Independent particle model for nucleon transfer in heavy ion collisions

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We formulate a three-dimensional quantum-mechanical model for the study of nucleon transfer in heavy ion collisions. This model describes the independent particle motion of nucleons initially separated in two adjacent cubic boxes and brought into contact through a window created in the separating wall. We calculate the one-sided flux and the spread of the mass distribution and compare them with results derived from other models.

I. INTRODUCTION

The particle exchange mechanism plays an important role in explaining the dissipative phenomena observed in heavy ion collisions.¹ Within the one-body dissipation mechanism² the relevant quantity for the dissipation rate is the one-sided current of nucleons which are exchanged via a window³ opened between the interacting nuclei.

Recently attempts have been made to derive the onesided current on a microscopic basis, starting either from time-dependent⁴ or adiabatic time-dependent⁵ Hartree-Fock wave functions. In such calculations some care must be taken because the quantum-mechanical definition^{6,7} of the one-sided current does not always have a meaning.⁸ This is due to the fact that the Wigner transform of the one-body density may assume regions of negative values which prevent the extension of the classical definition of the one-sided current^{2,3} to a quantum system.

In the present work we formulate a simple threedimensional quantum-mechanical model in order to obtain more insight into the dynamics of particle exchange. This model describes the independent particle motion of nucleons initially separated in two adjacent cubic boxes and brought into contact through a window created in the separating wall. We calculate a time-dependent one-sided current in two different ways. One is according to the currently used definition⁴⁻⁸ where the flux is evaluated as in Eq. (3.1) from the Wigner distribution function $f(\vec{r}, \vec{k}, t)$ of the one-body density. The other is given by the initial conditions of our model and better visualizes the physics contained in it. We start at t = 0 with one box empty and the other full. At t > 0 the window (suddenly) opens, the wave functions spread out of the filled box, and the matter flows into the empty box. As the box has a finite size, the waves travel up to the opposite wall where they are reflected. We study the correspondence between this matter flow and the result obtained from the currently used definition of the one-sided current. We also make a comparison between the presently obtained time averaged and the static one-sided current calculated in Ref. 9, at the same time indicating the value of the purely geometrical flux.

As we deal with an entirely quantum-mechanical system which simulates the mechanism of particle exchange we can also calculate the fluctuations σ^2 in the number of particles. The result is compared to that obtained by Buck and Feldmeier⁷ in a one-dimensional model without two-body collisions. The above authors indicated that such a term has a negligible effect on the order of magnitude of σ^2 .

In Sec. II we give a brief account of the static model and introduce the time-dependent wave functions. In Sec. III we present results for the matter flow calculated according to various definitions. Section IV deals with the calculation of the width of the mass distribution, and Sec. V is devoted to the conclusion.

II. THE MODEL

Our model consists of two adjacent hard-walled cubes having a length L on each edge and communicating through a window initially closed. Up to time t=0 the boxes are equally filled with A/2 particles each. At t=0the window opens and the particles move freely from one box to the other through the window. The geometry is indicated in Fig. 1. The window is situated in the plane z=0 and extends from x=-w to x=w and from y=-L/2 to y=L/2. The Hamiltonian is symmetric under reflections through each coordinate. We denote by π_i the parity for i=x, y, or z and choose the basis functions

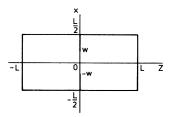


FIG. 1. Adjacent cubic boxes communicating through a rectangular window of size x = 2w, y = L in the z = 0 plane.

$$u_{n_{i}}^{\pi^{i}}(i) = \left[\frac{2}{L_{i}}\right]^{1/2} \begin{cases} \cos\frac{n_{i}\pi}{L_{i}}i & \pi_{i} = +1, n_{i} \text{ odd} \\ \sin\frac{n_{i}\pi}{L_{i}}i & \pi_{i} = -1, n_{i} \text{ even }, \end{cases}$$
(2.1)

which vanish on the planes $L_i/2$, where $L_x = L_y = L$ and $L_z = 2L$.

The system is described by functions defined inside the volume occupied by the two cubes. They have total parity $\pi = \pi_x \pi_y \pi_z$ and due to the geometry of the window they can be factorized as

$$\phi_{\alpha,n_y}^{\pi}(\vec{\mathbf{r}}) = u_{n_y}^{\pi_y}(y)\phi_{\alpha}^p(x,z) , \qquad (2.2)$$

with $p = \pi_x \pi_z$ and ϕ^p_α given by the expansion

$$\phi_{\alpha}^{p}(x,z) = \sum_{n} c_{\alpha n}^{p} u_{n_{x}}^{\pi_{x}}(x) u_{n_{z}}^{\pi_{z}}(z) , \qquad (2.3)$$

$$V_{nn'}^{p} = \left[\frac{2}{L}\right]^{2} \begin{cases} -\frac{L}{\pi} \left[\frac{\sin\frac{\pi}{L}(n_{x} + n'_{x})w}{n_{x} + n'_{x}} + p\frac{\sin\frac{\pi}{L}(n_{x} - n'_{x})w}{n_{x} - n'_{x}}\right] \\ \frac{L}{2} - w - p\frac{L}{2\pi n_{x}} \sin\frac{2\pi n_{x}w}{L} & \text{for } n_{x} = n'_{x} \end{cases}$$

for any n_z and n_z' odd, or equivalently $\pi_z = +1$, or else we have

$$V_{\rm nn'}^p = 0 \text{ for } \pi_z = -1$$
 (2.8)

as automatically implied by (2.1) and (2.4). Now let us call N the maximum value taken by n_z in the truncated space where the matrix equation (2.5) is solved. The case N+1 even reduces to N odd due to (2.8). At fixed N odd and for $\pi_z = 1$, n_x runs up to N for p = +1 and up to N-1 for p = -1. Hence the model has two parameters, N and λ . Adequate values of these parameters have been found in Ref. 9. Accordingly, in the present calculations we choose $\lambda = 10^4$ MeV fm and N = 19 whenever w < L/2.

Now, when the procedure of obtaining the eigenstates (2.2) of the communicating boxes has been settled, we can introduce the time-dependent wave functions. At t = 0 we assume that the window is closed and the particles fill one of the boxes, let us say for example the left box. For states initially localized in the left box (z < 0) we shall use the label L. Then at t > 0 the associated single particle wave function ψ_n^L will evolve in time according to

$$\psi_n^L(x,z,t) = \sum_{\alpha} d_{n\alpha}^{\pi_x p}(L) e^{-(i/\tilde{n})E_{\alpha}^p t} \phi_{\alpha}^p(x,z) .$$
(2.9)

The variable y is omitted for simplicity but taken into account in filling up the levels (see Table I). The states ψ_n^R initially localized in the right box and orthogonal to ψ_n^L are defined in a similar way with the associated expansion coefficients $d_{n\alpha}^{\pi_x p}(R)$. In both cases the coefficients are en-

where the summation runs through $n = (n_x, n_z)$ and the coefficients $c_{\alpha n}^p$ are determined from the boundary conditions imposed on the separating wall which requires the vanishing of the wave function at z = 0, $w \le |x| \le L/2$. Such a condition is automatically satisfied for $\pi_z = -1$ and achieved only approximately for $\pi_z = +1$ states. The procedure has been described in detail in Ref. 9 and it is equivalent with solving the Schrödinger equation for the following potential:

$$V = \lambda \delta(z) \theta(|x| - w); \quad \lambda \to \infty , \qquad (2.4)$$

in a truncated space, i.e., solving the matrix equation

$$\sum_{n'} \left[(\epsilon_n^p - E_{\alpha}^p) \delta_{nn'} + V_{nn'}^p \right] c_{\alpha n'}^p = 0 , \qquad (2.5)$$

where

$$\epsilon_n^p = \frac{\hbar^2}{2m} \frac{\pi^2}{L^2} (n_x^2 + \frac{1}{4} n_z^2)$$
(2.6)

and

$$\frac{\mathbf{n}_{\mathbf{x}}' \mathbf{n}_{\mathbf{x}}}{\mathbf{n}_{\mathbf{x}}'} \quad \text{for } \mathbf{n}_{\mathbf{x}} \neq \mathbf{n}_{\mathbf{x}}'$$

$$(2.7)$$

tirely determined by the initial conditions as the overlap integrals

$$d_{n\alpha}^{\pi_x p}(L) = \int u_n^L \phi_\alpha^p dv \text{ and } d_{n\alpha}^{\pi_x p}(R) = \int u_n^R \phi_\alpha^p dv , \quad (2.10)$$

TABLE I. The occupied states in a cubic box at t = 0 (closed window) for A = 32 and 80 particle systems. The energy eigenvalue is

$$\epsilon_n^{\pi} = \hbar^2/2m \left(\frac{\pi}{L}\right)^2 (n_x^2 + n_y^2 + \frac{1}{4}n_z^2) .$$

System	n _x	n _y	n _z	ϵ_n^{π} (MeV)	Occupation number Eq. (2.15)
A = 32	1	1	2	29.69	0.9998
	1	2	2	59.40	0.9998
	2	1	2	59.40	0.9998
	1	1	4	59.40	0.9994
A = 80	1	1	2	16.11	0.9998
	1	2	2	32.23	0.9998
	2	1	2	32.23	0.9998
	1	1	4	32.23	0.9994
	2	2	2	48.35	0.9998
	2	1	4	48.35	0.9994
	1	2	4	48.35	0.9994
	1	1	6	59.10	0.9985
	3	1	2	59.10	0.9998
	1	3	2	59.10	0.9998

(2.11)

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where

$$u_n^L = \frac{2}{L}\theta(-z)\sin\frac{\pi n_z}{2L}z \begin{cases} \sin\frac{\pi n_x}{L}x; & n_x \text{ even, } n_z \text{ even} \\ \cos\frac{\pi n_x}{L}x; & n_x \text{ odd, } n_z \text{ even} \end{cases}$$

and

$$u_n^R(z) = u_n^L(-z); \quad z > 0 \ . \tag{2.12}$$

We are reminded that p is the parity associated with the eigenstate ϕ_{α}^{p} under reflections in the x,z plane and π_{x} is the parity of the state (2.11) at reflection through the x axis. The product $\pi_{x}p = \pm 1$ serves as the upper index for the coefficients (2.10) and indicates that the function (2.9) has a given parity at reflections through the x axis but mixes symmetric and antisymmetric states ϕ_{α}^{p} under reflections through the z axis.

The integrals (2.10) give

$$d_{n\alpha}^{-1}(L) = d_{n\alpha}^{-1}(R) = \frac{\sqrt{2}}{2} \delta_{n\alpha} ,$$

$$d_{n\alpha}^{+1}(L) = -d_{n\alpha}^{+1}(R) = \frac{\sqrt{2}}{L} \sum_{n'} c_{\alpha n'}^{p} N_{n_{z}n'_{z}} \delta_{n_{x}n'_{x}} , \quad (2.13)$$

with

$$N_{n_z n'_z} = -\frac{L}{\pi} \left[\frac{1}{n_z + n'_z} + \frac{1}{n_z - n'_z} \right]; \quad n_z \text{ even, } n'_z \text{ odd }.$$
(2.14)

In the present work we consider two systems, A = 32and 80 particles, respectively. The occupied states at t = 0when the window is closed and each box contains A/2particles are indicated in Table I for both systems, each state being four times degenerate. They are given in increasing order of the energy

$$\frac{\hbar^2}{2m} \left[\frac{\pi}{L}\right]^2 (n_x^2 + n_y^2 + \frac{1}{4}n_z^2) .$$

We note that n_z is always even, indicating the cancellation of the wave function at z = 0 (closed window). The size L is chosen such as to have an average density $\rho = 0.17 \text{ fm}^{-3}$ in each box. This gives L=4.55 and 6.17 fm, for A=32and 80 particles, respectively. Now we want to approximate the eigenstates of the left (or right) box by the truncated expansion (2.9). As an example in Fig. 2 we plot ψ_n^L at t=0, x=y=0 as a function of z; n_x can take any odd value and the four curves correspond to $n_z = 2, 4, 6, \text{ and } 8,$ respectively. The truncation has been made at N = 15, i.e., the summation in Eq. (2.9) contains one state with p = -1 $(\alpha_z \text{ even})$ and eight states with p = +1 $[\alpha_z \pmod{\leq N}]$. We notice that the wave function ψ_n^L is very well localized. We found that a larger space, having N > 15, brings insignificant changes to the wave functions with $n_z \leq 8$ when w = L/2 and therefore limits the expansion at N = 15 for a fully open window. A quantitative measure of the localization is given by the probability of finding the particle

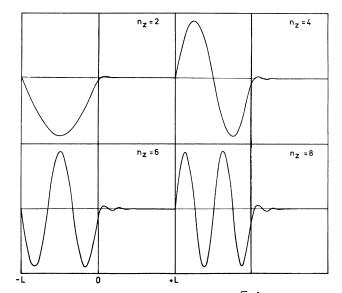


FIG. 2. Eigenstates of the left box $L/\sqrt{2}\psi_n^L$ at t=0, x=0. They are given by the truncated expansion (2.9) containing nine eigenstates of the big box (fully open window), i.e., N=15. The value of n_z is specified for each case; n_x can take any odd value.

within the volume of the left box

$$\int_{L} dv |\psi_{n}^{L}(t=0)|^{2} = \sum_{\alpha} \left[d_{n\alpha}^{\pi_{x}p}(L) \right]^{2}.$$
(2.15)

This is indicated in the last column of Table I for all initially occupied states. For $n_z = 8$ states, unoccupied at t = 0, it is about 0.997, i.e., close to 1 in all cases of interest. The latter will be used in Sec. IV for calculating the width of the mass distribution.

A comment is necessary about the truncation for the case w < L/2. A first truncation is made in search for the state ϕ_{α}^{p} . Not all these states are physically acceptable. The vanishing of the wave function on the intermediate wall is achieved to a very good approximation for the lowest states. The associated eigenvalue E^p_a becomes practically independent of λ . The highest excited states do not have such properties. But, in order to localize the wave function in the left or right box, it turned out that two or three of the lowest states carry most of the weight for all occupied and few unoccupied states close to the Fermi level. Hence the truncated space used in the preceding section to define ϕ_{α}^{p} can be even more restricted in (2.9). A practical alternative is to keep the same dimension, but instead of E^{p}_{α} in (2.9), we use the expectation values $\langle \phi_{\alpha}^{p} | T | \phi_{\alpha}^{p} \rangle$ of the kinetic energy operator which permits us to get rid of the region where E^p_{α} depends strongly on λ for the highest excited states. For all the states where the vanishing of the wave function on the intermediate wall is well achieved E^{p}_{α} and the expectation value of the kinetic energy are practically the same. By using $\langle \phi^p_{\alpha} \mid T \mid \phi^p_{\alpha} \rangle$ instead of E^{p}_{α} one can also ensure the satisfaction of the continuity equation. Details are discussed in Ref. 10.

III. MASS EXCHANGE

A measure of the number of nucleons exchanged through the window is the one-sided flux.^{2,3} In this sec-

tion we calculate the one-sided flux by using two different definitions and discuss the relation between the two calculations. We also compare the present results with the static one-sided flux.⁹

A. Definition I

This is an extension of the definition used in classical mechanics and it makes use of the Wigner distribution function $f(\vec{r}, \vec{k}, t)$ associated with the one-body density of the quantum system. According to this definition the one-sided flux in the z direction is given by^{4,5,8,9}

$$j_{+}(\vec{\mathbf{r}},t) = \frac{\hbar}{m} \int_{k_{z} \ge 0} d^{3}k \, k_{z} f(\vec{\mathbf{r}},\vec{\mathbf{k}},t) \,. \tag{3.1}$$

This also represents the expectation value of the quantum-mechanical operator introduced by Feldmeier.^{6,7} Classically the distribution function is positive, hence j_+ is also positive, but for quantum systems problems might arise because the Wigner function can have negative values. An example has been extensively discussed in Ref. 8.

Using definition (3.1), we calculate the one-sided current through the window, i.e., at z = 0,

$$J_{+}(t) = \int_{-w}^{w} dx \, \int_{-L/2}^{L/2} dy \, j_{+}(x,y,0,t) \tag{3.2}$$

in the following way. At t=0 the window is closed and each box contains A/2 particles. The occupied states ψ_n^L (t=0) of the left box are degenerate with ψ_n^R (t=0) of the right box. Adding together the contribution of each two initially degenerate states we define the one-sided flux associated with the state n,

$$j_{+,n}(\vec{\mathbf{r}},t) = j_{+,n}^{L}(\vec{\mathbf{r}},t) + j_{+,n}^{R}(\vec{\mathbf{r}},t) .$$
(3.3)

Then, the total flux reads

$$j_{+}(\vec{\mathbf{r}},t) = 4 \sum_{n \text{ (occ)}} j_{+,n}(\vec{\mathbf{r}},t) ,$$
 (3.4)

when the sum runs over all occupied states, listed in Table I and the factor 4 takes into account the spin-isospin degeneracy. Here by *n* we mean the set n_x, n_y, n_z . In filling up the levels in increasing order it is necessary for each state to add the contribution $(\hbar^2/2m)(\pi/L)^2 n_y^2$ to E_{α}^p . We are reminded that the y-dependent part of the wave function appearing as the factor $u_{n_y}^{\pi_y}(y)$ remains invariant in the process of building up the wave function ψ_n^L or ψ_n^R and was therefore omitted for simplicity.

From the Wigner function,

$$f_n(\vec{\mathbf{r}},\vec{\mathbf{k}},t) = \frac{1}{(2\pi)^3} \int d^3s \, e^{-i\,\vec{\mathbf{k}}\cdot\vec{\mathbf{s}}} \psi_n^* \left[\vec{\mathbf{r}}+\frac{\vec{\mathbf{s}}}{2}\right] \psi_n \left[\vec{\mathbf{r}}-\frac{\vec{\mathbf{s}}}{2}\right], \qquad (3.5)$$

of each occupied state we calculate $j_{+,n}^L$ and $j_{+,n}^R$ according to definition (3.1), and making the operations (3.2)–(3.4), we obtain

$$J_{+}(t) = \frac{\hbar}{m} \frac{1}{L^{2}} \sum_{n \text{ (occ)}} \sum_{\alpha,\alpha'} d_{n\alpha}^{\pi_{x}p}(L) d_{n\alpha'}^{\pi_{x}p}(L) F_{\alpha\alpha'}^{p} e^{(i/\hbar)(E_{\alpha}^{p} - E_{\alpha'}^{p})t}, \qquad (3.6)$$

with

$$F_{\alpha'\alpha''}^{p} = \sum_{n',n''} c_{\alpha'n'}^{p} c_{\alpha''n''}^{p} B_{n'_{x}n''_{x}}^{\pi} \left\{ (n'_{z} - n''_{z}) \operatorname{Si}\left[(n'_{z} - n''_{z}) \frac{\pi}{2} \right] + (n'_{z} + n''_{z}) \operatorname{Si}\left[(n'_{z} + n''_{z}) \frac{\pi}{2} \right] \right\},$$
(3.7)

where Si(x) is the sine integral¹¹ and

$$B_{n'_{x}n''_{x}}^{\pi_{x}} = \delta_{n'_{x}n''_{x}} - \frac{L}{2\lambda} V_{n'_{x}n''_{x}}^{\pi_{x}}, \qquad (3.8)$$

with $V_{n_x n'_x}^{\pi_x}$ given by an expression like (2.7), where *p* is replaced by π_x and *n* by n_x . We have $F_{\alpha\alpha'}^p = F_{\alpha'\alpha}^p$ which implies that $J_+(t)$ of Eq. (3.6) is in fact real. This quantity has been expressed in terms of $d_{n\alpha}^{\pi_x p}(L)$ only due to the relations (2.13).

In Figs. 3(a) and (b) the dotted line shows $J_{+}(t)$ obtained at w = L/2 for A = 32 and 80 particle systems, respectively. The result has been given over a full period which can be easily calculated. The one-sided current $J_{+}(t)$ returns to its initial zero value whenever

$$e^{(i/\hbar)(E^p_\alpha - E^p_{\alpha'})t} = 1 \text{ for all } \alpha, \alpha' .$$
(3.9)

For the first two modes E_1^+ $(n_x = n_y = n_z = 1)$ and E_3^+ $(n_x = 1, n_y = 1, n_z = 3)$ of the same parity the Eq. (3.9) is satisfied if

$$T = \frac{2L^2 m}{\pi \hbar} . \tag{3.10}$$

Relation (3.10) gives a period of $J_+(t)$ because any other difference $E^p_{\alpha} - E^p_{\alpha'}$ is a multiple of $E^+_1 - E^+_3$ when w = L/2, and this makes condition (3.9) fulfilled for any α whenever t = T. Relation (3.10) gives $T = 2.09 \times 10^{-22}$ s for A = 32 and $T = 3.85 \times 10^{-22}$ s for A = 80.

Contrary to the periodical behavior at the fully open window Eqs. (3.9) have no unique solution at $w \neq L/2$ because E^p_{α} are no more simply related and the current becomes aperiodical. This is illustrated in Figs. 4 and 5 for the A = 32 system at w = 1.6 and 1 fm. One can see that at w = 1.6 fm the deviation from periodicity is still small, but at w = 1 fm the periodicity has been entirely lost.

An interesting quantity is the time average \overline{J}_+ of the current. Its analytic expression is

$$\bar{J}_{+} = \frac{\hbar}{m} \frac{1}{L^2} \sum_{n \text{ (occ)}} \sum_{\alpha} \left[d_{n\alpha}^{\pi_{\chi} p}(L) \right]^2 F_{\alpha \alpha}^p .$$
(3.11)

(a)

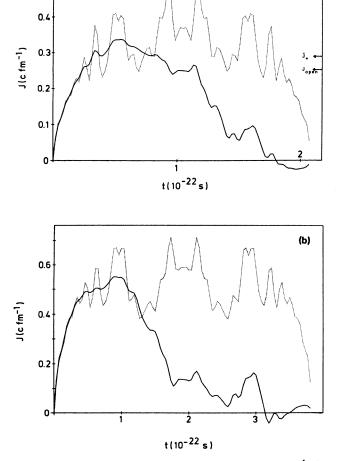


FIG. 3. The one-sided current J_+ (dotted line) and J^L (full line) as a function of time at w = L/2. The parameters of the model are $\lambda = 0, N = 15$. (a) A = 32 particles, (b) A = 80 particles. The arrows indicate \overline{J}_+ of Eq. (3.11) and $J_{\text{open}} = j_{\text{open}} \times 2wL$ with j_{open} from (3.12).

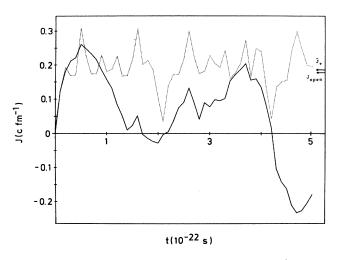


FIG. 4. Same as Fig. 3(a) but for w = 1.6 fm, $\lambda = 10^4$ MeV fm, N = 19.

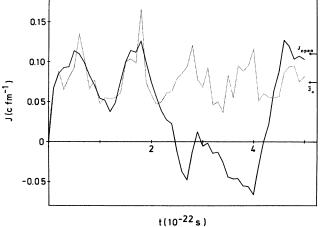


FIG. 5. Same as Fig. 3(a) but for w = 1 fm, $\lambda = 10^4$ MeV fm, N = 19.

As an example, values of \overline{J}_+ are given in Table II for w = 1 and 1.6 fm and L/2 for the A = 32 particle system. In Table II we also reproduce the values of the static onesided current obtained in Ref. 9 for the same values of w, λ , and N. It seems that there is a good agreement between the two results, the static value being somewhat smaller in all cases. The difference between the two results can be explained by the rearrangement in the eigenvector space of dimension N through definition (2.9) when each component ϕ_{α}^{p} has a weight $|d_{n\alpha}^{\pi_{x}p}|^{2} < 1$, while in the static calculations the weight is either 1 or 0.

For completeness we also reproduce the purely geometrical current obtained by multiplying the window area with the static one-sided flux j_{open} obtained at maximum opening

$$j_{\rm open} = \frac{\hbar}{m} \frac{1}{L^4} \sum_{n_z \, (\rm occ)} n_z {\rm Si}(\pi n_z) \,. \tag{3.12}$$

We note that a given n_z can correspond to several sets n_x, n_y and its multiplicity must be taken into account in (3.12).

B. Definition II

Now we consider one box full with A/2 particles and the other empty at t = 0. Let us choose the full left box. The window opens at t > 0 and the particles flow into the other box. The current $J^{L}(t)$ is obtained as in Eq. (3.2), but from the flux $j^{L}(\vec{r},t)|_{z=0}$ given by the contribution of

TABLE II. The one-sided current for the A = 32 particle system at three different window sizes w; \overline{J}_+ [Eq. (3.11)], J_{static} (Ref. 9), $J_{\text{open}} = j_{\text{open}} \times 2wL$ with j_{open} from Eq. (3.12).

w (fm)	\overline{J}_+ (c fm ⁻¹)	J_{static} (c fm ⁻¹)	J_{open} (c fm ⁻¹)
1	0.075	0.063	0.112
1.6	0.187	0.162	0.179
L/2	0.288	0.254	0.254

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all initially occupied states of the left box only

$$j^{L}(\vec{\mathbf{r}},t) \mid_{z=0} = 4 \sum_{n \text{ (occ)}} j^{L}_{n}(\vec{\mathbf{r}},t) \mid_{z=0} , \qquad (3.13)$$

where for j_n^L we use the quantum-mechanical definition

$$J^{L}(t) = \frac{\hbar}{m} \frac{\pi\sqrt{2}}{L^{2}} \sum_{n \text{ (occ)}} \sum_{\alpha \neq n} n_{z} d_{n\alpha}^{+1}(L) \sin \frac{E_{\alpha}^{p} - E_{n}^{-p}}{\hbar} t \sum_{n''} c_{\alpha n''}^{p} B_{n_{x}n_{x}''}^{p} , \qquad (3.15)$$

where n_z is even and n_z'' odd. Although by definition $d_{n\alpha}^{+1}$ implies $\alpha \neq n$ we specify it in the summation in order to avoid any confusion here or below.

In Figs. 3–5 the result for $J^{L}(t)$ is represented by a full line. In all cases one can see that there is an interval starting at t = 0 where J_+ and J^L are very similar. This interval increases as w decreases. Its length depends on the moment when the reflection on the opposite wall starts to dominate. In a large time interval $\tilde{J}^{\hat{L}}$ oscillates between positive and negative values showing the emptying and filling up of the left box, while J_+ remains always positive. The region where J_+ and J^L are close to each other plays an important role. By construction J^L has a one-sided nature before the reflection sets in, i.e., the particles leave the left box with $k_z > 0$ and travel into the right box. The practical coincidence between J_+ and J^L shows that J^R_+ , i.e., the contribution to J_+ coming from the particles initially present in the right box is negligible or vice versa whenever such a contribution is negligible, the exclusion of the interval $\int_{-\infty}^{0} dk_z$ in the calculation of the flux, as it is done in Eq. (3.1), is entirely justified. These considerations are illustrated in Figs. 6 and 7 for the system with A = 32 particles with a window of size w = 1 fm. In Fig. A = 32 particles with a window of size w = 1 fm. In Fig. 6 we compare J^L with J^L_+ , i.e., the one-sided flux (3.2) ob-tained by filling the left box only with A/2 particles at t = 0. Comparing $J_+ = J^L_+ + J^R_+$ from Fig. 5 with J^L_+ from Fig. 6, we see that J^R_+ becomes appreciable only beyond $t > 2 \times 10^{-22}$ s, while J^L_+ remains dominant from t = 0 up to $t \sim 2 \times 10^{-22}$ s. In the Appendix we show that

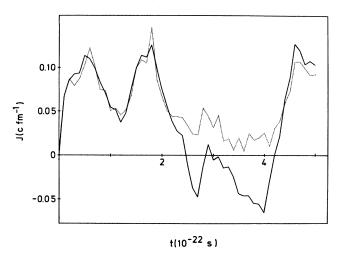


FIG. 6. The one-sided current for the full left box with A/2particles, empty right box: J_{+}^{L} (dotted line) and J^{L} (full line) as a function of time at w = 1 fm and A = 32 particles. The parameters are $\lambda = 10^4$ MeV fm, N = 19.

$$j_{n}^{L}(\vec{r},t) |_{z=0} = \frac{n}{m} \operatorname{Im}[\psi_{n}^{L}(\vec{r},t)^{*} \frac{d}{dz} \psi_{n}^{L}(\vec{r},t)]_{z=0}.$$
 (3.14)
ing expansion (2.9) and definitions (2.3) and (2.13), we

Usi obtain

 $i^{L}(\vec{r},t) = -\frac{\hbar}{2} Im [\eta^{L}(\vec{r},t)^{*} \frac{d}{d} \eta^{L}(\vec{r},t)]$

$$\frac{2}{n}\sum_{n \text{ (occ)}}\sum_{\alpha \neq n} n_z d_{n\alpha}^{+1}(L) \sin \frac{E_{\alpha} - E_n^{-1}}{\hbar} t \sum_{n''} c_{\alpha n''}^p B_{n_x n_x''}^p, \qquad (3.1)$$

$$J_{+}^{L} = \frac{1}{2}(J_{+} + J^{L}) ,$$

$$J_{+}^{R} = \frac{1}{2}(J_{+} - J^{L}) ,$$
(3.16)

where J_+ and J^L are given by Eqs. (3.6) and (3.15), respectively. When the two boxes are equally filled with A/4 particles J^L becomes zero and only at this moment J_{+}^{R} and J_{+}^{L} become equal. In Fig. 7 we compare J_{+}^{L} with a quantity called J_{-}^{L} defined as the current (3.2), but obtained from

$$j_{-}^{L} = \frac{\hbar}{m} \int_{k_{z} < 0} d^{3}\vec{k} \, k_{z} f^{L}(\vec{r}, \vec{k}, t) , \qquad (3.17)$$

where the Wigner function f^L is associated with the full left box and the empty right box. One can see that over the interval where J^R_+ is negligible, J^L_- is also negligible as we have previously mentioned. In fact, through the use of some algebra like in the Appendix and the fact that the Wigner function is symmetric under the reflections $\vec{k} \rightarrow -\vec{k}$, one can show that J_{+}^{R} and J_{-}^{L} are equal but of opposite sign.

IV. THE WIDTH OF THE MASS DISTRIBUTION

In the preceding section we have shown results for the current which gives a measure of the nucleon exchange between nuclei. In heavy ions another quantity of interest is the width of the mass or charge distribution. While the current is related to the one-body density via the continuity equation, the width of the mass distribution is expressed in terms of the two-body density matrix. Within the present model the variance σ^2 of the particle number

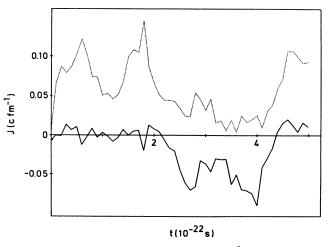


FIG. 7. Same as Fig. 6 but full line gives J_{-}^{L} defined according to (3.2) and (3.17).

(4.4)

in one box can be easily calculated.

We define the particle number operator in one box like in Refs. 7 and 12. If we arbitrarily choose the left box this operator is

$$N_L = \sum_{i=1}^{A} n_L(i) = \sum_{i=1}^{A} [1 - \theta(z_i)], \qquad (4.1)$$

with A the total number of particles and $\theta(z_i)$ the step function

$$\theta(z_i) = \begin{cases} 1 \text{ for } z_i > 0 \\ 0 \text{ for } z_i < 0 \end{cases}.$$
 (4.2)

The variance is

$$\sigma_L^2 = \langle N_L^2 \rangle - \langle N_L \rangle^2 , \qquad (4.3) \qquad \text{Sec. II. Then the matrix element } \langle o \mid n_L \mid u \rangle \text{ is given by}$$

$$\langle o \mid n_L \mid u \rangle = \frac{1}{2} \sum_{\alpha \neq o, u} d_{o\alpha}^{+1}(L) d_{u\alpha}^{+1}(L) [1 \pm (e^{(i/\hbar)(E_{\alpha}^p - E_u^{-P})t} + e^{-(i/\hbar)(E_{\alpha}^p - E_o^{-P})t})] \text{ for } o \neq u , \qquad (4.5)$$

where the plus (minus) sign is related to the case where both states $|o\rangle$ and $|u\rangle$ are initially localized in the left (right) box. If $|o\rangle$ and $|u\rangle$ are initially in different boxes the expression is

$$\langle o \mid n_L \mid u \rangle = \frac{1}{2} \sum_{\alpha \neq o, u} d_{o\alpha}^{+1}(L) d_{u\alpha}^{+1}(L) [-1 \pm (e^{(i/\hbar)(E_{\alpha}^p - E_u^{-p})t} - e^{-(i/\hbar)(E_{\alpha}^p - E_o^{-p})t})] \text{ for } o \neq u , \qquad (4.6)$$

with the + (-) sign for $|o\rangle$ in the left (right) box, respectively. If the number of states α is large enough the orthogonality between $|u\rangle$ and $|o\rangle$ is well satisfied, which implies

$$\sum_{\alpha \neq o, u} d_{o\alpha}^{+1}(L) d_{u\alpha}^{+1}(L) \simeq 0 \text{ for } o \neq u .$$
(4.7)

Using (4.7) and adding up the four contributions given by (4.5) and (4.6), we obtain

$$\sigma_L^2 = 2 \sum_{o \neq u} \left| \sum_{\alpha \neq o, u} d_{o\alpha}^{+1}(L) d_{u\alpha}^{+1}(L) e^{i/\hbar E_{\alpha}^{pt}} \right|^2.$$
(4.8)

The calculated value of σ_L^2 is shown in Figs. 8(a) and (b) for the systems A = 32 and 80, respectively, at w = L/2. One can notice a periodical structure again, like for the current J_+ . Each time the condition

$$e^{(i/\hbar)E^{p}_{\alpha}t} = 1$$
 for all α (4.9)

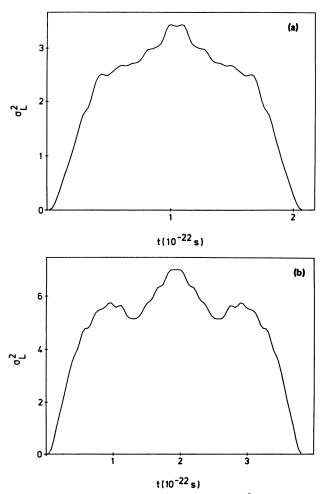
is fulfilled, σ_L^2 goes to zero. This takes place at the same period T given by (3.10), which can be easily seen if we use instead of (4.9) the equation

$$e^{(i/\hbar)(E_{\alpha}^{p}-E_{1}^{p})t}=1$$

also consistent with (4.8). As already mentioned the period T for A = 80 is about twice as large than that for A = 32. If we average over a period, the formula (4.8) gives

$$\overline{\sigma_L^2} = 2 \sum_{o \neq u} \sum_{\alpha \neq o, u} \left[d_{o\alpha}^{+1}(L) d_{u\alpha}^{+1}(L) \right]^2$$
(4.10)

because the coefficients $d_{n\alpha}^{\pi_x p}$ are real. This quantity should be compared to the classical value^{7,13} of σ_L^2 which at large *t* tends towards A/4, where A is the total number of particles. The ratio between σ_L^2 and the asymptotic classical value is 2/8 for A = 32 and 4.88/20 for A = 80,



where the expectation value $\langle \rangle$ is taken with respect to

the ground state. In the present model this state is a

where o and u represent occupied and unoccupied states, respectively. We discuss the case of two equally filled

boxes. Then the states o, u can either be chosen to have a

definite parity or be represented by wave functions initially localized in the left or right box as given by Eq. (2.9). In the first version the procedure would be close to a twocenter shell model description.¹² In the second version the parities are mixed in the wave functions, but for continui-

ty we maintain this basis such as it has been defined in

Slater determinant. Hence,

 $\sigma_L^2 = \sum_{o,u} |\langle o | n_L | u \rangle|^2,$

FIG. 8. The width of the mass distribution σ_L^2 as a function of time at w = L/2 and N = 15. (a) A = 32 particles; (b) A = 80 particles.

i.e., it has the same order of magnitude as that found in Ref. 7 from a one-dimensional model. This confirms the conclusion given in Ref. 7 that when the particles both at the left and right are forced to move in the same potential the mass dispersion is very small, as in time-dependent Hartree-Fock calculations. It would be interesting to extend the present study to the case where the two boxes have different sizes in order to see whether σ^2 increases in such a case.

V. CONCLUSIONS

We have formulated a simple three-dimensional quantum-mechanical independent particle model in order to get more insight into the quantities describing the exchange of nucleons between heavy ions. In this model the colliding nuclei are represented by two adjacent cubic boxes communicating through a rectangular window. The window suddenly opens with a given size and we calculate the time evolution of the one-sided current and of the width of the mass distribution at constant opening.

We have used two different definitions to calculate the one-sided flux of the window formula. One is the currently used definition introduced as an extension of the classical definition and based on the Wigner distribution function obtained from the single particle wave functions of the communicating system. The other is the quantummechanical definition applied in the case where one box is filled and the other box is empty. The calculations show that both definitions give similar results in the time interval before the reflection on the opposite wall sets in. We consider that such similarities bring support in using the extended classical definition where the integration over positive momenta only [see Eq. (3.1)] is not *a priori* justified in quantum mechanics.

The calculated width of the mass distribution σ_L^2 has a periodical structure, like J_+ , due to the finite size of the system. The ratio between the time average of σ_L^2 and its asymptotic classical value has the same order of magnitude as that obtained by Buck and Feldmeier⁷ in a onedimensional system. Our calculations therefore support the conclusion of Ref. 7 according to which the smallness of σ_L^2 is the combined effect of the Pauli principle and the use of an identical mean potential for both nuclei.

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APPENDIX

By using definition (2.9) and relations (2.13) one can write the density matrix associated with a state *n* initially occupied in the left box as

$$\rho_{n}^{L}(\vec{r},\vec{r}',t) = \sum_{\alpha} d_{n\alpha}^{\pi_{x}p}(L) d_{n\alpha'}^{\pi_{x}p}(L) e^{(i/\hbar)(E_{\alpha}^{p} - E_{\alpha'}^{p})t} \phi_{\alpha}^{p}(\vec{r}) \phi_{\alpha'}^{p}(\vec{r}') + \frac{\sqrt{2}}{2} \sum_{\alpha \neq n} d_{n\alpha}^{+1}(L) [e^{(i/\hbar)E_{\alpha}^{p} - E_{n}^{-p})t} \phi_{\alpha}^{p}(\vec{r}) \phi_{n}^{-p}(\vec{r}') + e^{-(i/\hbar)(E_{\alpha}^{p} - E_{n}^{-p})t} \phi_{n}^{-p}(\vec{r}) \phi_{n}^{p}(\vec{r}')], \qquad (A1)$$

i.e., as a sum of two distinct terms: the first term contains products of states ϕ_{α}^{p} and $\phi_{\alpha'}^{p}$ of the same parity, while in the second term one wave function carries the parity -pof the initially occupied state and the other has the opposite parity p.

By using the same relations one finds that the density matrix ρ_n^R associated with a state *n* initially occupied in the right box is given by the difference between the same two terms.

In calculating $J_{+,n}^L$ or $J_{+,n}^R$ according to (3.5), (3.1), and (3.2) one can treat the two terms separately. Making the

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summation over all occupied states, each four times degenerate, one obtains

$$J_{+}^{L} = 4 \sum_{n \text{ (occ)}} J_{+,n}^{L} = \frac{1}{2} (J_{+} + J^{L}) , \qquad (A2)$$

$$J_{+}^{R} = 4 \sum_{n \text{ (occ)}} J_{+,n}^{R} = \frac{1}{2} (J_{+} - J^{L}) .$$
 (A3)

In (A2) J_{+} is given by (3.6) and originates from the first term in (A1) and J^{L} is given by (3.15) and originates from the second term of (A1). Similar considerations are valid for (A3) derived from ρ_{n}^{R} .

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