

Self-Energy and Interaction Effects in Graphene from a Quantum Field Theory Approach

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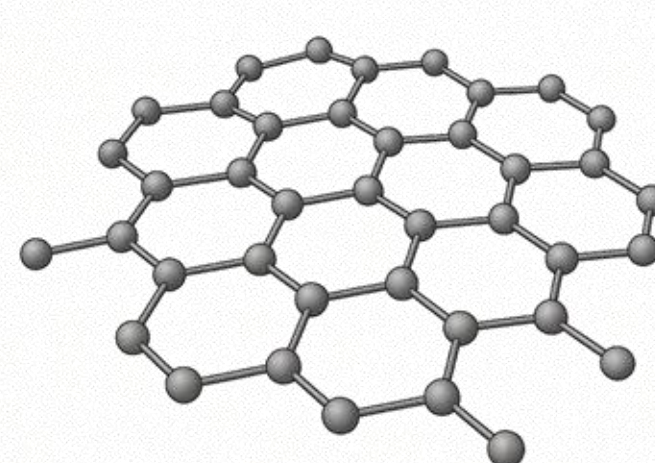
Abstract

- We perform a **Quantum Field theoretical study** of **Dirac fermions in graphene**, focusing on:
 - their electromagnetic interactions
 - emergence of a mass gap via symmetry breaking
- Using an **effective Lagrangian and Dimensional Regularization (DR)**, we compute one-loop corrections to the **electron propagator**.
 - A mass term is introduced to model **deviations from graphene's gapless behavior**.
- Our results show momentum-dependent corrections to the **Fermi velocity** and demonstrate conditions under which a **finite energy gap** forms.

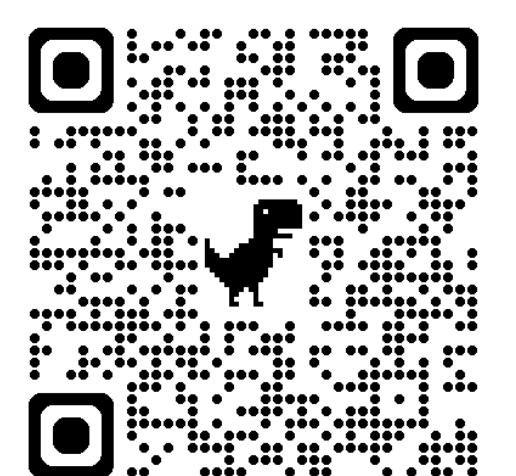
Motivation

- Our goal is to develop a comprehensive framework that applies **Quantum Field Theory (QFT) on a spacetime lattice** to model, renormalize, simulate, and validate **quantum phenomena in graphene-based and other condensed matter systems**.*
- We have employed our software **SUPER_QUANTA** to explore applications in lower-dimensional QFTs, most notably the study of graphene within the framework of **Quantum Electrodynamics (QED) in 2+1 dimensions**. For information on our previous software, **FEDILA**, scan the QR code below:

GRAPHENE



Graphene is a one-atom-thick sheet of carbon atoms that behaves as a two-dimensional material in many applications.



Theoretical QFT Model

- We extend **Quantum Electrodynamics (QED)** models to explore gap-related phenomena in condensed matter systems (see, e.g., *Phys. Rev. D* **109** (2024) 125014 by M. Bordag et al., *Nucl. Phys. B* **424** (1994) 595 by J. González et al., and references therein). A representative continuum formulation involves the following **Lagrangian density in Euclidean space** (in units $\hbar = c = 1$):

$$\mathcal{L}_{eff} = \bar{\psi}(i\tilde{\gamma}_\mu D_\mu - m v_F^2)\psi \delta(z) - \frac{1}{4} F_{\mu\nu} F_{\mu\nu},$$

- $\psi(x, y, z, t)$: Dirac spinor field, $\tilde{\gamma}_\mu$: modified gamma matrices to incorporate the Fermi velocity v_F , with spatial $\tilde{\gamma}_i = v_F \sigma_i$ ($i = 1, 2$) and temporal $\tilde{\gamma}_t = \gamma_t = \sigma_3$, $D_\mu = \partial_\mu - ieA_\mu$: covariant derivative including the interaction with the photon field A_μ , e : electric charge, m : mass term introduced to explore symmetry breaking and potential gap formation, $F_{\mu\nu} = \partial_\mu A_\nu - \partial_\nu A_\mu$: electromagnetic field strength tensor and $\delta(z)$: Dirac delta function, ensuring fermions are confined to the 2D graphene sheet at $z = 0$. In the limit $v_F = c$, the momentum integrals appearing in the fermion self-energy, $\Sigma_\psi(q, m)$, are automatically regularized in $D = 3 - 2\varepsilon$ dimensions (ε is the regulator).

Preliminary Calculations ($v_F = c = 1$) and Future Plans

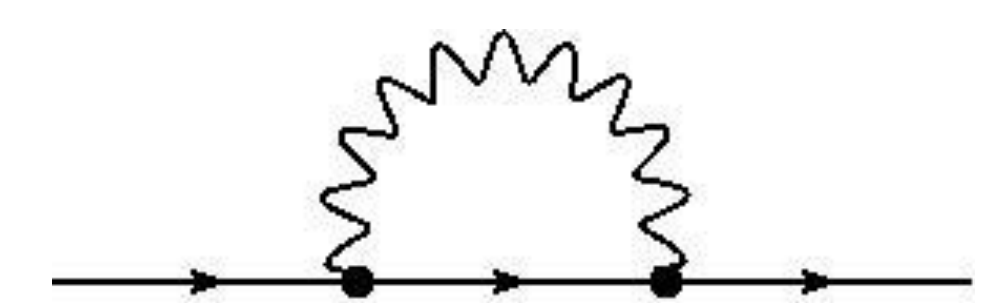
- The **electron self-energy is computed from the Feynman diagram of Fig. 1** in the naive approximation $v_F = c$ and is given by (ξ is the gauge parameter and $q_\mu = (E, q_x, q_y)$ is the external three-vector energy-momentum in Fig. 1):

$$\Sigma_\psi(q, m) = \frac{e^2}{18\pi q^3} \left[2(\xi + 2) m q^2 \arcsin\left(\frac{q}{\sqrt{m^2 + q^2}}\right) + i \gamma_\mu q_\mu \xi \left(m q + (q^2 - m^2) \arcsin\left(\frac{q}{\sqrt{m^2 + q^2}}\right) \right) \right] + O(e^4).$$

- In pristine graphene, the excitations exhibit a linear dispersion relation given by: $E = v_F |q|$, and the quantum corrections to $-i \Sigma_\psi(q, 0) = \frac{e^2}{16} \xi \begin{pmatrix} E & q_x + iq_y \\ q_x - iq_y & -E \end{pmatrix} \frac{1}{\sqrt{E^2 + k_x^2 + k_y^2}}$, modify the **Fermi velocity**: $v_F \rightarrow v_F \pm \frac{e^2}{16} \xi \frac{1}{|q|}$.

- The **energy gap ΔE** is proportional to the part of the self-energy which is not proportional to $i \gamma_\mu q_\mu$ at zero momentum:

$$\lim_{q \rightarrow 0} \Sigma_\psi(q, m) \Big|_{\gamma_\mu q_\mu = 0} = \frac{e^2}{4\pi} (2 + \xi) \Rightarrow \Delta E \approx \frac{e^2}{4\pi} (2 + \xi).$$



One-loop Feynman diagram contributing to the $\Sigma_\psi(q, m)$. A wavy (solid) line represents photons (electrons).

- **Future work** regards the computation of quantum corrections in the more realistic case where $v_F \neq 1$, and calculations within lattice regularization to simulate the theory.