Simulation of tunneling in the quantum tomography approach

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The new method for the simulation of nonstationary quantum processes is proposed. The method is based on the tomography representation of quantum mechanics, i.e., the state of the system is described by the nonnegative function (quantum tomogram). In the framework of the method one uses the ensemble of trajectories in the tomographic space to represent evolution of the system (therefore direct calculation of the quantum tomogram is avoided). To illustrate the method we consider the problem of nonstationary tunneling of a wave packet. Different characteristics of tunneling, such as tunneling time, evolution of spatial and momentum distributions, tunneling probability are calculated in the quantum tomography approach. Tunneling of a wave packet of composite particle, exciton, is also considered; exciton ionization due to the scattering on the barrier is analyzed.

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I. INTRODUCTION

Nowadays simulation of quantum systems is developed to a high extent (see, e.g., reviews [1, 2]). However, common simulation methods, for example, Path Integral Monte Carlo or Wigner dynamics, use non-positively defined functions (wave function, the Wigner function, etc.) to describe a quantum state. This leads to the difficulties with convergence of corresponding integrals, especially harmful for the simulation of Fermi systems (sign problem). There is a hope that using a real nonnegative function, describing the quantum state, one can avoid these difficulties.

The real nonnegative function in the phase space, completely describing the quantum state, was proposed 60 years ago [3, see also 1,4,5]. During the last decade another very interesting representation has been actively developed: the quantum tomography, operating with the ensemble of scaled and rotated reference frames, instead of the phase space [4,5,6,8,10,11]. In the framework of this formalism the state-describing function (called marginal distribution or quantum tomogram) is real and nonnegative. The advantage is that the quantum tomogram is a probability distribution shown to completely describe the quantum state [12,13]. It is one of the reasons why the quantum tomography has become so popular.

In this paper we propose a new method for computer simulation of nonstationary quantum processes based on the tomography representation of quantum mechanics and illustrate it considering the problem of nonstationary tunneling of a wave packet. Many simulation approaches in nonstationary quantum mechanics are based on the numerical solution of the time-dependent Schrödinger equation. There are also methods using the ensembles of classical trajectories to simulate quantum evolution. For example, the method of ”Wigner trajectories”, based on the Wigner representation [14], is well known (see, e.g., review [1] for details) and was recently successfully applied to investigate the tunneling of a wave packet [15, 16]. Using such trajectories, one does not have to store the large arrays representing, say, the wave function (contrary to grid methods), therefore there is a gain in computer memory.

The quantum tomogram \( w \) depends on the variables \( \{X, \mu, \nu\} \), where \( X = \mu q + \nu p \), and \( q, p \) are the coordinates and momenta of the system, respectively, \( \mu, \nu \) are the parameters of scaling and rotation of reference frame in the phase space. The quantum tomogram is nonnegative and normalized in \( X \) direction, therefore it can be interpreted as a distribution function of the value \( X \). In our method the ensemble of trajectories in space \( \{X, \mu, \nu\} \) is introduced to describe the quantum evolution. The trajectories are governed by the dynamical equations obtained from the evolution equation for the quantum tomogram.

We demonstrate the method considering the nonstationary tunneling of a wave packet through the potential barrier. For this problem we calculated tunneling times, which are of interest nowadays, both for basic science (e.g., what is the time spent by an atom to tunnel from the trap?) and for applications (electronic tunneling time is connected with the operation rate of some nanostructure-based devices). We also analyzed the evolution of the wave packet in coordinate and momentum spaces in details. Another demonstration was designed to show that our method is not restricted to one-particle simulations. Namely, we investigated the tunneling of a wave packet of composite quasiparticle, an exciton (coupled electron and hole in semiconductor), in a one-dimensional nanostructure (quantum wire). In this case we had two degrees of freedom. For this problem, in addition to the probability density evolution, we determined the probability of ionization due to electron and hole scattering on the barrier in different directions.
In Sec. [III] we present description of the method, in Sec. [III] describe the model problem and main results for a wave packet tunneling. Exciton tunneling is considered in Sec. [V] and the work is summarized in Sec. [V].

II. THE METHOD OF SIMULATION

The quantum tomogram \( w(X, \mu, \nu) \) is connected with the density matrix \( \rho(q, q') \) as [17, 18]:

\[
\rho(q, q') = \int w(X, \mu, q - q') e^{i(X - \mu q'q')/2} \frac{d\mu dX}{2\pi} \quad (1)
\]

\[
w(X, \mu, \nu) = \int e^{-i(k(X - \mu q - \nu p) + \mu u)} \rho(q + \frac{u}{2}, q - \frac{u}{2}) d\mu dq du \quad (2)
\]

Consider the case of the particle with mass \( m \) in one-dimensional space. If the Hamiltonian of the system is

\[
H = \frac{p^2}{2m} + V(q),
\]

then the integral transformation [29] applied to the time-dependent evolution equation for the density matrix gives [6]

\[
\dot{w} - \frac{\mu}{m} \frac{\partial w}{\partial \nu} - 2 \frac{\partial V(q)}{\partial q} \left( \nu \frac{\partial}{2 \partial X} \right) w + \sum_{n=1}^{\infty} \frac{(-1)^{n+1}}{(2n+1)!} \frac{\partial^{2n+1} V(q)}{\partial q^{2n+1}} \left( \nu \frac{\partial}{2 \partial X} \right)^{2n+1} \hat{w} = 0,
\]

where we use \( \hat{w} = 1 \) and \( \hat{q} \) is given by

\[
\hat{q} = \left( \frac{\partial}{\partial X} \right)^{-1} \frac{\partial}{\partial \mu}
\]

Eq. (1) can be rewritten as

\[
\frac{\partial w}{\partial t} = \frac{\partial w}{\partial X} G_X(X, \mu, \nu) + \frac{\partial w}{\partial \mu} G_\mu(X, \mu, \nu) + \frac{\partial w}{\partial \nu} G_\nu(X, \mu, \nu) = 0,
\]

where functions \( G \) depend on quantum tomogram, its derivatives and antiderivatives (the latter corresponding to terms with \((\partial/\partial X)^{-1}\) in Eq. (1)). Generalization for the case of more variables is straightforward because the form of the equations does not change. Functions \( G \) for the problem under investigation are given in Sec. [III]. The evolution equation rewritten as (9) has the form of continuity equation for the quantum tomogram

\[
\frac{dw}{dt} = \frac{\partial w}{\partial t} + \frac{\partial w}{\partial X} \dot{X} + \frac{\partial w}{\partial \mu} \dot{\mu} + \frac{\partial w}{\partial \nu} \dot{\nu} = 0
\]

This equation is analogous to the continuity equation for classical distribution function and Liouville equation. As known, the characteristics of Liouville equation are the classical trajectories in phase space and they obey Hamilton equations of motion. The quantum tomogram is nonnegative and we use it as a distribution function for trajectories in the space \( \{X, \mu, \nu\} \), obeying the equations analogous to Hamilton equations for the classical trajectories. From the comparison of Eq. (1) with Eq. (4) it is obvious that the trajectories are governed by the equations

\[
\dot{X} = G_X(X, \mu, \nu), \quad \dot{\mu} = G_\mu(X, \mu, \nu), \quad \dot{\nu} = G_\nu(X, \mu, \nu)
\]

The trajectories are used to avoid the direct calculations of the distribution function (contrary to grid methods where the wave function is calculated at every point to solve numerically Schrödinger equation). Hence it is necessary to use some approximation for the quantum tomogram and we use local exponential approximation (as in Ref. [17] for the Wigner function):

\[
w(X, \mu, \nu) \approx \int \Delta \delta_{\nu a}(t) A_a(t)(y - y_a(t) + b_a(t)(y - y_a(t))) e^{-i(y_a(t) - y)} d\mu dX \quad (9)
\]

where \( \nu = \{X, \mu, \nu\} \) and \( y_a \) is the point under consideration. Parameters of this approximation are matrix \( A_a \) and vector \( b_a \), and some combinations of these parameters enter the evolution equation (11) instead of the derivatives and antiderivatives of the quantum tomogram. Calculation of average \( X, \mu, \nu \) and their average products allows to obtain \( A_a \) and \( b_a \). After that, functions \( G \) are known and dynamical equations (3) can be solved numerically.

We would like to emphasize that we use the local approximation (9) only for the calculation of r.h.s. in the equations of motion (8). The use of ensemble of trajectories to represent the quantum tomogram means the approximation of quantum tomogram as a set of delta-functions, each delta-function corresponds to one trajectory. If the number of trajectories approaches infinity, the quantum tomogram can be approximated by the set of delta-functions with arbitrary precision. This is analogous to what is conventionally done in classical statistical mechanics for the distribution function in phase space. But unlike the classical statistical mechanics now the trajectories are not independent: the approximation (9) holds for this case

\[
\hat{w}(X, \mu, \nu) \approx \int \Delta \delta_{\nu a}(t) A_a(t)(y - y_a(t) + b_a(t)(y - y_a(t))) e^{-i(y_a(t) - y)} d\mu dX
\]

where \( \nu = \{X, \mu, \nu\} \) and \( y_a \) is the point under consideration. Parameters of this approximation are matrix \( A_a \) and vector \( b_a \), and some combinations of these parameters enter the evolution equation (11) instead of the derivatives and antiderivatives of the quantum tomogram. Calculation of average \( X, \mu, \nu \) and their average products allows to obtain \( A_a \) and \( b_a \). After that, functions \( G \) are known and dynamical equations (3) can be solved numerically.

The validity of this approximation holds if the quantum tomogram is smooth and the trajectories are close to each other. For example, this approximation can fail if one tries to consider a plain wave with wave vector \( k \); in this case \( w(X, \mu = 0, \nu = 1) = \delta(X - k) \). Approximation (9) also works not well for unbounded motion, because the trajectories scatter with time. If there are few trajectories in the region around the given point, then the approximation (9) will not reconstruct the quantum tomogram well, due to lack of statistics.

We consider the tunneling of wave packets through the barrier, and comparing our results with exact quantum computation we see that approximation (9) holds for this...
problem (see Secs. III, IV). For example, considering the tunneling through the potential barrier from the well, we deal with both the region of bounded motion (in the well) and unbounded motion (beyond the barrier), still the approximation works not bad (Sec. III). For higher initial energy the penetration through the barrier is higher, and the evolution of most trajectories corresponds to unbounded motion, then the validity of Eq. (9) becomes poorer (see the end of Sec. III), in agreement with the discussion above. But in general local approximation (9) works satisfactory even for quite long time intervals (see Fig. 1).

To obtain any information about the system, we have to calculate some average values. Consider an arbitrary operator $A(\hat{q}, \hat{p})$. Average value $\langle A \rangle$ of corresponding physical quantity is calculated in the tomographic representation of quantum mechanics as 

$$\langle A \rangle = \int A(\mu, \nu) e^{iX} w(X, \mu, \nu) dX d\mu d\nu, \quad (10)$$

where $A(\mu, \nu)$ is the Fourier component of the Weyl symbol $A^W(q, p)$ of operator $A(\hat{q}, \hat{p})$ (see, e.g., [1]):

$$A(\mu, \nu) = \int A^W(q, p) \exp(-i(\mu q + \nu p)) \frac{dq dp}{4\pi^2} \quad (11)$$

For the calculation of average values we use the following approximation of quantum tomogram:

$$w(X, \mu, \nu, t) = \sum_{j=1}^{J} \delta(X - X_j(t)) \delta(\mu - \mu_j(t)) \delta(\nu - \nu_j(t)), \quad (12)$$

where the summation is made over all $J$ trajectories; $X_j(t), \mu_j(t), \nu_j(t)$ are the coordinates of the $j$-th trajectory in $\{X, \mu, \nu\}$ space at time $t$. Such approximation corresponds to use of the ensemble of trajectories. In the regions, where $w(X, \mu, \nu)$ is small, trajectories are rare, and where it is large, trajectories are accumulated. The more trajectories are used, the better the approximation works. If during the simulation the wave function has the form of a compact wave packet, even consisting of several distinct parts, approximation (12) holds, because in this case one has the compact sets of trajectories providing good statistics. This is the case for problems considered, and therefore the use of this approximation does not change results essentially, in comparison with the exact quantum computation (Secs. III, IV).

For the operators $A(\hat{q})$, depending on $\hat{q}$ only, expression for $\langle A \rangle$ takes the form:

$$\langle A \rangle = \int A(X) w(X, \mu = 1, \nu = 0) dX, \quad (13)$$

where $A(X)$ is the function corresponding to the operator $A(\hat{q})$ in coordinate representation, $A(X) = A(q = X)$. The method of calculation of an average $\langle A(\hat{q}) \rangle$ at arbitrary time $t$, with the approximation (12), is quite simple. One just takes into account the trajectories with any $X$ and with $\mu(t), \nu(t)$ from the small region near $\mu = 1, \nu = 0$ only, and performs a summation of $A(X)$ over all such trajectories.

The developed method is similar to well-known method of Wigner trajectories (see [1] for review), where the ensemble of trajectories is introduced in the phase space, with the Wigner function used as a quasi-distribution function. The quantum tomogram is defined in the space $\{X, \mu, \nu\}$, that is not as simple for understanding as the phase space used in the Wigner approach. On the other hand, quantum tomogram is true distribution function, while the Wigner function can be both positive and negative. In spite of these differences the two approaches are quite close in general, and there are some difficulties common for both methods. The discussion (see above) of the approximation for the tomogram (such as (9)) concerns, in principle, the method of Wigner trajectories either. Another important example is the discontinuity of the Wigner trajectories discussed in Ref. [20], that is due to the fact that the trajectories are not independent as in classical statistical mechanics, their evolution depends on their local distribution. The same is applicable for the quantum tomography approach, where the trajectories are also not independent for the same reason. In Ref. [21] a possible alternative to Wigner function was proposed. Namely, in that work the authors proposed to use the Weyl transforms of some operators (the Wigner function, up to the constant, is the Weyl transform of the density operator) instead of the Wigner function to generate the ensemble of trajectories. The same can be introduced for the quantum tomography. Applying the transform $A$ to the matrix elements of an operator $A$ we obtain the symbol $w_A(X, \mu, \nu)$. This function is not nonnegative in general, but in analogy with the Wigner trajectories one can use $w_A(X, \mu, \nu)$ to develop some new ensemble of trajectories. Probably, as with the Weyl transforms trajectories [21], it will be more convenient to use the trajectories corresponding to $w_A$ in certain cases. Of course, this problem needs further investigation.

III. SIMULATION OF TUNNELING OF A WAVE PACKET

A. The model and calculated average values

We choose the external potential to coincide with the potential used in [12], for comparison of the results of simulation in quantum tomography approach with those obtained by other methods. We consider behavior of the wave packet in one-dimensional space in external potential

$$V(q) = \frac{m a^2 q^2}{2} - \frac{b q^3}{3} \quad (14)$$

As the potential has only the second and third powers of coordinate, all its derivatives of order more than the third vanish. Evolution equation in this case has the form
Except the initial mean coordinate packet. We consider the problem, where all parameters, these processes depend on the initial energy of the wave zero. The particle can oscillate in the potential well and situated to the left from q maximum at q = 0.

This potential has the minimum at q = 0 (V(q) = 0) and maximum at q = 0.6709 (V(0.6709) = 0.015), therefore here we consider the motion of a particle in the potential well with infinite left wall and the barrier of height 0.015 at q = 0.6709. This model problem roughly describes nonstationary tunneling of an atom from the trap.

Initially the particle represented by the wave packet is situated to the left from q = 0, its mean momentum is zero. The particle can oscillate in the potential well and can tunnel or pass above the barrier. The probabilities of these processes depend on the initial energy of the wave packet. We consider the problem, where all parameters, except the initial mean coordinate q0 of the wave packet, are fixed (initial mean momentum equals zero, dispersions of the wave packet in coordinate and momentum spaces are ≈ 0.3 and ≈ 1.6, respectively).

We solve the equations numerically. As in we consider three values of q0: −0.2, −0.3 and −0.4. The most interesting quantities characterizing tunneling are reaction probability and tunneling time. Reaction probability is defined as

$$\int_{q_0}^{\infty} |\psi(x, t)|^2 dx,$$

where q0 = 0.6709 (the point where potential has the maximum), the maximum value of reaction probability is unity. Reaction probability shows what part of the wave packet is currently beyond the barrier.

Tunneling time of the wave packet is also an important feature of tunneling. There are a lot of methods to determine tunneling time. We use the approach where tunneling time is calculated as the difference of presence times (see for review) at point x_a and x_b, located on the opposite sides of the barrier:

$$t_R(x_a, x_b) = \langle t(x_b) \rangle - \langle t(x_a) \rangle$$

The presence time at arbitrary point x_0 is

$$\langle t(x_0) \rangle = \frac{\int_0^\infty t |\psi(x_0, t)|^2 dt}{\int_0^\infty |\psi(x_0, t)|^2 dt}$$

B. Reaction probability

Here we present the results obtained by means of our method and compare them with the exact numerical solution of Schrödinger equation. In Fig. 4 we present the dependence of reaction probability on time for three values of initial mean coordinate of the wave packet q_0 = −0.2, −0.3 and −0.4, corresponding mean energies of the wave packet are ≈ 0.75V_0, ≈ 1.25V_0 and ≈ 2.0V_0, respectively. Solid lines represent the results of simulation in the quantum tomography approach (QT) and dashed lines correspond to the numerical solution of Schrödinger equation (exact quantum computation). Due to the increase of initial mean energy with the increase of |q_0|, the portion of high energy components in the wave packet grows. This leads to the higher portion of components, which pass through the barrier, either because their energy is greater than the height of the barrier, or due to tunneling. Therefore, with the growth of |q_0|, reaction probability becomes larger and one can see that the curves corresponding to different q_0 are above each other in Fig. 4. The time evolution of reaction probability is qualitatively the same for every q_0. The components, which have passed through the barrier, can not return, because for q > 0.6709 potential diminishes with the growth of coordinate, and so reaction probability can not decrease with time. At first it grows rapidly due to transmission of components with the energy higher than the height of the barrier, or due to tunneling. Then reaction probability continues to grow slowly, because of the tunneling. All these features present both for QT simulation and for exact quantum computation.

In comparison with the exact computation, reaction probability for the QT simulation is slightly higher. Note also some difference in the character of increase of the reaction probability for QT simulation and exact solution: in the former case the curves are not so smooth. These differences are due to the finite number of trajectories used in QT simulation: for smaller number of
FIG. 1: The dimensionless reaction probabilities for three values of initial mean coordinate of the wave packet: \( q_0 = -0.2, -0.3, \) and \(-0.4\) a.u. Solid lines are for the simulation in quantum tomography approach, dashed lines are for the exact numerical solution.

trajectories (not shown) reaction probability curves resemble staircase more evidently (this is connected with the overestimation of the role of wave packet oscillations in the well for the finite number of trajectories), and quantitative deviation from exact result is stronger. But in general, for quite large number of trajectories, as for the case shown in Fig. 1, QT simulation results on reaction probability are quite close to those obtained through the exact quantum computation (compare also with the method of ”Wigner trajectories” in the work by Donoso and Martens).

C. Evolution of the wave packet and tunneling times

Besides the reaction probability we also obtained a number of new qualitative and quantitative results, which described in details the behavior of the wave packet during tunneling. We also calculated tunneling times using the concept of presence time (see below).

The following discussion concerns tunneling of the wave packet with initial mean coordinate \( q_0 = -0.2 \). Note that the maximum of the potential is at the point \( x = 0.6709 \). We present the normalized probability density \( |\psi(x)|^2 \) in coordinate space (Figs. 2-4) and \( |\psi(p)|^2 \) in momentum space (Figs. 5, 6) for several successive time moments. In these figures smooth lines show the shape of the wave packet obtained by means of exact quantum computation. Histograms represent the result of single QT run. One can consider many runs with the same number of trajectories and average the probability density over all these runs to obtain smoother picture. But here we would like to show what QT simulation can give for one run, in comparison with the exact quantum computation. Therefore the histograms (QT) fit the smooth solid lines in Figs. 2-4 (exact solution) not ideally, still the resemblance is obvious.

First, consider the probability density \( |\psi(x)|^2 \) in coordinate space (Figs. 2-4). One can see (Fig. 2) that initially the wave packet has Gaussian form. It begins to move as a whole towards the potential minimum at \( x = 0 \) (initial mean momentum is zero, but the potential inclines in that direction), passes that point, is acceler-
ated, and collides with the barrier. During the motion
the wave packet broadens (due to dispersion in momentum
space, compare right and left plots in Fig. 2) but
the interaction with the barrier changes its form more
substantially ($t = 300$ and $t = 400$, Figs. 3 and 4). The wave
packet shrinks a little, some components pass through
the barrier and transmitted part can be seen beyond the
barrier ($x = 0.6709$). As the transmitted part can not
return and accelerates (potential diminishes with the dis-
tance for $x > 0.6709$), the enriching of the wave packet
by high-energy components must present (see below).

All features described in the previous paragraph are
common for both the exact solution and QT simulation.
The histograms in Figs. 2 - 4 fit the smooth lines rep-re-
senting exact solution better for earlier times, but even
after the interaction with the barrier (Fig. 4), resem-
blance is quite close. This shows that approximations
(9) and (12) work not bad for the problem under consid-
eration.

Proceed now to the evolution of the wave packet in
momentum space. To confirm our analysis on the ac-
celeration of the transmitted part of the wave packet,
we present the probability density $|\psi(p)|^2$ in momentum
space in Figs. 5 and 6 at times $t = 0$ and $t = 400$, respec-
tively. As the wave function is initially the Gaussian
wave packet, the initial distributions both in coordinate
and momentum space are Gaussian (compare Fig. 2 and
Fig. 4). But after the wave packet has interacted with
the barrier, the distribution in momentum space changes
substantially (Fig. 5). The higher the energy of incident
particle, the greater the tunneling probability. There-
therefore the barrier transmits mainly the wave packet com-
ponents with relatively high energy, serving as an energy
selector. Components with high momentum arise, as the

transmitted part of the wave packet is accelerated in the
region of lowering potential beyond the barrier, and we
observe the enriching of the wave packet by high energy
components. The resemblance between the histograms
(QT simulation) and smooth solid lines (exact solution)
is somewhat poorer for momentum distribution at large
times ($t = 400$, Fig. 5) than for coordinate distribution
(Fig. 4). This is due to the fact that one deals with fi-
nite number of trajectories and has to sample quite large
interval in momentum space, because the transmitted
part is permanently accelerated. Therefore, with time,
momentum distribution spreads and there are not many
trajectories with $\mu, \nu$ close enough to $\mu = 0, \nu = 1$ for

FIG. 4: Probability density in coordinate space for QT sim-
ulation (histogram) and exact solution (smooth line), at time
$t = 400$ a.u. The barrier is at the point 0.6709 a.u., $q_0 = -0.2$
a.u.

FIG. 5: Initial probability density in momentum space for
QT simulation (histogram) and exact solution (smooth line).
$t = 0$ a.u.

FIG. 6: Probability density in momentum space for QT simu-
lation (histogram) and exact solution (smooth line) at
$t = 400$ a.u.
IV. SIMULATION OF THE EXCITON TUNNELING

The method described in Sec. III can be used to simulate the evolution of systems with more than one degree of freedom. In this section we demonstrate such a possibility, considering nonstationary tunneling of the composite particle, exciton, through the potential barrier in one-dimensional (1D) semiconductor structure (quantum wire). Exciton is a bound state of electron and hole in semiconductor, therefore we deal with two degrees of freedom in contrast to Sec. III.

Possible experimental realization is as follows. Consider semiconductor quasi-one-dimensional nanostructure where the motion is allowed only in one direction (quantum wire). Transverse motion is restricted due to strong confining barriers. Potential barrier in the direction of allowed motion can be located at some point of the quantum wire either using the semiconductor heterojunction or by the gate. Using femtosecond laser pulses, we can form an excitonic wave packet, either by the quasi-resonance pumping, or exciting an electron from valence band with formation of a hole and subsequent binding of two particles into exciton. Then the excitonic wave packet can move to the barrier and with the help of some detectors one can investigate scattering of the exciton.

Keeping this in mind let us construct the model for the simulation. We use the constants corresponding to GaAs for reference (dielectric constant $\varepsilon = 12.5$, effective masses of electron and hole are $m_e(0) = 0.07m_e^{*}$ and $m_h(0) = 0.15m_e^{*}$, respectively, where $m_e^{*}$ is the electron mass in vacuum). 3D exciton in bulk GaAs is characterized by effective Bohr radius $a^* \approx 10nm$ and binding energy $E_b \approx 4meV$. We use unit of length $a^*$, unit of mass $m_e$, and $\hbar = 1$. Corresponding units of energy and time are $E_0 = \hbar^2/(m_e a^*) \approx 10meV$ and $t_0 = m_e a^*/\hbar \approx 100fs$.

The energy spectrum and wave functions of relative electron and hole motion in 3D exciton are analogous to those of hydrogen atom. But this is not the case for 1D exciton. First, electron-hole effective interaction potential in quasi-1D structure is not Coulomb. Indeed, if the exciton size in the direction of allowed motion is much greater than the width of the quantum wire (in the transverse direction), then the adiabatic approximation is applicable and 3D interaction potential must be averaged over the transverse degrees of freedom. Resulting 1D effective potential substantially differs from Coulomb (see [35] for the discussion of similar model). Second, corresponding energy spectrum and wave functions of electron and hole relative motion also change in comparison with the hydrogen-like states. We choose the wave function of the exciton ground state in Gaussian form.

Excitonic wave packet can be represented as a Gaussian wave packet in center-of-mass coordinates:

$$\Psi(x_e, x_h, t = 0) = \frac{e^{-r^2/(2\sigma)} e^{-(R-x_0)^2/(2S)+iR_0}}{(\pi\sigma)^{1/4} (\pi S)^{1/4}},$$  \hspace{1cm} (21)
where \( R = (m_e x_e + m_h x_h) / (m_e + m_h) \), \( r = |x_e - x_h| \), \( x_e \) and \( x_h \) are electron and hole coordinates, \( x_0, p_0 \) and \( S \) are parameters; for them we used the following values: \( x_0 = -10 \), \( p_0 = 3 \), \( S = 2 \) and \( \sigma = 1 \).

External potential is assumed to be zero everywhere except the region of barrier; we use the barrier of thickness equal to 5\( \text{nm} \), or 0.5 in accepted units. For simplicity we set the barriers for electron and hole to be the same and use both external and interaction potentials in quadratic form, cut at some distance. Then external potential is given by

\[
V_{\text{ext}}(x) = \begin{cases} C - D x^2, & \text{if } |x| < \sqrt{\frac{C}{D}}, \\ 0, & \text{if } |x| \geq \sqrt{\frac{C}{D}}. \end{cases} \tag{22}
\]

\( C \) is the height of the barrier, its width is \( \sqrt{C/D} = 0.5 \).

Interaction potential \( V_{\text{int}} \) is also assumed to be quadratic:

\[
V_{\text{int}}(r) = \begin{cases} B r^2 - A, & \text{if } r < \sqrt{\frac{A}{B}}, \\ 0, & \text{if } r \geq \sqrt{\frac{A}{B}}. \end{cases} \tag{23}
\]

where \( r = |x_e - x_h| \). Potential \( V_{\text{int}} \) can describe, e.g., e-h interaction in spatially indirect exciton, for example in coupled quantum wires with large interwire separation \[30\]. Initial wave function of relative motion, chosen to be Gaussian with unity dispersion, is negligible within one percent accuracy at \( r = 3 \). Thus we choose the radius of electron-hole interaction to be \( \sqrt{A/B} = 3 \).

We assume that we deal with a quasi-1D exciton with binding energy \( E_C = 1/8 \). In fact, for an exciton in quantum wire, the wave function, binding energy, etc., are essentially influenced by the properties of quantum wire. We also neglect the possibility of electron and hole recombination at the time scales studied.

Here we consider just an example, therefore use relatively simple model. Still this model contains the main features of exciton tunneling, such as the possibility of ionization, barrier and interaction with realistic strength and size, the fact that composite particle is the bounded state of two particles.

For stationary state the binding energy is \( -E_C = \int \Psi_{\text{int}}^*(r) H_{\text{int}}(r) \Psi_{\text{int}}(r) dr \), where \( \Psi_{\text{int}}(r) \) is the wave function of relative motion. Then, from Eq. \[23\] and condition \( \sqrt{A/B} = 3 \) we have \( A \approx 18E_C/17 \).

We do not use the variables of relative and center-of-mass motion, for QT simulation it is easier to deal with the initial conditions and evolution equation in coordinates of electron and hole \( x_e \) and \( x_h \). Potentials \[22\] and \[23\] are quadratic, that makes tomographic consideration of the problem easier (see Sec. \[11\]). Discontinuity is neglected. On the other hand, in coordinates \( x_e \) and \( x_h \) the evolution equations depend on trajectory distribution (see Sec. \[11\]), therefore the problem considered allows to employ all techniques, developed for one degree of freedom, in this case of two degrees of freedom.

Results of exciton tunneling simulation are presented in Figs. \[8\]-\[10\]. All parameters are fixed (see above), except the barrier height \( C \) in Fig. \[10\]. In Figs. \[8\] and \[9\] we depict the evolution of probability density for electron and hole, barrier height \( C = 1 \). Solid lines in Figs. \[8\] and \[9\] and circles in Fig. \[10\] correspond to the simulation in quantum tomography approach, while dashed lines and
squares represent the exact numerical computation. Unlike in Figs. 2, 3 here we show the results of several combined QT simulation runs, therefore corresponding lines are relatively smooth. Coincidence of QT and exact computation results, initially very good, becomes poorer with time (Figs. 8 and 9), but QT simulation reproduces main properties of exciton tunneling: wave packets broadening with time and due to interaction with the barrier, shrinking near the barrier, dividing into two parts (reflected and transmitted). Note that QT results are quite close to exact ones even for long times ($t = 30$), despite the fact that the motion is unbounded (see Sec. IV). Integral values (in Fig. 10) obtained in QT approach also agree with the exact results. Larger discrepancies correspond to higher barriers, probably due to larger inaccuracies, introduced by neglecting the potential discontinuity in the case of stronger interaction with the external potential.

The electron and hole wave packets begin the motion from the point $x = -10$ and, shrinking near the barrier, are partially reflected and transmitted. For the case presented in Figs. 8 - 9 about the half of wave packets is transmitted. Interesting is the question about the ionization probability of exciton, induced by interaction with the barrier. If electron and hole are scattered in different directions on the barrier, the distance between them can become quite large, but, in principle, there is a possibility that exciton is not ionized after such scattering, because one of the particles can be ‘pulled’ beyond the barrier, to the other particle, due to electron-hole attraction. On the other hand, the electron-hole interaction is cut at the distance $\sqrt{A/B}$ in our model. After the interaction with the barrier the wave packet divides into reflected and transmitted parts moving in opposite directions. For the time large enough, these two parts are well separated, the separation between them grows and the leakage through the barrier in both directions is negligible. Denote the probability of ionization due to electron and hole scattering in different directions as $P_{\text{Ion}}$. Then, the probability to find electron and hole in different directions in respect to the barrier, with e-h distance being larger than $\sqrt{A/B}$, approaches $P_{\text{Ion}}$ in the limit $t \to \infty$.

The probability of ionization due to electron and hole scattering in different directions on the barrier $P_{\text{Ion}}$ is presented in Fig. 10 depending on the barrier height $C$. For very high and very low barriers $P_{\text{Ion}}$ must approach zero, because in former case both particles are reflected and in the latter they both are transmitted. This trend is seen in Fig. 10 and $P_{\text{Ion}}$ depending on $C$ is maximal (other parameters are fixed, see above) at $C \approx 1$. Note that these features are obvious for curves representing both QT simulation (circles) and exact computation (squares), and in general two curves are quite close to each other.

V. CONCLUSION

We have developed the new method of numerical simulation of quantum nonstationary processes and applied it to the problem of tunneling of the wave packet through the potential barrier. The method is based on tomographic representation of quantum mechanics. The quantum tomogram is used in a sense as the distribution function for the ensemble of trajectories in space $X, \mu, \nu$, where $X = \mu q + \nu p$ is the coordinate measured in rotated and scaled reference frame, $q, p$ are coordinate and momentum of the system, respectively. The trajectories are governed by the equations, resembling the Hamilton equations of motion, therefore, some analogue of molecular dynamics can be used. The Gaussian approximation allows to avoid the direct calculation of quantum tomogram. Instead of quantum tomogram, the parameters of the approximation are used in the equations of motion. Those parameters can be obtained if one calculates the local moments of the ensemble of the trajectories.

The problem of nonstationary tunneling of the wave packet was considered. Our method gave the results in agreement with those obtained by the method of "Wigner trajectories" and by exact quantum computation.

Of course, we made only the first step in the development of this simulation method, having considered one-dimensional problem. But it is obvious that the generalization for the case of several dimensions is straightforward. In this case the method possesses the additional advantage of more rapid convergence due to the fact the quantum tomogram is nonnegative. This may help to overcome the problem of sign for fermionic systems. In the next work we intend to consider the many-body problem for fermionic and bosonic systems by means of the quantum tomography method. The similar approach (in
the framework of "Wigner trajectories") has been already applied for the investigation of the tunneling of two identical particles.\footnote{10}

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