Wiener–Hopf matrix factorization using ordinary differential equations in the commutative case

A.V. Shanin

Abstract

A matrix factorization problem is considered. The matrix to be factorized is algebraic, has dimension $2 \times 2$ and belongs to Moiseev’s class. A new method of factorization is proposed. First, the matrix factorization problem is reduced to a Riemann–Hilbert problem using the Hurd’s method. Secondly, the Riemann–Hilbert problem is embedded into a family of Riemann–Hilbert problems indexed by a variable $b$ taking values on a half–line. A linear ordinary differential equation (ODE1) with respect to $b$ is derived. The coefficient of this equation remains unknown at this step. Finally, the coefficient of the ODE1 is computed. For this, it is proven that it obeys a non-linear ordinary differential equation (ODE2) on a half–line. Thus, the numerical procedure of matrix factorization becomes reduced to two runs of solving of ordinary differential equations on a half–line: first ODE2 for the coefficient of ODE1, and then ODE1 for the unknown function. The efficiency of the new method is demonstrated on some examples.

1 Introduction

Many diffraction problems can be transformed into matrix factorization problems \[\tag{1}\] . Typically, these diffraction problems are 2D problems with different boundaries occupying positive and negative parts of the $x$-axis. There emerges a known matrix $G$ analytical in a thin strip going along the real axis of a complex variable $k$, and it is necessary to represent it as a product

$$G(k) = U^{-1}(k)W(k),$$

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where $W$ and $U$ are matrices analytical and having no zeros of the determinant in the upper and lower half-plane, respectively. Also, both matrices should have algebraic growth in corresponding half-planes.

In the scalar case, which can be considered as a degenerate case of $1 \times 1$ matrix, the solution can be readily achieved by taking the logarithm of the matrix and performing the additive decomposition by means of Cauchy’s integral and Sokhotsky’s formula [1]. Returning to the matrices of order $N > 1$, this approach can be generalized for Moiseev’s matrices [2] having form

$$G = \sum_{n=0}^{N-1} g_n(k) A^n(k),$$

(2)

where $g_n(k)$ are scalar function, and $A$ is a polynomial matrix. In the simplest case of matrix $A$ having distinct eigenvalues almost everywhere, $A$ can be decomposed as

$$A = T D T^{-1},$$

(3)

where $T$ is the matrix of the eigenvectors and $D$ is a diagonal matrix composed of the eigenvalues. Both $T$ and $D$ are algebraic matrices, and one can introduce the Riemann surface $\mathcal{R}$ on which $T$ and $D$ are single-valued. Further, the matrix factorization problem becomes reduced to a scalar Riemann–Hilbert problem on $\mathcal{R}$. This problem can be solved in terms of Abelian integrals with the help of Jacobi’s inversion problem [3]. So, the solution of the problem of factorization of (2) is known at least formally, and it possibly can be used for practical needs. Some examples can be found e.g. in [4]. Simpler, but more popular cases [5, 6] can be described as particular cases of (2). Khrapkov’s method [5] is rather simple and leads to straightforward computations, but for a broad class of matrices it produces non-algebraic growth at infinity. Moiseev’s method can be considered as a remedy enabling one to avoid this growth. Another technique to avoid the non-algebraic growth has been proposed in [7]. This technique also includes some numerical stages. A review of the commutative factorization and a development of ideas of [6] can be found in [8].

If $G$ cannot be represented as (2) then some numerical [9] or approximate (e.g. [10]) methods can be applied.

In the current work we consider matrices of Moiseev’s class (2) and develop a new technique which is arguably simpler in practical realization than the Moiseev–Zverovich or Daniele procedure. The new technique can be
applied only when the \( g_n(k) \) in (2) are algebraic functions. This is an important restriction, however in much of the practical situation this restriction is fulfilled. The new procedure comprises three steps. Firstly, the matrix factorization problem is reduced to a Riemann–Hilbert problem using the Hurd’s method [11]. Namely, instead of studying the factors \( W \) and \( U \) we are studying only the factor \( U \) continued into the upper half–plane of \( k \). The \( k \)–plane is cut along half–lines connecting the branch points of \( G \) located in the upper half–plane, namely the points \( k_j \), with \( k = +i\infty \). A Hilbert problem is formulated on the half–lines \((k_j, k_j + i\infty)\). As Hurd mentioned, the new problem can be simpler than the initial matrix factorization problem.

At the second step the Riemann–Hilbert problem is embedded into a family of Riemann–Hilbert problems indexed by a variable \( b \) taking values on a half–line. Namely, for the whole family the coefficients \( H_j(k) \) remain the same, but the contours on which the functional equations should be fulfilled are changed from \((k_j, k_j + i\infty)\) to \((k_j + b, k_j + i\infty)\), where \( b \) is an imaginary number taking values from 0 to \(+i\infty\). Thus, we can define the family of solutions \( U(b, k) \). The solution of the initial problem is denoted by \( U(0, k) \).

A linear ordinary differential equation (ODE1) with respect to \( b \) is derived for \( U(b, k) \). The coefficient of this equation remains unknown on this step.

Finally, step the coefficient of the ODE1 is computed. For this, it is proved that it obeys a non-linear ordinary differential equation (ODE2) on a half–line. Thus, the numerical procedure of matrix factorization becomes reduced to solving two ordinary differential equations on a half–line: first ODE2 for the coefficient of ODE1, and then ODE1 for the unknown function.

Some numerical results are presented. Namely, we demonstrate that the new procedure applied to a matrix belonging to the Khrapkov’s class is factorized exactly the same way as by the traditional Khrapkov’s procedure. Moreover, we apply our method to the matrix emerging in [4].

2 Problem formulation and Hurd’s procedure

Let \( G(k) \) be an algebraic matrix \( N \times N \) having no singularities and no zeros of determinant on the real axis and tending to the unit matrix \( I \) of dimension \( N \times N \) as \(|k| \to \infty \). Our aim is to find the decomposition (1) valid in some strip \(|\text{Im}[k]| < \epsilon \) with \( U \) having no singularities or zeros of the determinant in the lower half–plane and on the real axis, and \( W \) having no singularities or zeros of the determinant in the upper half–plane, maybe except several
points, where poles or zeros are allowed. We demand that the unknown functions $U$ and $W$ tend to $I$ as $|k| \to \infty$. Some restrictions on $G$ will be imposed below.

Apply Hurd’s procedure [11] as follows. Let $k_j, j = 1, \ldots, p$ be branch points of matrix $G$ in the upper half-plane. Connect the points $k_j$ with $i\infty$ by the cuts

$$\Gamma_j = (k_j, k_j + i\infty)$$

parallel to the imaginary axis. Let contours $\Gamma_j$ do not pass through other branch points, poles or zeros of the determinant of $G$. Continue function $U(k)$ into the upper half-plane cut along the lines $\Gamma_j$ by the relation

$$U(k) \equiv W(k)G^{-1}(k).$$  \hfill (4)

Note that $W$ is defined and regular in the upper half-plane and $G$ is defined in the upper half-plane with the cuts $\Gamma_j$. Define by $U(k^+)$ and $U(k^-)$ for $k \in \Gamma_j$ the values of $U$ taken on the right and on the left shore of $\Gamma_j$, respectively (see Fig. 1). Similarly, define the values $G(k^+)$ and $G(k^-)$. For some $k \in \Gamma_j$

$$U(k^+) \equiv W(k)G^{-1}(k^+),$$  

$$U(k^-) \equiv W(k)G^{-1}(k^-)$$

(note that $W(k^-) = W(k^+) = W(k)$). Then,

$$U(k^+) = U(k^-)H_j(k),$$ \hfill (5)

$$H_j(k) \equiv G(k^-)G^{-1}(k^+), \quad k \in \Gamma_j$$ \hfill (6)

The set of equations (6) taken for $j = 1, \ldots, p$ constitute the Riemann–Hilbert problem in Hurd’d formulation. It was Hurd’s observation that this problem can be simpler than the initial matrix factorization problem.

In our case each coefficient $H_j(k)$ of the Riemann–Hilbert problem can be continued analytically into some half-strip $\Omega + k_j$

$$\Omega = \{\text{Re}[k] < \epsilon, \text{Im}[k] > 0\}.$$  

Note that the point $k_j$ does not belong to $\Omega + k_j$. Also $H_j(k)$ are algebraic functions, so each of them can be continued onto some Riemann surface.

Let all $H_j(k)$ tend to $I$ $\text{Im}[k] \to \infty$ (this restriction is fulfilled if $G(k) \to I$ on all sheets). We forget about $W(k)$ and look for $U(k)$ on the complex
plane $k$ cut along the contours $\Gamma_j$ having no singularities and no zeros of the determinant on the cut plane, obeying the problem (5). We recall that $U(k) \to I$ as $|k| \to \infty$.

The behavior of $U(k)$ at the points $k_j$ will be specified below in such a way that the problem possesses a unique solution.

We assume also that matrices $H_j$ have distinct eigenvalues almost everywhere.

All restrictions described above correspond to a quite general matrix factorization problem and are easy to fulfil by, e.g. slight change of the contour position. If matrices $H_j$ do not tend to $I$ the method can be easily modified also. Here we are going to pose the most strong restriction: we assume that all branches of all matrices $H_j(k)$ taken for arbitrary affix $k$ commute with each other, i.e. for each $k$

$$H_{j_1}(k)H_{j_2}(k) = H_{j_2}(k)H_{j_1}(k),$$

(7)

where $H_{j_1}(k)$ and $H_{j_2}(k)$ are any possible continuations of $H_{j_1}$ and $H_{j_2}$ to $k$. The meaning of this restriction is discussed in the next section.

3 Functional–commutative and branch–commutative matrices

All existing analytical approaches to matrix factorization are available only for matrices admitting a commutative factorization, i.e. a representation of the form

$$G(k) = U^{-1}(k)W(k) = W(k)U^{-1}(k).$$

(8)
The theory of commutative matrix factorization starts from [12] where a concept of functional–commutative matrix has been introduced. A functional–commutative matrix is a matrix commuting with its singular integral. The property of functional–commutativity is not easy to check for an arbitrary matrix. That is why, in [13] we introduced branch–commutative matrices. Namely, an algebraic matrix $G(k)$ is called branch-commutative if for any $k$ the values $G_j(k)$ corresponding to different branches of $G$ commute with each other.

To formulate the main result of [13] we need one more definition. A Riemann surface of an algebraic matrix is called balanced if each sheet of it can be reached from any fixed sheet only by bypassing the branch points located in the upper half–plane and only bypassing the branch points lying in the negative half–plane. Most of the known matrices arising in practical problems have balanced Riemann surfaces.

The main result of [13] is as follows. If an algebraic matrix $G$ with balanced Riemann surface admits commutative factorization then it is branch–commutative. Vice versa, a branch–commutative matrix can be represented in the form (2), and thus the Moiseev’s method can be applied to it.

Note that if matrix $G$ is branch–commutative then the property of (7) for the matrices $H_j$ defined by (6) is valid. Thus, the matrices to which the method described here can be applied are (with some unimportant restrictions) the same as the matrices, to which the Moiseev’s method is applicable.

Let us formulate one important consequence of the property (7).

**Proposition 1** If property (7) is fulfilled then there exists rational matrix $B(k)$ commuting with all matrices $H_j(k)$.

The proof of the proposition is as follows. Represent $H_1$ in the form

$$H_1(k) = P(k)F_1(k)P^{-1}(k),$$

(9)

where $P(k)$ is the matrix composed of the eigenvectors of $H_1$ normalized, say, by making the first component of each vector equal to 1. Respectively, $F_1(k)$ is a diagonal matrix composed of scalar functions $f_1(k), \ldots, f_N(k)$.

It is known that if two matrices commute then normalized eigenvectors of the matrices coincide [14]. Matrix $P$ is algebraic, so it is single–valued on some Riemann surface. Since the values of $H_j(k)$ taken on different sheets (with the same $k$) commute, we can conclude that when a branch point of $P$ is bypassed the columns of $P$ are just permuted.
Construct matrix $B$ in the form

$$B(k) = P(k)D(k)P^{-1}(k), \quad (10)$$

where $D(k)$ is a diagonal matrix with the scalar functions $h_1(k), \ldots, h_N(k)$ on the diagonal. Let the functions $h_m$ be branches of some algebraic function $h$ having the same branch points as $P$. Moreover, let the values $h_m$ be permuted the same way as the columns of $P$ when the branch points are bypassed. Then the function $B(k)$ is single–valued, and therefore rational.

A proper choice of the functions $h_m$ is as follows:

$$h_m = \sum_{n=1}^{N} \beta_n(k)P_{n,m}(k), \quad (11)$$

where $\beta_n(k)$ is an arbitrary set of scalar rational functions (provided none of $h_m$ is identically zero). Values of $P_{n,m}$ are elements of matrix $P$.

Since $B$ commutes with $H_1$ and all matrices $H_j$ have distinct eigenvalues almost everywhere, matrix $B$ commutes with every $H_j$. Note that all other matrices $H_j$ can be represented as

$$H_j(k) = P(k)F_j(k)P^{-1}(k). \quad (12)$$

Due to arbitrariness of the choice of $\beta_n(k)$ one can make matrix $B$ having simple poles and tending to $I$ as $|k| \to \infty$. The matrix $B$ possessing all these properties plays an important role below.

## 4 A family of Riemann–Hilbert problems and derivation of ODE1

### 4.1 Family of Riemann–Hilbert problems

We have reduced the matrix factorization problem to finding the function $U(k)$ obeying equations (5) on the cuts $\Gamma_j$. To solve this problem we use the idea described in detail in [15]. Namely, we are fixing the functions $H_j(k)$, defined and continuous on contours $\Gamma_j$ (and regular in the strips $\Omega + k_j$), and introduce truncated contours

$$\Gamma_j(b) = (k_j + b, k_j + i\infty),$$

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where $b$ is an imaginary number $b \in (0, i\infty)$. Consider a family of problems for the function $U(b, k)$ set by the relations

$$U(b, k^+) = U(b, k^-)H_j(k), \quad k \in \Gamma_j(b).$$

(13)

We assume that for each $b$ the matrix function $U(b, k)$ is single-valued, continuous, and free of zeros of determinant on the plane of $k$ cut along the contours $\Gamma_j(b)$. It tends to $I$ as $|k| \to \infty$. We assume also that the behavior of $U(b, k)$ at the points $k_j + b$ is derived by continuity from the conditions formulated for large $\text{Im}[b]$ (see below). Obviously,

$$U(k) = U(0, k).$$

(14)

The main idea of the method is to study the behavior of $U(b, k)$ as a function of $b$.

Embedding of $U(k)$ into a family $U(b, k)$ enables us to define behavior of $U(k)$ at $k = k_j$ in the most natural way. Expand matrices $H_j$ in the form (12). The leading term of $U(b, k)$ near the point $k = k_j + b$ can be written in the form

$$U(b, k) \approx K_j(k_j + b)(k - (k_j + b))^{\log(F_j(k_j + b))/\nu}K^{-1}_j(k_j + b)$$

(15)

where $F_j$ are taken from (9), and $K_j$ are some non-singular matrices. The branch of the logarithm should be fixed as follows. For $b \to i\infty$ the matrices $F_j(k_j + b)$ tend to $I$. For these values we choose the branch of logarithm close to zero matrix. Then, for other values of $b$ we choose the branch of logarithm by continuity. Such a choice enables us to avoid discussing partial indices of the initial Riemann–Hilbert problem.

### 4.2 Form of ODE1 for a single cut

Let the number of cuts $p$ be equal to 1, i.e. let there exists only one cut $\Gamma_j$. This corresponds to a function $G(k)$ having a single branch point in the positive half-plane. This case has been studied in [15]. Here we formulate the main theorem of [15] with a short proof.

For this, it is necessary to introduce a notation of the ordered exponential (the term comes from quantum mechanics). Namely, let $\gamma$ be a contour (directed one) connecting the points $\tau_1$ and $\tau_2$ ($\tau_1$ is a starting point), and let $C(\tau)$ be a $N \times N$ matrix defined on $\gamma$. Consider a matrix equation

$$\frac{d}{d\tau}X(\tau) = C(\tau)X(\tau)$$

(16)
taken with the initial condition $X(\tau_1) = I$. Solve this equation along contour $\gamma$ and define the value $X(\tau_2)$. By definition,

$$\text{OE}_\gamma[C(\tau)d\tau] \equiv X(\tau_2).$$

This notation is just a convenient way to refer to a solution of an ordinary differential equation.

**Theorem 1 a)** There exists $N \times N$ matrix $s_1(b)$ analytical in the strip $\Omega$, such that $U(b,k)$ obeys an ordinary differential equation (ODE1)

$$\frac{\partial}{\partial b} U(b,k) = \frac{s_1(b)}{k - (k_1 + b)} U(b,k).$$

The initial condition for this equation is as follows:

$$U(i\infty,k) = I.$$  

**b)** Let there exist a $N \times N$ matrix $s_1(\tau)$ analytic in $\Omega$ and such that

$$\text{OE}_\gamma \left[ \frac{s_1(\tau)}{k - (\tau + k_1)} d\tau \right] = H_1(k)$$

for $k \in (k_1, k_1 + i\infty)$. Contour $\gamma$ is a concatenation of $\gamma^+$ and $\gamma^-$ (see Fig. 2). Then solution $U(b,k)$ is given by the formula

$$U(b,k) = \text{OE}_{\gamma_b} \left[ \frac{s_1(\tau)}{k - (\tau + k_1)} d\tau \right]$$

where contour $\gamma_b$ goes from $i\infty$ to $b$ along $\Gamma_1$.

![Figure 2: Contours $\gamma^+$ and $\gamma^-$](image-url)
Let us outline the proof. Consider part a). Consider the function

\[ S(b, k) = \frac{\partial U(b, k)}{\partial b} U^{-1}(b, k). \]  

(22)

This function is analytic in the plane cut along \((k_j + b, k_j + i\infty)\). Consider its behavior at the cut. Note that the coefficient \(H_1\) does not depend on \(b\). Thus,

\[ \frac{\partial U(b, k^+)}{\partial b} = \frac{\partial U(b, k^-)}{\partial b} H_1(k), \quad k \in \Gamma_1 + b. \]

Using this relation with (13) we conclude that \(S(b, k^+) = S(b, k^-)\), and therefore the function is single–valued. According to the condition at infinity for \(U\), \(S\) should decay as \(|k| \to \infty\). The only singularity of \(S\) in the finite part of the \(k\)–plane is \(k = k_1 + b\). The leading term of the singularity is given by (15). According to this, \(S\) has a simple pole at \(k = k_1 + b\), and

\[ S(b, k) = -\frac{1}{2\pi i(k - (k_1 + b))} K_1(k_1 + b) \log(F_1(k_1 + b)) K_1^{-1}(k_1 + b). \]  

(23)

Finally,

\[ s_1(b) = -\frac{1}{2\pi i} K_1(k_1 + b) \log(F_1(k_1 + b)) K_1^{-1}(k_1 + b). \]  

(24)

The choice of branch of the logarithm has been discussed in the previous subsection.

Analyticity of the coefficient \(s_1\) in \(\Omega + k_1\) follows from the fact that the cut on which functional equation (13) is set can be deformed (without changing its starting point) arbitrarily within \(\Omega + k_1\), and the solution remains the same while the contour changes.

The initial condition (19) follows from general properties of the Riemann–Hilbert problem [9].

Consider part b) of the theorem. Let us show that (21) is a solution of the problem (13). Note that due to analyticity of \(s_1\) the contour \(\gamma\) in (20) can be deformed provided that it does not cross the singularity \(\tau = k - k_1\) of the coefficient. Let be \(k \in \Gamma_1(b)\), and thus \(\text{Im}[k] > \text{Im}[k_1 + b]\). Deform contour \(\gamma\) into \(\gamma + b\):

\[ \text{OE}_{\gamma+b} \left[ \frac{s_1(\tau)}{k - (\tau + k_1)} d\tau \right] = H_1(k) \]  

(25)
According to general properties of the ordinary differential equations and the ordered exponential notations \cite{20},
\[
\text{OE}_{\gamma+b} \left[ \frac{s_1(\tau)}{k-(\tau+k_1)} d\tau \right] =
\]
\[
\left( \text{OE}_{b+\gamma} \left[ \frac{s_1(\tau)}{k-(\tau+k_1)} d\tau \right] \right)^{-1} \text{OE}_{b+\gamma} \left[ \frac{s_1(\tau)}{k-(\tau+k_1)} d\tau \right].
\] (26)

Note that according to (21)
\[
\text{OE}_{b+\gamma} \left[ \frac{s_1(\tau)}{k-(\tau+k_1)} d\tau \right] = U(b, k^+),
\]
\[
\text{OE}_{b+\gamma} \left[ \frac{s_1(\tau)}{k-(\tau+k_1)} d\tau \right] = U(b, k^-),
\]

Thus, (26) is equivalent to (13).

4.3 Form of ODE1 for several cuts

If there are \( p > 1 \) branch points \( k_j \) (and, thus, several cuts \( \Gamma_j \)) Theorem 1 can be modified, while the reasoning remains basically the same. Coefficient \( S \) from (22) can be proven to be single-valued and decaying, but it should have \( p \) simple poles \( k = k_j + b \). Therefore equation (18) has form
\[
\frac{\partial}{\partial b} U(b, k) = \left( \sum_{j=1}^{p} \frac{s_j(b)}{k_j^+ + b} \right) U(b, k).
\] (27)

with \( p \) unknown matrices \( s_j(b) \) analytical in \( \Omega \). The initial conditions are the same as for one cut (i.e. (19)). The form of the solution follows from (27) and (19):
\[
U(b, k) = \text{OE}_{\gamma_b} \left[ \sum_{j=1}^{p} \frac{s_j(\tau)}{k_j^+ - (\tau + k_j^+)} d\tau \right].
\] (28)

A generalization of (24) has form
\[
s_j(b) = -\frac{1}{2\pi i} K_j(k_j^+ + b) \log[F_j(k_j^+ + b)] K_j^{-1}(k_j^+ + b).
\] (29)
for some unknown matrices $K_j$ and known (up to transmutations) diagonal matrices $F_j$.

Finally, condition (20) should be rewritten as

$$
\text{OE}_\gamma \left[ \sum_{j=1}^p \frac{s_j(\tau)}{k - (\tau + \lambda_j)} d\tau \right] = H_m(k), \quad k \in \Gamma_m + b, \quad m = 1 \ldots p. \quad (30)
$$

5 ODE2

5.1 Derivation of ODE2

Formula (21) (or (28)) cannot be used immediately to find solution $U(k) = U(0,k)$ since matrix functions $s_j(b)$ are unknown. Thus, before finding $U$ one should find $s_j$ somehow. In [15] it has been proposed to use equation (20) to find $s_1$. A numerical procedure has been proposed and tested. Application of this procedure does not require branch–commutativeness, so the method is potentially applicable to a much wider class of problem than Moiseev's class. However, this procedure is rather sophisticated and it does not reveal the mathematical nature of the solution. Here we are proposing another technique reducing the determination of $s_j$ to solving a (nonlinear) ordinary differential equation. Unfortunately, the new technique is applicable only to Riemann–Hilbert problems obeying relations (7).

The key idea of the new method is to use matrix $B(k)$ defined by (10). Namely, we construct a rational matrix $B(k)$ commuting with $H_j(k)$ on all their sheets, behaving as $B(k) \to I$ as $|k| \to \infty$ and having only simple poles. Obviously, such matrix can be constructed by using the arbitrariness of the rational functions $\beta_n(k)$. Let the poles of $B(k)$ be located at the points $k = \rho_l$, $l = 1 \ldots d$.

Consider function

$$
V(b,k) = U(b,k)B(k).\quad (31)
$$

Note that

$$
\frac{\partial V(b,k)}{\partial b} V^{-1}(b,k) = \frac{\partial U(b,k)}{\partial b} U^{-1}(b,k) \equiv S(b,k). \quad (32)
$$

Thus, $V$ obeys ODE1 (27) with the same coefficient as $U$.

The key property of function $V$ is expressed by the following proposition.
Proposition 2  There exists function \( R(b, k) \), which is rational as a function of \( k \) for each \( b \), such that

\[
V(b, k) = R(b, k) U(b, k).
\] (33)

Construct function \( R \) as follows:

\[
R(b, k) = V(b, k) U^{-1}(b, k) = U(b, k) B(k) U^{-1}(b, k)
\] (34)

Consider the behavior of \( V(b, k) \) on the cuts \( \Gamma_j(b) \). Since \( B \) commutes with all \( H_j \),

\[
R(b, k^+) = U(b, k^-) H_j(k) B(k) H_j^{-1}(k) U^{-1}(b, k^-) = U(b, k^-) B(k) U^{-1}(b, k^-) = R(b, k^-), \quad k \in \Gamma_j(k).
\]

Thus, for each \( b \) function \( R(b, k) \) is a single–valued function of \( k \). At infinity \( R(b, k) \to I \). Obviously, \( R \) can only have simple poles at \( k = \rho_l \). Due to Liouville’s theorem, \( R(b, k) \) should be a rational function of \( k \). Moreover, one can conclude that \( R \) has form

\[
R(b, k) = I + \sum_{l=1}^{d} \frac{r_l(b)}{k - \rho_l},
\] (35)

where \( r_l(b) \) are some \( N \times N \) matrix functions of \( b \) defined in \( \Omega \).

Construct the coefficient of ODE1 for \( V \) using representation (33):

\[
\frac{\partial V}{\partial b} V^{-1} = R \frac{\partial U}{\partial b} U^{-1} R^{-1} + \frac{\partial R}{\partial b} R^{-1} = R S R^{-1} + \frac{\partial R}{\partial b} R^{-1}.
\] (36)

Comparing (36) with (32), conclude that

\[
\frac{\partial R(b, k)}{\partial b} = S R - R S \equiv [S, R].
\] (37)

Equation (37) is the global form of ODE2. One can easily see that this equation describes the evolution of \( R \) but from the first glance it is not clear how it can describe the evolution of \( S \). However, we possess some additional information about \( R \) and \( S \) (namely, both functions are rational with respect to \( k \)). This information is enough to transform (37) into a local form, which is a closed set of ordinary differential equations describing the evolution of \( R \) and \( S \).
Substitute (37) and
\[ S(b, k) = \sum_{j=1}^{p} \frac{s_j(b)}{k - (k_j + b)} \] (38)
into (37). Expand the right-hand side and left-hand side of (37) as a sum of simple fractions. Taking into account that
\[ \frac{1}{k - (k_j + b)} - \frac{1}{k - \rho_l} = \frac{1}{(k_j + b) - \rho_l} \left( \frac{1}{k - (k_j + b)} - \frac{1}{k - \rho_l} \right) \]
and considering the terms with each denominator separately, obtain equations
\[ \frac{d r_l(b)}{db} = \sum_{j=1}^{p} \frac{[s_j(b), r_l(b)]}{\rho_l - (k_j + b)}, \quad l = 1 \ldots d \] (39)
and
\[ \sum_{l=1}^{d} \frac{[s_j(b), r_l(b)]}{\rho_l - (k_j + b)} = 0, \quad j = 1 \ldots p \] (40)

System (39), (40) does not form a closed system of ordinary differential equations for finding the unknown matrices \( s_j(b), r_j(b) \). To make the system closed, consider (40) together with (29). Formulate the problem of finding matrices \( s_j \) provided that matrices \( r_l \) are known. Equations (29) provide information about the eigenvalues of \( s_j \), while (40) provide information about the eigenvectors of \( s_j \). Namely, the eigenvalues of \( s_j(b) \) are equal to the diagonal elements of
\[ \tilde{F}_j(b) = -\frac{1}{2\pi i} \log(F_j(k_j + b)). \]
where \( F_j(k_j + b) \) is a (known) diagonal matrix composed of the eigenvalues of \( H_j(k_j + b) \) (see (12)). According to (40), the eigenvectors of \( s_j(b) \) coincide with the eigenvectors of the matrix
\[ R(b, k_j + b) = \sum_{l=1}^{d} \frac{r_l(b)}{(k_j + b) - \rho_l}. \]

Define function \( F(X, Y) \) producing a matrix, whose eigenvalues coincide with the eigenvalues of \( X \), and the eigenvectors coincide with the eigenvectors of \( Y \) (provided all eigenvalues of \( Y \) are distinct). The function \( F \) is
defined ambiguously since the mapping between the eigenvalues of $X$ and eigenvectors of $Y$ is not defined. I.e. $\mathcal{F}$ is defined up to a permutation of order $N$. If this ambiguity is eliminated in a correct way,

$$s_j = \mathcal{F}(\tilde{F}_j(b), R(b, k_j + b)). \quad (41)$$

Equations (39) together with (41) form a closed system of equations for finding $r_l$ and $s_j$. This system is non-linear. The system (39), (41) will be called the ODE2 (in the local form). Derivation of the ODE2 is the main result of this paper. This result can be formulated in the form of the following theorem.

**Theorem 2** Let there be a family of Riemann–Hilbert problems (5) obeying the restrictions posed above, including the commutativity restrictions (7). The ODE1 for this family has the notation of (27). Then there exist such matrices $r_l(b)$ and such a choice of the function $\mathcal{F}$ that matrices $s_j(b)$, $r_l(b)$ obey the system (39), (41).

### 5.2 Initial conditions for ODE2 and choice of function $\mathcal{F}$

To make a numerical solution of ODE2 possible one should define the initial conditions and eliminate the ambiguity of defining the function $\mathcal{F}$. Since $U(b, k) \to I$ as $b \to i\infty$, one can conclude that

$$R(i\infty, k) = B(k), \quad (42)$$

where of course

$$R(i\infty, k) \equiv \lim_{b \to i\infty} R(b, k).$$

Thus, if

$$B(k) = I + \sum_{l=1}^{d} \frac{t_l}{b - \rho_l} \quad (43)$$

for some matrices $t_l$ (which are assumed to be known) then

$$r_l(i\infty) = t_l. \quad (44)$$

These relations play the role of initial conditions for the ODE2.
To eliminate the ambiguity of definition of function $F$, we need to establish a correspondence between the eigenvectors of the matrix $R(b, k_j + b)$ and the diagonal element of the (diagonal) matrix $\tilde{F}_j(b)$. Again, consider large values of $\text{Im}[b]$. For large imaginary $b$ the values of the coefficients $H_j(k)$ approximately commute with the common factor $B(b)$. Therefore, the solution $U(b, k)$ near the points $k = k_j + b$ approximately commutes with $B(b)$ or (which is the same in asymptotic sense) with $B(k_j + b)$. Thus, according to (34),

$$R(b, k_j + b) \approx B(k_j + b).$$

Using this relation, one can establish a natural correspondence between the eigenvectors of $B(k_j + b)$ and $R(b, k_j + b)$. Then, the eigenvectors of $B(k_j + b)$ are by construction the eigenvectors of $H_j(k_j + b)$. Thus, it is possible to establish a natural correspondence between the eigenvectors of $B(k_j + b)$ and $H_j(k_j + b)$. Finally, this gives correspondence between the diagonal elements of $F_j(k_j + b)$ and the eigenvectors of $R(b, k_j + b)$.

Thus, function $F$ can be defined without ambiguity for large $\text{Im}[b]$ and for other $b$ it can be defined by continuity.

5.3 Invariance of ODE2 with respect to the choice of $B(k)$

The choice of the factor $B(k)$ is not unique. Namely, if $B(k)$ obeys all restrictions then a combination

$$B'(k) = \sum_{m=0}^{N-1} g_m(k) B^m(k) \quad (45)$$

with rational scalar functions $g_m(k)$ also can be used as $B$, provided that $B'(k) \to I$ as $|k| \to \infty$ and $B'$ has only simple poles. The form of ODE2 changes when $B$ is substituted by $B'$. Let us show that this substitution does not change the solution $s_j(b)$. For this we remind that the system (39), (40) is equivalent to (37). The invariance of $s_j$ is established by the following proposition.

Proposition 3 Let $R$ be defined by (34), and

$$R'(b, k) = U(b, k) B'(k) U^{-1}(b, k), \quad (46)$$

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where $B'$ is defined by (45). Let (37) be valid for some matrix $S$. Then

$$\frac{\partial R'(b, k)}{\partial b} = [S, R'].$$ (47)

First, note that it follows from (46) that

$$R'(k) = \sum_{m=0}^{N-1} g_m(k) R^m(k)$$ (48)

Due to formal linearity of (47), it is sufficient to prove that

$$\frac{\partial R^m(b, k)}{\partial b} = [S, R^m].$$ (49)

This can be easily proved by induction.

6 Examples

6.1 Description of the numerical procedure

The numerical procedure straightforwardly follows from Theorem 1 and Theorem 2. Assume that matrices $H_j(k)$ are known explicitly, and let the matrix $B(k)$ be constructed and represented in the form (43).

First, ODE2 is solved along the positive imaginary axis of $b$ from $i\infty$ to 0. In practice, ODE2 is solved not from $b = i\infty$, but from $b = iL$, where $L$ is a large number playing the role of infinity. At the “infinite” point $b = iL$ initial condition for ODE2 are set in the form of

$$r_l(iL) = t_l.$$ (50)

At the point $b = iL$ function $F$ is constructed without ambiguity as follows. According to the argument above and according to (50), for numerical solution

$$R(iL, k_j + iL) = B(k_j + iL).$$ (51)

Represent $B(k_j + iL)$ in the form (10), i.e.

$$B(k_j + iL) = P_* D_* P_*^{-1},$$ (52)
where $D_*$ is a diagonal matrix. Function $H_j(k_j + iL)$ can be represented in the form (12), i.e.

$$P_*^{-1}H_j(k_j + ib)P_*$$

should be a diagonal matrix. According to (41),

$$s_j(iL) = -\frac{1}{2\pi i}P_* \log(P_*^{-1}H_j(k_j + ib)P_*)P_*^{-1},$$

(53)

where the branch of logarithm close to zero is taken. This procedure defines $s_j(iL)$ in a unique way.

Then ODE2, i.e. the system (39), (41) is solved numerically, say by Runge–Kutta method, from $b = iL$ to $b = 0$. On each step function $F$ is chosen such that new values $s_j(b - i\delta)$ are close to old values $s_j(b)$, i.e. such that $s_j(b)$ are continuous. As the result of this procedure, the matrices $s_j$ are found at points covering the segment $(iL, 0)$ densely enough.

Next, ODE1 is solved to find $U(b, k)$. A set of points $k = z_n$ at which the solution $U(k)$ will be found is selected. The initial conditions have form

$$U(iL, z_n) = I.$$  

(54)

Equation (28) is solved from $b = iL$ to $b = 0$ for the values $U(b, z_n)$ along the imaginary axis of $b$ say by Runge–Kutta method. As the result, the solution $U(k) = U(0, k)$ becomes known at the points $k = z_n$.

One can see that the numerical procedure is rather simple. It consists of two solutions of ordinary differential equations. If the segment $(iL, 0)$ is split into $N_b$ steps, and if there are $N_k$ points in the set $z_n$, then the first step takes $\sim N_b$ operations, and the second step takes $\sim N_k N_p$ operations. This makes difference with results of [15] where the first step takes $\sim N_p^2$ operations.

### 6.2 Khrapkov’s case

It is important to show that the proposed technique is equivalent to the known method in the simplest commutative case, namely in Khrapkov’s case [5]. Consider as an example a family of Riemann–Hilbert problems set on $\Gamma_1(b) = (k_1 + b, k_1 + i\infty)$ with the coefficient

$$H_1(k) = g_0(k)I + g_1(k)\Lambda(k),$$

(55)
where

$$\Lambda(k) = \begin{pmatrix} 1 & k \\ k & -1 \end{pmatrix},$$

$g_0$ and $g_1$ are some algebraic functions such that $g_0(k) \to 1$ as $|k| \to \infty$, $g_1$ tends to zero as $|k| \to \infty$ not slower than $1/|k|^2$.

A traditional solution of this problem is as follows. First, a solution $\tilde{U}(b,k)$ is constructed by the formula \cite{5}

$$\tilde{U}(b,k) = \exp(\bar{\xi}) \left( \cosh \left( \sqrt{\phi(k)}\bar{\eta} \right) I + \sinh \left( \sqrt{\phi(k)}\bar{\eta} \right) \frac{\Lambda(k)}{\sqrt{\phi(k)}} \right),$$

(56)

$$\phi(k) = k^2 + 1,$$

$$\bar{\xi}(b,k) = - \int_{k_1+b}^{k_1+i\infty} \frac{\xi(\tau)}{k - \tau} d\tau,$$

$$\bar{\eta}(b,k) = - \int_{k_1+b}^{k_1+i\infty} \frac{\eta(\tau)}{k - \tau} d\tau,$$

(57)

$$\xi(k) = - \frac{1}{4\pi i} \log \left( g_0^2(k) - \phi(k)g_1^2(k) \right),$$

$$\eta(k) = - \frac{1}{4\pi i \sqrt{\phi(k)}} \log \left( \frac{g_0(k) + g_1(k)\sqrt{\phi(k)}}{g_0(k) - g_1(k)\sqrt{\phi(k)}} \right),$$

(58)

(59)

Solution $\tilde{U}$ obeys all conditions except the condition $U \to I$ at infinity. Instead, for a fixed $b$

$$\tilde{U}(b,k) \to \cosh(\zeta(b))I + \sinh(\zeta(b)) \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} \equiv Q(b) \quad \text{as } |k| \to \infty;$$

(60)

$$\zeta(b) = - \int_{k_1+b}^{k_1+i\infty} \eta(\tau) d\tau.$$

(61)

Thus, one has to "correct" the behavior of $\tilde{U}$ by a left multiplication:

$$U(b,k) = Q^{-1}(b)\tilde{U}(b,k).$$

(62)

Let us consider the same problem from the point of view of the proposed method. One can check directly \cite{15} that the auxiliary solution obeys ODE1 in a slightly modified form:

$$\frac{\partial U(b,k)}{\partial b} = \tilde{S}(b,k)\tilde{U}(b,k), \quad \tilde{S}(b,k) = \frac{\xi(b+k_1)}{k-(b+k_1)}I + \frac{\eta(b+k_1)}{k-(b+k_1)}\Lambda(k).$$

(63)
Similarly, it can be checked that $Q(b)$ obeys equation

$$\frac{dQ}{db} = \eta(b + k_1) \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} Q(b). \tag{64}$$

Construct ODE1 for $U$. According to (62) the coefficient of this equation is equal to

$$S(b, k) = \frac{\partial \tilde{U}}{\partial b} = Q^{-1} \left( \tilde{S} - \frac{dQ}{db} Q^{-1} \right) Q = Q^{-1} \left( \tilde{S} - \eta(b + k_1) \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} \right) Q. \tag{65}$$

One can see that the coefficient has form of (18), i.e. for a fixed $b$ it is a rational function of $k$ having a simple pole at $k = b + k_1$ and decaying at infinity.

Now consider ODE2. Select a function $B(k)$ commuting with both branches of $H_1(k)$, having only simple poles and tending to $I$ at infinity. For example one can choose

$$B(k) = I + \frac{1}{k^2 - 1} \Lambda(k) \tag{66}$$

with simple poles at $k = \pm 1$. Define $\tilde{R} = \tilde{U}B\tilde{U}^{-1}$. One can see that $B$ commutes with $\tilde{U}$, and thus $\tilde{R}(b, k) = B(k)$. Note that

$$[\tilde{R}, \tilde{S}] = 0, \quad \frac{\partial \tilde{R}}{\partial b} = 0. \tag{67}$$

Define $R$ as (34). It can be expressed as

$$R(b, k) = Q^{-1}(b)\tilde{R}(b, k)Q(b) = Q^{-1}(b)B(k)Q(b) \tag{68}$$

Taking into account (67) and (65) it is easy to show that equation (37), i.e. ODE2 in the global form is valid for Khrapkov’s matrix.

Let us write down ODE2 in the local form (for demonstration purposes). Represent $\Lambda(k)$ in the form

$$\Lambda(k) = \sqrt{k^2 + 1} P(k) \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix} P^{-1}(k), \tag{69}$$

$$P(k) = \begin{pmatrix} \frac{1}{\sqrt{k^2 + 1}} & \frac{1}{\sqrt{k^2 + 1}} \\ \frac{1}{k} & -\frac{1}{k} \end{pmatrix}. \tag{70}$$
Similarly,

\[ H_1(k) = P(k)F_1(k)P^{-1}(k), \]

\[ F_1(k) = \begin{pmatrix} g_0(k) + \sqrt{k^2 + 1} g_1(k) & 0 \\ 0 & g_0(k) - \sqrt{k^2 + 1} g_1(k) \end{pmatrix} \] (71)

Since \( R \) has two poles (\( \rho_1 = 1, \rho_2 = -1 \)), we need a system of equations describing evolution of three matrices: \( r_1(b), r_2(b), \) and \( s_1(b) \). According to (39), first two equations have form

\[
\frac{dr_1(b)}{db} = \left[ s_1(b), r_1(b) \right]_{1 - (k_1 + b)}, \quad \frac{dr_2(b)}{db} = \left[ s_1(b), r_2(b) \right]_{1 - (k_1 + b)},
\] (72)

The third equation has form of (41):

\[
s_1(b) = -\frac{1}{2\pi i} F \left( \log(F_1(b + k_1)), \frac{r_1(b)}{k_1 + b - 1} + \frac{r_2(b)}{k_1 + b + 1} \right). \] (73)

Initial conditions for (72) should be taken in the form (44). For this, matrix \( B \) should be represented as a sum of simple fractions. As the result, we get

\[
r_1(i\infty) = \frac{1}{2} \begin{pmatrix} 1 & 1 \\ 1 & -1 \end{pmatrix}, \quad r_2(i\infty) = \frac{1}{2} \begin{pmatrix} -1 & 1 \\ 1 & 1 \end{pmatrix} \] (74)

Function \( F(X, Y) \) is implemented as follows. Let \( X \) be a diagonal matrix. Matrix \( Y \) is represented in the form \( Y = Y_1 Y_2 Y_1^{-1} \) numerically or analytically (\( Y_2 \) should be a diagonal matrix). The result is formed as

\[
F(X, Y) = Y_1 X Y_1^{-1}, \] (75)

or

\[
F(X, Y) = Y_1 X' Y_2^{-1} \] (76)

where \( X' \) is a matrix, whose diagonal elements are interchanged. The choice between these two forms is made by the following rule. For the point \( b = iL \) where conditions (74) are set matrix \( Y \) has two eigenvectors, one of which is close to

\[
a \begin{pmatrix} 1 \\ 1 \end{pmatrix},
\]

and another one is close to

\[
a \begin{pmatrix} 1 \\ -1 \end{pmatrix}
\]
These eigenvectors are columns of $Y$. If the first column corresponds to the vector of the first type, then form (75) is chosen at this point. Otherwise, form (76) should be chosen. At each new step function $\mathcal{F}$ is chosen to be approximately continuous.

### 6.3 Factorization of Antipov’s matrix

Here we consider a more sophisticated (but also commutative) case previously addressed in [4]. Matrix $G(k)$ is as follows:

$$G(k) = g_0(k)I + g_1(k)\Lambda(k),$$

where

$$\Lambda(k) = \begin{pmatrix} k^4 - \mu^4 & \alpha\mu^4/\tau \\ \alpha\mu^4/\tau & -k^4 + \mu^4 \end{pmatrix},$$ (78)

$$g_0(k) = \frac{(\psi(k) - \tau)(k^4 - \mu^4) - \alpha\mu^4}{\psi(k)(k^4 - \mu^4)}, \quad g_1(k) = \frac{\tau}{\psi(k)(k^4 - \mu^4)},$$ (79)

$$\psi(k) = \sqrt{k^2 - (1 + 0i)^2}$$ (80)

$\mu, \tau, \alpha$ are some scalar constant physical parameters. Notation (80) means that the only branch point in the upper half-plane is $k_1 = 1$.

Matrix (77) is related to a problem of scattering by a screen composed of a rigid half-plane and a flexible perforated sandwich half-plane. The boundary conditions for this problem were derived in [16]. The problem was reduced to the Wiener–Hopf problem in [4].

The problem belongs to the Khrapkov’s class. The most important function for such problem is

$$\phi(k) = \Lambda_{12}\Lambda_{21} - \Lambda_{11}\Lambda_{22},$$

having the property

$$\Lambda^2(k) = \phi(k)I.$$ (81)

In this case

$$\phi(k) = k^8 - 2\mu^4k^4 + \mu^8(1 + \alpha^2/\tau^2).$$ (82)

This function is a polynomial of degree 8. If a direct Khrapkov’s method [5] is applied then a solution grows rapidly (faster than algebraically) at infinity. Therefore, the Moiseev’s method should be applied. The method has been outlined in [4], however no numerical results have been presented. An

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(see (70)).
application of this method requires finding zeros of Riemann’s theta function and Weierstrass’ kernel quadratures.

We apply the method developed above to this problem. The following values of parameters are taken for computations:
\[ \mu = 2, \quad \tau = 0.25, \quad \alpha = 0.3. \]

Apply the Hurd’s method. There a single cut \( \Gamma_1 \) in the upper half-plane going from \( k_1 = 1 \) to \( 1 + i \infty \). Denote by \( g_0^+(k), g_1^+(k) \) the values of \( g_0(k), g_1(k) \) on the right shore of the cut, and by \( g_0^-(k), g_1^-(k) \) the values on the left shore of the cut. Note that these values are different due to the presence of the square root \( \psi \). The coefficient \( H_1(k) \) describing the multiplicative jump on \( \Gamma_1 \) is equal to
\[
H_1(k) = G(k^-)G^{-1}(k^+) = \frac{(g_0^- g_0^+ - \phi g_1^- g_1^+)I + (g_1^- g_0^+ - g_1^- g_0^+) \Lambda}{(g_0^+)^2 - (g_1^+)^2 \phi} \tag{83}
\]

Then, the function \( B(k) \) is chosen. We can take it in the form
\[
B(k) = I + \frac{1}{\xi(k)} \Lambda(k), \tag{84}
\]
where \( \xi(k) \) is a rational function. Since \( \lambda(k) \) grows as \( k^4 \), we can take \( \xi(k) \) as a polynomial of 5th order, namely
\[
\xi(k) = \prod_{l=1}^{5} (k - \rho_l). \tag{85}
\]

The choice of \( \rho_l \) can be done quite arbitrarily. We use the values
\[ \rho_1 = 2 + i, \quad \rho_2 = 2 - i, \quad \rho_3 = -i, \quad \rho_4 = -1 + i, \quad \rho_5 = -1 - i. \]

The scheme outlined above is implemented. The set of the points of interest \( k = z_n \) belong to the real segment \( k \in (-1, 1) \) (see Fig. 3). To determine them, it is necessary to find the values \( s_1(b) \) for \( b \in \Gamma_1 \). These values are found by solving ODE2. The result (i.e. the components of the matrix \( U(0, z_n) \)) is shown in Fig. 3.

Besides finding the values \( U(0, z_n) \) we perform a simple control of the whole procedure. For this, we find the values \( U(0, (z'_n)^+) \) and \( U(0, (z'_n)^-) \), where the points \( z'_n \) belong to the cut \( \Gamma_1 \) (see Fig. 3), the values \( U(0, (z'_n)^+) \)
represent the right shore of the cut, and the values $U(0, (z'_n)^-)$ represent the left shore of the cut. To determine the values on the shores we change the contour for solving ODE2 slightly. Namely, for the values $U(0, (z'_n)^+)$ we chose contour $\gamma^+$ in Fig. 3 and for values $U(0, (z'_n)^-)$ we chose contour $\gamma^-$. After that, we compute the combination $U^{-1}(0, (z'_n)^-)U(0, (z'_n)^+)M^{-1}(z'_n)$. In the ideal case this matrix should be equal to $I$, therefore its deviation from $I$ can be taken as a measure of relative accuracy of the computation. It has been found that the relative accuracy of the computation used to be of order $10^{-4}$.

![Figure 4: Solution $U(k)$ on the segment $k \in (-1, 1)$](image-url)
7 Conclusion

A new method for matrix factorization in the commutative (Moiseev’s) case is developed. The method is numerical, but it is based on two analytical properties of the factorization problem. It is applicable to algebraic matrices having the property of branch-commutativity, i.e. the matrices, whose values corresponding to different sheet over the same affix commute.

The factorization problem is transformed by Hurd’s procedure into a Riemann–Hilbert problem on a set of cuts. Then the Riemann–Hilbert problem is embedded into a family of Riemann–Hilbert problems indexed by a variable \( b \). The solution as a function of \( b \) is described by ordinary differential equation (ODE1) with an unknown coefficient \( S \). This coefficient is found by solving another ordinary differential equation, ODE2. Initial conditions for ODE1 and ODE2 are formulated. It is shown that the proposed procedure in the Khrapkov’s case is equivalent to the standard solution. Moreover, it is shown that the new procedure is applicable to Antipov’s matrix, and it does not lead to Jacobi’s inversion problem, which is not easy to implement.

In more general (non-commutative) cases ODE2 should be replaced by an OE-equation described in [15].

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