Description of the FORTRAN program KANTBP 3.0 for computing energy levels, reflection and transmission matrices, and corresponding wave functions in the coupled-channel adiabatic approach

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Abstract. Description of the program KANTBP 3.0 implemented in FORTRAN for calculating energy values, reflection and transmission matrices, and corresponding wave functions in a coupled-channel approximation of the adiabatic approach is presented. In this approach, a multidimensional Schrödinger equation is reduced to a system of the coupled second-order ordinary differential equations (SOODEs) on a finite interval with the homogeneous boundary conditions of the third type at the left- and right-boundary points for continuous spectrum problem, or a set of first, second and third type boundary conditions for discrete spectrum problem. The resulting system of these equations containing the potential matrix elements and first-derivative coupling terms is solved using high-order accuracy approximations of the finite element method. Efficiency of the schemes proposed is demonstrated on an example of solution of quantum transmittance two-dimensional problem for a pair of coupled ions connected by the harmonic oscillator interaction through the repulsive Coulomb-like barriers. As a test desk, the program is applied to the calculation of the reflection and transmission matrices and corresponding wave functions of the boundary-value problem for a set of N coupled-channel SOODEs which follows from the above two-dimensional problem.

Keywords: Scattering and tunneling problems, partial differential equations, finite-element method, high-order accuracy approximations

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1. Introduction

In the adiabatic approach [1] or the Kantorovich method [2], a multidimensional Schrödinger equation for quantum reflection [3], tunneling of a diatomic molecule incident upon a potential barrier [4, 5], fission model of collision of heavy ions [6], tunneling of a composite system thought barriers [7, 8, 9], the photoionization and decay of a hydrogen atom in magnetic field [10, 11, 12, 13, 14] is reduced by separating the longitudinal coordinate, labeled as \( z \in (-\infty, +\infty) \), from the transversal variables to a system of the second-order ordinary differential equations containing the potential matrix elements and first derivative coupling terms.

The purpose of this paper is to present a description the KANTBP 3.0 program [15] based on the use of the finite element method of high-order accuracy approximations [16, 17] for calculating reflection and transmission matrices and wave functions for such systems of coupled differential equations on finite intervals of the variable \( z \in [z_{\min}, z_{\max}] \) with homogeneous boundary conditions of the third-type at the left- and right-boundary points following from the above scattering problems. The third-type boundary conditions are formulated for the continuous problems under consideration by using known asymptotes for a set of linear independent asymptotic regular and irregular solutions in the open channels and a set of linear independent regular asymptotic solutions in the closed channels, respectively.

As a benchmark calculations the program is applied to the computation of the penetration coefficient for 2D-model of the pair charged particles connected by the harmonic oscillator interaction throughout symmetric or nonsymmetric as well as the Coulomb-like barriers [8].

As a test desk, the program is applied to the calculation of the reflection and transmission matrices and corresponding wave functions of the boundary-value problem for a set of \( N \) coupled-channel ordinary second order differential equations which follows from the above two-dimensional problem.

The paper is organized as follows. In Section 2 we give a brief overview of the problem. Description of the program KANTBP 3.0 is given in Section 3. Benchmark calculations are given in Section 4. Test desk is discussed in Section 5.

2. Statement of the problem

In the Kantorovich approach [2, 16], the multidimensional Schrödinger equation is reduced to a finite set of \( N \) ordinary second-order differential equations on the finite interval \([z_{\min}, z_{\max}]\) for the partial solution \( \chi^{(j)}(z) = (\chi_1^{(j)}(z), \ldots, \chi_N^{(j)}(z))^T \)

\[
(L - 2EI)\chi^{(j)}(z) \equiv \left(-I\frac{1}{z^{d-1}}\frac{d}{dz}z^{d-1}\frac{d}{dz} + V(z)\right)
+ Q(z)\frac{d}{dz} + \frac{1}{z^{d-1}}\frac{d}{dz}Q(z) - 2EI\chi^{(j)}(z) = 0. \tag{1}
\]
Here $I$, $V(z)$ and $Q(z)$ are the unit, symmetric and antisymmetric $N \times N$ matrices, respectively. We assume that $d = 1$, and $V(z)$ and $Q(z)$ matrices have the following asymptotic behaviour at large $z = z_\pm \rightarrow \pm \infty$

$$V_{ij}(z_\pm) = \left( \epsilon_j + \frac{2Z_j^\pm}{z_\pm} \right) \delta_{ij} + \sum_{l=2}^{\infty} \frac{v_{ij}^{(l,\pm)}}{z_\pm^l}, \quad Q_{ij}(z_\pm) = \sum_{l=1}^{\infty} \frac{q_{ij}^{(l,\pm)}}{z_\pm^l}, \quad (2)$$

where $\epsilon_1 \leq \ldots \leq \epsilon_N$ are the threshold values, and $\delta_{ij}$ is the Kronecker symbol.

In the present work, scattering problem is solved using the homogeneous third-type boundary conditions at $z = z_{\text{min}} \ll 0$ and $z = z_{\text{max}} \gg 0$:

$$\left. \frac{d\Phi(z)}{dz} \right|_{z=z_{\text{min}}} = R(z_{\text{min}})\Phi(z_{\text{min}}), \quad \left. \frac{d\Phi(z)}{dz} \right|_{z=z_{\text{max}}} = R(z_{\text{max}})\Phi(z_{\text{max}}), \quad (3)$$

where $R(z)$ is a unknown $N \times N$ matrix-function, $\Phi(z) = \{\chi^{(j)}(z)\}_{j=1}^N$ is the required $N \times N_o$ matrix-solution and $N_o$ is the number of open channels, $N_o = \max_{2E \geq \epsilon_j, j \leq N}$. From this we obtain the quadratic functional at $d = 1$ (similar to Eq. (23) in [16] and Eq. (5) in [17])

$$\Xi(\Phi, E, z_{\text{min}}, z_{\text{max}}) \equiv \int_{z_{\text{min}}}^{z_{\text{max}}} \Phi^\dagger(z) (L - 2EI) \Phi(z) dz = \Pi(\Phi, E, z_{\text{min}}, z_{\text{max}})$$

$$-\Phi^\dagger(z_{\text{max}})G(z_{\text{max}})\Phi(z_{\text{max}}) + \Phi^\dagger(z_{\text{min}})G(z_{\text{min}})\Phi(z_{\text{min}}), \quad (4)$$

where $\Pi(\Phi, E, z_{\text{min}}, z_{\text{max}})$ is the symmetric functional

$$\Pi(\Phi, E, z_{\text{min}}, z_{\text{max}}) = \int_{z_{\text{min}}}^{z_{\text{max}}} \left[ \frac{d\Phi^\dagger(z)}{dz} \frac{d\Phi(z)}{dz} - \Phi^\dagger(z)Q(z)\Phi(z) - 2E\Phi^\dagger(z)\Phi(z) \right] dz,$$

and $G(z) = R(z) - Q(z)$ is the $N \times N$ matrix-function which should be symmetric according to the conventual R-matrix theory [18]. Here the symbol $\dagger$ denotes the conjugate of a matrix.

### 2.1 The physical scattering asymptotic forms of solutions in longitudinal coordinates and scattering matrix

Matrix-solution $\Phi_v(z) = \Phi(z)$ describing the incidence of the particle and its scattering, which has the asymptotic form “incident wave + outgoing waves” (see Figure 1a), is

$$\Phi_v(z \rightarrow \pm \infty) = \begin{cases} \begin{align*} X^{(+)}(z)T_v, & z > 0, \quad v = \rightarrow, \\
X^{(+)}(z) + X^{(-)}(z)R_v, & z < 0, \quad v = \rightarrow, \\
X^{(-)}(z) + X^{(+)}(z)R_v, & z > 0, \quad v = \leftarrow, \\
X^{(-)}(z)T_v, & z < 0, \quad v = \leftarrow. \end{align*} \end{cases} \quad (6)$$
where the scattering matrix $S$ is composed of the reflection and transmission matrices

$$S = \begin{pmatrix} R_\to & T_\leftarrow \\ T_\to & R_\leftarrow \end{pmatrix}. \quad (10)$$

In addition, it should be noted that functions $X^{(\pm)}(z)$ satisfy relations

$$\text{Wr}(Q(z); X^{(\mp)}(z), X^{(\pm)}(z)) = \pm 2i\mathbf{I}_{oo}, \quad \text{Wr}(Q(z); X^{(\pm)}(z), X^{(\mp)}(z)) = 0, \quad (11)$$

where $\text{Wr}(\bullet; a(z), b(z))$ is a generalized Wronskian with a long derivative defined as

$$\text{Wr}(\bullet; a(z), b(z)) = a^T(z) \left( \frac{db(z)}{dz} - \bullet b(z) \right) - \left( \frac{da(z)}{dz} - \bullet a(z) \right)^T b(z). \quad (12)$$
This Wronskian will be used to estimate a desirable accuracy of the asymptotic expansions (7) and (17). Here the symbol $T$ denotes the transpose of a matrix.

Let us show that the scattering matrix (10) is symmetric and unitary. Using Eqs. (6) and (11), we have following relations

$$
\text{Wr}(Q(z); \Phi^\dagger_{\rightarrow}(z), \Phi_{\leftarrow}(z)) = \begin{cases} 
+2iT_{\rightarrow}^\dagger T_{\rightarrow}, & z > 0, \\
+2i(I_{oo} - R_{\rightarrow}^\dagger R_{\rightarrow}), & z < 0,
\end{cases}
$$

$$
\text{Wr}(Q(z); \Phi^*_{\rightarrow}(z), \Phi_{\leftarrow}(z)) = \begin{cases} 
-2iT_{\rightarrow}^\dagger T_{\rightarrow}, & z < 0, \\
-2i(I_{oo} - R_{\rightarrow}^\dagger R_{\rightarrow}), & z > 0,
\end{cases}
$$

$$
\text{Wr}(Q(z); \Phi_{\rightarrow}(z), \Phi^*_{\leftarrow}(z)) = \begin{cases} 
+2iR_{\rightarrow}^\dagger T_{\rightarrow}, & z > 0, \\
-2iR_{\rightarrow}^\dagger T_{\rightarrow}, & z < 0,
\end{cases}
$$

$$
\text{Wr}(Q(z); \Phi_{\rightarrow}(z), \Phi_{\leftarrow}(z)) = \begin{cases} 
+2iR_{\rightarrow}^\dagger R_{\rightarrow}, & z > 0, \\
-2iT_{\rightarrow}, & z < 0,
\end{cases}
$$

$$
\text{Wr}(Q(z); \Phi_{\leftarrow}(z), \Phi_{\rightarrow}(z)) = \begin{cases} 
+0, & z > 0, \\
+2i(R_{\rightarrow}^T - R_{\rightarrow}), & z < 0,
\end{cases}
$$

$$
\text{Wr}(Q(z); \Phi_{\rightarrow}(z), \Phi_{\leftarrow}(z)) = \begin{cases} 
-2i(R_{\rightarrow}^T - R_{\rightarrow}), & z > 0, \\
+0, & z < 0,
\end{cases}
$$

where the asterisk denotes the conjugate of a matrix. From here, we obtain the following properties of the reflection and transmission matrices:

$$
T_{\rightarrow}^\dagger T_{\rightarrow} + R_{\rightarrow}^\dagger R_{\rightarrow} = I_{oo} = T_{\leftarrow}^\dagger T_{\leftarrow} + R_{\leftarrow}^\dagger R_{\leftarrow},
$$

$$
T_{\rightarrow}^\dagger R_{\leftarrow} + R_{\rightarrow}^\dagger T_{\rightarrow} = 0 = T_{\rightarrow}^\dagger R_{\rightarrow} + T_{\leftarrow}^\dagger R_{\rightarrow},
$$

$$
T_{\rightarrow} = T_{\leftarrow}, \quad R_{\rightarrow} = R_{\rightarrow}, \quad R_{\rightarrow}^T = R_{\leftarrow}.
$$

This means that the scattering matrix (10) is symmetric and unitary.

Also matrix-solution $\hat{\Phi}_v(z) = \Phi(z)$ describing the incidence of the particle and its scattering, which has the inverse asymptotic form “incident waves + ingoing wave” (see Figure 1b), is

$$
\hat{\Phi}_v(z \to \pm \infty) = \begin{cases} 
X^{(+)}(z) + X^{(-)}(z) \hat{R}_v^\dagger, & z > 0, \quad v = \rightarrow, \\
X^{(+)}(z) \hat{T}_v^\dagger, & z < 0, \\
X^{(-)}(z) \hat{T}_v, & z > 0, \quad v = \leftarrow,
X^{(-)}(z) + X^{(+)}(z) \hat{R}_v^\dagger, & z < 0,
\end{cases}
$$

Note, that an equality $\hat{\Phi}^*_{\leftarrow}(z) = \Phi_{\leftarrow}(z)$ should be fulfilled from which we obtain $\hat{R}_{\rightarrow} = R_{\rightarrow}, \hat{R}_{\leftarrow} = R_{\leftarrow}, \hat{T}_v = T_v$. Therefore we consider below only matrix-solution $\Phi_v(z)$. 

2.2 Calculation of matrices $G(z_{\text{min}})$ at $v = \leftarrow$ and $G(z_{\text{max}})$ at $v = \rightarrow$

Suppose that a set of linear independent regular solutions $\Phi^{\text{reg}}_v(z) = \{X^{a}_j(z)\}_{j=1}^{N}$ for a problem under consideration is known at $z < 0$, $v = \leftarrow$, and at $z > 0$, $v = \rightarrow$, i.e.,

\[
\Phi^a_j(z) = X^{-}(z), \quad z < 0, \quad v = \leftarrow, \quad \text{and} \quad \Phi^a_j(z) = X^{+}(z), \quad z > 0, \quad v = \rightarrow,
\]

In the case of some channels are closed, we use additional linear independent regular asymptotic functions:

\[
\begin{align*}
\tilde{X}^{-}_ij(z) & \rightarrow q_j^{-1/2} \exp \left( + \left( q_j z + \sum_{j} \ln(2q_j|z|) \right) \delta_{ij}, \quad z < 0, \quad v = \leftarrow, \\
\tilde{X}^{+}_ij(z) & \rightarrow q_j^{-1/2} \exp \left( - \left( q_j z + \sum_{j} \ln(2q_j|z|) \right) \delta_{ij}, \quad z > 0, \quad v = \rightarrow, \\
q_j & = \sqrt{\epsilon_j - 2E}, \quad i = 1, \ldots, N, \quad j = N_o + 1, \ldots, N.
\end{align*}
\]

Then as shown in [17], the $G(z)$ matrix at $z = z_{\text{min}} < 0$, $v = \leftarrow$ and at $z = z_{\text{max}} > 0$, $v = \rightarrow$ can be found via the known set of linear independent regular solutions $\Phi^{\text{reg}}_v(z)$:

\[
G(z) = R(z) - Q(z) = \frac{d\Phi^{\text{reg}}_v(z)}{dz} (\Phi^{\text{reg}}_v(z))^{-1} - Q(z) \equiv (R(z) + R^T(z))/2. \tag{18}
\]

2.3 Calculation of matrix-solution $\Phi_v(z)$

After using the high-order accuracy approximations of the finite element method [16, 17, 19], the solution of a multichannel scattering problem at a fixed value of energy $E$ in open channels is reduced to a solution of the following algebraic problem with respect to matrix-solution $\Phi^h \equiv ((\chi^{(1)})^h, \ldots, (\chi^{(N)})^h)$

\[
G^p \Phi^h = (A^p - 2E B^p) \Phi^h = (M^p_{\text{max}} - M^p_{\text{min}}) \Phi^h, \tag{19}
\]

\[
\frac{d\Phi^h(z)}{dz} = (G(z) + Q(z))\Phi^h(z), \quad z = z_{\text{min}}, \quad z = z_{\text{max}}, \tag{20}
\]

where $A^p$ and $B^p$ are the symmetric $(LN) \times (LN)$ matrices, $L$ is the number of the nodes of the finite element grid on interval $[z_{\text{min}}, z_{\text{max}}]$, $M^p_{\text{max}}$ and $M^p_{\text{min}}$ are matrices with zero elements except the right-lower and left-upper $N \times N$ matrices equal to $G(z_{\text{max}})$ and $G(z_{\text{min}})$, respectively.

First, we consider the numerical algorithm for the calculation of matrix-solution $\Phi^h = \Phi^h_{\rightarrow}$. In this case Eq. (19) can be rewritten in the following form

\[
(G^p + M^p_{\text{min}}) \begin{pmatrix} \Phi^a_{\rightarrow} \\ \Phi^b_{\rightarrow} \end{pmatrix} = \begin{pmatrix} G^{aa}_{\rightarrow} & G^{ab}_{\rightarrow} \\ G^{ba}_{\rightarrow} & G^{bb}_{\rightarrow} \end{pmatrix} \begin{pmatrix} \Phi^a_{\rightarrow} \\ \Phi^b_{\rightarrow} \end{pmatrix} = \begin{pmatrix} 0 & 0 \\ 0 & G(z_{\text{max}}) \end{pmatrix} \begin{pmatrix} \Phi^a_{\rightarrow} \\ \Phi^b_{\rightarrow} \end{pmatrix}, \tag{21}
\]

where $\Phi^a_{\rightarrow}$ and $\Phi^b_{\rightarrow} \equiv \Phi^a_{\rightarrow}(z_{\text{max}})$ are the matrix-solutions of dimension $(LN - N) \times N_o$ and $N \times N_o$, respectively. From here, we obtain explicit expressions

\[
\Phi^a_{\rightarrow} = -(G^{aa}_{\rightarrow})^{-1}G^{ab}_{\rightarrow} \Phi^b_{\rightarrow}, \quad G(z_{\text{max}}) = G^{bb}_{\rightarrow} - G^{ba}_{\rightarrow}(G^{aa}_{\rightarrow})^{-1}G^{ab}_{\rightarrow}. \tag{22}
\]
From Eqs. (20) and (22) we can obtain the relation between $\Phi^b_\leftarrow$ and its derivative
\[
\frac{d\Phi^b_\leftarrow}{dz} = R(z_{\text{max}})\Phi^b_\leftarrow, \quad R(z_{\text{max}}) = G(z_{\text{max}}) + Q(z_{\text{max}}).
\] (23)

Note, that matrix $G(z_{\text{max}})$ is determined via the inverse of submatrix $G_{\text{zz}}^a$ calculation of which requires substantial computer resources. For evaluating Eq. (23) without such calculation of the inverse of submatrix $G_{\text{zz}}^a$, let’s consider the following auxiliary system of algebraic equations
\[
\begin{pmatrix}
G_{\text{aa}}^a & G_{\text{ab}}^a \\
G_{\text{ba}}^a & G_{\text{bb}}^a
\end{pmatrix}
\begin{pmatrix}
F^a_\leftarrow \\
F^b_\leftarrow
\end{pmatrix}
= \begin{pmatrix} 0 \\ 1 \end{pmatrix}.
\] (24)

As the determinant of the matrix $G^b + M_{\text{min}}^b$ is nonzero, the above equation has a unique solution
\[
F^a_\leftarrow = -(G_{\text{aa}}^a)^{-1}G_{\text{ab}}^a F^b_\leftarrow, \quad F^b_\leftarrow = (G_{\text{bb}}^b - G_{\text{ba}}^b(G_{\text{aa}}^a)^{-1}G_{\text{ab}}^a)^{-1}.
\] (25)

Taking this into account, the required $R(z_{\text{max}})$ matrix is equal to
\[
R(z_{\text{max}}) = (F^b_\leftarrow)^{-1} + Q(z_{\text{max}}).
\] (26)

Using Eqs. (23) and (6), we obtain the following matrix equation for the reflection $R_\leftarrow$ matrix:
\[
Y^{(+)}(z_{\text{max}})R_\leftarrow = -Y^{(-)}(z_{\text{max}}), \quad Y^{(+)}(z) = \frac{dX^{(\pm)}(z)}{dz} - R(z)X^{(\pm)}(z).
\] (27)

Then the required solution $\Phi^h_\leftarrow$ is calculated by formulae (6), (22) and (25)
\[
\Phi^a_\leftarrow = X^{(-)}(z_{\text{max}}) + X^{(+)}(z_{\text{max}})R_\leftarrow, \quad \Phi^a_\leftarrow = F^a_\leftarrow (F^b_\leftarrow)^{-1} \Phi^h_\leftarrow.
\] (28)

The transmission $T_\leftarrow$ matrix is determined from the matrix equation
\[
X^{(-)}(z_{\text{min}})T_\leftarrow = \Phi^h_\leftarrow(z_{\text{min}}).
\]

Note that, when some channels are closed, the $Y^{(\pm)}(z)$ and $X^{(-)}(z)$ are rectangular $N \times N_o$ matrices. Therefore, using the pseudoinverse matrices of $Y^{(\pm)}(z)$ and $X^{(-)}(z)$, we obtain the following formulae:
\[
R_\leftarrow = - \left( (Y^{(+)}(z_{\text{max}}))^T Y^{(+)}(z_{\text{max}}) \right)^{-1} (Y^{(+)}(z_{\text{max}}))^T Y^{(-)}(z_{\text{max}}),
\] (29)
\[
T_\leftarrow = \left( (X^{(-)}(z_{\text{min}}))^T X^{(-)}(z_{\text{min}}) \right)^{-1} (X^{(-)}(z_{\text{min}}))^T \Phi^h_\leftarrow(z_{\text{min}}).
\]

Now we will describe briefly a calculational scheme for matrix-solution $\Phi^h_\rightarrow = \Phi^h_\leftarrow$. The required $R(z_{\text{min}})$ matrix is equal to
\[
R(z_{\text{min}}) = (F^a_\rightarrow)^{-1} + Q(z_{\text{min}}),
\] (30)
and required solution $\Phi^b_{\rightarrow}$ is calculated as

$$\Phi^b_{\rightarrow} = F^b_{\rightarrow} (F^a_{\rightarrow})^{-1} \Phi^a_{\rightarrow}, \quad \Phi^{a}_{\rightarrow} = X^{(+)}(z_{\text{min}}) + X^{(-)}(z_{\text{min}}) R_{\rightarrow}. \quad (31)$$

Here $\Phi^a_{\rightarrow} \equiv \Phi_{\rightarrow}(z_{\text{min}})$ and $\Phi^b_{\rightarrow}$ are the matrix-solutions of dimension $N \times N_o$ and $(LN-N) \times N_o$. $F^a_{\rightarrow}$ and $F^b_{\rightarrow}$ are the matrices of dimension $N \times N$ and $(LN-N) \times N$ which are the solutions of the auxiliary system of algebraic equations

$$(G^p - M^p_{\text{max}}) \begin{pmatrix} F^a_{\rightarrow} \\ F^b_{\rightarrow} \end{pmatrix} = - \begin{pmatrix} G^a_{\rightarrow} \\ G^b_{\rightarrow} \end{pmatrix} \begin{pmatrix} F^a_{\rightarrow} \\ F^b_{\rightarrow} \end{pmatrix} = - \begin{pmatrix} I \\ 0 \end{pmatrix}. \quad (32)$$

Finally, we obtain the following matrix equations for the reflection $R_{\rightarrow}$ and transmission $T_{\rightarrow}$ matrices:

$$Y^{(-)}_{\rightarrow}(z_{\text{min}}) R_{\rightarrow} = - Y^{(+)}_{\rightarrow}(z_{\text{min}}), \quad Y^{(+)}_{\rightarrow}(z) = \frac{dX^{(+)}(z)}{dz} - R(z) X^{(+)}(z), \quad (33)$$

$$X^{(+)}_{\rightarrow}(z_{\text{max}}) T_{\rightarrow} = \Phi^b_{\rightarrow}(z_{\text{max}}).$$

The reflection $R_{\rightarrow}$ and transmission $T_{\rightarrow}$ matrices are evaluated using the pseudoinverse matrices of $Y^{(-)}_{\rightarrow}(z_{\text{min}})$ and $X^{(+)}_{\rightarrow}(z_{\text{max}}):

$$R_{\rightarrow} = - \left( (Y^{(-)}_{\rightarrow}(z_{\text{min}}))^T Y^{(-)}_{\rightarrow}(z_{\text{min}}) \right)^{-1} (Y^{(-)}_{\rightarrow}(z_{\text{min}}))^T Y^{(+)}_{\rightarrow}(z_{\text{min}}), \quad (34)$$

$$T_{\rightarrow} = \left( (X^{(+)}_{\rightarrow}(z_{\text{max}}))^T X^{(+)}_{\rightarrow}(z_{\text{max}}) \right)^{-1} (X^{(+)}_{\rightarrow}(z_{\text{max}}))^T \Phi^b_{\rightarrow}(z_{\text{max}}).$$

### 2.4 Asymptotic forms of regular and irregular solutions in the longitudinal coordinates

We calculate the asymptotic solution to a set of $N$ coupled ordinary differential equations (ODE) at large values of independent variable $|z| \gg 1$

$$\left( -\frac{1}{z^{d-1}} \frac{d}{dz} z^{d-1} \frac{d}{dz} + V_{ii}(z) - 2E \right) \chi_{ii}(z) = - \sum_{j=1, j \neq i}^{N} \left( V_{ij}(z) + Q_{ij}(z) \frac{d}{dz} \right) \frac{d}{dz} + \frac{1}{z^{d-1}} \frac{d}{dz} z^{d-1} Q_{ij}(z) \chi_{jj}(z). \quad (35)$$

Here $d \geq 1$ is the dimension of configuration space of a general scattering problem [17]. For the considered case, we put $d = 1$ and calculate asymptotic solution on two intervals $-\infty < z \leq z_{\text{min}}$ and $z_{\text{max}} \leq z < +\infty$. We assume that coefficients of Eqs. (35) can be represented in the general asymptotic form as

$$V_{ij}(z) = \left( \epsilon^{(0)}_j + \epsilon^{(1)}_j \frac{1}{z^l} \right) \delta_{ij} + \sum_{l=2}^{k_{\text{max}}+1} \frac{V^{(l)}_{ij}}{z^l}, \quad Q_{ij}(z) = \sum_{l=1}^{k_{\text{max}}+1} \frac{Q^{(l)}_{ij}}{z^l}. \quad (36)$$

Note that in general case coefficients $\epsilon^{(1)}_j, V^{(l)}_{ij}$ and $Q^{(l)}_{ij}$ are different for $z > 0$ and $z < 0$. Below we will consider only case of $z > 0$. 


**Step 1.** We construct the solution of Eqs. (35) in the form:

\[ \chi_{jiv}(z) = \phi_{jiv}(z)R_{i'}(z) + \psi_{jiv}(z)\frac{dR_{i'}(z)}{dz}, \quad (37) \]

where \( \phi_{jiv}(z) \) and \( \psi_{jiv}(z) \) are unknown functions, \( R_{i'}(z) \) is a known function. We choose \( R_{i'}(z) \) as solutions of the auxiliary problem treated like etalon equation \((Z_{i'}^{(k<1)} = Z_{i'}^{(k>k_{\text{max}})} = 0):\)

\[ \left(-\frac{1}{z^{d-1}}\frac{dz}{d}z^{d-1} + \sum_{k=1}^{k_{\text{max}}} \frac{Z_{i'}^{(k)}}{z^k} - p_{i'}^2 \right) R_{i'}(z) = 0. \quad (38) \]

**Remark 1.** If \( Z_{i'}^{(k\geq3)} = 0 \) then solutions to the last equation are presented via the hypergeometric functions, exponential, trigonometric, Bessel, Coulomb functions, etc. For example, if the leading terms of the asymptotic solutions are given by formula

\[ R_{i'}(z) = \frac{1}{\sqrt{p_{i'}z^{d-1}}} \exp \left( \pm i \left( p_{i'}z - \frac{Z_{i'}}{p_{i'}} \ln(2p_{i'}|z|) \right) \right), \quad (39) \]

the coefficients of potential in the etalon equation (38) have the form:

\[ Z_{i'}^{(1)} = 2Z_{i'}, \quad Z_{i'}^{(2)} = -\frac{(d-3)(d-1)}{4} \pm i \frac{Z_{i'}}{p_{i'}} - \frac{Z_{i'}^2}{p_{i'}^2}. \quad (40) \]

**Step 2.** At this step we compute the coefficients \( \phi_{jiv}(z) \) and \( \psi_{jiv}(z) \) of the expansion (37) in the form of series over inverse powers of \( z:\)

\[ \phi_{jiv}(z) = \sum_{k'=0}^{k_{\text{max}}} \phi_{jiv'}(z), \quad \psi_{jiv}(z) = \sum_{k'=0}^{k_{\text{max}}} \psi_{jiv'}(z). \quad (41) \]

After substitution of Eq. (37), (41) into (35) with the use of Eq. (38) and equating the coefficients at \( z^{-k'}R_{i'}(z) \) and \( z^{-k'}\frac{dR_{i'}(z)}{dz} \), we arrive to a set of recurrent relations at \( k' \leq k_{\text{max}}:\)

\[ \left( \epsilon_i^{(0)} - 2E + p_{i'}^2 \right) \phi_{iiv'}^{(k')} + \left( \epsilon_i^{(1)} - Z_{i'}^{(1)} \right) \phi_{iiv'}^{(k'-1)} - 2p_{i'}^2(k'-1)\psi_{iiv'}^{(k'-1)} = -f_{iiv'}^{(k')}, \quad (42) \]

\[ \left( \epsilon_i^{(0)} - 2E + p_{i'}^2 \right) \psi_{iiv'}^{(k')} + 2(k'-1)\phi_{iiv'}^{(k'-1)} + \left( \epsilon_i^{(1)} - Z_{i'}^{(1)} \right) \psi_{iiv'}^{(k'-1)} = -g_{iiv'}^{(k')}, \quad (43) \]
where the right hand sides \( f_{i'i'}^{(k')} \) and \( g_{i'i'}^{(k')} \) are defined by relations

\[
f_{i'i'}^{(k')} = -(k'-2)(k'-d)\phi_{i'i'}^{(k'-2)} + \sum_{k=2}^{k'} \left( V_{ii}^{(k)} - Z_{i'i'}^{(k)} \right) \phi_{i'i'}^{(k'-k)}
+ \sum_{k=1}^{k'} \left( Z_{i'i'}^{(k)}(2k' - 2 - k)\psi_{i'i'}^{(k'-k-1)} + \sum_{j=1,j \neq i}^{N} \left( \sum_{k''=1}^{k'} 2Q_{ij}^{(k)} Z_{i'i'}^{(k''-k-k'')} \right)
- 2p_{i'i'}^{2} Q_{ij}^{(k)} \psi_{i'i'}^{(k'-k)} + Q_{ij}^{(k)} (-2k' + k + d + 1) \phi_{i'i'}^{(k'-k-1)} + V_{ij}^{(k)} \phi_{i'i'}^{(k'-k)} \right);
\]

\[
g_{i'i'}^{(k')} = -(k'-1)(k'-3 + d)\psi_{i'i'}^{(k'-2)} + \sum_{k=2}^{k'} \left( V_{ii}^{(k)} - Z_{i'i'}^{(k)} \right) \psi_{i'i'}^{(k'-k)}
+ \sum_{j=1,j \neq i}^{N} \sum_{k=1}^{k'} \left( 2Q_{ij}^{(k)} \phi_{i'i'}^{(k'-k)} - Q_{ij}^{(k)} (2k' + d - 3 - k) \psi_{i'i'}^{(k'-k-1)} + V_{ij}^{(k)} \psi_{i'i'}^{(k'-k)} \right)
\]

with initial conditions \( p_{i'i'}^{2} = 2E - \epsilon_{i'i'}^{(0)}, \phi_{i'i'}^{(0)} = \delta_{ii}, \psi_{i'i'}^{(0)} = 0 \), at \( i' = i \) span over the open channels \( i = 1, \ldots, N \) and \( p_{i'i'} = u_{i'i'}, q_{i'i'} > 0, q_{i'i'}^{2} = \epsilon_{i'i'}^{(0)} - 2E \) at \( i' = i \) span over the closed channels \( i = N + 1, \ldots, N \) that followed from (7) and (17). Also from Eq. (42) at \( k' = 1 \) and \( i = i' \),

\[
\left( \epsilon_{i'i'}^{(1)} - Z_{i'i'}^{(1)} \right) \phi_{i'i'}^{(0)} = 0, \quad \left( \epsilon_{i'i'}^{(1)} - Z_{i'i'}^{(1)} \right) \psi_{i'i'}^{(0)} = 0,
\]

we obtain condition \( Z_{i'i'}^{(1)} = \epsilon_{i'i'}^{(1)} \).

**Step 3.** Here we perform calculation of the coefficients \( \phi_{i'i'}^{(k')} \) and \( \psi_{i'i'}^{(k')} \) by a step–by–step procedure of solving Eqs. (42) for \( 2E \neq \epsilon_{i'i'}^{(0)}, i \neq i' \) and \( k' = 1, \ldots, k_{\text{max}} \):

\[
\phi_{i'i'}^{(k')} = \left[ \epsilon_{i'i'}^{(0)} - \epsilon_{i'i'}^{(0)} \right]^{-1} \left[ -f_{i'i'}^{(k')} - \left( \epsilon_{i'i'}^{(1)} - Z_{i'i'}^{(1)} \right) \phi_{i'i'}^{(k'-1)} + 2p_{i'i'}^{2} (k' - 1) \psi_{i'i'}^{(k'-1)} \right],
\]

\[
\psi_{i'i'}^{(k')} = \left[ \epsilon_{i'i'}^{(0)} - \epsilon_{i'i'}^{(0)} \right]^{-1} \left[ -g_{i'i'}^{(k')} - 2(k' - 1) \phi_{i'i'}^{(k'-1)} - \left( \epsilon_{i'i'}^{(1)} - Z_{i'i'}^{(1)} \right) \psi_{i'i'}^{(k'-1)} \right],
\]

and for \( 2E \neq \epsilon_{i'i'}^{(0)}, i = i' \) and \( k' = 2, \ldots, k_{\text{max}} \):

\[
\phi_{i'i'}^{(k'-1)} = - \left[ 2(k' - 1) \right]^{-1} g_{i'i'}^{(k')} - \epsilon_{i'i'}^{(0)} \left( \epsilon_{i'i'}^{(0)} - 2E \right) \left( \epsilon_{i'i'}^{(0)} - 2E \right)^{-1} f_{i'i'}^{(k')}.
\]

Algorithm described above has been implemented in the MAPLE and FORTRAN (see description of SLAS program in section 3.2). Resulting output provided evaluation of the \( \chi_{i'i'}(z) \) and \( \chi_{i'i'}^{(0)}(z) \frac{dz}{dz} \). This algorithm has been examined with the results from [20].

**Remark 2.** The choice of appropriate values \( z_{\text{min}} \) and \( z_{\text{max}} \) for the constructed expansions of the linearly independent solutions for \( p_{c} > 0 \) is controlled by the fulfillment of the Wronskian condition (11), (12)

\[
\text{Wr}(Q(z); \chi^{*}(z), \chi(z)) = \pm 2i\text{Im}
\]

up to the prescribed precision \( \varepsilon_{W} \).
Figure 2: Flow diagram of the new version of the KANTBP 3.0 program.
3. Description of the program

Figure 2 presents a flow diagram for the program. The KANTBP 3.0 program is called from the main routine (supplied by a user) which sets dimensions of the arrays and is responsible for the input data. The KANTBP 3.0 program needs no installation. The description of all subroutines can be found in comments in the program source code. Also users can find instructions on how to compile the KANTBP 3.0 in the README file.

The calling sequence for the subroutine KANTBP is:

```
CALL KANTBP(TITLE, IPTYPE, ISC, NROOT, MDIM, IDIM, NPOL, RTOL, NITEM,
1       SHIFT, IPRINT, IPRSTP, NMESS, RMESH, NDIR, NDIL, NMIDIL,
2       THRSHL, IBOUND, FNOUT, IOUT, POTEN, Ioup, FMATR, IouM,
3       EWfWN, IOUF, TOT, ITOT, ZTOT, MTOT, MITOT, MZTOT)
```

In the present code each array declarator is written in terms of the symbolic names of constants. These constants are defined in the following PARAMETER statement in the main routine:

```
PARAMETER (MTOT=10000, MITOT=30000,
1       MZTOT=90000, NMESH=7, MDIM=4)
```

Here

- MTOT is the dimension of the working DOUBLE PRECISION array TOT. The last address ILAST of array TOT is calculated and then compared with the given value of MTOT. If ILAST > MTOT the message about an error is printed and the execution of the program is aborted. In the last case, in order to carry out the required calculation it is necessary to increase the dimension MTOT of array TOT to the quantity ILAST taken from the message.

- MITOT is the dimension of the working INTEGER array ITOT. The last address ILAST of array ITOT is calculated and then compared with the given value of MITOT. If ILAST > MITOT the message about an error is printed and the execution of the program is aborted. In the last case, in order to carry out the required calculation it is necessary to increase the dimension MITOT of array ITOT to the quantity ILAST taken from the message.

- MZTOT is the dimension of the DOUBLE COMPLEX working array ZTOT. The last address ILAST of array ZTOT is calculated and then compared with the given value of MZTOT. If ILAST > MZTOT the message about an error is printed and the execution of the program is aborted. In the last case, in order to carry out the required calculation it is necessary to increase the dimension MZTOT of array ZTOT to the quantity ILAST taken from the message.

- NMESH is the dimension of the DOUBLE PRECISION array RMESH containing the information about the subdivision of the longitudinal interval $[z_{min}, z_{max}]$ on subintervals and number of elements on each one of them. NMESH is always odd and $\geq 3$. 
MDIM is the dimension of the DOUBLE PRECISION array THRSHL and INTEGER array NDIL containing information about a set of threshold values and numbers of coupled differential equations, respectively.

In order to change the dimensions of the code, all one has to do is to modify the single PARAMETER statement defined above in the main program unit.

We have added a new flag ISC for performing the calculation of the reflection and transmission matrices:

- $= 1$ – calculation of the reflection and transmission matrices is carried out only with direction $v = \leftarrow$;
- $= 2$ – calculation of the reflection and transmission matrices is carried out only with direction $v = \rightarrow$;
- $= 3$ – calculation of the reflection and transmission matrices is carried out with both directions $v = \leftarrow$ and $v = \rightarrow$. Also the properties (2.4) of the reflection and transmission matrices are verified.

The meaning of the all arguments except above are presented in [16].

New output data

The results of the calculation of the reflection and transmission matrices and corresponding wave functions are written using unformatted segmented records into file EVWFN, according to the following operator:

```
WRITE(IOUF) NDIM,NN,NOPEN,NGRID,((RR(I,J),I=1,NOPEN),J=1,NOPEN),
 1 ((TT(I,J),I=1,NOPEN),J=1,NOPEN),
 1 (XGRID(I),I=1,NGRID),((R(I,J),I=1,NN),J=1,NOPEN)
```

In the above, parameters presented in the WRITE statement have the following meaning:

- NDIM is the number of coupled equations,
- NGRID is the number of finite-element grid points,
- $NN = NGRID \times NDIM$,
- NOPEN is the number of open channels,
- Arrays RR and TT contain the reflection and transmission matrices values calculated,
- Array XGRID contains the values of the finite-element grid points,
- Array R contains NOPEN eigenfunctions each per NN elements in length stored (see the scheme in [16]).

New user-supplied subroutines
ASYMSL is the name of the new user-supplied subroutine for the scattering problem, $\Phi_-(z)$, which calculates the regular $X^-(z)$, irregular $X^+(z)$ asymptotic rectangle-solutions and their derivatives at $z = z_{\text{max}}$, and regular $X^-(z)$ asymptotic square-solution and its derivative at $z = z_{\text{min}}$. It should be written as follows:

```fortran
SUBROUTINE ASYMSL(ZMIN,ZMAX,NDIM,NOPEN,QR,SHIFT,THRSHL,
1       PREGL,DREGL,PREGR,PIRRR,DREGR,DIRRR,IOUT)
C . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . .
C ..
C . P R O G R A M
C . TO CALCULATE THE REGULAR, IRREGULAR
C . ASYMPTOTIC MATRIX SOLUTIONS PREGR, PIRRR
C . AND THEIR DERIVATIVES DREGR, DIRRR AT ZMAX,
C . THE REGULAR MATRIX SOLUTION PREGL AND ITS
C . DERIVATIVE DREGL AT ZMIN
C .
C 
IMPLICIT REAL*8 (A-H,O-Z)
DIMENSION QR(NOPEN),THRSHL(NDIM)
COMPLEX*16 PREGL(NDIM,NDIM),DREGL(NDIM,NDIM),
1       PREGR(NDIM,NOPEN),PIRRR(NDIM,NOPEN),
1       DREGR(NDIM,NOPEN),DIRRR(NDIM,NOPEN)
RETURN
END
```

ASYMSR is the name of the new user-supplied subroutine for the scattering problem, $\Phi_+(z)$, which calculates the regular $X^+(z)$, irregular $X^-(z)$ asymptotic rectangle-solutions and their derivatives at $z = z_{\text{min}}$, and regular $X^+(z)$ asymptotic square-solution and its derivative at $z = z_{\text{max}}$. It should be written as follows:

```fortran
SUBROUTINE ASYMSR(ZMIN,ZMAX,NDIM,NOPEN,QR,SHIFT,THRSHL,
1       PREGL,DREGL,PREGR,PIRRL,DREGL,DIRRL,IOUT)
C . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . .
C ..
C . P R O G R A M
C . TO CALCULATE THE REGULAR, IRREGULAR
C . ASYMPTOTIC MATRIX SOLUTIONS PREGR, PIRRL
C . AND THEIR DERIVATIVES DREGL, DIRRL AT ZMIN,
C . THE REGULAR MATRIX SOLUTION PREGL AND ITS
C . DERIVATIVE DREGL AT ZMAX
C .
C 
```
Here arrays QR and THRSHL contain a set of momentum and threshold values, respectively; SHIFT contains the given double energy spectrum value; NDIM is the number of coupled equations; NOPEN is the number of open channels; IOUT is number of the output logical device for printing out the results of the calculation.

To set the third-type boundary conditions at both points \( z_{\text{min}} < 0 \) and \( z_{\text{max}} > 0 \), flags IBOUND and IDIM always should be 8 and 1. Here IBOUND is parameter defining the type of boundary conditions, and IDIM is dimension of the envelope space [16].

3.1 Description of new subprogram units

The function of each new subroutine is briefly described below. Additional details may be found in COMMENT cards within the program.

- Subroutine ADDVEK assembles the element into the corresponding global complex vector using a compact storage form. This is modified version of subroutine ADDVEC [16] for complex arithmetics.

- Subroutine ASSMBC controls the calculation of element complex stiffness matrix and assembles them into the corresponding global complex matrix. This is modified version of subroutine ASSMBL [16] for complex arithmetics.

- Subroutine CHECRT controls the properties (2.4) of calculated reflection \( R_{\leftarrow} \), \( R_{\rightarrow} \) and transmission \( T_{\leftarrow} \), \( T_{\rightarrow} \) matrices.

- Subroutine CHECKN prints error messages when input data are incorrect and stops the execution of program KANTBP.

- Subroutine DECOMC calculates \( LDL^T \) factorization of matrix. This factorization is used in subroutine REDBAC to reduce and back-substitute the iteration vectors. They are modified versions of subroutines DECOMP and REDBAK [19] for complex arithmetics.

- Subroutine GAUSSC calculates linear equation solution by the Gauss-Jordan matrix inversion method. This is a modified version of subroutine GAUSSJ [21] for complex arithmetics.
• Subroutine HQPOTN calculates potential matrix elements of coupling in the Gaussian nodes of the finite-element mesh. This is a modified version of subroutine HQPOT [16] which calculates potential matrix elements of coupling in the \( z = z_{\text{min}} \).

• Subroutine SCSOLC calculates the reflection and transmission matrices and corresponding wave functions, and writes them into file EVWFN, if necessary.

Note, that the function of the other subroutines has been described in [16, 17].

### 3.2 Description of the auxiliary SLAS program

The calling sequence for the subroutine SLAS is:

```fortran
CALL SLAS(XS,SHIFT,THRSHL,IDIM,NDIM,NOPEN,NMAX,KMAX,
1 MAXZ,ZPREGR,ZDREGR,FXSAS,ZAS0)
```

where arguments have the following type and meaning:

**Input data**

<table>
<thead>
<tr>
<th>Variable</th>
<th>Type</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>XS</td>
<td>REAL*8</td>
<td>value of ( z_{\text{min}} ) or ( z_{\text{max}} ).</td>
</tr>
<tr>
<td>SHIFT</td>
<td>REAL*8</td>
<td>the given double energy spectrum.</td>
</tr>
<tr>
<td>THRSHL</td>
<td>REAL*8</td>
<td>array THRSHL of dimension NDIM containing values of the thresholds ( \epsilon_i^{(0)} ).</td>
</tr>
<tr>
<td>IDIM</td>
<td>INTEGER</td>
<td>dimension of the envelope space.</td>
</tr>
<tr>
<td>NDIM</td>
<td>INTEGER</td>
<td>number of coupled differential equations.</td>
</tr>
<tr>
<td>NOPEN</td>
<td>INTEGER</td>
<td>number of open channels.</td>
</tr>
<tr>
<td>NMAX</td>
<td>INTEGER</td>
<td>number of required linear independent solutions and always ( NMAX \leq NDIM ).</td>
</tr>
<tr>
<td>KMAX</td>
<td>INTEGER</td>
<td>maximal order of asymptotic expansions of matrix elements ( V(z) ) and ( Q(z) ) at large (</td>
</tr>
<tr>
<td>MAXZ</td>
<td>INTEGER</td>
<td>value of ( k'_{\text{max}} ) in etalon equation (38).</td>
</tr>
<tr>
<td>ZPREGR</td>
<td>COMPLEX*16</td>
<td>array ZPREGR of dimension NDIM × NMAX. In input ZPREGR(I,I) contains value of regular/irregular solution of I-th etalon equation (38), while on output ZPREGR contains asymptotic regular/irregular matrix-solution of Eq. (35).</td>
</tr>
<tr>
<td>ZDREGR</td>
<td>COMPLEX*16</td>
<td>array ZDREGR of dimension NDIM × NMAX. On input ZDREGR(I,I) contains first derivative of regular/irregular solution of I-th etalon equation (38), while on output ZDREGR contains first derivative of asymptotic regular/irregular matrix-solution of Eq. (35).</td>
</tr>
</tbody>
</table>
FXSAS LOGICAL*8 if FXSAS = .TRUE. then calculates only recommended value of boundary points. In othercase calculates ZPREGR and ZDREGR.

• ZAS0 is the name of the external user-supplied subroutine for evaluating coefficients $Z_i^{(k)}$ of the etalon equation (38), and should be written as follows:

```fortran
SUBROUTINE ZAS0(ZAS,MAXZ,SHIFT,THRSHL,NDIM,NOPEN,ABSB)
C . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . .
C ..
C . P R O G R A M
C . TO CALCULATE THE COEFFICIENTS OF ETALON
C . EQUATION
C .
C .
C . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . .
IMPLICIT REAL*8 (A-H,O-Z)
COMPLEX*16 ZAS(NDIM,MAXZ)
RETURN
END
```

• VQAS0 is the name of the new user-supplied subroutine for the evaluate the coefficients $V_{ij}^{(l)}$ ($2 \leq l \leq k_{\text{max}}$) and $Q_{ij}^{(l)}$ ($1 \leq l \leq k_{\text{max}}$) of the asymptotic expansion (36) of matrix element $V(z)$ and $Q(z)$ at large $|z|$, and should be written as follows:

```fortran
SUBROUTINE VQAS0(VAS,QAS,KMAX,NDIM,ABSB)
C . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . .
C .
C . P R O G R A M
C . TO CALCULATE THE COEFFICIENTS OF THE ASYMPTOTIC
C . MATRIX ELEMENTS $V(z)$ and $Q(z)$ WITH ORDER KMAX
C . AT LARGE $|z|$.
C .
C .
C . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . .
IMPLICIT REAL*8 (A-H,O-Z)
DIMENSION VAS(NDIM,NDIM,2:KMAX),QAS(NDIM,NDIM,KMAX)
RETURN
END
```

Here as followed from (44), on output ZAS(I,1) should be equal $\epsilon_I^{(1)}$, and on input the parameter ABSB = sign(XS).

Subroutine SLAS in program KANTBP, is called as needed via subroutines ASYMSL and ASYMSR.
4. Benchmark calculation of penetration coefficient

Wave function $\tilde{\Psi}(\tilde{x}, \tilde{y})$ of two particles (or ions) labeled by $i = 1, 2$ connected with oscillator potential penetration through repulsive (Coulomb) barriers $\tilde{U}(\tilde{x}_i)$ in the center-mass-system satisfies the two-dimensional Schrödinger equation [8]:

$$
\left( -\frac{\hbar^2}{2M} \frac{\partial^2}{\partial \tilde{y}^2} - \frac{\hbar^2}{2\mu} \frac{\partial^2}{\partial \tilde{x}^2} + \frac{\mu}{2} \tilde{\omega}^2 \tilde{x}^2 + \tilde{U}_1(\tilde{x}_1) + \tilde{U}_2(\tilde{x}_2) - \tilde{E} \right) \tilde{\Psi}(\tilde{x}, \tilde{y}) = 0,
$$

(48)

where $\tilde{\omega}$ is the oscillator frequency, $\tilde{E}$ is the energy, $\tilde{x}_1 = \tilde{y} + s_1 \tilde{x}$, $\tilde{x}_2 = \tilde{y} - s_3 \tilde{x}$ are variables in the laboratory system of coordinates. The parameters $s_1 = \frac{m_1}{M}$, $s_3 = \frac{m_2}{M}$ are defined via masses of particles $m_1$ and $m_2$, and their total $M = m_1 + m_2$ and reduced $\mu = \frac{m_1 m_2}{M}$ masses.

Using the transformation of variables

$$
x = x^{-1}_{osc} \tilde{x}, \quad y = \sqrt{\frac{M}{\mu}} x^{-1}_{osc} \tilde{y},
$$

(49)

with the oscillator units of length $x_{osc} = \sqrt{\frac{\hbar}{\mu}}$, the corresponding Eq. (48) leads to the following dimensionless equation

$$
\left( -\frac{\partial^2}{\partial y^2} - \frac{\partial^2}{\partial x^2} + x^2 + V(x, y) - \mathcal{E} \right) \Psi(x, y) = 0,
$$

(50)

where $\mathcal{E} \equiv 2E = \frac{E}{E_{osc}}$ and $V(x, y)$ are the dimensionless energy and barrier potential in units of energy $E_{osc} = \frac{\hbar \tilde{\omega}^2}{2}$

$$
V(x, y) = U_1(x_1) + U_2(x_2) \equiv \frac{1}{E_{osc}} \left( \tilde{U}_1(\tilde{x}_1) + \tilde{U}_2(\tilde{x}_2) \right).
$$

(51)

where $x_1 = s_2 y + s_1 x$ and $x_2 = s_2 y - s_3 x$ with $s_2 = \sqrt{s_1} M$.

**Model A.** We choose barrier potentials $U_i(x_i)$ with effective charges $\tilde{Z}_i > 0$ in the form of the repulsive truncated Coulomb potential cut off on small $0 < \bar{x}_{min} < 1$ and large $\bar{x}_{max} > 1$ distances from $x_i = 0$ as [7]

$$
U_i(x_i) = \begin{cases} 
\frac{2\tilde{Z}_i}{\bar{x}_{min}} - \frac{2\tilde{Z}_i}{\bar{x}_{max}}, & |x_i| \leq \bar{x}_{min}, \\
\frac{2\tilde{Z}_i}{|x_i|} - \frac{2\tilde{Z}_i}{\bar{x}_{max}}, & \bar{x}_{min} < |x_i| \leq \bar{x}_{max}, \\
0, & |x_i| > \bar{x}_{max}.
\end{cases}
$$

(52)

**Model B.** We define the Coulomb-like potentials $U_i(x_i)$ that depend on the integer parameter $s \geq 2$ and truncation parameter $\bar{x}_{min} > 0$ [8]:

$$
U_i(x_i) = \frac{2\tilde{Z}_i}{\sqrt{|x_i|^s + \bar{x}_{min}^s}}.
$$

(53)
The asymptotic boundary conditions for the solution \( \Psi(y, x) = \{\Psi_i(y, x)\}_{i=1}^{N_o} \) with direction \( v = \rightarrow \) can be written in the obvious form

\[
\Psi_i(y \to -\infty, x) \to B_i^{(0)}(x) \frac{\exp \left( \frac{\imath}{\sqrt{p_{i_o}}} \left( p_{i_o} y - \text{sign}(y) \frac{Z_{i_o}}{p_{i_o}} \ln(2|p_{i_o}|) \right) \right)}{\sqrt{p_{i_o}}} \\
+ \sum_{j=1}^{N_o} B_j^{(0)}(x) \frac{\exp \left( -\frac{\imath}{\sqrt{p_j}} \left( p_j y - \text{sign}(y) \frac{Z_{j_o}}{p_j} \ln(2|p_j|) \right) \right)}{\sqrt{p_j}} R_{ji_o}.
\]

\[
\Psi_i(y \to +\infty, x) \to \sum_{j=1}^{N_o} B_j^{(0)}(x) \frac{\exp \left( \frac{\imath}{\sqrt{p_j}} \left( p_j y - \text{sign}(y) \frac{Z_{j_o}}{p_j} \ln(2|p_j|) \right) \right)}{\sqrt{p_j}} T_{ji_o},
\]

\[
\Psi_i(y, x \to \pm \infty) \to 0.
\]

Here \( N_o \) is the number of open channels at fixed energy \( 2E = p^2 + \varepsilon_{i_o}^{(0)} > 0; Z_{12} = 0 \) for model A and \( Z_{12} = (\hat{Z}_1 + \hat{Z}_2)/s_2 \) for model B; \( R_{ji_o} \) and \( T_{ji_o} \) are unknown reflection and transmission amplitudes; \( B_j^{(0)}(x) \) are the basis functions of oscillator corresponding to energy \( \varepsilon_j^{(0)} = 2j - 1 \) at \( j \geq 1 \)

\[
\left( -\frac{\partial^2}{\partial x^2} + x^2 - \varepsilon_j^{(0)} \right) B_j^{(0)}(x) = 0, \quad \int_{-\infty}^{+\infty} B_i^{(0)}(x) B_j^{(0)}(x) dx = \delta_{ij}.
\]

### 4.1 Kantorovich expansion

We construct a desired solution of the boundary-value problem (BVP) \((50), (54)\) in the form of Kantorovich expansion:

\[
\Psi'(x, y) = \sum_{j=1}^{N} B_j(x; y) \chi_{ji}(y).
\]

The basis functions \( B_j(x; y) \) in the fast variable \( x \) and the potential curves \( \varepsilon_j(y) \) that depend continuously on slow variable \( y \) as a parameter are chosen as solutions of the BVPs for the equation on grid \( \Omega_x \{x_{\min}(y), x_{\max}(y)\} \)

\[
\left( -\frac{d^2}{dx^2} + x^2 + V(x, y) - \varepsilon_j(y) \right) B_j(x; y) = 0,
\]

which are subject to the boundary, normalization and orthogonality conditions

\[
B_j(x_{\min}(y); y) = B_j(x_{\max}(y); y) = 0,
\]

\[
\int_{x_{\min}(y)}^{x_{\max}(y)} B_i(x; y) B_j(x; y) dx = \delta_{ij}.
\]
By substituting (56) into (50), (54) and by taking average over (58), we obtain the BVP for a set of $N$ coupled ODEs that describes the slow subsystem for the partial solutions $\chi^{(j)}(y) = \left(\chi^{(j)}_1(y), \ldots, \chi^{(j)}_N(y)\right)^T$:

\[
\left(-I\frac{d^2}{dy^2} + V(y) + Q(y)\frac{d}{dy} + \frac{dQ(y)}{dy} - 2EI\right)\chi^{(j)}(y) = 0. \tag{59}
\]

Here $I$ is the unit $N \times N$ matrix, $V(y)$ and $Q(y)$ are the effective potential $N \times N$ matrices:

\[
V_{ij}(y) = \varepsilon_j(y)\delta_{ij} + \int_{y_{\min}}^{y_{\max}} \frac{\partial B_i(x; y)}{\partial y} \frac{\partial B_j(x; y)}{\partial y} dx, \tag{60}
\]

\[
Q_{ij}(y) = -\int_{y_{\min}}^{y_{\max}} B_i(x; y) \frac{\partial B_j(x; y)}{\partial y} dx.
\]

The eigenvalue problem (57), (58) was solved by the ODPEVP program [22] for $y \in [y_{\min}, y_{\max}]$ good separated eigenvalues $|\varepsilon_i(y) - \varepsilon_{i-1}(y)| > \epsilon > 0$ where $\epsilon \sim 0.05$ for the double precision arithmetic. This condition is valid for accepted values of parameters of considered models. In the case of non-good separated eigenvalues, i.e. if $0 < |\varepsilon_i(y_*) - \varepsilon_{i-1}(y_*)| \leq \epsilon$, one should generate a more dense grid in vicinity $v_y = |y - y_*| < \epsilon_*$ of avoided crossing points $y_*$ and/or use multi-precision arithmetic. For long-range potentials one should construct appropriate asymptotic expansion for eigenvalues and corresponding eigenfunctions $y \in (-\infty, +\infty) \setminus [y_{\min}, y_{\max}]$ to build up asymptotic effective potentials with leading terms

\[
V_{ij}(y) = \left(\varepsilon_j^{(0)} + \text{sign}(y)\frac{2Z12}{y}\right)\delta_{ij} + \sum_{k=3}^{k_{\max}} \frac{V_{ij}^{(k, \pm)}}{y^k} + O(y^{-k_{\max}-1}),
\]

\[
Q_{ij}(y) = \sum_{k=3}^{k_{\max}} \frac{Q_{ij}^{(k, \pm)}}{y^k} + O(y^{-k_{\max}-1}), \tag{61}
\]

where the sign “+” is for $y > 0$ and “−” is for $y < 0$.

For given number $N$ of Eq. (59), the values $x_{\min}$ and $x_{\max}$ of grid $\Omega_x \{x_{\min}, x_{\max}\}$ were chosen in the region $|x| > x_0 = \sqrt{2N - 1}$, where the Hermite polynomial [24] (or basis function $B_i(x; y)$ in a general case) has no zeros. These values are computed with prescribed precision $eps > 0$ from the condition

\[
\exp\left(-\int_{x_0}^{x} dx \sqrt{x^2 - x_0^2}\right) \leq eps, \tag{62}
\]

which in the given case leads to inequality

\[
\exp\left(-x\sqrt{x^2 - x_0^2}/2\right) \left(x + \sqrt{x^2 - x_0^2}\right)^{x_0^2/2} x_0^{-x_0^2/2} \leq eps. \tag{63}
\]
To find an approximate solution, at the first step we choose initial approximation \( x_{\text{max}} = x_0 \), after which it is increased with step 1 until condition (63) is satisfied. Values \( y_{\text{min}} < x_{\text{min}} \) and \( y_{\text{max}} > x_{\text{max}} \) were chosen from the condition that potential (53) is negligible on the interval \( x_{\text{min}} < x < x_{\text{max}} \).

The matching points \( y_1^{\text{match}} \) and \( y_2^{\text{match}} \) of the numerical (60) and asymptotic (61) effective potential were calculated as follows:

\[
y_1^{\text{match}} = \min \left( y_-, y^V \right), \quad y_2^{\text{match}} = \max \left( y_+, y^V \right),
\]

\[
y^Q = \pm k_{\text{max}} \sqrt{|Q_{NN-1}(k_{\text{max}}, \pm)}| \eps, \quad y^V = \pm k_{\text{max}} \sqrt{|V_{NN}(k_{\text{max}}, \pm)}| \eps,
\]

since \( |Q_{ij}(k_{\text{max}}, \pm)}| < |Q_{NN-1}(k_{\text{max}}, \pm)}|, |V_{ij}(k_{\text{max}}, \pm)}| < |V_{NN}(k_{\text{max}}, \pm)}| \). The values \( y_{\text{min}} \) and \( y_{\text{max}} \) were satisfied by inequalities \( y_{\text{min}} < y_1^{\text{match}} < x_{\text{min}} \) and \( y_{\text{max}} > y_2^{\text{match}} > x_{\text{max}} \) that should be calculated from conditions

\[
y_{\text{min}} = \min \left[ y_1^{\text{match}}, \min_{j,\delta} \frac{|\phi_j^{\text{max}}(-\delta)}{\eps}, \min_{j,\delta} \frac{|\psi_j^{\text{max}}(-\delta)}{\eps} \right],
\]

\[
y_{\text{max}} = \max \left[ y_2^{\text{match}}, \max_{j,\delta} \frac{|\phi_j^{\text{max}}(\delta)}{\eps}, \max_{j,\delta} \frac{|\psi_j^{\text{max}}(\delta)}{\eps} \right].
\]

For the calculation of asymptotic solutions of model B, we have used etalon equation (38) at \( d = 1, k'_{\text{max}} = 1 \) and \( Z_{\nu}^{(1)} = 2 \sign(y)Z_{12} \), which corresponds to known solutions in the open channels

\[
R_{t,0}(p, y) = p^{-1/2} \left\{ \begin{align*}
(G_0(p, +) \pm i F_0(p, +)) \exp(\mp \iota \omega)/2, & \quad y > 0, \\
(G_0(p, -) \mp i F_0(p, -)) \exp(\pm \iota \omega)/2, & \quad y < 0,
\end{align*} \right.
\]

and in the closed channels

\[
R_{t}(q, y) = q t \exp(-t/2) U(1 + Z_{12}/q, 2, t), \quad t = 2q_y |y|.
\]

Here \( F_0(p, y) \) and \( G_0(p, y) \) are the regular and irregular continuum zero order Coulomb functions calculated by subroutine RCWFNN [20] which is a modified version of the subroutine RCWF [23] for the DOUBLE PRECISION accuracy, \( \sigma_t = \exp(\Gamma(1 + Z_{12}/p)) \) is the Coulomb phase shift [24], and \( U(a, b, c) \) is the confluent hypergeometric function of second kind calculated by subroutine CHGU [25]. Note that, for the numerical calculation we have neglected the exponentially small factor \( \exp(-t/2) \) in \( R_{t}(q, y) \) and its first derivative, since this factor is canceled during evaluation of \( R(y) \) matrix in Eq. (18). The coefficients \( V_{ij}^{(k, \pm)} \), \( \phi_j^{(k, \pm)} \) and \( \psi_j^{(k, \pm)} \) have been implemented in MAPLE and FORTRAN up to order \( k_{\text{max}} = 11 \) using an algorithm described in [8] and Section 2.4.
Figure 3: The total probabilities $T \equiv |T|_1^2$ of penetration vs energy $\mathcal{E} = 2E$ through truncated Coulomb (52) (upper panel) and Coulomb-like (53) (lower panel) potential barriers.

Below we have used values of parameters: $m_1 = m_2 = 1$, $\bar{x}_{\text{min}} = 0.1$, $\bar{Z}_1 = \bar{Z}_2 = 0.5$ and $\bar{Z}_1 = \bar{Z}_2 = 1$. Also $\bar{x}_{\text{max}} = 5$ for model $A$, and $s = 8$ for model $B$. In the considered examples we used grids $\Omega_x \{x_{\text{min}}, x_{\text{max}}\} = \{-10(768)10\}$ and $\Omega_y \{y_{\text{min}}, y_{\text{max}}\} = \{-125(200) - 25(100) - 6(200)6(100)25(200)125\}$ with the Lagrange elements of the order $p = 4$ between the nodes. In the above, the number of grid elements for grids $\Omega_x$ and $\Omega_y$ is shown in the parentheses. At the boundary points $y_{\text{min}}$ and $y_{\text{max}}$ the absolute accuracy $\varepsilon_{\text{Wr}}$ of calculated Wronskian (12) was less than $10^{-11}$.

The total probabilities $T \equiv |T|_1^2 = \sum_{j=1}^{N_0} |T_{1j}|^2$ of penetration through truncated Coulomb (52) and Coulomb-like (53) potential barriers of models $A$ and $B$ are shown in Figure 3. These pictures illustrate the important peculiarity that a more realistic nontruncated Coulomb-like barrier, being more wide than truncated one, leads to a set of the probability maximums having a bigger half-width. It can be used for verification of the models of type $A$ and $B$ and quantum transparency effect. Positions of peaks of transmission coefficient demonstrated the quantum transparency effect correspond to the real part of energies of metastable states.
Figure 4: The profiles of the absolute values of the wave function $|\Psi_{\leftarrow}(x, z)|$ of the model B at $m_1 = m_2 = 1$, $\bar{x}_{\text{min}} = 0.1$, $s = 8$ and $Z_1 = Z_2 = 0.5$. Left panel: the resonance transmission at $2E = 8.1403$ and $|T|_{11}^2 = 0.9259$. Right panel: the total reflection at $2E = 9.4748$ and $|T|_{11}^2 = 0.0161$.

imbedded in continuum while the half width corresponds to imaginary part. The profiles of wave functions corresponding to the resonance transmission and total reflection are shown in Figure 4. One can see that in resonance transmission case the probability density have a maximum in a vicinity of barrier due to metastable states imbedded into continuum while in the case of total reflection in a vicinity of barrier it is suppressed.

5. Test desk

We consider the BVP (50) with parameters $m_1 = 1$, $m_2 = 3$, $\bar{x}_{\text{min}} = 0.1$, $Z_1 = Z_2 = 0.1$, $s = 8$ and $N = 4$. The corresponding BVP (1)–(3) and the matrix elements $V(z)$, $Q(z)$ have been solved on grid $\Omega_x \{x_{\text{min}}, x_{\text{max}}\} = \{-x_{\text{min}}(64), x_{\text{max}}\}$ with accuracy $10^{-10}$. Boundary points $x_{\text{max}} = -x_{\text{min}} \approx 8.8$ were defined by the inequality (63). All calculation details of this problem were written into file ODPEVP.LPR.

The following values of numerical parameters and characters have been used in the test run via the supplied input file SQRTBT.INP:

```plaintext
&PARAS TITLE=' REFLECTION AND TRANSMISSION MATRICES ',
   IPTYPE=1, ISC=3, IDIM=1, NPOL=4,
   SHIFT= 4D0, IPRINT=1, IPRSTP=120,
   RMESH=-25D0, 100D0, -6D0, 100D0, 6D0, 100D0, 25D0,
   NDIR=1, NDIL=4, NMDIL=1, THRSHL= 1.D0, 3D0, 5D0, 7D0, IBOUND=8,
```
Description of the FORTRAN program KANTBP 3.0

FNOUT='KANTBP.LPR',IOUT=7,POTEN='ODPEVP.PTN',IOUP=10,
FMAIR='KANTBP.MAT',IOUM=11,EVWFN='KANTBP.WFN',IOUF=0

&END

TEST RUN OUTPUT

PROBLEM: REFLECTION AND TRANSMISSION MATRICES

********

CONTROL INFORMATION

-------------------------------------------------------
NUMBER OF DIFFERENTIAL EQUATIONS . . . . (MDIM ) = 4
NUMBER OF FINITE ELEMENTS . . . . . . . . (NELEM ) = 300
NUMBER OF GRID POINTS . . . . . . . . . . . (NGRID ) = 1201
ORDER OF SHAPE FUNCTIONS . . . . . . . . (NPOL ) = 4
ORDER OF GAUSS-LEGENDRE QUADRATURE . . . (NGQ ) = 5
DIMENSION OF ENVELOPE SPACE . . . . . . . (IDIM ) = 1
BOUNDARY CONDITION CODE . . . . . . . . (IBOUND) = 8
DOUBLE ENERGY SPECTRUM . . . . . . . . . (SHIFT ) = 4.00000

SUBDIVISION OF RHO-REGION ON THE FINITE-ELEMENT GROUPS:

*******************************************************************************
NO OF BEGIN OF LENGTH OF GRID END OF
GROUP ELEMENTS INTERVAL ELEMENT STEP INTERVAL
----- --------- -------- --------- -------- --------
1  100  -25.000  0.19000  0.04750  -6.000
2  100  -6.000  0.12000  0.03000  6.000
3  100   6.000  0.19000  0.04750  25.000
********************************************************************************

NDIM, MDIM= 4 4

TOTAL SYSTEM DATA

---------------------------------------------------------------
TOTAL NUMBER OF ALGEBRAIC EQUATIONS . . . . (NN ) = 4804
TOTAL NUMBER OF MATRIX ELEMENTS . . . . . . (NWK) = 60010
MAXIMUM HALF BANDWIDTH . . . . . . . . . . (MK ) = 20
MEAN HALF BANDWIDTH . . . . . . . . . . (MMK) = 12

*******************************************************************************

CALCULATION OF WAVE FUNCTION WITH DIRECTION <---

NUMBER OF OPEN CHANNELS . . . . . . . . (NOPEN) = 2
VALUE OF I-TH MOMENTUM . . . . . . . . . (I,QR ) = 1 0.1732E+01
VALUE OF I-TH MOMENTUM . . . . . . . . . (I,QR ) = 2 0.1100E+01
IM PART OF WRONSKIANS

-2.00000  -1.79667E-08
-1.79667E-08  -2.00000

RE PART OF RR MATRIX

-1.94759  -5.91176E-03
-5.91176E-03  -4.85403E-01

IM PART OF RR MATRIX

-1.24681  0.172716
0.172716  0.931470

RE PART OF TT MATRIX

0.600469  -3.17926E-01
-3.17926E-01  -2.76469

IM PART OF TT MATRIX
Description of the FORTRAN program KANTBP 3.0

| RR_<-|^2 + |TT_<-|^2 | 1.00000 | 0.353866E-09 |
| M A X I M A L A B S O L U T E E R R O R =0.457613E-09 |

| RR_>|^2 + |TT_>|^2 | 1.00000 | -.353867E-09 |
| M A X I M A L A B S O L U T E E R R O R =0.457614E-09 |

| RE P A R T: TT_>^1 * RR_<- + RR_>^1 * TT_<- | 0.179243E-09 | 0.475463E-09 |
| IM P A R T: TT_>^1 * RR_<- + RR_>^1 * TT_<- | -0.475400E-09 | 0.169614E-09 |
| M A X I M A L A B S O L U T E E R R O R =0.515210E-09 |

| RR_<-^T - RR_<- | 0.00000 | -.128480E-09 |
| M A X I M A L A B S O L U T E E R R O R =0.489364E-09 |

| RR_>^T - RR_> | 0.00000 | 0.128480E-09 |
| M A X I M A L A B S O L U T E E R R O R =0.489364E-09 |

| TT_>^T - TT_<- | 0.258571E-12 | -.318056E-11 |
| M A X I M A L A B S O L U T E E R R O R =0.511799E-09 |

********************************************************************************
References


Description of the FORTRAN program KANTBP 3.0


