ap-
lPLIED to very large overlap. A suitable continuation into re-
 gions of compact shapes would be the diabatic energies of
the entrance-channel configuration \([15,16]\). Furthermore,
this frozen-density potential should be regarded as a suit-
able reference for a distribution of diabatic potentials (and
barriers) which are due to the mixture of configurations in
the approaching nuclei \([40]\).

2.2 Frozen-density potential

We consider the interaction potential \( V(R, \Theta) \) between a
spherical nucleus and an axially symmetric nucleus as a
function of the center-to-center distance \( R \) between the
nuclei and the angle \( \Theta \) between the symmetry axis of the
deformed nucleus and the line connecting the centers of
the nuclei. Denoting the energies of the interacting nuclei by
\( E_{12}(R, \Theta) \) and the energies of the non-interacting nuclei
by \( E_1, E_2 \), we define the interaction potential by

\[
V(R, \Theta) = E_{12}(R, \Theta) - E_1 - E_2.
\]

In the frozen-density approximation these energies are de-
termined by the energy-density functional \( \mathcal{E} = \mathcal{E}[\rho_p(r), \rho_n(r)] \),
i.e. \([39]\)

\[
E_{12}(R, \Theta) = \int \mathcal{E}[\rho_{1p}(r) + \rho_{2p}(R, \Theta, r), \rho_{1n}(r) + \rho_{2n}(R, \Theta, r)] \, dr,
\]

\[
E_1 = \int \mathcal{E}[\rho_{1p}(r), \rho_{1n}(r)] \, dr,
\]

\[
E_2 = \int \mathcal{E}[\rho_{2p}(r), \rho_{2n}(r)] \, dr,
\]

where \( \rho_{1p}, \rho_{2p}, \rho_{1n} \text{ and } \rho_{2n} \) are the frozen proton and
neutron densities of the spherical nucleus (index 1) and the
deformed nucleus (index 2), respectively.

2.3 Energy-density functional

For an accurate calculation of the interaction potential
between two nuclei we need an energy-density functional
which well describes both the bulk and surface prop-
ties of the nuclei. Suitable semi-classical expressions have
been obtained for Skyrme interactions and by an extended
Thomas-Fermi (ETF) approximation to the intrinsic ki-
netic energies, which includes all terms up to second order
in the spatial derivatives.

According to \([10]\) the following expression for the energy-
density functional

\[
\mathcal{E}[\rho_p(r), \rho_n(r)] = \frac{\hbar^2}{2m}[\tau_p(r) + \tau_n(r)] + \mathcal{V}_{sk}(r) + \mathcal{V}_c(r)
\]

has been deduced. The kinetic parts for protons \((i = p)\) and
neutrons \((i = n)\) are given by

\[
\tau_i(r) = \frac{3}{5}(3 \pi^2)^{5/3} \rho_i^{2/3} + \frac{1}{36} \left( \frac{\nabla \rho_i}{\rho_i} \right)^2 + \frac{1}{3} \Delta \rho_i + \frac{1}{6} \left( \frac{\nabla \rho_i}{\rho_i} \right)^2 + \frac{1}{12} \rho_i \left( \frac{\nabla \rho_i}{\rho_i} \right)^2 \right)^2,
\]

where \( \mathcal{V}_{sk}(r) \) and \( \mathcal{V}_c(r) \) are the Skyrme and core
potentials, and \( \rho_i \) is the density of the \( i \)-type nucleons.

\[
\mathcal{V}_{sk}(r) = \frac{3}{5}(3 \pi^2)^{5/3} \rho_i^{2/3} + \frac{1}{36} \left( \frac{\nabla \rho_i}{\rho_i} \right)^2 + \frac{1}{3} \Delta \rho_i + \frac{1}{6} \left( \frac{\nabla \rho_i}{\rho_i} \right)^2 + \frac{1}{12} \rho_i \left( \frac{\nabla \rho_i}{\rho_i} \right)^2 \right)^2,
\]

where \( \mathcal{V}_c(r) \) is the core potential, and \( \rho_i \) is the density of the \( i \)-type nucleons.

\[
\mathcal{V}_c(r) = \frac{3}{5}(3 \pi^2)^{5/3} \rho_i^{2/3} + \frac{1}{36} \left( \frac{\nabla \rho_i}{\rho_i} \right)^2 + \frac{1}{3} \Delta \rho_i + \frac{1}{6} \left( \frac{\nabla \rho_i}{\rho_i} \right)^2 + \frac{1}{12} \rho_i \left( \frac{\nabla \rho_i}{\rho_i} \right)^2 \right)^2,
\]
where \( W_0 \) denotes the strength of the Skyrme spin-orbit interaction, while \( \rho = \rho_p + \rho_n \) and

\[
f_i(r) = 1 + \frac{2m}{\hbar^2} \left( \frac{3t_1 + 5t_2}{16} + \frac{t_2x_2}{4} \right) \rho_i(r). \tag{7}\]

The nuclear interaction part \( V_{sk} \) results from the Skyrme force and reads

\[
V_{sk}(r) = \frac{t_0}{2} \left[ (1 + \frac{1}{2}x_0)\rho^2 - (x_0 + \frac{1}{2})(\rho_p^2 + \rho_n^2) \right] + \frac{1}{12}t_3\rho^3 \left[ (1 + \frac{1}{2}x_3)\rho^2 - (x_3 + \frac{1}{2})(\rho_p^2 + \rho_n^2) \right] + \frac{1}{4}[t_1(1 + \frac{1}{2}x_1) + t_2(1 + \frac{1}{2}x_2)]\rho^2 \\
+ \frac{1}{4}[t_2(x_2 + \frac{1}{2}) - t_1(x_1 + \frac{1}{2})]((\tau_p\rho_p + \tau_n\rho_n) \\
+ \frac{1}{16}[3t_1(1 + \frac{1}{2}x_1) - t_2(1 + \frac{1}{2}x_2)](\nabla\rho)^2 \\
- \frac{1}{16}[3t_1(1 + \frac{1}{2}x_1 + \frac{1}{2}x_2) + t_2(1 + \frac{1}{2}x_2)](\nabla\rho_0)^2 + (\nabla\rho_0)^2) \\
- \frac{W_0^2}{4}\frac{2m}{\hbar^2} \left[ \frac{\rho_p}{\hbar^2} (2\nabla\rho_p + \nabla\rho_n)^2 + \frac{\rho_n}{\hbar^2} (2\nabla\rho_p + \nabla\rho_n)^2 \right],
\]

where \( t_0, t_1, t_2, x_0, x_1, x_2, \alpha \) and \( W_0 \) are Skyrme-force parameters. The Coulomb-energy density is determined by

\[
V_c(r) = \frac{e^2}{2\rho_p(r)} \int \frac{\rho_p(r')}{|r - r'|} dr' - \frac{3e^2}{4} \left( \frac{3}{\pi} \right)^{1/3} (\rho_p(r))^{4/3}, \tag{9}
\]

where the last term is the local approximation to the exchange contribution.

Thus, if the proton and neutron density distributions in both nuclei are known, the interaction potential can be calculated from the semi-classical expressions (1)–(9).

### 2.4 Determination of density distributions

The charge densities of nuclei are well described in Hartree-Fock, Hartree-Fock-Bogoliubov (HFB) and semi-classical approaches [11,12,28]. However, HFB describes best various other ground-state properties of nuclei [11,12,28] and therefore has been chosen here for calculating the density distributions.

For spherical nuclei we have used the HFB code [47] with very small radial mesh-point intervals of 0.025 fm. The gradients and laplacians of the corresponding densities are evaluated numerically by using Lagrange formulas. The density distributions of deformed nuclei are also obtained in the HFB approximation by using the code HFODD (v. 1.75r) [13]. The densities, gradients and laplacians at any point of space are found by Lagrange interpolation based on output from the HFODD code. This code calculates the proton and neutron densities and their gradients and laplacians at special points which are related to the Gauss-Hermite integration. For an accurate evaluation of the densities and their derivatives we use 48 points for the Gauss-Hermite integration for each of the three space directions.

Note that the semi-classical approach based on microscopic HFB densities is quite accurate for the calculation of binding energies. For example, the differences of the binding energies between HFB and the semi-classical expression are only –0.9 MeV for \(^{48}\)Ca and –0.3 MeV for \(^{208}\)Pb, when the Skyrme force SkM* is used. Moreover, differences in the energies cancel to a large extent in the expression (1), such that the accuracy is expected to be even higher for the nucleus-nucleus potential SMP.

### 2.5 Relation to similar calculations

The frozen-density approximation has been frequently applied [11,12,28,47] in determining the nucleus-nucleus potentials for the synthesis of SHEs. In contradistinction to our approach, however, there are several differences. For example in [11,12,28,47] the Coulomb interaction potential is approximated by that of two point charges or of homogeneous charge distributions, respectively. In [11,12,28] the nuclear part of the nucleus-nucleus potential is calculated from a folding procedure with Landau-Migdal interactions [17]. However, since this effective interaction is tailored to describe the force between quasiparticles in the Fermi-liquid, the saturation properties of nuclear matter cannot be obtained. Furthermore, the nuclear density distributions, used for the folding procedure in [11,28], are chosen as Fermi distributions with fitted values of the radius and the diffuseness parameters. The kinetic energy density was limited in [50] to the Thomas-Fermi contribution which is the first term in (7). Similar calculations for light nuclei have been performed in [51].

### 3 Entrance-channel potentials

As described in section 2, the nucleus-nucleus potential is evaluated by numerical integration of the semi-classical energy-density functional with the (frozen) HFB nucleon densities of the separated nuclei. The integrals in Eqs. (2)-(4) are evaluated by using the Gauss-Legendre method with a suitable number of dots. The numerical integration is performed in cylindrical coordinates. The system of two spherical nuclei is axially symmetric, and hence the nuclear part of the potential is reduced to the calculation of 2-dimensional integrals, while the Coulomb part is obtained numerically from 3-dimensional integrals. Since axial symmetry is lost for the system of a spherical nucleus and a deformed nucleus, the nuclear and Coulomb parts have to be calculated from 3- and 5-dimensional integrals, respectively.

In Fig. 1a we present the SMPs for \(^{58}\)Fe + \(^{64}\)Ni evaluated for different Skyrme forces [12,28,51]. The potentials obtained for SkM*, SkP and SLy4 are very close to each other at all distances down to \( R = 6 \) fm. Considerable repulsion is observed when the nuclei overlap and
the density is doubling. Due to the large value of the compression modulus (stiff equation of state), this repulsion is particularly large for the SIII force and causes the discrepancy with respect to the other Skyrme forces. In the following sections 3.1 to 3.3 either SkP or SkM\(^*\) is used, which practically give the same interaction potentials.

The ground-state Q-values \((Q = E_{\text{CN}} - E_1 - E_2)\) is obtained from the experimental energies \(E_1, E_2\) of projectile and target taken from [56]. The compound-nucleus energy \(E_{\text{CN}}\) is also obtained from this table or, if experimental values are absent, from the Thomas-Fermi approach to nuclear masses [53]. The neutron separation energies, which mark the thresholds for neutron emission in Fig. 1, are also deduced from [56,57].

Due to the deep pocket inside the barrier, light ions easily fuse after tunneling through or passing over the barrier. Both, the barrier height and the potential pocket are well above the ground-state energy, such that the adiabatic potential surface exhibits large gradients in the fusion direction driving the system into the compound-nucleus shape.

The barriers obtained from different analytical expressions for the nucleus-nucleus potential introduced by Bass in 1974 [33] and in 1980 [34], by Swiatecki et al. (proximity 77) [35] and by Krappe-Nix-Sierk (KNS) [36] are spread over a large interval as shown in Fig. 1b. The difference between the barriers is around 3 MeV in energy and 1 fm in position. The KNS barrier is closest to the SMP barrier. Since the KNS potential depends on the shape parametrization for distances smaller than touching, results for this potential is presented only up to the touching point, which is given by \(R = r_0(A_1^{1/3} + A_1^{1/3})\) with \(r_0 = 1.2\) fm.

### 3.1 Cold-fusion systems

In this section we present results on entrance-channel potentials of spherical projectiles and targets, which have been used in the synthesis of SHEs by cold-fusion [1,2,3,4], and compare these potentials with those of symmetric systems with almost equal projectiles and targets.

#### 3.1.1 Cold-fusion systems with \(^{208}\text{Pb}\) targets

Cold-fusion reactions with light projectiles \((Z \approx 20...36)\) on lead or bismuth targets have been used in the synthesis of SHEs [1,2,3,4]. In Fig. 2 the interaction potentials are presented for the systems with lead as target and 8 projectiles from \(^{48}\text{Ca}\) till \(^{96}\text{Kr}\). The following features are observed.

- The interaction potentials, which are obtained from different standard expressions [34,35,36], are spread over even larger intervals for heavier systems as compared to \(^{58}\text{Fe}+^{64}\text{Ni}\) (cf. Fig. 1b). As for the light system the KNS potential is closest to our SMP around the barrier followed by the Bass potential from 1980 [34].
- The potential pockets are much shallower than for \(^{58}\text{Fe}+^{64}\text{Ni}\) and tend to vanish with increasing size of the projectile. For \(^{96}\text{Zr}+^{208}\text{Pb}\) (not shown in Fig. 2) no pocket exists anymore.
- Since we consider the depth of the pockets to be important for the fusion probability, we attribute the observed reduction of SHE formation with increasing size of the projectile, at least partially, to the decreasing pocket depth.
- The observed fusion windows (vertical thick bars in Fig. 2) lie systematically about 5 to 10 MeV below our barriers. This subbarrier fusion leading to SHEs is probably related to the distribution of barriers and pockets due to the structure of the approaching nuclei [15] and/or to transfer and capture in transfer-channel pockets, and possibly to other degrees of freedom.
- A correlation is indicated between the width of the observed fusion window and the depth of potential pocket (see cases \(^{50}\text{Ti}+^{208}\text{Pb},\ \ ^{58}\text{Fe}+^{208}\text{Pb}\) and \(^{64}\text{Ni}+^{208}\text{Pb}\) in Fig. 2).
The difference between the barrier position and the ground-state $Q$-value for fusion decreases with increasing charge of the projectile, cf. \cite{4,19}. Due to the small differences between the barrier heights and the ground-state energies large gradients towards the compound nucleus are missing in the adiabatic potential. Therefore, the shape evolution from the touching configuration to the compound nucleus becomes extremely sensitive to details of the potential landscape in cold-fusion reaction for SHEs with $Z \gtrsim 110$ \cite{17,18}. Note that the observed fusion windows in cold-fusion reactions lie about $(10 – 15)$ MeV above the ground-state $Q$-value \cite{1,4}. 

**Fig. 2.** SMPs for the collision systems $^{48}$Ca+$^{208}$Pb, $^{50}$Ti+$^{208}$Pb, $^{54}$Cr+$^{208}$Pb, $^{58}$Fe+$^{208}$Pb, $^{64}$Ni+$^{208}$Pb, $^{70}$Zr+$^{208}$Pb, $^{76}$Ge+$^{208}$Pb, and $^{86}$Kr+$^{208}$Pb obtained with the SkP Skyrme force as compared to the proximity potential (1977), the Bass potentials (1974, 1980) and the KNS potential. The ground-state $Q$-values are indicated by the lowest triangles at the left vertical axes. The other 6 triangles mark, respectively, the thresholds for the emission of 1, 2, 3, 4, 5 and 6 neutrons. The observed fusion windows [1–4] are indicated by the vertical bars on the right-hand sides of the individual diagrams.
3.1.2 Symmetric systems

SHEs might also be formed by the fusion of similar nuclei, such as Sn+Xe, Xe+Xe, Xe+Nd and others. The potentials for the collision systems 124Sn+130Xe, 124Sn+136Xe, 130Xe+136Xe, 136Xe+136Xe, 138Xe+150Nd and 144Ce+150Nd are presented in Fig. 3. As compared to the cold-fusion systems with 208Pb targets (Fig. 2) the differences between the barrier heights and the ground-state Q-values are smaller for symmetric systems leading to the same SHE. Also the potential pockets are shallower and at somewhat larger distances. Therefore, as compared to the asymmetric systems of cold-fusion, the symmetric systems seem to be less favorable because of the following reasons.

- The capture process is suppressed by the shallowness of the potential pocket.
- The shape of the system at capture is less compact, and hence a longer shape evolution is needed to reach the compound-nucleus configuration, such that the formation probability of the compound nucleus is reduced by the larger competition of other decays.
- The capture windows are expected to lie 5 to 10 MeV below the barriers, and hence are, apart from the lightest system 124Sn+130Xe, below the In fusion threshold.

3.2 Hot-fusion systems

The fusion of light nuclei with heavy deformed nuclei (U, Pu, Cm) leads already for collisions at barrier energies to large excitation energies of about (30 – 50) MeV in the compound nucleus. Therefore, in such hot-fusion reactions typically 3 to 5 neutrons have to be emitted in order to reach the compound-nucleus ground state. The projectiles are spherical and the targets are well-deformed prolate nuclei in these hot fusion reactions.

3.2.1 Ca projectiles

Successful syntheses of SHEs with Z = 112, 114 and 116 have been reported for 48Ca projectiles on the targets 238U, 244Pu and 248Cm. In this section we discuss the entrance-channel potentials for the systems 48Ca+238U, 244Pu, 248Cm and 252Cf and compare these with the corresponding potentials for 40,42Ca.

The interaction potentials as obtained from our semi-microscopic method are shown in Fig. 4 for various orientations of the deformed nuclei. For these systems the lowest barriers are obtained for $\Theta = 0^\circ$, i.e. when 48Ca touches the side of the deformed nucleus, while the barrier is highest for $\Theta = 90^\circ$, when 48Ca touches the side (cf. 40,42Ca for lighter systems).

Outside the range of nuclear interactions the Coulomb interaction tends to rotate the deformed nucleus into the $\Theta = 90^\circ$ position (side position). However, the time for such a rotation is typically

$$\tau_{\text{rot}} \approx \frac{\pi}{2} \frac{1}{\omega_{\text{rot}}} = 2 \cdot 10^{-20} \text{ s},$$

where a rotational energy $\hbar \omega_{\text{rot}} \approx 50$ keV has been inserted. Characteristic collision times on the approaching part of the Coulomb trajectory are of the order $2 \cdot 10^{-21}$ s, and hence the angle of rotation during the approach is negligible.

As seen in Fig. 4 the positions of barriers and pockets depend quite strongly on the orientation of the deformed nucleus, the shift in distance and energy reaching 2 fm and 20 MeV, respectively, between tip and side positions. In general, the difference between the barrier and the ground-state energy of the compound system is considerably higher in these very asymmetric (hot-fusion) systems as compared to the cold-fusion systems of section 3.1. Therefore, SHEs can be formed only in 3n to 4n reactions, which reduces the survival probability strongly due to the small branching ratio $\Gamma_n/\Gamma_f$ of neutron emission to fission.

The bombarding energies in the fusion reactions 48Ca on 238U, 244Pu, 248Cm have been determined from the maximum cross-section for symmetric fission events which indicate the formation of a compact relatively long-living compound nucleus. These energies are indicated by vertical bars in Fig. 4. For the cold-fusion reactions with 208Pb targets (cf. Fig. 2) the fusion windows lie about 5 to 10 MeV below our SMP barriers. If we assume that the fusion mechanisms are similar also for the deformed targets – and we do not see any reason to doubt this – we have to conclude from Fig. 4 that the side orientation ($\Theta \approx 90^\circ$) is the relevant fusion channel for the formation of the SHEs. This conclusion is supported by the experimental analysis of fusion reaction between lighter nuclei which show that fusion through the tip orientation ($\Theta = 0^\circ$) is strongly suppressed. Moreover, narrow fusion windows for the synthesis are expected also for hot-fusion reactions. On the basis of these considerations we expect the fusion window for the synthesis of SHE 118 in the collision 48Ca on 252Cf around 206 MeV.

Since the $\alpha$-decay chain of the superheavy isotopes, formed in the 48Ca induced reactions, do not reach the area of known isotopes, it is difficult to decide on the isotope formed in the synthesis. Lighter Ca isotopes ($^{40,42}$Ca) would lead to isotopes formed earlier in cold-fusion reactions, if the number of evaporated neutrons in 48Ca and 40,42Ca induced reactions are the same. Although the pockets have the same or even slightly larger depths, the excitation energies are considerably larger for 40,42Ca as compared to the 48Ca induced reactions. Therefore SHE formation is probably too much suppressed for reactions with 40,42Ca projectiles.

3.2.2 Systems with 238U and 252Cf targets

As discussed in the preceding section 3.2.1, hot-fusion systems with uranium and transuranium targets may be good candidates for the synthesis of SHEs heavier than those reached with 208Pb targets. Like for 208Pb we study in this section a series of projectiles with increasing mass for 238U and 252Cf targets. 238U has been chosen as one of the most convenient targets with respect to radioactivity,
Fig. 3. SMPs for the collision systems $^{124}\text{Sn}+^{130}\text{Xe}$, $^{124}\text{Sn}+^{136}\text{Xe}$, $^{130}\text{Xe}+^{136}\text{Xe}$, $^{136}\text{Xe}+^{136}\text{Xe}$, $^{136}\text{Xe}+^{150}\text{Nd}$ and $^{144}\text{Ce}+^{150}\text{Nd}$ obtained with the SkP Skyrme force and the proximity potentials (1977), the Bass potentials (1974, 1980) and the KNS potential. The notations are the same as in Fig. 2.

availability and target properties. Moreover it is one of the largest-bound isotopes in this mass region, and hence is in this respect similar to $^{208}\text{Pb}$ yielding relatively small compound-nucleus excitation energies. $^{252}\text{Cf}$ has been depicted as the heaviest target available [61].

The entrance-channel potentials are shown in Figs. 6 and 7 for $^{238}\text{U}$ and $^{252}\text{Cf}$ targets, respectively. The dependence on the orientation of the deformed nuclei is similar to that of the deformed targets with $^{48}\text{Ca}$ and $^{40,42}\text{Ca}$ in Figs. 4 and 5. Again, for the side orientation ($\Theta = 90^{\circ}$) the barrier and pocket shapes are considerably more compact than for $^{208}\text{Pb}$ and by about 25 MeV higher in energy. The general decrease of the depth and width of the pockets with increasing charge of the projectile is similar to that for the $^{208}\text{Pb}$ target in Fig. 3.
Some interesting points of Fig. 6 for 238U are summarized as follows.

- The potential pocket vanishes for the 78Ge projectiles leading to the SHE 316124, whereas this situation happens already for 96Zr on 208Pb forming SHE 304122.
- The potential pocket for 58Fe+238U is also somewhat larger than the one for 86Kr+208Pb, both leading to SHE 118. However, higher excitation energies are encountered for bombarding energies at the barrier of 58Fe+238U.
- Comparing the entrance-channel potentials for different nickel isotopes on 238U, we notice no difference in the shape; but the height of the barrier, measured with respect to the compound-nucleus ground-state energy, reduces considerably with increasing neutron number. This effect has been recognized already by comparing 40,42Ca with 48Ca in Figs. 4 and 5. Therefore the use of more neutron-rich projectiles may be in general favorable for SHE formation, because the compound nucleus is formed with less excitation energy.

For the series of projectiles on 252Cf we recognize from Fig. 7 that the potential pockets are considerably deeper as compared to those with 238U for the same compound system, and even somewhat deeper for the same projectile. The pocket for the side orientation vanishes for 78Ge+252Cf = 330130 while this happens already for the compound system 316124 if 252Cf is replaced by 238U. In addition, also the compound-nucleus excitation energies are lower for the 252Cf target. From all these features we expect larger cross-sections for the synthesis of the SHEs with 252Cf than with 238U targets. If we consider the pocket of the side orientation to be decisive, the limit for fusion is reached around 70Zn+252Cf=322128.

3.3 Warm-fusion systems

Recently 233,234Cm and Fm have been synthesized in the collisions 40Ar, 50Ti+198Pt [12]. The cross-section for 50Ti+198Pt is comparable with the one for the cold-fusion reaction 40Ar+208Pb. 198Pt is oblate with deformation constant $\beta_2 = -0.10$ [13] and offers a qualitatively new entrance channel for the synthesis of SHEs.

In Fig. 8 we present the entrance-channel potentials for 40Ar till 100Mo + 198Pt. In contradistinction to the
Fig. 5. The SMPs for the collision $^{40,42}\text{Ca}$ on $^{238}\text{U},^{244}\text{Pu},^{248}\text{Cm}$ and $^{252}\text{Cf}$ evaluated with the SkM* Skyrme force. The SMPs are evaluated for different angular orientations of the deformed nuclei. The notations are the same as in Fig. 4.

prolate targets $^{238}\text{U},^{244}\text{Pu},^{248}\text{Cm}$ and $^{252}\text{Cf}$ the dependence of the barrier on the orientation is opposite with the lowest barrier for $\Theta = 90^\circ$ (side position) and the highest barrier for $\Theta = 0^\circ$ (tip position), the difference being about 10 MeV. The difference between the highest and lowest barriers for $^{198}\text{Pt}$ is smaller than the one for uranium and transuranium cases due to the smaller $|\beta_2|$. Furthermore, the pockets are considerably deeper for the tip position than for the side position.

The larger tip-position pockets for the oblate shape as compared to the spherical and prolate shapes is essentially due to the small curvature of its surface at the tip position. Indeed, the interaction potential in the proximity approach around touching is proportional to the reduced radii of the surfaces, i.e.

$$V_N \propto \frac{R_1R_2}{R_1 + R_2}$$

with $R_1, R_2$ the radii of the near touching surfaces. If nucleus 2 is deformed, the curvature $C_2 = 1/R_2$ is replaced by the mean curvature $\overline{C}_2 = 1/\overline{R}_2$. We consider a spheroid with half axes $a = b = R_2(1 - \delta)$ and $c = R_2(1 + 2\delta)$. For the same mass and absolute value of the deformation parameter $\delta$ we have

$$\overline{R}_2^{ps} = R_2(1 + 2\delta)$$

with $\delta > 0$ for the side position of the prolate nucleus and

$$\overline{R}_2^{ot} = R_2(1 + 4|\delta|)$$

with $\delta < 0$ for the tip position of the oblate nucleus. Therefore as compared to spherical nuclei, the pockets, which are caused by the nuclear attraction between the surfaces of the nuclei, are deeper for the side position of prolate nuclei and even more for the tip position of oblate nuclei. The Coulomb interaction favors the depths of the pockets for the oblate tip position as well as for the prolate side position.

In general for the same compound nucleus, the barrier energies measured with respect to the ground-state compound energy lie for $^{198}\text{Pt}$ in between the barriers for the cold-fusion target $^{208}\text{Pb}$ and the hot-fusion targets $^{238}\text{U},^{244}\text{Pu},^{248}\text{Cm}$ and $^{252}\text{Cf}$, and hence we refer to $^{198}\text{Pt}$ as the warm-fusion target. Here we imply that the SHE fusion window is again located about 5 to 10 MeV below the SMP barriers of the most compact pockets in the tip position.
Fig. 6. The SMPs for the systems $^{50}\text{Ti}$, $^{54}\text{Cr}$, $^{58}\text{Fe}$, $^{64}\text{Ni}$, $^{68}\text{Ni}$, $^{70}\text{Zn}$, $^{78}\text{Ge}$, $^{74}\text{Se}$, $^{82}\text{Se}$ + $^{238}\text{U}$ evaluated with the SkM* Skyrme force. The SMPs are evaluated for different angular orientations of $^{238}\text{U}$. The notations are the same as in Fig. 4.

4 Conclusion

As pointed out already in the introduction, the capture of the colliding nuclei within the pocket of the interaction potential is a decisive first step in the fusion process. For the heavy systems under consideration these pockets are rather shallow and vanish for too large systems. The population of long-living quasi-bound states in the pocket is expected to be limited in bombarding energy from below by the barrier penetrability (including transfer channels and polarization of the nuclei). For incident energies above the barrier, reflection of the system by the large repulsive core of the SMP potential prevents the capture in the pocket. The resulting window in bombarding energy for capture should increase with increasing depths of the potential pockets. These arguments are consistent
Fig. 7. The SMPs for the systems $^{50}$Ti, $^{54}$Cr, $^{58}$Fe, $^{58}$Ni, $^{64}$Ni, $^{64}$Ni, $^{70}$Zn, $^{78}$Ge, $^{74}$Se, $^{82}$Se + $^{252}$Cf evaluated with the SkM* Skyrme force. The SMPs are evaluated for different angular orientation of $^{252}$Cf. The notations are the same as in Fig. 4.

with the observed fusion window in the collision energy which is typically 5 MeV wide and lies about (5 – 10) MeV below our SMP barrier. Note, however, that the observed fusion window for SHE formation may differ from the capture window by the additional limitation towards higher energies due to $\Gamma_n/\Gamma_f \ll 1$.

There is a second limitation of the fusion window from below, because the capture has to be at high enough energies, such that the system can reach the compound-nucleus shape and fall into the potential minimum by emitting one neutron. The experimentally observed squeezing of the fusion window for $^{64}$Ni + $^{208}$Pb and $^{70}$Zn + $^{208}$Pb may be ascribed to this effect. From this point of view the synthesis of $^{86}$Kr + $^{208}$Pb should be suppressed, because the energies necessary for the 1n fusion is above the capture window (cf. Fig. 2).
Fig. 8. The SMPs for the systems $^{40}\text{Ar}$, $^{48}\text{Ca}$, $^{50}\text{Ti}$, $^{54}\text{Cr}$, $^{58}\text{Fe}$, $^{64}\text{Ni}$, $^{70}\text{Zn}$, $^{78}\text{Ge}$, $^{82}\text{Se}$, $^{86}\text{Kr}$, $^{96}\text{Zr}$, $^{100}\text{Mo}$ + $^{198}\text{Pt}$ evaluated with SkM* Skyrme force. The SMPs are evaluated for different angular orientation of $^{198}\text{Pt}$. The notations are the same as in
On the basis of these arguments we define the following rules for the determination of the best candidates for the synthesis of SHEs.

- The SMP barrier should lie about 10 to 15 MeV above the 1n fusion threshold, but not above the 2n fusion threshold to avoid the reduction of the fusion cross-section by an additional factor $I_n/I_f$. This condition yields an optimum fusion window (5 to 10 MeV below the barrier) for the formation of a compound nucleus with excitation energy 5 MeV above the 1n-emission threshold.
- The pocket depth should be as large as possible, because the deeper the pocket is, the larger the capture window becomes, and hence the better is the chance of fusion.
- For the subsequent formation of a compound nucleus it is best to have a most compact capture configuration.

We illustrate these rules for the synthesis of SHE 118 with hot-, cold- and warm-fusion systems studied in section 3.

- As mentioned already, the cold-fusion system $^{86}$Kr+$^{208}$Pb (cf. Fig. 2) has its capture window below the 1n-channel, and hence is not expected to be a good candidate.
- The symmetric system $^{144}$Ce+$^{150}$Nd (cf. Fig. 3) has no pocket, and hence no capture window at all.
- The hot-fusion system $^{48}$Ca+$^{252}$Cf (cf. Fig. 4) has nice capture properties, however needs to emit about 2 to 3 neutrons, which reduce the survival probability by several orders due to factors $I_n/I_f \ll 1$.
- The hot-fusion system $^{40}$Ca+$^{252}$Cf (cf. Fig. 5) has less attractive capture properties (as compared to the $^{48}$Ca case) and needs to emit even 5 to 6 neutrons.
- The system $^{58}$Fe+$^{238}$U (cf. Fig. 6) has only a tiny pocket and needs to emit about 3 to 4 neutrons.
- The warm-fusion system $^{96}$Zr+$^{198}$Pt has also a tiny tip-positioned pocket, but needs to emit only one neutron.

From this we conclude that among the studied systems the most attractive projectile-target combinations for the synthesis of SHE 118 are $^{48}$Ca+$^{252}$Cf at $E_{\text{coll}} \approx 206$ MeV (side collision) and $^{96}$Zr+$^{198}$Pt at $E_{\text{coll}} \approx 330$ MeV (tip collision). While $^{48}$Ca+$^{252}$Cf is more compact, $^{96}$Zr+$^{198}$Pt needs to emit only 1 neutron instead of 2 to 3. It is hard to judge, which of these features are more important for the synthesis of SHE 118. Generally for deformed target nuclei, the experiments on the synthesis of SHEs may profit from using polarized targets by aligning prolate targets for side collisions and oblate targets for tip collisions.

One should be aware, that other effects not due to capture, compactness and neutron emission may influence the final formation of a SHE. In particular special features of the potential landscape, e.g. with respect to decay by quasi-fission, are of importance. We hope to obtain an improved guidance for experiments from a scattering model which treats capture, collective evolution and neutron decay on the same level and fully quantum-mechanically.

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