Renormalization group procedure for effective particles: 
elementary example of exact solution with 
finite mass corrections and no involvement of vacuum *†

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Renormalization group procedure for effective particles in the front form of Hamiltonian dynamics 
is applied to an elementary quantum field theory for two species of particles mixed through a mass-
like interaction term. The model interaction generates only finite terms and the procedure yields 
a whole family of equivalent effective theories. The exact solution for the family is found without 
involvement of the vacuum state in the dynamics. Physical spectrum is obtained at the end of 
the procedure in the form of free particles with definite masses. Since the procedure is designed 
general terms, it could be used for the purpose of constructing effective dynamics also in other 
theories than the elementary model.

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I. INTRODUCTION

Renormalization group procedure for effective particles (RGPEP) [1] in the front form (FF) of Hamiltonian dy-
namics [2] is designed for application in solving relativistic quantum field theories. In realistic theories, where non-
perturbative solutions of the RGPEP equations are hard to find, the equations can be initially solved only order-by-
order in a perturbative expansion [3] or after making other drastic simplifications of unknown accuracy. This article 
describes instead an application of the RGPEP to an elementary but instructive model of a quantum field theory in 
four dimensions which is soluble exactly. The model exhibits a non-perturbative scale-evolution of a mass matrix for 
effective particles and shows how the RGPEP can deal with the quantum vacuum problem [4]. The final result of 
the RGPEP in the model is a free theory of the particles whose masses appear in the exact eigenvalues of the initial 
Hamiltonian.

Since the RGPEP involves basic elements of the canonical field quantization [5, 6] and renormalization of Hamiltoni-
ans by techniques other than integrating out high-energy degrees of freedom [7, 8], the elementary model application 
is described including all details needed to make the presentation self-contained. Besides the FF of Hamiltonian 
dynamics, the description often refers to the commonly used form of dynamics, designated the instant form (IF) by 
Dirac [2].

Section II briefly introduces the RGPEP. The elementary example is defined in Section III. Section IV describes 
solution of the RGPEP equations. The vacuum problem is discussed in Section V. Section VI concludes the article. 
Appendix A describes derivation of the same physical solution but obtained using an alternative RGPEP generator 
to the one used in the main text.

II. SUMMARY OF RGPEP

The concept of effective particles as degrees of freedom in a relativistic quantum field theory is introduced through 
a transformation [2]

\[
\psi_s = U_s \psi_0 U_s^\dagger .
\]

(1)

\(\psi_s\) is a quantum field operator built from creation and annihilation operators for effective particles of size \(s\). These 
creation and annihilation operators are commonly denoted by \(q_s\). The operator \(\psi_0\) is the field operator built from 
the particle operators \(q_0\) that correspond to the bare, point-like particles, and \(s = 0\). By definition, all kinematical 
quantum numbers that label operators \(q\) on both sides of Eq. II, such as a three-momentum, charge, spin, isospin, 
flavor, color, and the like, are not altered by \(U_s\).

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The intuitive interpretation of parameter $s$ in terms of a size of the effective particles in the elementary model will be explained later. It is based on the general RGPEP feature that effective interactions contain the form factors that limit how far off energy shell the interactions can extend. The corresponding energy-width of the form factors is determined by $1/s$ (see below). The value $s = 0$ corresponds to absence of form factors. For a finite $s$, the effective Hamiltonian is band-diagonal on the energy scale and the band width is $\sim 1/s$. The principle of using the band-diagonal structure for the purpose of renormalization is formulated in [9]. It is convenient to use the parameter $t = s^4$ and label operators with $t$ rather than $s$ itself.

A canonical Hamiltonian density is built from products of fields $\psi_0$ and their derivatives. A corresponding Hamiltonian is obtained by integrating the density over a space-time hyper-surface. The result is a polynomial $\mathcal{H}_0(q_0)$ with coefficients $c_0$. If a term in $\mathcal{H}_0(q_0)$ contains a product of $n$ operators $q_0$, the coefficient has $n$ arguments. Each argument is a set of quantum numbers carried by a corresponding particle. Similarly, an effective-particle Hamiltonian $\mathcal{H}_t(q_t)$ is defined through its coefficients $c_t$.

The RGPEP employs the equality

$$\mathcal{H}_t(q_t) = \mathcal{H}_0(q_0),$$

which means that the same dynamics is expressed in terms of different operators. The change of $q_0$ to $q_t$ is accompanied with the change of coefficients $c_0$ to $c_t$ so that the physics is not changed. For example, the expansion of eigenstates of the Hamiltonian into the $t$-dependent Fock components involves the wave functions that depend on $t$, but the states as a whole do not depend on $t$ at all.

Variation of the coefficients $c_t$ with $t$ is described by the equation obtained by differentiating both sides of

$$\mathcal{H}_t(q_0) = \mathcal{U}_t^\dagger \mathcal{H}_0(q_0) \mathcal{U}_t,$$

with respect to $t$. One obtains

$$\mathcal{H}_t'(q_0) = [\mathcal{G}_t(q_0), \mathcal{H}_t(q_0)],$$

where $\mathcal{G}_t = -\mathcal{U}_t^\dagger \mathcal{U}_t$ is called a generator. Correspondingly,

$$\mathcal{U}_t = T \exp \left( - \int_0^t d\tau \mathcal{G}_\tau \right),$$

where $T$ orders operators from left to right in the order from a smallest to largest $t$.

In the RGPEP, the generator is defined by

$$\mathcal{G}_t = [\mathcal{H}_f, \mathcal{H}_{P_t}].$$

The operator $\mathcal{H}_f$, called the free Hamiltonian, is the part of $\mathcal{H}_0(q_0)$ that does not depend on the coupling constants,

$$\mathcal{H}_f = \sum_i p_i^- q_{0i}^\dagger q_{0i}.$$

The sum over subscript $i$ extends over all particle species and their quantum numbers, including integration over momenta, and

$$p_i^- = \frac{p_i^+ 2 + m_i^2}{p_i^2}.$$  

This is the FF free-energy of a particle with mass $m_i$ and kinematical momentum components $p_i^+$ and $p_i^-$. The operator $\mathcal{H}_{P_t}$ is defined using the Hamiltonian $\mathcal{H}_t$. Namely, if $\mathcal{H}_t(q_0)$ is of the form

$$\mathcal{H}_t(q_0) = \sum_{n=2}^\infty \sum_{i_1, i_2, \ldots, i_n} c_t(i_1, \ldots, i_n) q_{0i_1}^\dagger \cdots q_{0i_n},$$

where the coefficients $c_t(i_1, \ldots, i_n)$ are to be found using RGPEP, the operator $\mathcal{H}_{P_t}(q_0)$ is defined by

$$\mathcal{H}_{P_t}(q_0) = \sum_{n=2}^\infty \sum_{i_1, i_2, \ldots, i_n} c_t(i_1, \ldots, i_n) \left( \frac{1}{2} \sum_{k=1}^n p_{ik}^+ \right)^2 q_{0i_1}^\dagger \cdots q_{0i_n}. $$
This means that $\mathcal{H}_{P\ell}$ differs from $\mathcal{H}_t$ by multiplication of each and every term by a square of a total $+$ momentum involved in a term. In summary, the coefficients $c_t$ of products of operators $q_i$ in the effective Hamiltonians $\mathcal{H}_t(q_t)$, are solutions of the equation

$$\mathcal{H}'_t = [[\mathcal{H}_f, \mathcal{H}_{P\ell}], \mathcal{H}_t],$$

(11)

where all operators are written as polynomials in $q_0$ and the initial condition is provided by a regulated canonical Hamiltonian with counterterms.

The counterterms are calculated in the RGPEP using a condition that for finite $t$ the coefficients $c_t$ with finite arguments do not depend on the regularization parameters used in the canonical Hamiltonian [9]. The difficulty of satisfying the cutoff-independence condition for $c_t$ originates in the fact that the coefficients appear in the solution for $\mathcal{H}_t(q_t)$ while the counterterms are inserted in the initial condition $\mathcal{H}_0(q_0)$ and in-between there is a solution of the RGPEP that spans the range from 0 to $t$. However, there is no special difficulty associated here with the counterterms because the coefficients $c_t$ with finite arguments do not develop any dependence on regularization in the model and solutions for them are known exactly. Therefore, the adjustment of counterterms in the model only amounts to specifying their finite parts. These parts form the initial mass matrix. The only regularization dependence in the example appears in one overall constant term in $\mathcal{H}_t$, which is a pure number and drop out from Eq. (11).

The generic feature of narrowness of $\mathcal{H}_t$ as $t$ increases can be seen by introducing a projector $R$ on a subspace in the Fock space. Let $\mathcal{H}_R = R \mathcal{H}_t R$. The corresponding projected equation reads (for details, see Appendix C in [1])

$$\mathcal{H}'_R = [[\mathcal{H}_f, \mathcal{H}_{PR}], \mathcal{H}_R].$$

(12)

The free Hamiltonian $\mathcal{H}_f$ commutes with $R$. The matrix version of Eq. (12) resembles the Wegner flow equation introduced in the IP of dynamics for Hamiltonians in condensed matter physics [10–12]. Similar equations are also successfully used in nuclear physics [13–15]. In relativistic quantum field theories, a narrow matrix must be obtained introduced in the IF of dynamics for Hamiltonians in condensed matter physics [10–12]. Similar equations are also introduced in the IF of dynamics for Hamiltonians in condensed matter physics [10–12].

III. MODEL HAMILTONIAN

Let a theory of two real scalar fields $\phi$ and $\chi$ have a classical Lagrangian density

$$\mathcal{L} = \frac{1}{2} [(\partial \phi)^2 - \mu^2 \phi^2] + \frac{1}{2} [(\partial \chi)^2 - \nu^2 \chi^2] - m^2 \phi \chi.$$  

(14)

The last term is called the mass mixing term.

A. Classical Hamiltonian

In terms of the variables $x^\pm = x^0 \pm x^3$ and $x^\perp = (x^1, x^2)$ used to label points in space-time, so that $\partial^\pm = 2\partial/\partial x^\mp$, the Lagrangian density reads

$$\mathcal{L} = \frac{1}{2} [\partial^+ \phi \partial^- \phi - (\partial^\perp \phi)^2 - \mu^2 \phi^2] + \frac{1}{2} [\partial^+ \chi \partial^- \chi - (\partial^\perp \chi)^2 - \nu^2 \chi^2] - m^2 \phi \chi.$$  

(15)

The FF of dynamics involves the four-momentum [16, 17]

$$P^\mu = \frac{1}{2} \int dx^- d^2 x^\perp \mathcal{T}^{\mu \nu}(x),$$  

(16)
where the energy-momentum tensor density component relevant for constructing the model Hamiltonian is
\[ T^{+-}(x) = \partial^+ \phi \partial^- \phi + \partial^+ \chi \partial^- \chi - 2\mathcal{L}. \]  
(17)

Hence,
\[ P^- = \frac{1}{2} \int dx^- d^2 x^\perp \left[ (\partial^\perp \phi)^2 + \mu^2 \phi^2 + (\partial^\perp \chi)^2 + \nu^2 \chi^2 + 2m^2 \phi \chi \right]. \]  
(18)

B. Quantization

Let the fields \( \phi \) and \( \chi \) at \( x^+ = 0 \) have the Fourier decompositions \[ \phi(x^-, x^\perp) = \int [p] a_p e^{-ipx}, \]
(19)
\[ \chi(x^-, x^\perp) = \int [p] b_p e^{-ipx}, \]
(20)
where \([p]\) denotes the measure \( d^+ p d^2 p^\perp / [2|p^+|(2\pi)^3] \) of integration over momentum variables \( p^+ = p^0 + p^3 \) and \( p^\perp = (p^1, p^2) \). In this notation, the integration over momentum variables extends from \(-\infty\) to \(+\infty\) for all three components of \( p \) (a need for a cutoff on the range of \( p \) is still ignored at this point). Quantum theory is obtained by imposing commutation relations
\[ [a_p, a_q] = [b_p, b_q] = 2p^+ (2\pi)^3 \delta^3 (p + q). \]  
(21)

The absence of \( \dagger \) in the commutation relations is intended, since it is the sign of \( p^+ \) that distinguishes the operators that create field quanta from operators that annihilate them. Such kinematical distinction between creation and annihilation processes is not available in the standard, IF approaches. The “annihilation” operators with negative \( p^+ \) correspond to creation operators and one has
\[ a_{-p} = a_p^\dagger, \]  
(22)
\[ b_{-p} = b_p^\dagger. \]  
(23)

Note that these relations involve the change of sign of \( p^\perp \).

After quantization, the classical fields \( \phi \) and \( \chi \) are turned into operators that create and annihilate quanta on the front hyper-plane, \( \hat{\phi} \) and \( \hat{\chi} \). The commutation relations of Eq. (21) correspond to the spatial commutation relations
\[ [\hat{\phi}(x), \partial^+ \hat{\phi}(y)] = [\hat{\chi}(x), \partial^+ \hat{\chi}(y)] = i\delta^3 (x - y). \]  
(24)

The inverse relations are
\[ a_p = |p^+| \int d^3 x e^{+ipx} \hat{\phi}(x), \]
(25)
\[ b_p = |p^+| \int d^3 x e^{+ipx} \hat{\chi}(x), \]
(26)
where \( d^3 x = dx^- d^2 x^\perp \) and integrals extend from \(-\infty\) to \(+\infty\) on the \( x^+ = 0 \) hyper-plane (the behavior of fields in spatial infinity remains unspecified at this point).

C. Quantum Hamiltonian

The quantum Hamiltonian is obtained from Eq. (18) by inserting operator versions of Eqs. (19) and (20) for \( \hat{\phi} \) and \( \hat{\chi} \), respectively, and by normal ordering,
\[ P^- = \frac{1}{2} \int dx^- d^2 x^\perp : \left[ (\partial^\perp \phi)^2 + \mu^2 \phi^2 + (\partial^\perp \chi)^2 + \nu^2 \chi^2 + 2m^2 \phi \chi \right] : . \]  
(27)
The normal ordering is defined using Feynman’s convention 19 with the ordering parameter set equal to \( p^+ \). In this convention, it is understood that operators \( a_p \) are ordered in products according to the value of \( p^+ \) so that the greater \( p^+ \) the further to the right the operator.

All terms in the Hamiltonian are bilinear in fields and all of them contain one and the same integral

\[
\int dx d^2x \int [q \rho] e^{-i q^+ x} = \int [q \rho] 2(2\pi)^2 \delta^3(q + p).
\]

According to Eq. (27),

\[
\mathcal{P}^- = \frac{1}{2} \int [p] \frac{\rho}{|p|^3} = \left[ \left( p^{\perp 2} + \mu^2 \right) a_p a_{-p} + \left( p^{\perp 2} + \nu^2 \right) b_p b_{-p} + 2m^2 a_p b_{-p} \right].
\]

The normal ordering produces the operator that properly counts the FF energy of field quanta,

\[
\mathcal{P}^- = \int [p] \theta(p^+) \left[ \left( \frac{p^{\perp 2} + \mu^2}{p^+} \right) a_p a_p + \left( \frac{p^{\perp 2} + \nu^2}{p^+} \right) b_p b_p + \frac{m^2}{p^+} (a_p b_p + b_p a_p) \right].
\]

The last term describes the mixing of bare particles of type \( a \) associated with field \( \phi \) and of type \( b \) associated with field \( \chi \). From now on, the function \( \theta(p^+) \) is included in the integration measure \([p]\).

The diverging number that is removed by the FF normal ordering in \( \mathcal{P}^- \),

\[
\Omega^- = 2(2\pi)^3 \delta^3(0) \int [p] \left( p^{\perp 2} + \mu^2/2 + \nu^2/2 \right),
\]

involves factors \( V_F = 2(2\pi)^3 \delta^3(0) \) and \( \rho\Omega = \int [p] \left( p^{\perp 2} + \mu^2/2 + \nu^2/2 \right) \). Factor \( V_F \) has an interpretation of a volume of the front that a Hamiltonian density is integrated over. Factor \( \rho\Omega \) is associated with a ground-state energy, cf. 20 [24]. As a number, \( \Omega^- \) does not contribute to the commutators in Eq. (11) and it is not included in the RGEPEP discussion in the next section. However, regarding application of the RGEPEP to more complex theories, one should remember that the vacuum issue is not limited in them to a constant such as \( \Omega^- \), cf. 25 [26].

### IV. SOLUTION OF THE RGPEP EQUATION

According to Sec. I, Eq. (11) has the form,

\[
\mathcal{P}^- = \left[ [\mathcal{P}^-, \mathcal{P}^-], \mathcal{P}^- \right],
\]

and should be solved using Eq. (30) as the initial condition,

\[
\mathcal{P}^-_0(a_0, b_0) = \int [p] \left[ \left( \frac{p^{\perp 2} + \mu^2}{p^+} \right) a_p a_p + \left( \frac{p^{\perp 2} + \nu^2}{p^+} \right) b_p b_p + \frac{m^2}{p^+} (a_p b_p + b_p a_p) \right].
\]

#### A. Equations for coefficients \( c_i \)

On the basis of hindsight, the relevant operators can be written as

\[
\mathcal{P}^-_0(a_0, b_0) = \int [p] \left[ A_{tp} a_{tp}^\dagger a_{tp} + B_{tp} b_{tp}^\dagger b_{tp} + C_{tp} (a_{tp}^\dagger b_{tp} + b_{tp}^\dagger a_{tp}) \right],
\]

\[
\mathcal{P}^-_0(a_0, b_0) = \int [p] \left( A_{0p} a_{0p}^\dagger a_{0p} + B_{0p} b_{0p}^\dagger b_{0p} \right),
\]

\[
\mathcal{P}^-_0(a_0, b_0) = \int [p] \left( p^{\perp 2} A_{tp} a_{tp}^\dagger a_{tp} + B_{tp} b_{tp}^\dagger b_{tp} + C_{tp} (a_{tp}^\dagger b_{tp} + b_{tp}^\dagger a_{tp}) \right),
\]

where the coefficients generically denoted by \( c_i \) in Sec. II read

\[
A_{tp} = \frac{p^{\perp 2} + \mu^2}{p^+},
\]

\[
B_{tp} = \frac{p^{\perp 2} + \nu^2}{p^+},
\]

\[
C_{tp} = \frac{m^2}{p^+}.
\]
The initial conditions for these coefficients, denoted by $c_0$ in Sec. II, are set by fixing the mass-squared parameters at $t = 0$,

$$\mu_0 = \mu,$$  \hspace{1cm} (40)  
$$\nu_0 = \nu,$$  \hspace{1cm} (41)  
$$m_0 = m,$$  \hspace{1cm} (42)

with constants $\mu$, $\nu$, and $m$, taken from $\mathcal{P}_0^-$ in Eq. (33). These parameters include the finite parts of mass-squared counterterms as discussed in Section II.

In this notation, the generator has the form

$$[\mathcal{P}_\tau^+, \mathcal{P}_\tau^-] = \int [p] (A_{0p} - B_{0p}) p^{+2} C_{tp} \left( a_{0p}^\dagger b_{0p} - b_{0p}^\dagger a_{0p} \right).$$  \hspace{1cm} (43)

Eq. (32) reads

$$\mathcal{P}_t^{-,\dagger}(a_0, b_0) = \int [p] \left[ A_{tp} a_{0p}^\dagger a_{0p} + B_{tp} b_{0p}^\dagger b_{0p} + C_{tp} \left( a_{0p}^\dagger b_{0p} + b_{0p}^\dagger a_{0p} \right) \right]$$  \hspace{1cm} (44)

$$= \int [p] (p^{+2}) (A_{0p} - B_{0p}) (A_{tp} - B_{tp}) C_{tp} \left( a_{0p}^\dagger b_{0p} + b_{0p}^\dagger a_{0p} \right)$$  \hspace{1cm} (45)

$$+ \int [p] 2p^{+2} (A_{0p} - B_{0p}) C_{tp}^2 \left( a_{0p}^\dagger a_{0p} - b_{0p}^\dagger b_{0p} \right).$$  \hspace{1cm} (46)

By equating coefficients in front of the same bare particle operators (or evaluating matrix elements between bare one-particle states of types $a$ and $b$), one obtains a set of equations for the coefficients $A_{tp}$, $B_{tp}$, and $C_{tp}$ in $\mathcal{P}_t^{-,\dagger}(a_0, b_0)$. Namely,

$$A_{tp} = 2p^{+2} (A_{0p} - B_{0p}) C_{tp},$$  \hspace{1cm} (47)  
$$B_{tp} = -2p^{+2} (A_{0p} - B_{0p}) C_{tp}^2,$$  \hspace{1cm} (48)  
$$C_{tp} = (p^{+2}) (A_{0p} - B_{0p}) (A_{tp} - B_{tp}) C_{tp}.$$  \hspace{1cm} (49)

This set contains as many triplets of equations as there are different triplets of momentum labels $p$, which a priori is an infinite number when one does not regulate the field expansions into their Fourier components by imposing cutoffs on some discretized set of variables $p^+$ and $p^\perp$. However, it is clear that the modes with different values of $p$ are decoupled. They evolve in $t$ independently of each other. This simplification is a consequence of the bilinear nature of the initial Lagrangian. In addition, new generic simplifications occur thanks to the FF boost invariance of Eq. (11).

**B. Generic simplification due to boost invariance**

In full detail, Eqs. (17), (38), and (49), read

$$\left( \frac{p^{+2} + \mu^2}{p^+} \right)' = 2p^{+2} \left( \frac{p^{+2} + \mu^2}{p^+} - \frac{p^{+2} + \nu^2}{p^+} \right) \left( \frac{m_t^2}{p^+} \right)^2,$$  \hspace{1cm} (50)

$$\left( \frac{p^{+2} + \nu^2}{p^+} \right)' = -2p^{+2} \left( \frac{p^{+2} + \mu^2}{p^+} - \frac{p^{+2} + \nu^2}{p^+} \right) \left( \frac{m_t^2}{p^+} \right)^2,$$  \hspace{1cm} (51)

$$\left( \frac{m_t^2}{p^+} \right)' = (p^{+2}) \left( p^{+2} + \mu^2 - p^{+2} + \nu^2 \right) \left( \frac{p^{+2} + \mu_t^2}{p^+} - \frac{p^{+2} + \nu_t^2}{p^+} \right) \left( \frac{m_t^2}{p^+} \right)^2.$$  \hspace{1cm} (52)

It is visible that the kinematical variables $p^+$ and $p^{+\perp}$ drop out. This feature is special to the FF of dynamics. Thus, the a priori infinite set of different equations for infinitely many coefficients with different kinematical variables $p$, actually reduces to a single set of just 3 equations for 3 mass parameters that are independent of $p$,

$$\left( \mu_t^2 \right)' = 2 \delta \mu^2 \left( m_t^2 \right)^2,$$  \hspace{1cm} (53)  
$$\left( \nu_t^2 \right)' = -2 \delta \mu^2 \left( m_t^2 \right)^2,$$  \hspace{1cm} (54)  
$$\left( m_t^2 \right)' = -\delta \mu^2 \left( \mu_t^2 - \nu_t^2 \right) m_t^2,$$  \hspace{1cm} (55)
where
\[ \delta \mu^2 = \mu^2 - \nu^2. \]  

This set can be written as a differential matrix equation,
\[
\begin{bmatrix}
\mu_1^2 & \mu_2^2 \\
m_1^2 & m_2^2 \\
\nu_1^2 & \nu_2^2
\end{bmatrix}
\begin{bmatrix}
\beta'
\end{bmatrix}
= \begin{bmatrix}
\begin{bmatrix}
\mu_1^2 & 0 \\
0 & \nu_2^2
\end{bmatrix},
\begin{bmatrix}
0 & m_1^2 \\
m_1^2 & 0
\end{bmatrix},
\begin{bmatrix}
\mu_2^2 & m_2^2 \\
m_2^2 & \nu_2^2
\end{bmatrix}
\end{bmatrix}
\begin{bmatrix}
\alpha
\end{bmatrix},
\]

for a 2 × 2 matrix that will be called mass-squared matrix below.

Note that Eq. (57) would be a Wegner-like equation if the first matrix on the right-hand side contained \( \mu_t \) and \( \nu_t \) instead of the initial mass parameters \( \mu \) and \( \nu \). Such change corresponds to inserting \( \mu^2 \) and \( \nu^2 \) in place of \( \mu^2 \) and \( \nu^2 \), respectively, in \( P_f \) of Eq. (35). The resulting Wegner-like equation for the mass-squared matrix can be solved by proceeding in a way analogous to the one described below. This is shown in Appendix A. The explicit solution described in next sections is for constant masses in \( H_f \).

C. Analytic solution for masses

One can introduce a dimensionless variable
\[ u = \delta \mu^4 t, \]
and, denoting differentiation with respect to \( u \) with a prime, one obtains Eqs. (53), (54), and (55), in the form
\[
\begin{align*}
\alpha' &= 2 \gamma^2, \\
\beta' &= -2 \gamma^2, \\
\gamma' &= - (\alpha - \beta) \gamma,
\end{align*}
\]
where the dimensionless functions of \( u \) are
\[
\begin{align*}
\alpha &= \mu_1^2 / \delta \mu^2, \\
\beta &= \nu_1^2 / \delta \mu^2, \\
\gamma &= \mu_2^2 / \delta \mu^2.
\end{align*}
\]

If \( \mu^2 = \nu^2 \), so that \( \delta \mu^2 = 0 \), the mass parameters do not evolve with \( t \) irrespective of the initial value of mass-mixing parameter \( m \). It is assumed from now on that \( \mu^2 > \nu^2 \), so that \( \delta \mu^2 > 0 \).

Regarding the mass degeneracy in the initial theories, one should observe that in order to trigger an RGPEP evolution towards a solution when initially \( \mu = \nu \), one has to introduce an artificial splitting of masses in \( H_f \). For example, such splitting is needed in the case of local theories with massless bare particles and chiral symmetry. Two other physically important cases in which the mass degeneracy and its minimal lifting may play important roles as far as an application of RGPEP is concerned, are neutrinos in electroweak interactions and \( u \) and \( d \) quarks in QCD.

Eqs. (53) and (54) imply that the sum \( \alpha + \beta \) as a function of \( u \) is a constant. This constant, denoted by \( T = T / \delta \mu^2 \), results from the constancy of a trace of the mass-squared matrix, \( T = m_1^2 + m_2^2 \), where \( m_1^2 \) and \( m_2^2 \) denote its eigenvalues. The remaining coupled set of equations reads
\[
\begin{align*}
\delta' &= 4 \gamma^2, \\
\gamma' &= - \delta \gamma,
\end{align*}
\]
where \( \delta = \alpha - \beta \). Multiplying the first of these two equations by 2\( \delta \) and the second by 2\( \gamma \), one arrives at
\[
\begin{align*}
\delta'^2 &= 8 \delta \gamma^2, \\
\gamma'^2 &= -2 \delta \gamma^2,
\end{align*}
\]
and concludes that
\[
\epsilon^2 = \delta^2 + 4 \gamma^2 \]
do not depend on \( u \). In fact,
\[
\epsilon^2 = T^2 - 4D,
\]
where $D = D/\delta \mu^4$ and $D$ is the determinant of the mass-squared matrix. Hence, $\epsilon^2 = (m_1^2 - m_2^2)^2/\delta \mu^4$. Using the constant $\epsilon$, one can eliminate $\gamma^2$ from Eq. (67) to obtain

$$\delta' = \epsilon^2 - \delta^2,$$

(71)

which is an ordinary differential equation. Since the difference between eigenvalues of a hermitian $2 \times 2$ matrix is never smaller than the difference between its diagonal matrix elements, one always has $\delta' > 0$ except when $\delta = \epsilon$ and the mass-squared matrix is diagonalized. Without any loss of generality one can assume $\epsilon > 0$.

Integrations of Eqs. (71) and then (66) produce solutions for the elements of mass-squared matrix as functions of $t$,

$$\mu_t^2 = \frac{1}{2} (\mu^2 + \nu^2) + \frac{1}{2} \delta \mu_t^2,$$

(72)

$$\nu_t^2 = \frac{1}{2} (\mu^2 + \nu^2) - \frac{1}{2} \delta \mu_t^2,$$

(73)

$$\delta \mu_t^2 = \delta \mu^2 \frac{\cosh x_t + \epsilon \sinh x_t}{\cosh x_t + \epsilon^{-1} \sinh x_t},$$

(74)

$$m_t^2 = m^2 \frac{1}{\cosh x_t + \epsilon^{-1} \sinh x_t},$$

(75)

where $x_t = \delta \mu^2 \delta m^2 t$. Note that $\epsilon = \sqrt{1 + (2m^2/\delta \mu^2)^2}$. For $t \to \infty$, one obtains

$$\mu_\infty^2 = m_1^2,$$

(76)

$$\nu_\infty^2 = m_2^2,$$

(77)

$$m_\infty^2 = 0.$$  

(78)

These results mean that the RGPEP eventually produces a Hamiltonian for the two new species of particles of types 1 and 2 that are free, i.e., they no longer mix due to interactions, and their masses squared are given by the eigenvalues $m_1^2$ and $m_2^2$ of the initial mass-squared matrix.

D. Effective particles

The result of RGPEP is a family of Hamiltonians $P_t = P_t(a_t, b_t)$ for $t \geq 0$, which is obtained from $P_t(a, b)$ in Eq. (34) by replacement of $a_0$ and $b_0$ by $a_{tp}$ and $b_{tp}$, respectively. The effective particle operators are obtained from Eq. (1). Namely,

$$a_{tp} = U_t a_0 U_t^\dagger,$$

(79)

$$b_{tp} = U_t b_0 U_t^\dagger,$$

(80)

where $U_t$ is given in Eq. (40) as a solution of

$$U_t' = -U_t [P_t^+, P_t^-].$$

(81)

The generator, i.e., the commutator on the right-hand side of Eq. (81), is given in Eq. (43). Using results of the previous section, the generator can be written as

$$[P_t^+, P_t^-] = \delta \mu^2 m_t^2 \int [p] \left( a_{0p}^* b_0 - b_{0p}^* a_0 \right).$$

(82)

Boost invariance of the RGPEP thus yields the generator that is a product of a function of $t$ times a constant operator. The $t$-ordered exponential in Eq. (83) is

$$U_t = \exp(\varphi_t A),$$

(83)

where

$$\varphi_t = -\delta \mu^2 \int_0^t m_\tau^2 d\tau$$

(84)

$$= \arctan \sqrt{\frac{\epsilon + 1}{\epsilon - 1}} - \arctan e^{\epsilon t} \sqrt{\frac{\epsilon + 1}{\epsilon - 1}},$$

(85)

$$A = \int [k] \left( a_{0k}^* b_0 - b_{0k}^* a_0 \right).$$

(86)
The effective particle operators $a_{tp}$ and $b_{tp}$ are obtained from the formula

$$q_{tp} = e^{\varphi_i A} q_{0p} e^{-\varphi_i A}$$

(87)

with $q = a$ and $q = b$, respectively, using

$$[A, a_{0p}] = -b_{0p} ,$$

(88)

$$[A, b_{0p}] = a_{0p} .$$

(89)

Suppose there exists a combination

$$q_{0p} = a_{0p} + z_1 b_{0p},$$

(90)

for which one has

$$[A, q_{0p}] = z_2 q_{0p} ,$$

(91)

where $z_1$ and $z_2$ are some complex numbers. This is possible when $z_1 = z_2 = \pm i$ and

$$e^{\varphi_i A} q_{0p} \pm e^{-\varphi_i A} = e^{\pm \varphi_i} q_{0p} \pm .$$

(92)

where $q_{0p} \pm = a_{0p} \pm i b_{0p}$. Knowing that

$$a_{0p} = \frac{1}{2} (q_{0p} + q_{0p}^-) ,$$

(93)

$$b_{0p} = \frac{1}{2} (q_{0p} - q_{0p}^-) ,$$

(94)

one obtains

$$a_{tp} = \cos \varphi_t a_{0p} - \sin \varphi_t b_{0p} ,$$

(95)

$$b_{tp} = \sin \varphi_t a_{0p} + \cos \varphi_t b_{0p} ,$$

(96)

and the inverse relations

$$a_{0p} = \cos \varphi_t a_{tp} + \sin \varphi_t b_{tp} ,$$

(97)

$$b_{0p} = -\sin \varphi_t a_{tp} + \cos \varphi_t b_{tp} .$$

(98)

Eqs. (95) and (96) provide explicit definitions of annihilation operators for effective particles corresponding to the RGPEP parameter $t = s^4$. The corresponding relations for creation operators are obtained by hermitian conjugation.

E. Interpretation of $s$ as the effective particle size

The interpretation of parameter $s$ as a size of effective particles requires explanation in the context of our mass-mixing model because the mass-mixing interaction does not change any three-momentum that could be an argument of a form factor whose width might be related to a concept of a particle size. However, in more advanced theories, interactions change an invariant mass of the interacting particles when their relative momenta change. To be specific, consider a fermion of mass $m_f$ that emits a boson of mass $m_b$. The associated change of invariant mass squared is

$$\mathcal{M}_{fb,f}^2 = \left( \sqrt{m_f^2 + k^2} + \sqrt{m_b^2 + k^2} \right)^2 - m_f^2 .$$

(99)

The exponential factor of the type $\exp (-s^4 \mathcal{M}_{fb,f}^4)$ becomes $\exp [-\langle 2sk \rangle^4]$ for large $k$. This is the origin of interpreting the parameter $s$ as a size of effective particles in complex theories. Namely, only particles with small size $s$ can interact producing a large momentum $k$, cf. [18].

In the mass-mixing model, there is no change of relative three-momentum involved. Instead, the interaction strength $m_f^2$ in Eq. (96) is limited in strength roughly by $\exp (-\delta \mu^2 \delta m^2 s^4)$. The change of interaction strength comes solely from the change of a particle mass. Therefore, the role of the effective particle size parameter $s$ is reduced to taming changes in the mass. The point-like, bare particles at $s = 0$ can change mass through a mass-mixing interaction by arbitrary amounts that are introduced in the initial $\mathcal{P}$. But the effective particles of large $s$ can change mass only by amounts not exceeding $1/s$, as if the motion of their constituents could not involve a large excitation without breaking them apart. Thus, when $s$ is large, the effective particles can only change their masses by small amounts. Eventually, when $s \to \infty$, they cannot change mass at all, which means that they do not interact through a mass mixing term at all (see Section IV C below). In any case, the RGPEP suggests that mass mixing in low-energy effective theories should be small. Realistic effective theories appear to share this feature.
F. Constance of the Hamiltonian

The effective Hamiltonian, $\mathcal{P}_t^\text{e} = \mathcal{P}_t^\text{e}(a_t, b_t)$, is obtained from $\mathcal{P}_t^\text{e}(a_0, b_0)$ in Eq. (134) by replacing $a_{0p}$ and $b_{0p}$ in the latter by $a_{tp}$ and $b_{tp}$. The result is

$$\mathcal{P}_t^\text{e} = \int [p] \left[ A_{tp} a_{tp}^\dagger a_{tp} + B_{tp} b_{tp}^\dagger b_{tp} + C_{tp} (a_{tp}^\dagger b_{tp} + b_{tp}^\dagger a_{tp}) \right],$$

(100)

where the coefficients $A_{tp}$, $B_{tp}$, and $C_{tp}$ are given in Eqs. (37), (38), (39), respectively, and the mass parameters in them are given in Eqs. (72), (83), and (85). Using Eqs. (95) and (96), one obtains

$$\mathcal{P}_t^\text{e} = \mathcal{P}_0^\text{e} + \int [p] \left[ \Delta \mu^2 p^+ a_{op}^\dagger a_{op} + \Delta \nu^2 p^+ b_{op}^\dagger b_{op} + \Delta m^2 p^+ (a_{op}^\dagger b_{op} + b_{op}^\dagger a_{op}) \right],$$

(101)

where

$$\Delta \mu^2 = \mu_t^2 c^2 + \nu_t^2 s^2 + 2 m_t^2 cs - \mu^2, \quad (102)$$

$$\Delta \nu^2 = \mu_t^2 s^2 + \nu_t^2 c^2 - 2 m_t^2 cs - \nu^2, \quad (103)$$

$$\Delta m^2 = -(\mu_t^2 - \nu_t^2)cs + m_t^2 (c^2 - s^2) - m^2.$$

(104)

$s = \sin \varphi_t$, and $c = \cos \varphi_t$. Direct inspection demonstrates that $\Delta \mu^2 = \Delta \nu^2 = \Delta m^2 = 0$ for all values of $t$ in the range from 0 to $\infty$, which means that the operators $\mathcal{P}_t^\text{e} = \mathcal{P}_t^\text{e}(a_t, b_t)$ and $\mathcal{P}_0^\text{e} = \mathcal{P}_0^\text{e}(a_0, b_0)$ are the same for all values of $t$.

G. Spectrum of the theory

The initial Hamiltonian, $\mathcal{P}_0^\text{e}$ in Eq. (33), is transformed as a result of the RGPEP to $\mathcal{P}_t^\text{e}$ in Eq. (100). At the same time, the RGPEP secures equality $\mathcal{P}_t^\text{e} = \mathcal{P}_0^\text{e}$, as shown in Section IV. Since the eigenvalues and eigenstates of $\mathcal{P}_0^\text{e}$ and $\mathcal{P}_t^\text{e}$ are identical, one can derive them using any value of $t$ one wishes. The simplest to discuss is the case of $t \to \infty$, because in this case there is no mass mixing, $m_{\infty} = 0$. The mixing vanishes in the limit $t \to \infty$ provided that initially $\mu \neq \nu$. This is assumed in what follows. The case of $\mu = \nu$ is addressed near the end of this section.

The effective theory with $t = \infty$ is a free theory, with a correspondingly simple spectrum. Details of the spectrum are described below for two reasons. One reason is the completeness of the article. The other reason is a preparation for the discussion in Section V concerning the ground state, or vacuum. Simplicity of the RGPEP illustrated here is contrasted with complexity of other approaches there.

In the limit of $t \to \infty$,

$$\mathcal{P}_\infty^\text{e}(a_\infty, b_\infty) = \int [p] \left[ \frac{p^+}{p^+} a_{\infty p}^\dagger a_{\infty p} + \frac{p^+}{p^+} b_{\infty p}^\dagger b_{\infty p} \right],$$

(105)

where

$$a_{\infty p} = \cos \varphi_\infty a_{0p} - \sin \varphi_\infty b_{0p}, \quad (106)$$

$$b_{\infty p} = \sin \varphi_\infty a_{0p} + \cos \varphi_\infty b_{0p}, \quad (107)$$

and the angle $\varphi_\infty$ is

$$\varphi_\infty = - \arctan \sqrt{\frac{1 - \epsilon}{\epsilon + 1}}.$$  

(108)

Note that this angle is the same as the one in Eq. (A10) that results from solving RGPEP equations with a different generator in Appendix A.

Eigenvalues of the Hamiltonian in Eq. (105) are free energies of $n_{\infty 1}$ particles of mass $m_1$ and $n_{\infty 2}$ particles of mass $m_2$, each with some momentum components $p^+$ and $p^+$,

$$P_{\{p_{i1}, i=1, \ldots, n_{\infty 1}\}, \{p_{j2}, j=1, \ldots, n_{\infty 2}\}} = \sum_{i=1}^{n_{\infty 1}} \frac{p_{i1}^+ + m_1^2}{p_{i1}^+} + \sum_{j=1}^{n_{\infty 2}} \frac{p_{j2}^+ + m_2^2}{p_{j2}^+}.$$

(109)
The spectrum is degenerate. The eigenstates can be closely identified because the RGPEP provides expressions for the operators $a_\infty$ and $b_\infty$. A complete set of eigenstates (not normalized) is defined by writing

$$ |\{p_{1i}, i = 1, \ldots, n_{\infty1}\}, \{p_{2j}, j = 1, \ldots, n_{\infty2}\} \rangle = \prod_{i=1}^{n_{\infty1}} a_{0p_{1i}}^* \prod_{j=1}^{n_{\infty2}} b_{0p_{2j}}^* |0\rangle, \quad (110) $$

where $|0\rangle$ denotes the vacuum state. The vacuum state is annihilated by all annihilation operators of all particles for all values of $t$ and one can treat $|0\rangle$ as one and the same state for all values of the parameter $t = s^4$.

Since the creation operators $a_{0p_{1i}}^*$ and $b_{0p_{2j}}^*$ are given by linear combinations of $a_{0p}^*$ and $b_{0p}^*$ implied by Eqs. (106) and (107) through hermitian conjugation, the eigenstates defined in Eq. (110) can also be written as combinations of states created from the same vacuum state by products of the operators $a_{0p}^*$ and $b_{0p}^*$ with the corresponding momenta. The total number of particles in every resulting component of an eigenstate is the same. However, an eigenstate with definite numbers $n_{\infty1}$ and $n_{\infty2}$ of effective particles with $t = \infty$ corresponds to a combination of states with varying numbers of initial particles, $n_{01}$ and $n_{02}$, that satisfy the condition $n_{01} + n_{02} = n_{\infty1} + n_{\infty2}$. If the total number of particles is large, a simple state of effective particles with $t = \infty$ is a complex mixture of many states made of bare particles corresponding to $t = 0$.

When $\mu = \nu$, the RGPEP does not change the particle operators, since the generator is zero. On the other hand, it is clear that a non-zero mixing term $m^2$ causes the eigenvectors of mass-squared matrix in a classical Lagrangian to be definite combinations of the initial basis vectors. In the quantum theory, in order to generate a solution using the RGPEP, one may introduce a small artificial difference between the initial masses. When the initial mass degeneracy is changed by $m$, a non-zero mixing term in the Lagrangian of Eq. (14) is treated as a coupling constant. Setting $m = 0$, one obtains a Lagrangian of a free theory,

$$ \mathcal{L}_0 = \frac{1}{2} (\partial^2 \phi^2 - \mu^2 \phi^2) + \frac{1}{2} \nu^2 \phi^2 - \frac{1}{2} (\partial^2 \chi^2 - \mu^2 \chi^2) + \frac{1}{2} \nu^2 \chi^2. \quad (111) $$

The IF quantization of a free theory is well-known and nothing new is said here about it except for stressing one aspect that concerns the vacuum. Namely, when one evaluates $H_0 = \int d^3x \mathcal{H}_0$, where the Hamiltonian density $\mathcal{H}_0$ is canonically obtained from $\mathcal{L}_0$, the terms that involve products of two creation or two annihilation operators all cancel out, as desired. This happens because of the free energy formulae, $E_0^2(p) = \mu^2 + p^2$ and $E_0^2(p) = \nu^2 + p^2$, that are used in defining the time derivatives, or canonical momenta for the field variables. These energy formulae produce the desired cancellations in the sum of terms involving $\pi^2_\phi$, $\nabla \phi^2$, and $\mu^2 \phi^2$, and similarly for $\pi^2_\chi$, $\nabla \chi^2$, and $\nu^2 \chi^2$. The resulting $H_0$ in the IF of dynamics has the form

$$ H_0(a_0, b_0) = \int |p|_a \sqrt{\mu^2 + p^2} a_{0p}^* a_{0p} + \int |p|_b \sqrt{\nu^2 + p^2} b_{0p}^* b_{0p}. \quad (112) $$

V. THE VACUUM PROBLEM

The vacuum problem appears in the quantization of fields [4–6]. One starts with quantizing a free classical theory. This renders a quantum theory of non-interacting particles in terms of a free Hamiltonian $H_0$. Interaction terms are added to $H_0$ in the form of $H_I$. The latter can be constructed by starting from local products of classical fields multiplied by coupling constants and replacing the classical fields with the quantized ones. The vacuum problem becomes apparent when one attempts to solve the eigenvalue problem for $H = H_0 + H_I$. The problem is that $H_I$ takes eigenstates of $H_0$ out of the Hilbert space. In particular, the ground state of the free theory, denoted by $|0\rangle$, is changed by $H_I$ to a state with an infinite norm. The situation is further discussed below using the mass-mixing example, in which the vacuum problem appears in a similar way as in the model used by Dirac to discuss the vacuum problem [4].

A. Vacuum problem due to mass-mixing

The parameter $m^2$ in the mixing term in the Lagrangian of Eq. (14) is treated as a coupling constant. Setting $m = 0$, one obtains a Lagrangian of a free theory,

$$ \mathcal{L}_0 = \frac{1}{2} (\partial^2 \phi^2 - \mu^2 \phi^2) + \frac{1}{2} (\partial^2 \chi^2 - \mu^2 \chi^2). \quad (111) $$

The IF quantization of a free theory is well-known and nothing new is said here about it except for stressing one aspect that concerns the vacuum. Namely, when one evaluates $H_0 = \int d^3x \mathcal{H}_0$, where the Hamiltonian density $\mathcal{H}_0$ is canonically obtained from $\mathcal{L}_0$, the terms that involve products of two creation or two annihilation operators all cancel out, as desired. This happens because of the free energy formulae, $E_0^2(p) = \mu^2 + p^2$ and $E_0^2(p) = \nu^2 + p^2$, that are used in defining the time derivatives, or canonical momenta for the field variables. These energy formulae produce the desired cancellations in the sum of terms involving $\pi^2_\phi$, $\nabla \phi^2$, and $\mu^2 \phi^2$, and similarly for $\pi^2_\chi$, $\nabla \chi^2$, and $\nu^2 \chi^2$. The resulting $H_0$ in the IF of dynamics has the form

$$ H_0(a_0, b_0) = \int |p|_a \sqrt{\mu^2 + p^2} a_{0p}^* a_{0p} + \int |p|_b \sqrt{\nu^2 + p^2} b_{0p}^* b_{0p}. \quad (112) $$
The integration measures obtain the subscripts \( a \) and \( b \) because of the energies in their denominators. The non-zero commutation relations are \([a_{bp}, a_{bq}^+]=2E_a(p)\,(2\pi)^3\delta^3(\vec{p}-\vec{q})\) and \([b_{bp}, b_{bq}^+]=2E_b(p)\,(2\pi)^3\delta^3(\vec{p}-\vec{q})\). Possible additional quantum numbers can be ignored here.

An infinite constant \( \Omega^0 \) has been removed by the IF normal ordering, analogous to the constant \( \Omega^- \) removed from the FF Hamiltonian \( \mathcal{P}^- \), see Eqs. (40) and (41). The constant \( \Omega^0 \) can be subtracted this way \([4, 27]\), or it can also be included in variational estimates of the ground-state energy when interaction terms are taken into account \([20, 21]\).

In a theory set up this way, the IF vacuum problem emerges in the model due to the interaction term,

\[
H_I = \int d^3x \frac{m^2}{2} \phi \chi
\]

\[
= \int \frac{d^3p}{(2\pi)^3} \frac{m^2}{4E_aE_b} \left( a_{bp}^+ b_{b0-p}^0 + a_{bp}^+ b_{b0-p}^0 + b_{bp}^+ a_{b0} + a_{bp} + b_{bp}^+ \right).
\]

The result of action of \( H_I \) on the vacuum state is

\[
H_I |0\rangle = \int \frac{d^3p}{(2\pi)^3} \frac{m^2}{4E_aE_b} a_{bp} b_{b0-p}^0 |0\rangle.
\]

This state has an infinite norm. The infinity occurs through two factors. One factor is the volume of space in which the states with definite three-momentum are normalized. In the case of a state in Eq. (113), the three-momentum is zero. Another source of infinity is the integral over all momentum labels \( p \). This divergence results from the infinite number of momentum scales in the theory.

Acting on the state in Eq. (113) with \( H_I \) again also generates infinity. Multiple action of \( H_I \) creates further infinities. For example, the infinities appear in action of the evolution operator \( U = \exp(-iHt) \) on \( |0\rangle \), since \( U \) involves all powers of \( H_I \) \([4]\).

If states of the theory are built starting from \( |0\rangle \), the mixing operator \( H_I \) creates infinities in all of them. Removal of the infinities requires a cutoff on the range of momentum \( p \) in the Fourier expansions of fields \( \phi \) and \( \chi \). However, every cutoff on the momentum range in a theory violates the Lorentz symmetry \([4]\). One has to re-design the quantization procedure in the example in order to recover the quantum theory that was straightforwardly found as a solution using the RGPEP in previous sections.

On the one hand, it is known in the elementary model what needs to be done to solve it. On the other hand, one can look at the model as sharing some basic features with theories in which a more complex \( H_I \) is added to \( H_0 \) and it is not known how to deal with the vacuum problem in them beyond perturbation theory. Therefore, the model is of interest as a potential source of ideas about how to use the RGPEP to try to work around the vacuum problem in complex theories and attempt to break through the barriers that this problem poses in general.

### B. General scope of vacuum problems

Dirac pointed out that problems with vacuum may require a re-interpretation of quantum field theory \([4]\). He argued for such re-interpretation in the case of QED. Similar divergences occur in the vacuum problem of QCD but they cannot be as easily worked around as Dirac suggested for QED \([20]\).

One reason is that the coupling constant in QCD is much larger than in QED. The QED coupling constant is so small that one can use very large cutoffs in diverging terms in perturbation theory and still does not need to worry about the Lorentz-symmetry violation in practice. In QCD, where the coupling constant is much larger than in QED, the cutoffs would have to be much smaller than in QED in order to exclude large terms in perturbation theory. But much smaller cutoffs on \(|\vec{p}|\) could lead to effects that violate the Lorentz symmetry much stronger. Asymptotic freedom enables perturbative calculations in QCD but does not solve the vacuum problem. The other reason is the need for explaining spontaneous chiral symmetry breaking \([28, 29]\) for which a non-trivial vacuum structure is seen as the origin. The third reason is the desire to explain confinement. Confinement is often associated in the literature with a concept of a complex ground state. In any case, the ground state of QCD still awaits a construction. More generally, questions concerning a ground-state, spontaneous symmetry breaking, and mass generation, are of concern in the present standard model and theories trying to explain its origin. A famous ambiguity involved in the vacuum concept is the vacuum energy density, which can be seen as relevant to cosmology \([30, 31]\).

In the FF of Hamiltonian dynamics, the vacuum problem does not appear in the same way as in the IF. For example, the vacuum problem in the FF version of QCD can be formulated as a renormalization group problem for Hamiltonians \([20]\). Using the RGPEP, one can also envision a scenario for solving the canonical FF of QCD in which the effects commonly associated with a gluon condensate in vacuum \([32]\) may actually originate in an analogous expectation value but merely in the gluon medium that exists only inside the volume of a hadron, rather than in the...
entire space [1]. Discussions of the idea that condensate parameters may actually correspond to expectation values of operators in the medium present inside hadrons, instead of the entire space, are available in [33–36], including implications for cosmology.

The scope of vacuum problems is broad enough to suggest that the features that enable RGPEP to work around the vacuum problem and produce an exact quantum solution in the elementary example, should be identified. This is done in the next section.

C. RGPEP path around the vacuum

The general features that enable RGPEP to circumvent the vacuum problem and still produce a relativistic solution in the model stem from the FF of Hamiltonian dynamics. The key properties of the FF are the positivity of $p^+$ and boost invariance. The RGPEP takes advantage of these properties in the design of its generator.

The positivity of $p^+$ results from the assumption that for a free particle of an arbitrary mass $\mu > 0$ one can write for arbitrary three-momentum $\vec{p}$ that

$$ p^+ = \sqrt{\mu^2 + \vec{p}^2} + p^z \geq 0. \quad (116) $$

Thus, one assumes in the FF of quantum dynamics that a creation operator for a particle may only carry positive $p^+$ as a label. This feature is summarized in Eqs. (22) and (23) in Section IIIB.

Positivity of $p^+$ in Eq. (116) implies that the classical, translation-invariant mass-mixing interaction term in Eq. (18),

$$ P_{\vec{p}}^{-} = \int dx^+ d^3 x \phi \chi, \quad (117) $$

results in the quantum interaction operator in Eq. (33),

$$ P_{\vec{p}}^{-} = \int [p] \theta(p^+) \frac{m^2}{p^+} (a_{0\vec{p}}^+ b_{0\vec{p}} + b_{0\vec{p}}^+ a_{0\vec{p}}), \quad (118) $$

which does not contain any terms of the type $a_{\vec{p}}^+ b_{-\vec{p}}^+$ and $a_{\vec{p}} b_{-\vec{p}}$ that appear in Eq. (114) for $H_1$ in the IF of quantum dynamics. Such terms are excluded because both $p^+$ and $-p^+$ in them are required to be positive. This is not possible for particles of a finite mass in a presence of a cutoff on $p = |\vec{p}|$, as is visible in Eq. (116), no matter how large such cutoff is. Note also that the FF integration measure $[p]$ does not depend on the mass $\mu$ used in the condition (116).

When the cutoff on $|\vec{p}|$ can be made arbitrarily large, one can have boost invariance in practice in an arbitrarily large range of momenta provided that the theory respects the symmetry [29]. This is the case at $s=0$ in the RGPEP. In order to maintain the Lorentz symmetry in an effective theory, the sliding cutoff parameter $\lambda = 1/s$ emerges in the RGPEP through its equations. They are so designed that the sliding cutoff is not limiting $|\vec{p}|$ of individual particles. Instead, the effective-theory cutoff limits only the changes of invariant mass caused by interactions. The mass is invariant with respect to all 7 FF kinematical symmetries, including boost invariance.

The boost invariance is secured by design of the RGPEP generator in Eq. (19). The commutator guarantees that only connected interactions are generated. The total transverse momenta of interacting particles before and after an interaction cancel each other in the arguments of resulting vertex form factors. Spectators do not contribute to these arguments. The multiplication by a total + momentum squared of interacting particles in the definition of $H_{P_1}$, Eq. (10), results in the factor $p^+ 2$ in Eq. (36). In the absence of sensitivity to cutoffs on $p^+$, this factor removes $p^+$ from the RGPEP evolution equation entirely. Therefore, the arguments of resulting vertex form factors depend only on the change of invariant mass squared among the particles that are involved in the interaction.

These features, combined with the absence of divergences due to separation of momentum modes, reduce the RGPEP in the elementary mass-mixing model to solving an evolution equation for particle masses as functions of $t = s^4$. Quite generally, renormalized equations for coefficients $c_i$ in $H_{P_1}$ may involve only masses, relative momenta, coupling constants, and the parameter $t$. Thus, in the elementary model, the equations involve only $\mu_{1}^2$, $\nu_{1}^2$, mass-mixing parameter $m_1^2$, and $t$ itself. These equations are independent of the particle momentum $p$. As a result, the RGPEP equations render a different representation of the same relativistic quantum theory for every value of $t$.

Each and every one of the effective theories derived using the RGPEP, is defined in terms of a different basis in the space of operators acting in the Fock space. In the mass-mixing example, the effective representations tend in the limit of $t \to \infty$ to a relativistic theory of free particles with masses $m_1$ and $m_2$. No variation of the ground state $|0\rangle$ with $t$ is required in the procedure.
D. Standard, IF approach versus RGPEP

The comparison relies on a change of field variables in the classical Lagrangian of Eq. \((14)\). The new variables are determined by diagonalization of the mass-squared matrix. The mass terms,

\[
-2\mathcal{L}_{\text{mass}} = \mu^2 \phi^2 + \nu^2 \chi^2 + 2m^2 \phi \chi
\]

(119)
can be written in the form of a \(2 \times 2\) matrix sandwiched with a doublet of fields \(\Psi = [\phi, \chi]\). Namely,

\[
-2\mathcal{L}_{\text{mass}} = \Psi^\dagger M^2 \Psi
\]

(120)

\[
= [\phi, \chi] \begin{bmatrix} \mu^2 & m^2 \\ m^2 & \nu^2 \end{bmatrix} \begin{bmatrix} \phi \\ \chi \end{bmatrix}.
\]

(121)
The eigenvalues of matrix \(M^2\), denoted by \(m_1^2\) and \(m_2^2\) above Eq. \((65)\) in Section IV C, and the corresponding eigenvectors, are

\[
m^2_{1,2} = (\mu^2 + \nu^2)/2 \pm \sqrt{(\mu^2 - \nu^2)^2/4 + 4m^4},
\]

(122)

\[
v_1 = \begin{bmatrix} \cos \varphi_\infty \\ -\sin \varphi_\infty \end{bmatrix}, \quad v_2 = \begin{bmatrix} \sin \varphi_\infty \\ -\cos \varphi_\infty \end{bmatrix},
\]

(123)

where \(\varphi_\infty\) is given in Eq. \((108)\). Inverting the relation

\[
\Psi = \xi v_1 + \zeta v_2,
\]

(124)
one can define the fields

\[
\xi = \cos \varphi_\infty \phi - \sin \varphi_\infty \chi,
\]

(125)

\[
\zeta = \sin \varphi_\infty \phi + \cos \varphi_\infty \chi.
\]

(126)

This is a unitary change of field variables. Since the terms that involve derivatives of the fields \(\phi\) and \(\chi\) have equal coefficients in the classical Lagrangian density of Eq. \((14)\), the density can be written as

\[
\mathcal{L} = \frac{1}{2} [(\partial \xi)^2 - m_1^2 \xi^2] + \frac{1}{2} [(\partial \zeta)^2 - m_2^2 \zeta^2].
\]

(127)

This classical expression can now be quantized from scratch in the IF of dynamics.

The IF quantization involves definitions of the fields \(\xi\) and \(\zeta\) and their conjugated momenta. The quantization leads to a theory of particles with masses \(m_1\) and \(m_2\) when one defines the quantum fields \(\xi\) and \(\zeta\) and their conjugated momenta \(\pi_\xi\) and \(\pi_\zeta\) using energy expressions \(E_\xi^2(p) = m_1^2 + p^2\) and \(E_\zeta^2(p) = m_2^2 + p^2\) in defining the time derivatives of the fields, respectively. The new energy expressions guarantee that all terms of the type \(a_\xi^i a_\xi^{-i}\) or \(a_\xi^i a_\zeta^{-i}\) cancel out in the Hamiltonian. Eqs. \((125)\) and \((126)\) imply that

\[
a_{\xi p} = \cos \varphi_\infty a_{0 p} - \sin \varphi_\infty b_{0 p},
\]

(128)

\[
a_{\zeta p} = \sin \varphi_\infty a_{0 p} + \cos \varphi_\infty b_{0 p}.
\]

(129)

These relations match Eqs. \((106)\) and \((107)\). The matching shows that the IF quantization of fields \(\xi\) and \(\zeta\) produces the same result as the solution obtained entirely in one quantum theory using the RGPEP, in which there is no need to re-quantize the theory due to inclusion of the mass-mixing interaction term.

On the basis of knowing the full quantum implications of the mass-mixing interaction term in the FF of Hamiltonian dynamics, one can also write expressions for the initial IF quantum fields \(\hat{\phi}, \hat{\chi}, \hat{\pi}_\phi,\) and \(\hat{\pi}_\chi\), in terms of the fields \(\xi, \zeta, \hat{\pi}_\xi,\) and \(\hat{\pi}_\zeta,\) using Eq. \((124)\) and right energies for the time derivatives needed in \(\hat{\pi}_\xi\) and \(\hat{\pi}_\zeta\). Substituting these expressions into a classical IF Hamiltonian that canonically corresponds to the Lagrangian density of Eq. \((14)\), one obtains the IF quantum Hamiltonian that explicitly describes the same physics as the FF quantum Hamiltonian obtained from the RGPEP at \(t = \infty\). The vacuum-altering terms cancel out for all modes with a finite momentum.

However, when in a more complex theory than the elementary example some additional interaction terms cause divergences and other effects that are difficult to see through, the IF quantization approach may get stuck due to lack of a right guess for the time derivatives. In contrast, the RGPEP still indicates a direction for further studies in realistic cases. Namely, while the free theory that results from diagonalization of a bilinear part in a Lagrangian...
density is certainly not sufficient for establishing how to deal with the IF vacuum problem, the RGPEP promises some capability to work around the vacuum problem using the FF.

The FF Hamiltonian at \( t = s^4 = 0 \), \( \mathcal{P}_0^- \), involves fields \( \phi \) and \( \chi \). Their conjugate “momenta,” \( \pi_\phi = \partial^+ \phi \) and \( \pi_\chi = \partial^+ \chi \), do not involve FF time derivatives, i.e., they do not involve derivatives with respect to \( x^+ \). Instead, the “momenta” are expressed through gradients of the fields in the front hyper-plane. Rotation of the quantum fields automatically rotates the quantum “momenta.”

By the way, the fields \( \xi \) and \( \zeta \) can be used as initial variables also in the FF. The RGPEP provides no additional value in such setup, since there is no interaction between the free fields \( \xi \) and \( \zeta \). However, when more interactions are added, nothing prevents the RGPEP from application to the whole quantum theory using the effective particle operators associated with the fields \( \xi \) and \( \zeta \), instead of \( \phi \) and \( \chi \).

It should be mentioned that an interesting example of the IF application of a similarity renormalization group procedure in a fixed source model has been recently considered by Jones and Perry [37]. In the fixed source model, the interaction term is only linear in the quantum field variables, different momentum modes evolve separately, and one obtains the well-known solution in an elegant way. The fixed source model does not appear to suggest how to proceed in the IF when interaction terms involve more than one field and create a genuine vacuum problem.

VI. CONCLUSION

The case of a theory with two free fields with a mass-mixing interaction term can be generalized to theories with an arbitrary number \( n > 2 \) of fields and mass mixing terms. In such theories, the RGPEP equation describes the evolution of a mass matrix of dimension \( n \times n \) with \( s \). The solution tends for \( s \to \infty \) to a diagonal matrix, whose eigenvalues provide physical masses for \( n \) species of free particles.

Degeneracy of the mass matrix, which may correspond to a symmetry in a theory, prevents its full diagonalization via the RGPEP equation. In this case, an artificial infinitesimal breaking of the degeneracy can be introduced in order to enable RGPEP to identify a solution in the limit \( s \to \infty \), as the artificial breaking is being removed.

The diagonalization of the mass matrix does not correspond to a minimization of a classical potential in the IF. Instead, it corresponds to identification of the eigenmodes in classical field oscillations. One has to use the eigenmodes in the IF quantization procedure in order to solve a vacuum problem in the absence of interactions other than the mass mixing. However, when such additional interaction terms, involving products of more than two fields, are included in a theory, the IF vacuum problem can no longer be solved using the field combinations that correspond to eigenvectors of the mass matrix. The additional interactions typically contribute to particle masses, bound states may develop, and, as it would have to happen in the case of confinement, the full theory eigenmodes do not even correspond to the fields present in an initial Lagrangian.

The intriguing feature of the RGPEP, illustrated here in the elementary model with a mass-mixing interaction term, is that it applies to quantum theories via steps that are essentially independent of the type of interaction one grapples with, while the vacuum problem is treated in a new way. Namely, the vacuum stays simple while the interaction terms evolve towards expressions in terms of effective degrees of freedom. This feature makes the RGPEP a deserving candidate for application to more realistic theories than the elementary model discussed here. It is evident from the works referenced in this article that the RGPEP can be applied to realistic quantum field theories. The elementary example described here is thus of interest not only as an illustration of an exact non-perturbative solution of the RGPEP equations but also as the indicator of a difference between the options one has got left for treating vacuum problems in the IF and FF of Hamiltonian dynamics in relativistic quantum field theories.

Appendix A: Solution for \( \mathcal{H}_f \) dependent on \( t \)

Discussion of Eq. (57) in Section IVB included the case of \( \mathcal{H}_f \) containing masses dependent on \( t \), which yields a \( 2 \times 2 \) mass-squared matrix equation of the form

\[
\begin{bmatrix}
\mu_1^2 & m_1^2 \\
m_1^2 & \nu_1^2
\end{bmatrix} = \begin{bmatrix}
\mu_1^2 & \nu_1^2 \\
0 & \nu_1^2
\end{bmatrix} \cdot \begin{bmatrix}
\mu_1^2 & m_2^2 \\
m_2^2 & \nu_2^2
\end{bmatrix} \cdot \begin{bmatrix}
\mu_2^2 & m_2^2 \\
m_2^2 & \nu_2^2
\end{bmatrix}.
\]

(A1)

This equation matches the Wegner equation for a Hamiltonian matrix [10] of a two-level system. Its analytic solution is well-known but as far as the author knows it was never considered before in the context of particle masses in an exactly soluble quantum field theory in the FF of dynamics.

Proceeding as in Section IV.C one obtains

\[
\delta' = 4 \delta \gamma^2,
\]

(A2)
\[ \gamma' = -\delta^2 \gamma. \] (A3)

Multiplying the first of these two equations by \(2\delta\) and the second by \(2\gamma\), one arrives at
\[ \delta'^2 = 8\delta^2 \gamma^2, \] \[ \gamma'^2 = -2\delta^2 \gamma^2, \]
which implies the same constant \(\epsilon^2 = \delta^2 + 4\gamma^2\) as in Section [IVC]. After eliminating \(\gamma^2\) from Eq. (A3),
\[ \delta'^2 = 2\delta^2 (\epsilon^2 - \delta^2). \] (A6)

The solutions corresponding to Eqs. (74) and (75), are
\[ \delta_{\mu t}^2 = \delta_{\mu}^2 e^{\epsilon x t} \sqrt{\epsilon^2 - 1 + e^{2\epsilon x t}}, \] \[ m_{\mu t}^2 = m^2 e^{\epsilon \sqrt{\epsilon^2 - 1 + e^{2\epsilon x t}}}, \]
where \(x_t = (\delta m^2)^2 t\).

Since the generator given in Eqs. (43) and (82) is now altered to contain the varying \(\delta_{\mu t}^2\) instead of the constant \(\delta_{\mu}^2\), the angle \(\varphi_t\) given by Eq. (84) is replaced by
\[ \varphi_t = -\int_0^t \frac{\delta_{\mu t}^2 m_{\mu t}^2 d\tau}{\sqrt{\epsilon^2 - 1 + e^{2\epsilon x t}}}, \] (A9)
This result deviates from the result in Eq. (84) for finite values of \(t\). The difference in the angles of rotation, \(\varphi_t\), implies different combinations of operators \(a_{\mu 0}\) and \(b_{\mu 0}\) in Eqs. (84) and (A9) for the same \(t\). This means that the effective particle operators at any finite \(t > 0\) depend on the choice of the generator, although the Hamiltonians as operators are just one and the same operator for all values of \(t\) and both choices of the generator. When \(t \to \infty\), Eqs. (84) and (A9) produce the same result for \(\varphi_{\infty}\) for arbitrary values of \(\epsilon > 1\),
\[ \varphi_{\infty} = \arctan \frac{1 + \epsilon}{\epsilon - 1} - \pi/2 = \frac{1}{2} \arctan \frac{1}{\sqrt{\epsilon^2 - 1}} - \frac{\pi}{4}. \] (A10)
Thus, the change in the generator from a constant \(H_t\) to a \(t\)-dependent full free part of \(H_t\), does not lead to any change in the effective particles that one obtains for \(t \to \infty\) as a solution of the theory.

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