Instability of a uniform electric field in pure non-Abelian Yang-Mills theory

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We study the Schwinger process in a uniform non-Abelian electric field using a dynamical approach in which we evolve an initial quantum state for gluonic excitations. We evaluate the spectral energy density and number density in the excitations as functions of time. The total energy density has an ultraviolet divergence which we argue gets tamed due to asymptotic freedom, leading to \( g^4 E^4 t^4 \) growth, where \( g \) is the coupling and \( E \) the electric field strength. We also find an infrared divergence in the number density of excitations whose resolution requires an effect such as confinement.

I. INTRODUCTION

The Schwinger effect \(^1\), whereby non-perturbative quantum effects in a background electric field lead to electron-positron pair production, has received much attention (for example, see the reviews \(^2\)). Heuristically, the electric field pulls apart the electron-positron pairs that are fluctuating in and out of the vacuum. For weak electric field strength \( E \ll \pi m_e^2/e \), where \( e \) and \( m_e \) are the positron’s electric charge and mass, the Schwinger process can be thought of as a quantum tunneling event and is exponentially suppressed in the dimensionless combination of variables \( \pi m_e^2/eE \). The rate of creation of number density of electrons is,

\[
\dot{n} \propto e^2 E^2 \exp(-\pi m_e^2/eE)
\]  

(1)

Schwinger’s original computation deals with the probability of vacuum persistence\(^3\) for spin-0 and spin-1/2 matter fields and has since been generalized to numerous other systems besides electromagnetism. Pair creation has also been studied for massless charges, equivalently for supercritical electric fields \( (eE \gg \pi m_e^2) \) in 1+1 dimensions \(^8\)\(^9\)\(^10\)\(^11\)\(^12\). In these cases, the exponential suppression of the original Schwinger effect is not present, and other techniques have to be employed as pair creation is no longer a tunneling process that is exponentially suppressed. For example, the massless Schwinger model can be solved completely, including backreaction on the electric field, and results in a \( 1/\sqrt{t} \) decay of the electric field strength \(^1\).

In the present work we are interested in non-Abelian gauge theory in the background of a homogeneous (color) electric field and the consequent Schwinger pair creation of “gluons”\(^2\). The process has been investigated before using effective action techniques to calculate vacuum persistence amplitudes \(^13\)\(^22\) with the result that there is a constant rate of particle number density production, still given by (1) with \( m_e = 0 \). The result is surprising to us since the Schwinger process can be viewed as a tunneling process and one might expect that the Wentzel–Kramers–Brillouin (WKB) \(^23\)\(^24\) and other approximations used to obtain (1) would break down for massless gauge fields. For this reason we wish to re-examine the problem using a different approach.

We take a dynamical approach to the problem. (For a kinetic approach to QED, see \(^25\)\(^30\)\.) At the initial time, we consider a color electric field background and quantum excitations in their non-interacting ground state. We then evolve the system, just as one would do for small quantum excitations in a time-dependent background. We use the method of Bogoliubov transformations \(^31\)\(^32\), re-set as a “classical-quantum correspondence” (CQC) whereby quantum evolution is described in terms of the classical evolution of a related classical system \(^33\)\(^34\). The method is explained in Appendix A. Our approach simply evolves an initial state in contrast to other methods that compute (non-interacting) vacuum persistence amplitudes.

Our results indeed differ from (1) with \( m_e \) set to zero. We find that the energy density in excitations, \( \mathcal{E} \), grows with time as \( \mathcal{E} \propto g^4 E^4 t^4 \) where \( g \) is the non-Abelian coupling constant (see also \(^37\)\(^38\)). We also examine the number density of particles in the leading adiabatic approximation \(^39\) and find that it is not well-defined as there is a zero frequency mode at all times for which the particle number density diverges. The total number density, found by integrating over all excitation modes, also diverges.

In Sec. I, we set up the basic framework, identifying the background field and the small excitations. Here we also discuss some limitations of our setup. Sec. II diagonalizes the Hamiltonian by expanding the quantum fields in modes. We find it convenient to discretize the modes for numerical analysis (Sec. IV). The quantum analysis is reduced to a classical analysis in Sec. V following which we numerically evaluate the spectral energy density and the total energy in quantum excitations as a function of time in Sec. VI. The number density of quantum excitations is discussed in Sec. VII and shown to diverge for all times due to the presence of a zero frequency mode. Sec. VIII briefly considers the case of adiabatically turning on the external field. We conclude in Sec. IX.

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1 For a nice account on the relation between the vacuum persistence probability and rate of pair creation see \(^7\).

2 We will refer to the model as “pure QCD” even though we will consider the simpler SU(2) gauge group. The “pure” means that we will only consider gauge fields and not include any other fields.
II. SETUP

We will consider a pure $SU(2)$ gauge theory with Lagrangian density,

$$ L = -\frac{1}{4}(W^a_{\mu\nu})^2 , \; a = 1, 2, 3 \quad (2) $$

with the field strength defined in the usual way

$$ W^a_{\mu\nu} = \partial_\mu W^a_\nu - \partial_\nu W^a_\mu + g e^{a\kappa} W^\kappa_\mu W^c_\nu \quad (3) $$

We have to make certain approximations to proceed with our analysis. Our main approximation is that we expand the fields about a fixed background electric field to quadratic order in the Lagrangian and ignore higher order interactions.

Let us write

$$ W^a_\mu = A^a_\mu + Q^a_\mu \quad (4) $$

where $A^a_\mu$ is a classical background and $Q^a_\mu$ denotes quantum excitations on top of the classical background. We will work in temporal gauge, so $W^a_0 = 0$, and take

$$ A^a_\mu = -\dot{A}(t) \delta^a \delta_{\mu 3} \quad (5) $$

Then there is an externally imposed classical electric field but no magnetic field,

$$ E^a_\mu = -\partial_\mu A^a_\mu = \dot{A}(t) \delta^a \delta_{\mu 3}, \; B^a_\mu = 0. \quad (6) $$

We will evolve the quantum variables, $Q^a_\mu$, assuming that they are in their non-interacting ground state at $t = 0$. We insert (4) in (2), then use the background (5), and expand to quadratic order in the $Q^a_\mu$ to obtain,

$$ L = \frac{1}{2}(\dot{Q}^1)^2 - \frac{1}{4}(\partial_\mu Q^1_\mu - \partial_\mu Q^1_\mu - g A(Q^1_\mu \delta^3 - \delta^3 Q^2_\mu))^2 
+ \frac{1}{2}(\dot{Q}^2_\mu)^2 - \frac{1}{4}(\partial_\mu Q^2_\mu - \partial_\mu Q^2_\mu + g A(Q^1_\mu \delta^3 - \delta^3 Q^2_\mu))^2 
+ \frac{1}{2}(\dot{Q}^3_\mu - \dot{\hat{A}})^2 - \frac{1}{4}(\partial_\mu Q^3_\mu - \partial_\mu Q^3_\mu + O((Q^3)^3)) \quad (7) $$

The classical electric field is externally imposed, i.e. there are external sources that produce and maintain the electric field $E^3_\mu$ which is assumed to be constant. Therefore, we take

$$ A(t) = E t, \quad \dot{\hat{A}} = E \quad (8) $$

Then the variables $Q^1_\mu$ decouple from the other quantum variables. We can calculate the rate of particle production in a fixed external field by considering the truncated Lagrangian,

$$ L' = \frac{1}{2}(\dot{Q}^1)^2 + \frac{1}{2}(\dot{Q}^2)^2 - \frac{1}{4}(Q^1_{ij})^2 - \frac{1}{4}(Q^2_{ij})^2 \quad (9) $$

where,

\begin{align*}
Q^1_{ij} &\equiv \partial_i Q^i_j - \partial_j Q^j_i - g E t (Q^2_{ij} \delta^3 - \delta^3 Q^2_{ij}) \\
Q^2_{ij} &\equiv \partial_i Q^i_j - \partial_j Q^j_i + g E t (Q^1_{ij} \delta^3 - \delta^3 Q^1_{ij}) \quad (10)
\end{align*}

At this level of approximation, the $Q^1_i$ and $Q^2_i$ fluctuations do not backreact on the background electric field. The backreaction will only appear due to the cubic and higher order terms in the $Q^3_i$ in (7).

We will expand the excitations in momentum modes in the next section. There are quantum issues at both ends of the spectrum. For modes with low energy, the coupling constant is strong and confinement should play a role. The lowest energy excitations will be massive glueballs, not massless gluons.\(^4\) Modes with very high energy are in the regime of asymptotic freedom as the coupling constant becomes small. Inclusion of these effects in our calculations is beyond our reach and we shall proceed based on (9) as if it is the full story and see if there are any inconsistencies. Indeed we will encounter two inconsistencies in this approach. In Sec. \(\text{VII}\)A we will encounter an ultraviolet divergence that we argue will be resolved by properly accounting for asymptotic freedom. We will also encounter a divergence in the number density of excitations at low energy at all times whose interpretation will change radically once we take confinement into account.

III. EXPANSION IN MODES

In this section we expand the fields in modes and diagonalize the Lagrangian. The calculations are straightforward if tedious; the end result for the diagonalized Lagrangian is given in (34) and (35).

The gauge fields are expanded in the physical transverse modes as,

$$ Q^1_\mu = \int \frac{d^3p}{\sqrt{2E_p}} \sum_{r=0}^3 \left[ b^r_{p} e^{ip \cdot x} + \text{h.c.} \right] \quad (11) $$

where \(d^3p = d^3p/(2\pi)^3\), \(E_p = |\mathbf{p}|\), and the reality of $Q^1$ implies $a^r_{-p} = a^r_{p}$. Similarly,

$$ Q^2_\mu = \int \frac{d^3p}{\sqrt{2E_p}} \sum_{r=0}^3 \left[ b'^r_{p} e^{ip \cdot x} + \text{h.c.} \right] \quad (12) $$

with $b'^r_{-p} = b^r_{p}$.

The magnetic fields are

\begin{align*}
(B^1)_i &= \frac{1}{2} \epsilon_{ijk} Q^1_{jk} = \nabla \times Q^1 - g E t Q^2 \times \hat{z} \quad (13) \\
(B^2)_i &= \frac{1}{2} \epsilon_{ijk} Q^2_{jk} = \nabla \times Q^2 + g E t Q^1 \times \hat{z} \quad (14)
\end{align*}

\(^4\) Indeed the assumed background uniform electric field itself ignores confinement. The spirit of the present work is that we work as if there is no confinement and study the consequences.
Therefore

\[ \mathbf{B}^1 = \int \frac{d^3 p}{\sqrt{2E_p}} \sum_{r,s=0}^3 \left[ (a^*_p \mathbf{p} + gEt b^*_p \hat{r} \hat{z}) \times \mathbf{\epsilon}_p e^{i \mathbf{p} \mathbf{\times} \mathbf{r}} + h.c. \right] \]

\[ = \int \frac{d^3 p}{\sqrt{2E_p}} \sum_{r,s=0}^3 \left[ a^*_p \mathbf{p} \times \mathbf{\epsilon}_p e^{i \mathbf{p} \mathbf{\times} \mathbf{r}} + h.c. \right] \]

(15)

where \( a^*_p \equiv (a^*_p \mathbf{p} + gEt b^*_p \hat{r} \hat{z}) \times \mathbf{\epsilon}_p \), and

\[ \mathbf{B}^2 = \int \frac{d^3 p}{\sqrt{2E_p}} \sum_{r,s=0}^3 \left[ (b^*_p \mathbf{p} - gEt a^*_p \hat{r} \hat{z}) \times \mathbf{\epsilon}_p e^{i \mathbf{p} \mathbf{\times} \mathbf{r}} + h.c. \right] \]

\[ = \int \frac{d^3 p}{\sqrt{2E_p}} \sum_{r,s=0}^3 \left[ b^*_p \mathbf{p} \times \mathbf{\epsilon}_p e^{i \mathbf{p} \mathbf{\times} \mathbf{r}} + h.c. \right] \]

(16)

where \( b^*_p \equiv (b^*_p \mathbf{p} - gEt a^*_p \hat{r} \hat{z}) \times \mathbf{\epsilon}_p \).

These expressions give,

\[ \mathcal{E}_{B1} = \frac{1}{2} \int d^3x(B^1)^2 \]

\[ = \int \frac{d^3 p}{2E_p} \sum_{r,s=0}^3 \left[ a^*_p \mathbf{p} \cdot \mathbf{a}^*_p + h.c. \right] \]

(17)

Note that there are two polarizations and \( \{ \mathbf{\epsilon}_p^1, \mathbf{\epsilon}_p^2, \mathbf{\hat{p}} \} \) form a right-handed orthonormal basis. For example,

\[ \mathbf{\epsilon}_p^1 = (\cos \theta \cos \phi, -\cos \theta \sin \phi, \sin \theta) \]

(18)

\[ \mathbf{\epsilon}_p^2 = (-\sin \phi, \cos \phi, 0) \]

(19)

\[ \mathbf{\hat{p}} = (\sin \theta \cos \phi, \sin \theta \sin \phi, \cos \theta) \]

(20)

We check

\[ (\mathbf{\hat{z}} \times \mathbf{\epsilon}_p^r) \cdot (\mathbf{\hat{z}} \times \mathbf{\epsilon}_p^s) = p^2 \delta^{rs} \]

(21)

\[ (\mathbf{\hat{z}} \times \mathbf{\epsilon}_p^r) \cdot (\mathbf{\hat{z}} \times \mathbf{\epsilon}_p^s) = p \cos \theta \delta^{rs} \]

(22)

\[ (\mathbf{\hat{z}} \times \mathbf{\epsilon}_p^r) \cdot (\mathbf{\hat{z}} \times \mathbf{\epsilon}_p^s) = \delta^{rs} - (\mathbf{\hat{z}} \cdot \mathbf{\hat{z}}) (\mathbf{\hat{r}} \cdot \mathbf{\hat{z}}) \]

(23)

and with the choice of vectors in \([18],[20]\),

\[ (\mathbf{\hat{r}} \cdot \mathbf{\hat{z}}) (\mathbf{\hat{z}} \cdot \mathbf{\hat{z}}) = \sin^2 \theta \delta^{rs} \]

(24)

Therefore

\[ \mathcal{E}_{B1} = \int \frac{d^3 p}{E_p} \sum_{r,s=1}^2 \left[ p^2 |a^*_p|^2 + g^2 E^2 t^2 |b^*_p|^2 (1 - \sin^2 \theta \delta_{r1}) \right. \]

\[ -igEt p_z (a^*_p b^*_p - a^*_p b^*_p) \]

(25)

Similarly

\[ \mathcal{E}_{B2} = \int \frac{d^3 p}{E_p} \sum_{r,s=1}^2 \left[ p^2 |b^*_p|^2 + g^2 E^2 t^2 |a^*_p|^2 (1 - \sin^2 \theta \delta_{r1}) \right. \]

\[ -igEt p_z (a^*_p b^*_p - a^*_p b^*_p) \]

(26)

\[ \mathcal{E}_{B1+2} = \mathcal{E}_{B1} + \mathcal{E}_{B2} \]

\[ = \int \frac{d^3 p}{E_p} \left[ (p^2 + g^2 E^2 t^2 \cos^2 \theta) (|a^*_p|^2 + |b^*_p|^2) \right. \]

\[ -igEt p_z \cos \theta (a^*_p b^*_p - a^*_p b^*_p) \]

\[ + (p^2 + g^2 E^2 t^2) (|a^*_p|^2 + |b^*_p|^2) \right. \]

\[ -igEt p_z \cos \theta (a^*_p b^*_p - a^*_p b^*_p) \]

(27)

Next let

\[ \alpha^r_p = \alpha^r_p + i \beta^r_p, \quad \beta^r_p = \gamma^r_p + i \delta^r_p \]

(28)

where \( \alpha^r_p, \beta^r_p, \gamma^r_p \), and \( \delta^r_p \) are real. The reality conditions, \( \alpha^r_p = \alpha^r_p \) and \( \beta^r_p = \beta^r_p \) imply

\[ \alpha^r_p = \alpha^r_p, \quad \beta^r_p = -\beta^r_p, \quad \gamma^r_p = \gamma^r_p, \quad \delta^r_p = -\delta^r_p \]

(29)

For convenience, define

\[ P = gEt, \quad P_z = P \cos \theta \]

(30)

Then

\[ \mathcal{E}_{B1+2} = \int \frac{d^3 p}{E_p} \left[ (p^2 + P_z^2) (|\alpha_p^1|^2 + (|\delta_p^1|^2) + 4pP_z \alpha_p^1 \delta_p^1 \right. \]

\[ + (p^2 + P_z^2) (|\beta_p^1|^2 + (|\gamma_p^1|^2) + 4pP_z \beta_p^1 \gamma_p^1 \]

\[ + (p^2 + P_z^2) (|\beta_p^2|^2 + (|\gamma_p^2|^2) + 4pP_z \beta_p^2 \gamma_p^2 \]

\[ + (p^2 + P_z^2) (|\beta_p^2|^2 + (|\gamma_p^2|^2) + 4pP_z \beta_p^2 \gamma_p^2 \]

(31)

The energy has separated into four disjoint sectors: \( (\alpha_p^1, \delta_p^1), (\beta_p^1, \gamma_p^1), (\alpha_p^2, \delta_p^2), (\alpha_p^2, \delta_p^2) \). The \( (\alpha_p^1, \delta_p^1) \) sector is equivalent to the \( (\beta_p^1, \gamma_p^1) \) sector under \( P_z \rightarrow -P_z \) which is the same as \( g \rightarrow -g \). Similarly, for the \( (\alpha_p^2, \delta_p^2) \) and \( (\beta_p^2, \gamma_p^2) \) sectors. The \( (\alpha_p^1, \delta_p^1) \) and \( (\alpha_p^2, \delta_p^2) \) sectors look very similar but they differ in their first terms: the former has \( P_z \) while the latter has \( P \).

The energy in the electric field is

\[ \mathcal{E}_{E1+2} = \frac{1}{2} \int d^3x \left[ ( \partial_t \mathbf{W}^1)^2 + ( \partial_t \mathbf{W}^2)^2 \right] \]

\[ = \int \frac{d^3 p}{E_p} \sum_{r=1}^2 \left[ (\dot{\alpha}_p^r)^2 + (\dot{\beta}_p^r)^2 + (\dot{\gamma}_p^r)^2 + (\dot{\delta}_p^r)^2 \right] \]

and we only need to solve for the dynamics of the two quantum systems,

\[ L_1 = \int \frac{d^3 p}{E_p} \left[ (\dot{\alpha}_p^1)^2 + (\dot{\delta}_p^1)^2 - (\dot{p}^2 + P_z^2) (|\alpha_p^1|^2 + (|\delta_p^1|^2) \right. \]

\[ -4pP_z \alpha_p^1 \delta_p^1 \]

(31)

\[ L_2 = \int \frac{d^3 p}{E_p} \left[ (\dot{\alpha}_p^2)^2 + (\dot{\delta}_p^2)^2 - (\dot{p}^2 + P_z^2) (|\alpha_p^2|^2 + (|\delta_p^2|^2) \right. \]

\[ -4pP_z \alpha_p^2 \delta_p^2 \]

(32)
As noted above, the Lagrangians for \((\beta^p_p, \gamma^p_p)\) are related to \(L_1\) and \(L_2\) by \(g \rightarrow -g\). Also recall that there is time-dependence in these Lagrangians because \(P\) and \(P\) grow in proportion to \(t\) as defined in (30).

The Lagrangians can be diagonalized by using linear combinations,

\[
\phi_{p,\pm} = \frac{\alpha_{1,p}^1 \pm \delta_{1,p}^1}{\sqrt{2}}, \quad \psi_{p,\pm} = \frac{\alpha_{2,p}^2 \pm \delta_{2,p}^2}{\sqrt{2}}
\]

Then

\[
L_1 = \int \frac{d^3p}{E_p} \left[ \dot{\phi}_{p,\pm}^2 - (p + P)^2 \phi_{p,\pm}^2 + \dot{\psi}_{p,\pm}^2 - (p - P)^2 \psi_{p,\pm}^2 \right] \]

\[
L_2 = \int \frac{d^3p}{E_p} \left[ \dot{\phi}_{p,\pm}^2 - (p + P)^2 \phi_{p,\pm}^2 + \dot{\psi}_{p,\pm}^2 - (p - P)^2 \psi_{p,\pm}^2 \right] \]

where \(P = gEt\), \(p_\perp = p \sin \theta\) and \(p_\parallel = p \cos \theta\).

Similarly, we can easily write down \(L_3\) and \(L_4\) for the \((\beta^p_p, \gamma^p_p)\) sector since these are related to \(L_1\) and \(L_2\) by \(g \rightarrow -g\). The full Lagrangian is given by the sum of \(L_1, \ldots, L_4\).

### IV. DISCRETIZATION

We now discretize the integrations in (34) and (35). For example,

\[
L_1 = \sum_p (\Delta p)^3 \left[ \dot{\phi}_{p,\pm}^2 - (p + P)^2 \phi_{p,\pm}^2 \right] \equiv \sum_p L_{1p}
\]

and similarly for \(L_2\). The volume element in momentum space is

\[
(\Delta p)^3 = \left( \frac{2\pi}{L} \right)^3 = \left( \frac{2\pi}{V} \right)^3
\]

where \(L\) is the size of the (compactified) spatial domain and \(V\) is its volume. The Hamiltonian for each mode can now be written as that for a simple harmonic oscillator,

\[
H_{1p} = \frac{\pi^2_{p,\pm}}{2m_p} + \frac{m_p}{2} (p + P)^2 \vert \phi_{p,\pm} \vert^2
\]

\[
H_{2p} = \frac{\pi^2_{p,\pm}}{2m_p} + \frac{m_p}{2} (p - P)^2 \vert \phi_{p,\pm} \vert^2
\]

where \(\pi_{p,\pm}\) and \(\Pi_{p,\pm}\) are conjugate momenta to \(\phi_{p,\pm}\) and \(\psi_{p,\pm}\), and

\[
m_p = 2 (\Delta p)^3 = 2 \left( \frac{2\pi}{L} \right)^3
\]

Note that \(m_p\) has dimensions of mass squared (not mass).

As noted at the end of Sec. III, we will also have \(H_{3p}\) and \(H_{4p}\) corresponding to the \((\beta^p_p, \gamma^p_p)\) sector.

The Hamiltonians in (38) and (39) are those of simple harmonic oscillators with time dependent frequencies,

\[
\omega_{\phi_{p,\pm}} = + \sqrt{(p \pm P)^2 + \pi^2_{p,\pm}}
\]

\[
\omega_{\psi_{p,\pm}} = + \sqrt{(p \pm P)^2 + \pi^2_{p,\pm}}
\]

The + signs in the pre-factor are to emphasize that we are taking the positive square root. The structure of the frequencies is easy to understand because the \(p_\pm \pm P\) follows from the contravariant derivative acting on the excitations. The different forms of \(\omega_{\phi_{p,\pm}}\) and \(\omega_{\psi_{p,\pm}}\) arise since the variables \(\phi_{p,\pm}\) are associated with \(\tilde{a}_p^{-}\), while \(\psi_{p,\pm}\) are associated with \(\tilde{a}_p^{+}\).

An important point for us is that there are certain modes for which the frequency vanishes. For example, \(\omega_{\phi_{p,\pm}} = |p - P|\) and, at any time, the frequency vanishes for \(p = P = gEt \cos \theta\). (Similarly for \(\omega_{\psi_{p,\pm}}\)). The frequencies \(\omega_{\phi_{p,\pm}}\) and \(\omega_{\psi_{p,\pm}}\) do not vanish for \(t > 0\). We will return to this point in our discussion of the particle number density in Sec. VII.

### V. CLASSICAL-QUANTUM CORRESPONDENCE

To obtain particle production of the fields \(\phi_{p,\pm}\), \(\psi_{p,\pm}\) we will use the CQC. (See Appendix A for a summary of the CQC.) Then the variables \(\phi_{p,\pm}\) and \(\psi_{p,\pm}\) are complexified and we solve the classical equations of motion

\[
\ddot{\phi}_{p,\pm} + (P_\pm \pm p_\perp)^2 \phi_{p,\pm} = 0
\]

\[
\ddot{\psi}_{p,\pm} + (P_\pm \pm p_\perp)^2 \psi_{p,\pm} = 0
\]

with initial conditions,

\[
\phi_{p,\pm}(t = 0) = -\frac{i}{\sqrt{2m_p}} \psi_{p,\pm}(t = 0)
\]

\[
\dot{\phi}_{p,\pm}(t = 0) = \sqrt{\frac{p}{2m_p}} \psi_{p,\pm}(t = 0)
\]

(Note that the simple harmonic oscillator frequencies at \(t = 0\) are simply \(p\) because \(P(t = 0) = 0\).) Using (40) we write

\[
\phi_{p,\pm}(0) = -\frac{i}{2} \left( \frac{L}{2\pi} \right)^{3/2} = \psi_{p,\pm}(0)
\]

\[
\dot{\phi}_{p,\pm}(0) = \frac{p}{2} \left( \frac{L}{2\pi} \right)^{3/2} = \dot{\psi}_{p,\pm}(0)
\]
To calculate the energy density in excitations, we simply need to evaluate the classical energy in the complexified \( \phi_{p,\pm} \) and \( \psi_{p,\pm} \) as we describe next.

VI. ENERGY DENSITY PRODUCTION

The energy in the complexified variables \( \phi_{p,\pm} \) \( \psi_{p,\pm} \) follows from (34) and (35).

\[
\mathcal{E}_1 = \int \frac{d^3p}{E_p} \left[ |\dot{\phi}_{p,+}|^2 + (p + P_z)^2 |\phi_{p,+}|^2 + |\dot{\phi}_{p,-}|^2 + (p - P_z)^2 |\phi_{p,-}|^2 \right] (49)
\]

\[
\mathcal{E}_2 = \int \frac{d^3p}{E_p} \left[ |\dot{\psi}_{p,+}|^2 + ((p_z + P)^2 + p_z^2) |\psi_{p,+}|^2 + |\dot{\psi}_{p,-}|^2 + ((p_z - P)^2 + p_z^2) |\psi_{p,-}|^2 \right] (50)
\]

The energies in the \((\beta^*_p, \gamma^*_p)\) sector give identical expressions and we will include these in the end in the total energy by multiplying by a factor of two.

The energy expressions in (49) and (50) include the ground state energy – the \(\omega/2\) of the simple harmonic oscillator – whereas we are interested in the energy of the excitations only. As described in Appendix A the ground state energy can be discarded by writing the energies as,

\[
\mathcal{E}_1 = \int \frac{d^3p}{E_p} \left[ |\dot{\phi}_{p,+} - i\omega\phi_{p,+} + \phi_{p,+}|^2 + |\dot{\phi}_{p,-} - i\omega\phi_{p,-} + \phi_{p,-}|^2 \right] (51)
\]

\[
\mathcal{E}_2 = \int \frac{d^3p}{E_p} \left[ |\dot{\psi}_{p,+} - i\omega\psi_{p,+} + \psi_{p,+}|^2 + |\dot{\psi}_{p,-} - i\omega\psi_{p,-} + \psi_{p,-}|^2 \right] (52)
\]

The total energy in the excitations is

\[
\mathcal{E} = 2(\mathcal{E}_1 + \mathcal{E}_2) = 8\pi \int_0^1 dp \int dp \int_0^1 du \times \left[ |\dot{\phi}_{p,+} - i\omega\phi_{p,+} + \phi_{p,+}|^2 + |\dot{\phi}_{p,-} - i\omega\phi_{p,-} + \phi_{p,-}|^2 + |\dot{\psi}_{p,+} - i\omega\psi_{p,+} + \psi_{p,+}|^2 + |\dot{\psi}_{p,-} - i\omega\psi_{p,-} + \psi_{p,-}|^2 \right] (53)
\]

where the factor of two in the first line accounts for the excitations in the \((\beta^*_p, \gamma^*_p)\) sector, \(u = \cos \theta\), and we have used the symmetry under \(u \rightarrow -u\) to restrict \(u\) to the interval \([0, 1]\). We remark that the expressions occurring in the integrands of (51) and (52), such as \(\dot{\phi}_{p,+} - i\omega\phi_{p,+} + \phi_{p,+}\), are the usual Bogolyubov \(\beta\) coefficients up to a factor of \(1/\sqrt{\omega}\), where \(\omega\) stands for the frequency.

Now all that is required is to solve the equations of motion in (43) and (44), insert the solutions in the energy expressions above, and perform the integrations. The first step is formally accomplished since the solutions to the equations of motion can be written in terms of parabolic cylindrical functions. However, we have found it more practical to solve the differential equations numerically followed by numerical integration.

A. Numerical evaluation of the energy

We have numerically solved the differential equations in (43) and (44) for \(u \in [0, 1]\) and \(p \in [0, p_c]\) where \(p_c \gg gE t\) is a momentum cutoff. As we solve the differential equations, we also numerically evaluate the energy in (53). In the numerical computations we choose units so that \(gE = 1\).

We define the spectral energy density, \(\rho(p, t)\), via

\[
\frac{\mathcal{E}}{L^3} = \int_0^\infty dp \rho(p, t) (54)
\]

or explicitly,

\[
\rho(p) = \frac{8\pi p}{L^3} \int_0^1 du \sum_{s=\pm} \left[ |\dot{\phi}_{p,s} - i\omega\phi_{p,s} + \phi_{p,s}|^2 + |\dot{\psi}_{p,s} - i\omega\psi_{p,s} + \psi_{p,s}|^2 \right] (55)
\]

In Fig. 1, we plot the spectral energy density rescaled by \(t^{-3}\), i.e., \(t^{-3} \rho(p, t)\), as a function of \(p/t\) at three different times. It is clear that the peak of the spectrum \(\rho\) grows as \(t^3\) and the width grows as \(t\), implying that the total energy in excitations grows as \(t^4\).

The linear growth of the width of \(\rho(p, t)\) follows from the form of the oscillation frequencies in (41) and (42).
For very large values of $p$, the time dependence of the frequencies can be ignored. Then the solutions for $\phi_{p,\pm}$ and $\psi_{p,\pm}$ are simply trigonometric functions for which there is no contribution to the energy. Hence there is no particle production for $p \gg gEt$: there is only particle production for $p \lesssim gEt$ and so the width in $p$ contributing to particle production grows linearly in $t$. To understand the growth of the peak of $\rho(p,t)$ that goes as $t^3$, we note that the peak is located at $p \approx gEt/2$. Thus the $p$ prefactor in (55) contributes one factor of $t$. In the integrand, the initial conditions for the variables $\dot{\phi}_{p,\pm}$ and $\dot{\psi}_{p,\pm}$ are proportional to $p$ as in (48) and since the $p$ that contributes to the energy integral grows proportional to $t$, and the variables enter quadratically in the energy integral, the peak of the spectral energy density grows as $t^3$.

The $t^3$ growth of the total energy density is further confirmed in Fig. 2 where we plot the total energy density divided by $t^3$ vs. time.

To obtain the total energy density, $: E :$, we have integrated over $p \in [0,p_c]$ where $p_c \gg gEt$ is a cutoff. To study the dependence of our result on the cutoff, we zoom into the large $p$ behavior of $\rho(p,t)$, shown in Fig. 3. This gives $\rho(p,t) \propto 1/p$ at large $p$ and the total energy diverges logarithmically as the cutoff $p_c$ is taken to infinity. However this ultraviolet divergence will be controlled once asymptotic freedom is taken into account. To see this in more detail, note that $\rho(p,t)$ in Fig. 1 has a dominant peak structure followed by the $1/p$ fall off, which after integration over $p$, lead to an asymptotic contribution to the energy density given by $g^2 \log(p_c/M)$, where $M$ is a renormalization scale. However the asymptotic value of the coupling constant at the cut-off scale evolves from its value $g_M$ at the renormalization scale as [40],

$$ g^2 = \frac{g^2_M}{1 + g^2_M \log(p_c/M)} , \quad (56) $$

implying,

$$ \lim_{p_c \to \infty} g^4 \log(p_c/M) \to 0 . \quad (57) $$

Hence the $1/p$ tail contribution to the integration vanishes in the $p_c \to \infty$ limit once we take the running of $g^2$, i.e. asymptotic freedom, into account.

VII. NUMBER DENSITY PRODUCTION

The number density of produced particles, $n$, is ambiguous at intermediate times though there is no ambiguity in the infinite time limit if the electric field is switched off [39]. Here we adopt the definition that the number density is given by the energy density for each mode divided by the frequency of that mode. Then using (53) we have,

$$ n = \frac{8 \pi}{\int_0^\infty dp \int_0^1 du} \sum_{s=\pm} \left[ \frac{\left| \phi_{p,s} - i \omega \phi_{p,s} \phi_{p,s} \right|^2}{\omega \phi_{p,s}} + \left| \psi_{p,s} - i \omega \psi_{p,s} \psi_{p,s} \right|^2}{\omega \psi_{p,s}} \right] \quad (58) $$

The expressions for the frequencies are given in (41) and (42). The issue is that at any given time, there is a momentum mode for which the frequency vanishes. Thus the integrand in (58) is singular at all times.

To determine if the singularity is integrable, we focus our attention on the $\phi_{p,-}$ sector for which the number density is

$$ n_{\phi,-} = \frac{8 \pi}{\int_0^\infty dp \int_0^1 du} \left| \frac{\phi_{p,-} - i |p - gEt|}{|p - gEt|} \right|^2 \quad (59) $$

First consider the numerator of the integrand. For $p = gEt$ it is simply $|\phi_{p,-}|^2$. Non-vanishing energy
in the \( \phi_{p,-} \) excitations implies that \( |\dot{\phi}_{p,-}|^2 \neq 0 \). So the singularity structure of the integral is

\[
n_{\phi,-} = 8\pi \langle p|\dot{\phi}_{p,-}|^2 \rangle \int_0^\infty dq \int_0^1 du \frac{1}{|q-u|} \quad (60)
\]

where \( q \equiv p/gEt \) and \( \langle \cdot \rangle \) denotes an effective value along the singular curve, \( p = gEtu \) (or \( q = u \)), in the integration plane. By transforming integration variables to the singular curve, \( x_\pm = q \pm u \) it is clear that the integral is logarithmically divergent due to the singularity along \( x_- = 0 \).

We note that the divergence in particle number density arises at low energy where the frequencies vanish. It is worth emphasizing that this is a particular behavior of the massless theory and will not be present in massive cases such as QED. In the full interacting theory we can expect this infrared divergence to be resolved due to confinement effects. At such low frequencies, the “soft” gluons are confined and are only present in massive glueball states. In our analysis, as discussed in Sec. II, we are examining where the road goes when we ignore confinement effects. It is interesting that the calculation without confinement leads to a divergent number density of excitations, suggesting a self-inconsistency.

\section{VIII. ADIABATIC CASE}

Often in the Bogoliubov method, the background is taken to turn on adiabatically, survive for a certain time period, and then slowly turn off. Then the asymptotic vacua are unambiguously defined and the total energy density of particles produced is evaluated at \( t \to \infty \).

We have also treated the adiabatic case. Now our choice for \( A(t) \) is,

\[
A(t) = E\tau \left[ \tanh \left( \frac{t - t_E}{\tau} \right) - \tanh \left( \frac{-t_E}{\tau} \right) \right] \quad (61)
\]

where \( t_E \) is some large time at which the electric field is maximum and \( \tau \) is the duration for which the field is turned on. The electric field is given by,

\[
\dot{A} = E \text{sech}^2 \left( \frac{t - t_E}{\tau} \right) \quad (62)
\]

Our analysis from the previous sections remains unchanged except that the frequencies in (63) and (64) now become

\[
\omega_{\phi,p,\pm} = \pm \sqrt{(p \pm gA(t)u)^2} \quad (63)
\]

\[
\omega_{\psi,p,\pm} = \pm \sqrt{(p_\pm gA(t))^2 + p_\perp^2} \quad (64)
\]

We have repeated the numerical analysis in this case and show the results for the energy density versus \( \tau \) in Fig. 4.

Once again we find : \( \dot{E} \propto \tau^2 \).

We can also consider the number density of particles in the adiabatic case. From Eq. (63) we see that the mode with \( p = gA(\infty)u \) has vanishing frequency \( \omega_{\phi,p,-} \) as \( t \to \infty \). Similarly the mode with \( p_\perp = gA(\infty) \), \( p_\perp = 0 \) has vanishing frequency \( \omega_{\psi,p,-} \) as \( t \to \infty \). Therefore the singularity discussed in Sec. VII is present even with the short duration electric field of (62).

\section{IX. CONCLUSIONS}

We have considered the fate of a uniform color electric field in a pure gauge Yang-Mills theory with SU(2) gauge group\(^5\). Our approach assumes initial conditions with a uniform electric field and quantum excitations in their non-interacting ground state. We then truncate the excitations to quadratic order in the action and evolve the system using the CQC.

We find that there is rapid particle (gluon) production for which we are able to characterize the energy spectrum and its time dependence as shown in Fig. 1. Production occurs for modes in a range \( p \in [0, gEt] \) with peak production at \( p \approx gEt/2 \). The amount of energy produced in an interval \( dp \) grows as \( t^3 \). We can understand this growth in terms of phase space factors that give one power of \( t \), and the square of the amplitude of vacuum fluctuations that are proportional to \( p^2 \), and hence grow as \( t^2 \). The \( t^3 \) dependence we find is in sharp contrast to usual Schwinger pair production, as in (1), for which the energy grows linearly with time. However, in contrast to the usual Schwinger pair productions of massive particles, gauge excitations are massless and this may be sufficient to explain the different production rates.

To the order in which we perform our calculations, the coupling constant does not run with energy scale. However, we encounter an ultraviolet divergence in our calculation of the energy produced in excitations. We have argued that the divergence would get controlled if we were to properly account for the running of the coupling

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\(^5\) The gauge group is not important for our analysis since SU(N) models have SU(2) subgroups.
constant at high energies (asymptotic freedom). In calculating the particle number density we also encounter an infrared divergence that we argue will be absent in the confining theory. In other words, if we ignore confinement, our analysis implies that the particle number density diverges, which we interpret as an indication of a lack of self-consistency of the unconfined assumption.

Our analysis leaves open several directions of interest such as the backreaction of particle production on the background color electric field. If we simply use energy conservation as a guide, the background electric field has an energy density proportional to \(E^2\) whereas the energy density in excitations grows as \(g^4E^4t^4\). Equating these two gives the decay time for the electric field to be \(\tau \sim (g\sqrt{E})^{-1}\). However, we cannot exclude the possibility that the electric field will be antiscreened as argued for non-Abelian gauge theories (see for example Sec. 16.7 of [10]), in which case the electric field strength would actually increase with time. We hope to investigate this issue by a more detailed analysis in the future.

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**Appendix A: Classical-Quantum Correspondence**

The CQC is equivalent to the Bogoliubov transformation method for calculating particle production in the case of a time dependent but spatially homogeneous background. Once the model is discretized, each mode is time dependent and spatially homogeneous. So we only need to demonstrate the method for a simple harmonic oscillator.

Consider the Hamiltonian for a simple harmonic oscillator of mass \(m\) – which can also be time dependent in the general case – with an arbitrary time dependent frequency \(\omega(t)\),

\[
H = \frac{p^2}{2m} + \frac{m}{2} \omega^2(t)x^2
\]  

(A1)

Then the quantum operators \(x\) and \(p\) can be written as

\[
x = z^a a_0 + z a_0^\dagger, \quad p = m(\dot{z}^a a_0 + \dot{z} a_0^\dagger)
\]  

(A2)

where \(a_0\) and \(a_0^\dagger\) are the initial annihilation and creation operators defined in terms of the initial position operator \(x_0\) and momentum operator \(p_0\),

\[
a_0 = \frac{p_0 - im\omega_0 x_0}{\sqrt{2m\omega_0}}, \quad a_0^\dagger = \frac{p_0 + im\omega_0 x_0}{\sqrt{2m\omega_0}}
\]  

(A3)

where \(\omega_0\) is the frequency at the initial time, and \(z(t)\) is a complex-valued c-number function of time. Using the Heisenberg equations for \(x\) and \(p\) we find that \(z(t)\) must satisfy

\[
\ddot{z} + \omega^2 z = 0
\]  

(A4)

The initial conditions follow from the above relations and are,

\[
z_0 = \frac{-i}{\sqrt{2m\omega_0}}, \quad \dot{z}_0 = \sqrt{\frac{\omega_0}{2m}}
\]  

(A5)

The Wronskian is constant and fixed by the initial conditions,

\[
m(z^* \dot{z} - z \dot{z}^*) = i
\]  

(A6)

Other quantum operators can be re-written in terms of \(z(t), x_0\) and \(p_0\). Expectation values are written entirely in terms of \(z(t)\). In particular, the energy is

\[
\langle H \rangle = \frac{1}{2} m |\dot{z}|^2 + \frac{1}{2} m\omega^2 |z|^2
\]  

(A7)

This can also be written as

\[
\langle H \rangle = \frac{m}{2} |\dot{z} - i\omega z|^2 - \frac{i}{2} m\omega(z^* \dot{z} - z \dot{z}^*)
\]  

(A8)

The second term is the Wronskian and hence is a constant determined by the initial conditions. In the quantum simple harmonic oscillator this term corresponds to the ground state energy \(\omega/2\). So the energy in the excitations is given by

\[
\mathcal{E} := \frac{m}{2} |\dot{z} - i\omega z|^2
\]  

(A9)

which is what we use in [51] and [52].
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