Intrinsic dissipative Floquet superconductors beyond mean-field theory

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We study the intrinsic superconductivity in a dissipative Floquet electronic system in the presence of attractive interactions. Based on the functional Keldysh theory beyond the mean-field treatment, we find that the system shows a time-periodic bosonic condensation and reaches an intrinsic dissipative Floquet superconducting (SC) phase. Due to the interplay between dissipations and periodic modulations, the Floquet SC gap becomes “soft” and contains the diffusive fermionic modes with finite lifetimes. However, bosonic modes of the bosonic condensation are still propagating even in the presence of dissipations.

I. INTRODUCTION

Periodic driving schemes provide a simple way to study systems out of equilibrium \cite{1–3, 4–33}. An interesting example is the Floquet topological superconductor \cite{11, 13, 14, 23, 26, 34}. Non-equilibrium superconductivity, including the enhancement of superconductivity due to non-equilibrium electrons \cite{34–41} or non-equilibrium phonons \cite{42–47}, and dynamics of Cooper correlations due to time-dependent interactions \cite{48–57}, has been widely studied and obtained great exciting results. In this paper, however, we want to understand Floquet structures and their behaviors in the presence of dissipations, which is unavoidable and a sensitive factor to Floquet engineering. Floquet superconductivity can be induced in two different routes: proximity-induced SC and intrinsic SC. The proximitized SC provides not only Cooper correlations but also strong dissipations, which could significantly change the behavior of Floquet systems \cite{58, 59}. For the intrinsic case, SC is created due to the interaction instability near the Fermi surface or other strong interaction effects \cite{60, 61}. Those interaction instabilities and related dissipations could be significantly modified by the periodic driving potential in Floquet engineering; and so, the Floquet treatment, i.e. Floquet theorem for quadratic BCS mean-field Hamiltonian, could be unreliable in the Floquet engineering. Therefore, a careful self-consistent treatment of all critical factors, i.e. periodical driving force, interaction instabilities and dissipations, should be considered for Floquet engineering. Based on those motivations, we focus on the questions: can the interaction cause fermion-to-boson transition in the dissipative Floquet systems; and how well is the Floquet BCS mean-field treatment in describing periodically driven intrinsic SC?

In this paper, we study Cooper instability for a realistic periodically driven electronic system with interactions and dissipations. Based on the functional Keldysh field theory \cite{62, 63}, we consider both the stationary point analysis and the Gaussian fluctuation, which is beyond the mean-field theory. We show that the system develops a periodic bosonic condensation, and reaches a dissipative Floquet superconducting phase below a critical value $\gamma_c$ or $T_c$, where $\gamma$ is the system-bath coupling and $T$ is the temperature of the bath. The fermionic quasiparticle shows a “soft” energy gap, acquires a finite lifetime and becomes diffusive. However, bosonic modes of the condensation are still propagating even in the presence of dissipations. In addition, we also find that the oscillation amplitude of the order parameter is a non-monotonic function of dissipation; and therefore, a certain finite dissipation will be helpful for the Floquet SC.

Summary of the Treatment

In order to avoid bringing confusions to readers and make our manuscript easy to read, we summarize the treatment we used in this manuscript. Our treatment is a generalization of the method used in the equilibrium superconducting case, to the dissipative Floquet system. The first step is to write down the Hamiltonian Eq.(1) of the composite system(electronic system with periodic driving + normal metal bath) under consideration. We then apply a time-dependent unitary transformation shown in Eq.(2) to obtain an equivalent time-independent electronic system (then we can directly use the knowledge of the static electronic system). We focus on the possibility of dissipative Floquet superconductivity in the presence of the bath. We further apply the functional Keldysh field theory \cite{63} to obtain the action Eq.(3) from the Hamiltonian Eq.(2); and this treatment is a standard routine (see Appendix A). Next, we use the Hubbard-Stratonovich transformation to decouple the four-fermion attractive interaction by introducing an auxiliary bosonic field, and then integrate out fermionic degrees of freedom using Gaussian integrals, arriving at an effective bosonic field theory Eq.(6) (see

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Appendix A). Those are formulated in Sec.II. After that, based on the effective bosonic theory, we do the stationary point analysis in Sec.III and the Gaussian fluctuation approximation, which is beyond the mean-field theory, in Sec.IV. Generalized Feynman diagram rules are developed in Sec.IV to facilitate the analysis.

II. INTRINSIC DISSIPATIVE FLOQUET SC

We consider a single-band electronic system with a time-periodic chemical potential and attractive interactions, coupling to a normal fermionic bath. The Hamiltonian of the whole system can be written as

$$H(t) = H_D + H_{int} + H_T + H_B,$$

where $H_D = \sum_{k\sigma} [\epsilon_k - \mu_0 - \mu(t)] c_{k\sigma}^\dagger \text{c}_{k\sigma}$ describes the non-interacting electronic system with the time-periodic chemical potential $\mu_0 + \mu(t)$, where $\mu(t) = -K \cos(\Omega t)$ with $\Omega$ the driving frequency, and $c_{k}^\dagger$ ($c_k$) is the fermionic creation (annihilation) operator. $H_{int} = -g \sum_{k_1 \ldots k_n} c_{k_1}^\dagger a_{q_1}^\dagger \ldots c_{k_n}^\dagger a_{q_n}$ with $g > 0$ describes the attractive interaction. Here, $H_B = \sum_{kq} a_{q}^\dagger a_q$ is the fermionic bath Hamiltonian, which provides dissipations; and $a_{q}^\dagger$ is the creation (annihilation) operator of the bath. Such a bath is necessary for a driving interacting system to avoid the featureless infinite-temperature state [64] and thermalize to a non-trivial phase [3]. One can imagine either an unavoidable dissipation resources or a large engineered equilibrium system weakly coupled to the small driven part; and the bath is in equilibrium state with temperature $T$. The system-bath coupling term can be written as

$$H_T = W \sum_{kq} \left( c_{k\sigma}^\dagger a_{q\sigma} + h.c. \right)$$

with $W$ being the coupling strength.

For a periodically driven system, it is convenient to consider a rotating frame [45, 58, 65, 66] by using of a time-dependent unitary transformation $U_F = \exp(-i f(t) \sum_{k\sigma} c_{k\sigma}^\dagger \text{c}_{k\sigma})$ with $df/dt = -\mu(t)$, which results in an equivalent system in the rotating frame

$$H_F(t) = U_F^\dagger (H(t) - i \partial_t) U_F$$

$$= \sum_{k\sigma} (\epsilon_k - \mu_0) c_{k\sigma}^\dagger \text{c}_{k\sigma} + H_{int}$$

$$+ W \sum_{kq} \left( e^{i f(t)} c_{k\sigma}^\dagger a_{q\sigma} + h.c. \right) + H_B.$$  \hspace{1cm} (2)

In the absence of the bath, that is $W \to 0$, we reach a time-independent system with equilibrium superconductivity, because the interaction term fully commutes with our periodic driven chemical potential term (thus commutes with $U_F$). Note that the original electron-phonon interaction also commutes with the periodic driven term, thus the phonon-induced attractive interaction is unchanged under the periodic driving. In those cases, without coupling to a bath, the periodic driving is trivial, and can be removed using a time-dependent unitary transformation. It is also our purpose to consider such a simple model that the periodical driving potential cannot simply spoil the formation of the SC correlation for the analysis of Floquet SC. In the presence of the bath with finite $W$, the superconductivity can be modified by the time-dependent system-bath coupling as shown in Eq.(2).

Then, the standard procedure leads to the total action of the system in the closed time contour [63]:

$$S = \int_c \int_c^{t'} \sum_k \tilde{\Psi}_{sk}(t) \tilde{Q}_{0k}^{-1}(t - t') \tilde{\Psi}_{sk}(t') + S_{int}$$

$$+ \int_c \int_c^{t'} \sum_q \tilde{\Psi}_{sq}^{-1}(t) \tilde{Q}_{0q}^{-1}(t - t') \tilde{\Psi}_{sq}(t')$$

$$+ \int_c \int_c^{t'} \sum_{kq} \left[ \tilde{\Psi}_{sk}(t) \tilde{M}(t) \tilde{\Psi}_{sq}(t) + h.c. \right],$$

where $\tilde{\Psi}_{sk} = [\tilde{\psi}_{k\uparrow}, \tilde{\psi}_{-k\uparrow}]$; $\tilde{\Psi}_{sq} = [\tilde{\phi}_{q\uparrow}, \tilde{\phi}_{-q\uparrow}]$; $\tilde{Q}_{0k}$ and $\tilde{Q}_{0q}$ are free-fermion Green’s functions in Nambu space, $\tilde{M}(t) = \begin{bmatrix} We^{i f(t)} & 0 \\ 0 & -We^{-i f(t)} \end{bmatrix}$, and in $S_{int} = g \int_c \int_c \sum_{kk'} \tilde{\psi}_{k\uparrow} \tilde{\psi}_{-k\uparrow} \tilde{\psi}_{k'\uparrow} \tilde{\psi}_{-k'\uparrow}$, we only count in terms describing the interaction between electrons with opposite momenta and spins.

In order to study the non-equilibrium steady state in such a system, one can first integrate out the bath’s degrees of freedom with a Keldysh functional integral formalism [58, 63] to obtain an equivalent effective fermionic action:

$$S' = \int_c \int_c^{t'} \sum_k \tilde{\Psi}_{sk}^{-1}(t) \tilde{Q}_k^{-1}(t, t') \tilde{\Psi}_{sk}(t') + S_{int},$$  \hspace{1cm} (4)

where $\tilde{Q}_k(t, t')$ is the dressed Green’s function giving by the Dyson’s equation

$$\tilde{Q}_k(t, t') = \tilde{Q}_{0k}(t - t')$$

$$+ \int_{-\infty}^{+\infty} dt_1 dt_2 \tilde{Q}_{0k}(t - t_1) \tilde{\Sigma}_k(t_1, t_2) \tilde{Q}_k(t_2, t'),$$

with the self-energy from the bath being

$$\tilde{\Sigma}_k(t_1, t_2) = \sum_q \tilde{M}(t_1) \tilde{Q}_{0q}(t_1 - t_2) \tilde{M}(t_2).$$

These Green’s functions can be derived analytically through perturbative expansions in the small parameter $\kappa = K/\Omega$ (the driving potential reads as $\mu(t) = -K \cos(\Omega t)$) (see Appendix A for details).

For the four-fermion interaction, it is common to decouple them through the Hubbard-Stratonovich transformation [63, 67]. Such a procedure will introduce an auxiliary bosonic field, denoted as $\Delta$ here. Applying the Keldysh transformation for the bosonic fields and the Keldysh-Larkin-Ovchinnikov transformation for the fermionic fields [63], one turns the effective fermionic action—$S'$ into the Keldysh@Nambu space. Note that
now $S'$ is in a quadratic form with respect to the fermionic degrees of freedom, thus one can also integrate out fermionic fields and arrives at the effective bosonic action:

$$S_{\text{eff}} = -\frac{1}{g} \int dt \Delta^\alpha \sigma^\alpha_1 \Delta^{\beta}(t) - i \text{Tr} \left[ 1 - \bar{Q}_k (\gamma^\alpha \otimes \Delta^\alpha) \right], \tag{6}$$

where we have restricted the pairing between electrons with opposite momenta, thus $\Delta^\alpha$ here denotes the zero-momentum bosonic field; $\Delta^\alpha = (1/\sqrt{2}) (\Delta^\alpha \hat{r}_+ + \Delta^\alpha \hat{r}_-)$ with $\alpha \in \{c, \bar{c}\}$ introduced by the Keldysh rotation; $\hat{r}^\alpha = \hat{r}^\alpha_1 \hat{\bar{r}}_0$, $\hat{r}^\alpha = \hat{r}^\alpha_0 \hat{\bar{r}}_1$, and $\hat{r}_\pm = \hat{r}_1 \pm i \hat{r}_2$, with $\hat{r}_\mu (\mu = 0, 1, 2, 3)$ being the four Pauli matrices. One should not confuse the dissipation strength $\gamma$ with matrices $\hat{r}^\alpha$ in the Keldysh space and we use the hat symbol for $2 \times 2$ matrices acting in either Nambu or Keldysh spaces and the check symbol to denote $4 \times 4$ matrices acting in the Keldysh$\otimes$Nambu space.

In the non-equilibrium case, one will often encounter two-time functions, like $\bar{Q}_k(t, t')$. Here, we show how to transform them into the Floquet representation [58] widely used in our discussion.

Due to the periodic driving, those two-time functions will possess the discrete time-translational symmetry $Q(t, t') = Q(t + \tau, t' + \tau)$, where $\tau$ is the period. One can introduce two new variables $s \equiv t$ and $u \equiv t - t'$ and define the new function $Q(t, t') \rightarrow Q(s, u)$ which satisfies $Q(s + \tau, u) = Q(s, u)$ for all $u$. Thus, one can perform the Fourier transformation for $u$ and the Fourier series expansion for $s$:

$$Q(s, \omega) = \int_{-\infty}^{+\infty} du e^{i\omega u} Q(s, u), \tag{7}$$

$$Q(n, \omega) = \frac{1}{\tau} \int_{0}^{\tau} ds e^{-i\omega n s} Q(s, \omega). \tag{8}$$

Turning $Q(n, \omega)$ into a matrix form, known as the Floquet structure, one has

$$Q = \begin{bmatrix} \cdots & Q(0, \omega + \Omega) & Q(1, \omega) & Q(2, \omega - \Omega) \\ Q(-1, \omega + \Omega) & Q(0, \omega) & Q(1, \omega - \Omega) \\ Q(-2, \omega + \Omega) & Q(-1, \omega) & Q(0, \omega - \Omega) & \cdots \end{bmatrix}. \tag{8}$$

III. STATIONARY POINT ANALYSIS

In equilibrium, the stationary point analysis of the effective bosonic action is just the mean-field theory and provides the gap equation. Out of equilibrium, we don’t know what exactly the mean field is. However, we can always consider the stationary point, around which the action can be expanded perturbatively. A variation of the bosonic action Eq.(6) with respect to $\bar{\Delta}^{\alpha}(t)$ generates the stationary point equation:

$$\bar{\Delta}^{\alpha}(t) = i \frac{g}{\sqrt{2}} \text{Tr} \left[ (\hat{r}^\alpha \otimes \hat{r}^-_\alpha) \sum_k \bar{Q}_k \right], \tag{9}$$

where in the time domain,

$$\bar{Q}_k^{-1} (t, t') = \bar{Q}_k^{-1} (t + \tau, t' + \tau) - \hat{r}^\alpha \otimes \Delta^\alpha \delta(t - t'), \tag{10}$$

and $\bar{Q}_k(t, t')$ is the non-interacting fermionic Green’s function dressed by the self-energy of the bath in the Keldysh$\otimes$Nambu space as mentioned before. We want to ask: 1) if we have a stationary point solution for $\bar{\Delta}^{\alpha}(t)$; 2) if the solution describes the SC order parameter or “gap” of the dissipative Floquet superconductor. We will address the two questions below.

As the system is periodic in the time domain, it is natural to assume that $\bar{Q}_k \bar{\Delta}(t, t') = \bar{Q}_k \bar{\Delta}(t + \tau, t' + \tau)$, where $\tau$ is the period. Indeed, the validity of this ansatz will be confirmed later. Then, in the stationary point, the dominated field $\bar{\Delta}(t)$ is also periodic in time. The Fourier transformation and the Fourier series expansion shown in Eq.(7) lead to

$$\bar{\Delta}^{\alpha}_n = i \frac{g}{\sqrt{2}} \int \frac{d\omega}{2\pi} \text{Tr} \left[ (\hat{r}^\alpha \otimes \hat{r}^-) \sum_k \bar{Q}_k \Delta(n, \omega) \right], \tag{11}$$

where $\bar{Q}_k \Delta(n, \omega)$ is one matrix element of the Floquet matrix $-\bar{Q}_k$ which has an infinite-dimension structure,

![FIG. 1. These three figures are plotted under conditions: $\omega_D/\Omega = 10, \Omega = 100K, T = 0K, gp_F = 0.2$ and $\kappa = 0.4$. Here $\omega, \Delta_0, \gamma$, and the amplitude of $\Delta^\alpha(t)$ are all rescaled by being divided by $\Omega$. (a) The tendency of $\Delta_0$, i.e. the soft gap with the effective temperature $\gamma$. It scales as $\sqrt{\gamma_0/(\gamma_0 - \gamma)}$, where $\gamma_0$ is the transition temperature. This scaling suggests that the critical value is $1/2$ ;(b) The DOS of quasiparticles in our intrinsic dissipative Floquet superconductors; (c) The amplitude of the order parameter $\bar{\Delta}^{\alpha}(t)$ for different $\gamma$’s. Note that the amplitude of $\bar{\Delta}^{\alpha}(t)$ is non-monotonic, as the amplitude is proportional to $\gamma_0$, and $\Delta_0$ changes in the opposite direction with the changing of $\gamma$.](image)
where we define $\Delta^\alpha_n$ for the solution $\Delta^0_q(t)$ reaches a maximum. This comes from the competition of two effects resulted from the dissipation strength. On one hand, in the rotating frame, the periodicity of the order parameter comes from the system-bath coupling. Thus, we need $\gamma$ to be large in order that the fermionic system has a clear periodicity. On the other hand, we need $\gamma$ to be small in order that it will not kill the order parameter. Thus, the competition of these two effects results in the maximum value. From this physical interpretation, one can think that at this certain value of $\gamma$, the periodicity of the order parameter is most clear and stable, which makes it easier to detect this periodicity experimentally.

In fact, this stationary point analysis can be regarded as a Floquet BCS mean-field treatment in our self-consistent functional formalism. To obtain a more comprehensive understanding, let’s go beyond this mean-field treatment.

**IV. FORMATION OF PERIODIC BOSONIC CONDENSATION**

From the previous discussion, we find that the bosonic field in the stationary point has a non-zero value when $\gamma$ is below a critical value $\gamma_c$. One can think that the non-zero bosonic field results from the condensation of Cooper pairs, just as the equilibrium SC case. The fact that some bosonic condensation exists in a pure fermionic system implies a fermion-to-boson phase transition[70]. We now discuss how this bosonic condensation is formed.
in the dissipative Floquet systems.

In the vicinity of \( \gamma_c \), \( \Delta^{cl}(t) \) is small, which means one can expand \( S_{\text{eff}} \) around the critical point. Thus, we expand the Trln-term in \( S_{\text{eff}} \) in powers of \( \Delta^{cl}(t) \), and simply keep terms up to the second order in \( \Delta^{cl}(t) \). This can be easily achieved by using of the series expansion \( \ln(1+x) = -\sum_{n=1}^{\infty} \frac{1}{n} (-x)^n \). Then applying transformations shown in Eq.(7), one can obtain the Gaussian action in the frequency space

\[
S^{(2)} = \int_0^{\Omega} \frac{d\omega}{2\pi} \bar{\Delta}(\omega) \left[ -\frac{1}{2} \sigma^2 + \sum_k \bar{\Gamma}_k \right] \Delta(\omega), \tag{16}
\]

where to avoid redundancy, we have restricted the integration range in \([0, \Omega]\), as the first Floquet-Blochlouin zone(FBZ) \([59, 71]\), and the vector is defined as \( \bar{\Delta} = [\Delta^d, \Delta^a]^t \), and \( \Delta = [\Delta^d, \Delta^a] \). The Floquet matrix structure denoted by the underline has already defined in Eq.(8). Note that the identity in the Floquet space—\( \sum \) only has diagonal elements and all of them are \( \frac{1}{\gamma} \). The matrix elements \( \Gamma_k(n, \omega) \) of \( \Gamma_k \) comes from the Fourier transformation and the Fourier series expansion of \( \Gamma_k^{\alpha\beta}(t, t') \), which is defined as

\[
\Gamma_k^{\alpha\beta}(t, t') = \frac{i}{2} \text{Tr} \left[ \bar{Q}_{kp}(t, t') \gamma^\alpha \bar{Q} - k h(t, t') \gamma^\beta \right], \tag{17}
\]

with subscript \( p \) and \( h \) denoting particle and hole, respectively.

Then, the Green’s function of the bosonic field in the Floquet space can be defined as

\[
\mathcal{D} = -i \langle \Phi(\omega) \Phi(\omega) \rangle
\sim -i \left[ \begin{array}{cc}
\bar{\Delta}(\omega + \Omega) & \bar{\Delta}(\omega - \Omega)
\end{array} \right] \left[ \begin{array}{cc}
\bar{\Delta}(\omega + \Omega) & \bar{\Delta}(\omega - \Omega)
\end{array} \right],
\tag{18}
\]

where \( \langle \rangle \) describes the average with respect to the weight \( \exp(i\mathcal{S}^{(2)}) \). The function defined in the Floquet basis \( \Phi(\omega) = [\cdots, \bar{\Delta}^t(\omega + \Omega), \bar{\Delta}^t(\omega), \bar{\Delta}^t(\omega - \Omega), \cdots]^t \), and \( \bar{\Phi}(\omega) = [\cdots, \bar{\Delta}(\omega + \Omega), \bar{\Delta}(\omega), \bar{\Delta}(\omega - \Omega), \cdots] \). In the second equality, we keep terms up to the second order in \( \kappa \), and then truncate \( \mathcal{D} \) to a \( 3 \times 3 \) Floquet matrix(note that each matrix element of a Floquet matrix here is also a \( 4 \times 4 \) matrix in the Keldysh\( \otimes \)Nambu space). As will see in the following content, the transition temperature will be modified in this case, but not in the \( \mathcal{O}(\kappa) \) case discussed in Sec.III. The retarded part of one matrix element of the Floquet matrix \( \mathcal{D} \), defined as

\[
\mathcal{D}^{cl, q}(n, \omega) = -i \langle \Delta^{cl}(\omega + n\Omega) \Delta^{cl}(\omega) \rangle
\tag{19}
\]

following from Eq.(18) is enough for the following discussion. The \( (n, \omega) \) component of the Floquet matrix—\( \mathcal{D} \) has the following physical meaning: a bosonic excitation \( \Delta(\omega) \) originally has energy \( \omega \); due to the external driven field, it will absorb \( n \) parts of energy(\( n\Omega \)), and then becomes a bosonic mode with energy \( \omega + n\Omega \). If \( \mathcal{D}^{cl, q}(n, \omega) \) has a pole, then this process is inevitable, as poles of Green’s functions correspond to quasiparticle excitations. Owing to structures of the distribution function and Green’s functions \( Q_k(t, t') \), no poles exist when \( \omega \neq 0 \). At \( \omega = 0 \), this bosonic excitation is similar to the equilibrium counterpart. But, due to the periodic driving, two pairing electrons can form bosonic modes with energy \( n\Omega \), and generate the SC order parameter at higher harmonics. Then, the order parameter becomes periodic in time and has higher harmonics \( \Delta^{q}(t) = \sum_q \Delta_q e^{i\Omega Q} \) (i.e. the ansatz we made in Sec.III)

In order to confirm this structure, we still need to check that if the bosonic modes at different harmonics occur at the same phase transition.

We first consider the \( n = 0 \) component—\( \mathcal{D}^{cl, q}(0, 0) \), which will result in zero-energy modes as the equilibrium case. The pole equation for the bosonic Green’s function(see Appendix C) can be written as

\[
1 - g \sum_k \mathcal{D}^{cl, q}(0, 0) = 0, \tag{20}
\]

one will find the solution at zero temperature

\[
\gamma = \gamma_c \equiv \omega_D \frac{1}{\omega_D} e^{-\frac{\pi^2}{g_{P}F(1-2q)}}, \tag{21}
\]

which holds for \( \gamma \ll \Omega \ll \omega_D \) and terms are kept up to \( \mathcal{O}(\kappa^2) \) for deriving this. This result reduces to the SC transition temperature in equilibrium case for \( \kappa \to 0 \). For finite temperature, we numerically plot the phase diagram in Fig.2. Note that \( \mathcal{D}^{cl, q}(0, 0) \) can also be interpreted as the normalized attractive interaction constant, and thus the divergence suggests that two electrons with opposite momenta and spins will form a bound state, known as Cooper pair. Therefore, we can see that our dissipative Floquet system could still develop a clear fermion-to-boson phase transition with a modification in the transition temperature due to the driving field.

We now analyze Green’s functions \( \mathcal{D}^{cl, q}(\pm n, 0) \) for \( n \geq 0 \), which involves at most \( n \) order transition processes between different Floquet bands. For example, after absorbing \( m \) parts of energies(\( m\Omega \)), the bosonic mode

\[
\text{FIG. 2. Phase Diagram for } (\gamma, T). \text{ We choose } \omega_D/\Omega = 10, \Omega = 100K, \quad g_{P}F = 0.2 \text{ and } \kappa = 0.4. \text{ } T \text{ and } \gamma \text{ are rescaled by being divided by } \Omega.
\]
will transit from \(a\Omega\) to \((a + m)\Omega\) with \(|m| \leq n\). In that sense, we require perturbation calculations to keep terms up to \(O(\kappa^n)\).

Here, we develop diagram rules to facilitate this analysis. For simplicity, we will introduce diagram rules through an example.

When we study Cooper instability in the equilibrium case, the most important factor is the vertex of the two-electron correlation function, and under the random phase approximation (RPA), there is just one kind of vertices [67]. However, in our Floquet system, there will be more kinds of vertices, as particles can absorb or emit energies.

Comparing to the equilibrium case, one can observe that \(-\hat{D}^{cl,q}(n,\omega)\) is the so-called vertex. As for \(-\hat{D}^{cl,q}(0,0)\) shown in Appendix C, it has three terms and they represent different kinds of vertices. Briefly, all vertices can be classified into two classes. One represents the direct process without emitting or absorbing energies, known as the direct vertex, and the other describes the indirect process containing energy exchanges, known as the indirect vertex. For example, in \(-\hat{D}^{cl,q}(0,0)\), the first term represents the direct vertex, and the last two terms give indirect vertices.

As for direct vertices, they are just the same with those in the equilibrium case [67]. However, for indirect vertices, we should take absorbing and emitting processes into consideration, which only appears in the periodically driven system. They are described by \(\Gamma^{cl,q}(m,\omega)\) and we call them the energy exchange vertices. The index \(m\) denotes the number of energies absorbed or emitted by particles (the energy unit is \(\Omega\)), and they absorb energy if \(m > 0\) and emit energy if \(m < 0\). \(\omega\) denotes the sum of initial energies of the two scattering electrons. Through the analytical expression, one can find that indirect vertices are constructed from direct vertices and energy exchange vertices. As it should be, the total number of energy exchanges should be consistent with \(n\) in \(-\hat{D}^{cl,q}(n,\omega)\).

Following above descriptions, we now write down diagram rules for different vertices, which will significantly facilitate the calculation of \(\hat{D}^{cl,q}(n,\omega)\):

1. Attach \(g/[1 - g\hat{\Gamma}^{cl,q}(0,\omega)]\) to a direct vertex shown in Fig.3(a);
2. Attach \(\hat{\Gamma}^{cl,q}(m,\omega)(\sim \kappa^{|m|} + O(\kappa^{|m|+2}))\) to an energy exchange vertex shown in Fig.3(b);
3. The indirect vertex shown in Fig.3(c) is constructed from direct vertices and indirect vertices, thus one should attach

   \[
   \frac{g}{1 - g\hat{\Gamma}^{cl,q}(0,\omega)} \Gamma^{cl,q}(m_1,\omega) \frac{g}{1 - g\hat{\Gamma}^{cl,q}(0,\omega + m_1\Omega)} \times \Gamma^{cl,q}(m_1,\omega + m_2\Omega) \frac{g}{1 - g\hat{\Gamma}^{cl,q}(0,\omega + (m_1 + m_2)\Omega)}
   \]

   to an indirect vertex. Read from the left to the right.

4. As for the diagram of one complete process, we just need to sum up those relative direct vertices and indirect vertices. As an example, \(-\hat{D}^{cl,q}(0,0)\) can be expressed as Fig.3(d) with \(\omega = 0\), when we keep terms up to \(O(\kappa^2)\).

Having shown the diagram rules for Floquet vertices or the bosonic Green’s functions (In Appendix C, we also summarize the diagram rules and show more examples for clarity.), one can now turn to analyze \(\hat{D}^{cl,q}(n,0)\) used for demonstrating the periodicity of the bosonic condensation.

As we now want to study the general case, that is \(n\) can be arbitrary non-zero integers, terms should be kept up to \(O(\kappa^n)\). Therefore, suppose that we keep terms up to \(O(g^2\kappa^n)\) for simplicity. According to diagram rules discussed previously, we can observe that the number of direct vertices are no more than 2, as one direct vertex will contribute a factor \(g\). Thus, there is one energy exchange vertex at most. For \(\hat{D}^{cl,q}(\pm n,0)\) with \(n \neq 0\), there must be energy exchange vertices. Therefore, there is one energy exchange vertex and two direct vertices for

FIG. 3. Diagrams for different vertices. (a) Direct vertex; (b) Energy exchange vertex; (c) Indirect vertex; (d) Diagram for \(-\hat{D}^{cl,q}(0,\omega)\).
the diagram of $-\hat{D}^{cl,q}(\pm n, 0)$ in $\mathcal{O}(g^2 \kappa^n)$ case. Thus, it is easy to obtain

$$\hat{D}^{cl,q}(\pm n, 0) = -g^2 \hat{\Gamma}^{cl,q}(\pm n, 0) \left[ 1 - g \hat{\Gamma}^{cl,q}(0, \pm n \Omega) \right] \left[ 1 - g \hat{\Gamma}^{cl,q}(0, 0) \right],$$

where we have $\hat{\Gamma}^{cl,q}(\pm n, \omega) \equiv \sum_k \hat{\Gamma}^{cl,q}(\pm n, \omega)$. Since $\hat{\Gamma}^{cl,q}(\pm n, \omega) \sim \kappa^n + \mathcal{O}(\kappa^{n+2})$, $D^{cl,q}(\pm n, 0) \sim g^2 \kappa^n$. In the $\mathcal{O}(\kappa^n)$ case, terms that is proportional to $\kappa^{|m|}$ with $|m| < n$ are still under our consideration, and one can derive similar expressions for $\hat{D}^{cl,q}(m, 0)$ from the same procedure. Note that the pole of $\hat{D}^{cl,q}(\pm n, 0)$ is the same as that of $\hat{D}^{cl,q}(0, 0)$ shown in Eq. (20). We also note that the term $1 - g \hat{\Gamma}^{cl,q}(0, \pm n \Omega)$ in the denominator is always non-zero. For arbitrary higher order corrections of interaction constant $g$, the term $1 - g \hat{\Gamma}^{cl,q}(0, 0)$ always appears in their denominator as $\hat{D}^{cl,q}(\pm n, 0)$. Thus, all Green’s functions—$\hat{D}^{cl,q}(\pm n, 0)$—exhibit the same pole structure (effective transition temperature $\gamma_t$) for all different $n$’s. The discussion above suggests that the bosonic excitation is periodic in the time domain, and can be expanded as $\Delta(t) = \sum_n \Delta(n \Omega) e^{i n \Omega t}$. Due to the presence of the pole structure, the bosonic modes in the condensation are propagating and dissipillation even with system-bath coupling. This result also confirms the ansatz we made in Sec.III.

V. DISCUSSION AND SUMMARY

The DOS of quasiparticles in Fig.1(b) shows that in the finite $\gamma$, there still exists energy levels in the gap. Actually, this comes from the fact that the lifetime of quasi-particles are now finite and they are diffusive, which can be derived from the Green’s function of quasiparticles—$Q_k^B \Delta(0, \omega)$, Eq.(B6). Just like the broadening of the peak in the DOS of dissipative Floquet Majorana zero modes shown in Ref.[59]. One can think this is because in the presence of the bath, the fermionic degrees of freedom or electrons before the SC phase transition are diffusive, which can be found from the Green’s functions Eq.(B5). However, the surprising thing is the bosonic mode or condensation resulting from the superconducting phase transition is still a propagating mode in spite of the existence of dissipations, as the Green’s functions of these bosonic modes—$D^{cl,q}(\pm n, 0)$—do not contain imaginary parts in the denominator.

In summary, based on the functional Keldysh field theory with a self-consistent treatment of all building blocks of our system, we demonstrate that the BCS mean-field treatment in the dissipative Floquet case is equivalent to the stationary point analysis in the functional or path integral formalism, where one can always implement the stationary point analysis. Moreover, based on the Gaussian fluctuation approximation, which is beyond the mean-field theory, we also consolidate the validity of the BCS mean-field theory in the dissipative Floquet scenario. Note that this system possesses a more structured gauge $U(1)$ symmetry, which could be an interesting point to be discussed in the future.

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Appendix A: Obtaining the Bosonic Effective Action

In this section, we show some key steps that lead to the effective action Eq.(6).

We start from the total Hamiltonian in the rotating frame, shown in Eq.(2)

$$H_F(t) = \sum_{k\sigma} \left( \epsilon_k - \mu_0 \right) c_{k\sigma}^\dagger c_{k\sigma} + H_{int} + \sum_{kq\sigma} \left( \epsilon_{k+q}^F(t) c_{kq\sigma}^\dagger a_{q\sigma} + h.c. \right) + H_B,$$

where expressions of $H_{int}$ and $H_B$ can be found in the main text, below Eq.(1).

The next step is to write the action from this Hamiltonian, and the procedure is quite standard. For completeness, we start from the construction of the functional Keldysh field theory(see Ref.[63] for details). In the rotating frame, our system is governed by the Hamiltonian $H_F(t)$. The evolution of the density matrix can expressed as $\rho(t) = U_{t,-\infty} \rho(-\infty) U_{t,-\infty}^\dagger$, where $U_{t,t'} = \mathcal{T} \exp \left( -i \int_t^{t'} dt H_F(t) \right)$ is the evolution operator, and $\mathcal{T}$ is the time-ordering operator. The expectation value of some observable can be calculated through the generating function, defined as

$$Z[V] = \frac{\text{Tr} \left[ U_C[V] \rho(-\infty) \right]}{\text{Tr}[\rho(-\infty)]},$$

where expressions of $H_{int}$ and $H_B$ can be found in the main text, below Eq.(1).

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$$Z[V] = \frac{\text{Tr} \left[ U_C[V] \rho(-\infty) \right]}{\text{Tr}[\rho(-\infty)]},$$

where expressions of $H_{int}$ and $H_B$ can be found in the main text, below Eq.(1).
where \( U_C[V] \equiv U_{-\infty,+\infty}[V]U_{+\infty,-\infty}[V] \), and \( U_-[V] \) is generated by \( \hat{H}_F(t) \equiv \hat{H}_F(t) \pm \mathcal{O} V(t) \), where the plus(minus) sign refers to the forward(backward) part of the contour, Fig.4(a). \( V(t) \) is an auxiliary field, and will be set to zero after taking the derivative of \( Z[V] \) with respect to \( V(t) \).

Eq.(A2) implies that calculating the generating function is the key for deal with a many-body problem. The functional formalism(path integral formalism) is a useful method to rewrite the generating function. The standard procedure, which can be found in any quantum field theory textbook, is:

- Divide the closed time contour into \((2N - 2)\) intervals with length \( \delta_i \), Fig.4(b);
- Insert the revolution of unity in the coherent state basis(for fermions, we need the Grassmann number, denoted as \( \psi \) in the following, for help);
- Take the limit \( \delta_i \to 0 \), and then one will get the generating function \( Z = \int D[\bar{\Psi}, \Psi] \exp(iS) \) in the continuum limit. For our system, we have

\[
S = \int_c dt \int_c dt' \sum_k \bar{\Psi}_{sk}^\dagger(t) \hat{Q}_{sk}^{-1}(t-t') \bar{\Psi}_{sk}(t') + S_{int} \\
+ \int_c dt \int_c dt' \sum_q \bar{\Psi}_{aq}^\dagger(t) \hat{Q}_{aq}^{-1}(t-t') \bar{\Psi}_{aq}(t') \\
+ \int_c dt \sum_{kq} \left[ \bar{\Psi}_{sk}^\dagger(t) \hat{M}(t) \bar{\Psi}_{aq}(t) + h.c. \right].
\]

The definition of those quantities can be found in the main text. Here, we do not include \( V(t) \), as we will not use it.

Integrating out the degrees of freedom of the bath using Gaussian integrals, one can obtain the equivalent fermionic action:

\[
S' = \int_c dt \int_c dt' \sum_k \bar{\Psi}_{sk}^\dagger(t) \hat{Q}_{k}^{-1}(t,t') \bar{\Psi}_{sk}(t') + S_{int},
\]

where

\[
\hat{Q}_{k}(t,t') = \hat{Q}_{sk0k}(t-t') + \int_{-\infty}^{+\infty} dt_1 dt_2 \hat{Q}_{sk0k}(t-t_1) \hat{\Sigma}_{k}(t_1,t_2) \hat{Q}_{k}(t_2,t'),
\]

with the self-energy from the bath being \( \hat{\Sigma}_{k}(t_1,t_2) = \sum_q \hat{M}(t_1) \hat{Q}_{aq0k}(t_1-t_2) \hat{M}^\dagger(t_2) \), and

\[
\hat{M}(t) = \begin{bmatrix}
W e^{if(t)} & 0 \\
0 & -W e^{-if(t)}
\end{bmatrix},
\]

where \( f(t) = (K/\Omega) \sin(\Omega t) \). We now turn to discuss how to calculate the dressed Green’s function \( \hat{Q}_{k}(t,t') \) through the perturbative expansion with respect to \( \kappa = K/\Omega \).
Using the Fourier transformation Eq.(7), Eq.(A5) can be rewritten as

$$\tilde{Q}_k(n, \omega) = \delta_{n0} \tilde{Q}_{s0k} + \sum_{n_1} \tilde{Q}_{s0k}(\omega + n\Omega)\dot{\Sigma}_k[n_1, \omega + (n - n_1)\Omega] \tilde{Q}_k(n - n_1, \omega),$$  \hspace{1cm} (A6)

where $\dot{\Sigma}_k(n, \omega) = \sum_q \sum_{n_2} \dot{M}_{n+n_2} \tilde{Q}_{s0q}(\omega - n_2\Omega) \dot{M}_{n_2}^\dagger$, and

$$\dot{M}_n = \begin{bmatrix} WJ_n(\frac{K}{\Omega}) & 0 \\ -W(1)^nJ_n(\frac{K}{\Omega}) & 0 \end{bmatrix},$$  \hspace{1cm} (A7)

where $J_n(x)$ is the Bessel function of the first kind. For weak driving amplitude $\kappa$, i.e., $\kappa \ll 1$, one can expand the Bessel function to the lowest order in $\kappa$. Up to $O(\kappa^2)$, we have

$$\dot{M}_0 = \begin{bmatrix} W & 0 \\ 0 & -W \end{bmatrix}, \quad \dot{M}_1 = \begin{bmatrix} \frac{W}{2} \kappa & 0 \\ 0 & \frac{W}{2} \kappa \end{bmatrix}, \quad \dot{M}_{-1} = \begin{bmatrix} -\frac{W}{2} \kappa & 0 \\ 0 & -\frac{W}{2} \kappa \end{bmatrix}.$$  \hspace{1cm} (A8)

Then, according to the Dyson equation Eq.(A6), we can derive the dressed Green’s function in the frequency space.

Next, we need to deal with the interaction action $S_{int}$, which is four-fermion interaction. One can treat this term through the Hubbard-Stratonovich transformation [63]. The idea of this transformation is that we multiply the Bessel function to the lowest order in $\kappa$. Thus, we have

$$\tilde{S}_F = \int dt \int d\eta \int d\xi \int d\varphi \int d\psi \tilde{M}_0 \tilde{M}_1 \tilde{M}_{-1} \tilde{M}^\dagger_0 \tilde{M}^\dagger_1 \tilde{M}^\dagger_{-1} \tilde{M} \tilde{M}^\dagger,$$

where $\Delta$ is a complex bosonic field. Thus, we have

$$e^{iS_{int}} = \int D[\tilde{\Delta}, \Delta] \exp \left[ i \int dt \int g \sum_{k} \tilde{\psi}_{k\uparrow} \tilde{\psi}_{-k\downarrow} \psi_{-k\uparrow} \psi_{k\downarrow} - \frac{1}{g} \tilde{\Delta} \Delta \right],$$  \hspace{1cm} (A10)

then we make a variable shift

$$\tilde{\Delta} \rightarrow \tilde{\Delta} - g \sum_{k} \tilde{\psi}_{k\uparrow} \tilde{\psi}_{-k\downarrow},$$

$$\Delta \rightarrow \Delta - g \sum_{k} \psi_{-k\uparrow} \psi_{k\downarrow},$$  \hspace{1cm} (A11)

and $e^{iS_{int}}$ becomes

$$e^{iS_{int}} = \int D[\tilde{\Delta}, \Delta] \exp \left\{ -i \int dt \left[ \frac{1}{g} \tilde{\Delta} \Delta - \Delta \sum_{k} \tilde{\psi}_{-k\uparrow} \psi_{k\downarrow} - \tilde{\Delta} \sum_{k} \tilde{\psi}_{k\uparrow} \psi_{-k\downarrow} \right] \right\}.$$  \hspace{1cm} (A12)

One can find that the generating function $Z = \int D[\tilde{\Delta}, \Delta, \tilde{\Psi}, \Psi] \exp \left[ iS(\tilde{\Delta}, \Delta, \tilde{\Psi}, \Psi) \right]$ now is in a quadratic form with respect to the fermionic field, which means we can also integrate them out using Gaussian integrals. Finally, one will obtain the effective bosonic action Eq.(6).

**Appendix B: Solving the Stationary Point Equation**

The key to solve the stationary point equation is to derive $\tilde{Q}_{k,\Delta}(n, \omega)$. Since $\tilde{Q}^{-1}_{k,\Delta} = \tilde{Q}^{-1}_{s0k} - \dot{\Sigma}_k$ (derived from the Dyson equation Eq.(5)), where $\tilde{Q}^{-1}_{s0k}$ is the free fermionic Green function and $\dot{\Sigma}_k$ is the self-energy provided by the normal metal bath in the Keldysh+Nambu space. Note that $\tilde{Q}_{k,\Delta}^{-1} = (\tilde{Q}_{k,\Delta}^{-1})^{-1}$, we have the following expansion:

$$\tilde{Q}_{k,\Delta}^{-1} = \begin{bmatrix} \tilde{Q}_{k,\Delta}^{-1(0)} - \dot{\Sigma}_{k}^{(1)} + \frac{1}{\sqrt{2}} (\Delta^{\alpha(1)} \otimes \hat{\tau}^+ + \Delta^{\alpha(1)} \otimes \hat{\tau}^-) + O(\kappa^2) \end{bmatrix}^{-1}$$

$$\equiv \tilde{Q}_{k,\Delta}^{(0)} + \tilde{Q}_{k,\Delta}^{(1)} + O(\kappa^2),$$  \hspace{1cm} (B1)
where the superscript \((i)\) also stands for the \(i\)th order in \(\kappa\),

\[
\tilde{Q}^{-1}_{k,\Delta} = \hat{Q}_{slk}^{-1} - \tilde{\Sigma}^{(0)}_k - \frac{1}{\sqrt{2}} \left[ \Delta^{\alpha(0)} \otimes (\hat{\gamma}^\alpha \otimes \hat{\tau}_+^\alpha) + \Delta^{\alpha(0)} \otimes (\hat{\gamma}^\alpha \otimes \hat{\tau}_-^\alpha) \right], 
\]  

(B2)

and

\[
\tilde{Q}^{(1)}_{k,\Delta} = \hat{Q}^{(0)}_{k,\Delta} \tilde{\Sigma}^{(1)}_k + \frac{1}{\sqrt{2}} \left[ \Delta^{\alpha(1)} (\hat{\gamma}^\alpha \otimes \hat{\tau}_+) + \Delta^{\alpha(1)} (\hat{\gamma}^\alpha \otimes \hat{\tau}_-) \right] \tilde{Q}^{(0)}_{k,\Delta}. 
\]  

(B3)

Due to the feature of Floquet matrices, \(\tilde{Q}^{(0)}_{k,\Delta}\) is a diagonal matrix, and \(\tilde{Q}^{(1)}_{k,\Delta}\) is a secondary diagonal matrix, etc.

For \(\Delta^q_0\), we have

\[
\Delta^q_0 = i \frac{g}{\sqrt{2}} \int \frac{d\omega}{2\pi} \text{Tr} \left[ (\hat{\gamma}^q \otimes \hat{\tau}_-) \sum_k \tilde{Q}_{k,\Delta}(0, \omega) \right] = i \frac{g}{\sqrt{2}} \int \frac{d\omega}{2\pi} \sum_k \left[ \tilde{Q}^K_{k,\Delta}(0, \omega) \right]^{12},
\]  

(B4)

where the superscript ”12” stands for the matrix element in the first row and second column. When we keep terms up to \(O(\kappa)\), normal fermionic Green’s functions with dissipations are

\[
\tilde{Q}^{R/A}_{k,p}(0, \omega) = \frac{1}{\omega - \epsilon_k \pm i\gamma}, \quad \tilde{Q}^{R/A}_{k,h}(0, \omega) = \frac{1}{\omega + \epsilon_k \pm i\gamma}, \quad \tilde{Q}^K_{k,p/h}(0, \omega) = -2i\gamma \frac{\tanh(\frac{\pi \omega}{2T})}{(\omega + \epsilon_k)^2 + \gamma^2}.
\]  

(B5)

From \(\tilde{Q}^K_{k,p/h}(0, \omega)\), we find that when we keep terms up to \(O(\kappa)\), the fermionic distribution function is still the Fermi-Dirac distribution, which is different from that in \(O(\kappa^2)\) case [58]. Then Eq.(B2) leads to the Green’s functions of quasiparticles, dressed by the bosonic field \(\Delta\):

\[
\tilde{Q}^{R/A}_{k,\Delta}(0, \omega) = \frac{1}{(\omega \pm i\gamma)^2 - \epsilon_k^2 - \Delta^2} \left[ \frac{\omega + \epsilon_k \mp i\gamma}{\Delta} \right. 
\]  

\[
\left. \frac{\omega - \epsilon_k \pm i\gamma}{\Delta} \right], \quad \text{B6}
\]

and

\[
\tilde{Q}^K_{k,\Delta}(0, \omega) = \tilde{Q}^R_{k,\Delta}(0, \omega) \cdot F(\omega) - F(\omega) \cdot \tilde{Q}^K_{k,\Delta}(0, \omega),
\]  

(B7)

where we have assumed \(\Delta\) to be real, \(F(\omega) = [1 - 2n_F(\omega)] \mathbb{1}_N\) with \(\mathbb{1}_N\) being the identity in the Nambu space and \(n_F(\omega) = 1/(e^{\beta\omega} + 1)\) is the Fermi-Dirac distribution.

Substituting \(\sum_k\) with \(\int d\epsilon_k p_F\), one will get the quasi-classical Green’s function [58, 69, 72] defined as \(\tilde{Q}^K_{\Delta} = \sum_k \tilde{Q}^K_{k,\Delta}\):

\[
\left[ \tilde{Q}^K_{\Delta}(0, \omega) \right]^{12} = i\pi p_F \left[ \frac{\Delta \tanh(\frac{\pi \omega}{2T})}{\sqrt{(\omega - i\gamma)^2 - \Delta^2}} - \frac{\Delta \tanh(\frac{\pi \omega}{2T})}{\sqrt{(\omega + i\gamma)^2 - \Delta^2}} \right]. 
\]  

(B8)

Then according to Eq.(B4), one will get the gap equation in the main text.

Now we turn to consider \(\Delta^q_{\pm 1}\). Seeing as \(\tilde{Q}^{(1)}_{k,\Delta}\) is not a diagonal matrix, for convenience, we truncate the Floquet matrix to a \(3 \times 3\) one. Then, one can solve the matrix equation Eq.(B3) to get

\[
\tilde{Q}_{k,\Delta}(\pm 1, \omega) = \tilde{Q}^{(0)}_{k,\Delta}(0, \omega \pm \Omega) \left\{ \tilde{\Sigma}_k(\pm 1, \omega) + \frac{1}{\sqrt{2}} \left[ \Delta^q_{\pm 1} (\hat{\gamma}^\alpha \otimes \hat{\tau}_+^\alpha) + \Delta^q_{\pm 1} (\hat{\gamma}^\alpha \otimes \hat{\tau}_-^\alpha) \right] \right\} \tilde{Q}^{(0)}_{k,\Delta}(0, \omega).
\]  

(B9)

Solving this equation, one can derive \(\tilde{Q}^K_{k,\Delta}(\pm 1, \omega)\), then one can get equations that decide \(\Delta^q_{\pm 1}\). Since it is difficult to analytically solve them, we resort to numerical calculations.
Appendix C: Bosonic Green’s Functions and The Diagram Rules

In this section, we briefly discuss the derivation of the bosonic Green’s functions. Here, we will keep terms up to the second order in $\kappa$. As shown in the main text, the transition temperature will be modified in this case, but not in the $O(\kappa)$ case. The second order approximation in $\Delta^\alpha(\Delta^\sigma)$ with $\alpha \in \{cl, q\}$ of the effective bosonic action—$S_{\text{eff}}$ has been derived and shown in Eq.(16). Then, according to the feature of Gaussian integrals, one can get

$$
\hat{D} = -i\left[ \begin{array}{cc}
\hat{\Delta} (\omega + \Omega) \\
\hat{\Delta} (\omega) \\
\hat{\Delta} (\omega - \Omega)
\end{array} \right] \begin{bmatrix}
\hat{\Delta} (\omega + \Omega) & \hat{\Delta} (\omega) & \hat{\Delta} (\omega - \Omega)
\end{bmatrix}
$$

where we have truncated the Floquet matrix to a $3 \times 3$ one, $\hat{\Gamma} = \sum_k \hat{\Gamma}_k$, the superscript $(i)$ denotes terms kept up to the $i$th order in $\kappa$,

$$
\mathbb{1} - g \hat{\Gamma}^{(0)} \otimes \hat{\sigma}_1 = \begin{bmatrix}
\mathbb{1}_K - g \hat{\Gamma}(0, \omega + \Omega) \hat{\sigma}_1 \\
0 \\
0 \\
0
\end{bmatrix},
$$

$$
g \hat{\Gamma}^{(1)} \otimes \hat{\sigma}_1 = \begin{bmatrix}
0 \\
g \hat{\Gamma}(1, \omega) \hat{\sigma}_1 \\
g \hat{\Gamma}(1, \omega - \Omega) \hat{\sigma}_1 \\
g \hat{\Gamma}(1, -\omega) \hat{\sigma}_1
\end{bmatrix},
$$

$$
g \hat{\Gamma}^{(2)} \otimes \hat{\sigma}_1 = \begin{bmatrix}
0 \\
0 \\
g \hat{\Gamma}(2, \omega - \Omega) \hat{\sigma}_1 \\
g \hat{\Gamma}(2, -\omega) \hat{\sigma}_1
\end{bmatrix},
$$

where $\mathbb{1}_K$ is the identity matrix in the Keldysh space. Expanding $(\mathbb{1} - g \hat{\Gamma} \otimes \hat{\sigma}_1)^{-1}$ up to $\kappa^2$, one will get

$$
(\mathbb{1} - g \hat{\Gamma} \otimes \hat{\sigma}_1)^{-1} \approx (\mathbb{1} - g \hat{\Gamma}^{(0)} \otimes \hat{\sigma}_1)^{-1} + (\mathbb{1} - g \hat{\Gamma}^{(0)} \otimes \hat{\sigma}_1)^{-1} \cdot g \hat{\Gamma}^{(1)} \otimes \hat{\sigma}_1 \cdot (\mathbb{1} - g \hat{\Gamma}^{(0)} \otimes \hat{\sigma}_1)^{-1}
$$

Then one can get matrix elements of $\hat{D}^{cl,q}(n, \omega)$ up to the second order in $\kappa$: 

$$
\hat{D}^{cl,q}(0, \omega) = -\frac{g}{1 - g \hat{\Gamma}^{cl,q}(0, \omega)}
$$

$$
\frac{g^2 \hat{\Gamma}^{cl,q}(1, \omega - \Omega) \hat{\Gamma}^{cl,q}(-1, \omega)}{1 - g \hat{\Gamma}^{cl,q}(0, \omega)} \frac{1 - g \hat{\Gamma}^{cl,q}(0, \omega - \Omega)}{1 - g \hat{\Gamma}^{cl,q}(0, \omega)}
$$

$$
\frac{g^2 \hat{\Gamma}^{cl,q}(-1, \omega + \Omega) \hat{\Gamma}^{cl,q}(1, \omega)}{1 - g \hat{\Gamma}^{cl,q}(0, \omega)} \frac{1 - g \hat{\Gamma}^{cl,q}(0, \omega + \Omega)}{1 - g \hat{\Gamma}^{cl,q}(0, \omega)}
$$
\[
\dot{D}^{cl,q}(\pm 1, \omega) = -\frac{g^2 \hat{\Gamma}^{cl,q}(\pm 1, \omega)}{1 - g \hat{\Gamma}^{cl,q}(0, \omega)} \frac{1 - g \hat{\Gamma}^{cl,q}(0, \omega \pm \Omega)}{1 - g \hat{\Gamma}^{cl,q}(0, \omega \pm 2\Omega)}.
\]
\[
\dot{D}^{cl,q}(\pm 2, \omega) = -\frac{g^2 \hat{\Gamma}^{cl,q}(\pm 2, \omega)}{1 - g \hat{\Gamma}^{cl,q}(0, \omega)} \frac{1 - g \hat{\Gamma}^{cl,q}(0, \omega \pm \Omega)}{1 - g \hat{\Gamma}^{cl,q}(0, \omega \pm 2\Omega)} - \frac{g^3 \hat{\Gamma}^{cl,q}(\pm 1, \omega)}{1 - g \hat{\Gamma}^{cl,q}(0, \omega)} \frac{1 - g \hat{\Gamma}^{cl,q}(0, \omega \pm \Omega)}{1 - g \hat{\Gamma}^{cl,q}(0, \omega \pm 2\Omega)}.
\]

In order to make the diagram rules we show in the main text clearer, we summarize them again and give one more example.

The diagram rules read as:

1. Attach \(g/(1 - g\hat{\Gamma}^{cl,q}(0, \omega))\) to a direct vertex shown in Fig. 3(a);
2. Attach \(\hat{\Gamma}^{cl,q}(m, \omega)(\sim \kappa|m| + \mathcal{O}(\kappa|m|^{2}))\) to an energy exchange vertex shown in Fig. 3(b);
3. The indirect vertex shown in Fig. 3(c) is constructed from direct vertices and indirect vertices, thus one should attach

\[
\frac{g}{1 - g\hat{\Gamma}^{cl,q}(0, \omega)} \hat{\Gamma}^{cl,q}(m_1, \omega) \frac{g}{1 - g\hat{\Gamma}^{cl,q}(0, \omega + m_1 \Omega)} \hat{\Gamma}^{cl,q}(m_1, \omega + m_2 \Omega) \frac{g}{1 - g\hat{\Gamma}^{cl,q}(0, \omega + (m_1 + m_2) \Omega)}
\]

4. As for the diagram of one complete process, we just need to sum up those relative direct vertices and indirect vertices.

In the main text, we use the example of \(\dot{D}^{cl,q}(0, 0)\) to introduce the diagram rules, and in the following, we will use the diagram rules to derive \(\dot{D}^{cl,q}(\pm 2, \omega)\), and comparing it to Eq. (C8) to show the correctness of the diagram rules. Here, we also keep terms up to the second order in \(\kappa\).

For \(\dot{D}^{cl,q}(\pm 2, \omega)\), since there exist energy exchange processes \(n = \pm 2\), thus in the diagram, the energy exchange vertex must exist.

- The simplest process is that there are only one energy exchange vertex, which stands for one particle absorbing/emitting two parts of energies \((\pm 2\Omega)\), and two direct vertices. The diagram is shown in Fig. 5(a);
- One may find that a process that contains two energy exchange vertices, which stands for one particle first absorbing/emitting one part of energy \((\pm \Omega)\), then absorbing/emitting one part of energy as well, can also contribute to \(\dot{D}^{cl,q}(\pm 2, \Omega)\). It indeed does, and the diagram is shown in Fig. 5(b);
- Since the leading order term of the process, in which one particle absorbs/emits \(n\) parts of energies, is proportional to \(\kappa^n\)(that is \(\hat{\Gamma}^{cl,q}(\pm n, \omega) \sim \kappa^n + \mathcal{O}(\kappa^{n+2})\)), in the \(\mathcal{O}(\kappa^2)\) case, there are only the above two diagrams that contribute to \(\dot{D}^{cl,q}(\pm 2, \omega)\). The diagram for \(\dot{D}^{cl,q}(\pm 2, \omega)\) is just the sum of Fig. 5(a) and Fig. 5(b);
- Use the diagram rules shown previously, one can easily derive the expression of \(\dot{D}^{cl,q}(\pm 2, \omega)\):

\[
-\dot{D}^{cl,q}(\pm 2, \omega) = \frac{g^2 \hat{\Gamma}^{cl,q}(\pm 2, \omega)}{1 - g \hat{\Gamma}^{cl,q}(0, \omega)} \frac{1 - g \hat{\Gamma}^{cl,q}(0, \omega \pm \Omega)}{1 - g \hat{\Gamma}^{cl,q}(0, \omega \pm 2\Omega)}
\]

\[
+ \frac{g^3 \hat{\Gamma}^{cl,q}(\pm 1, \omega)}{1 - g \hat{\Gamma}^{cl,q}(0, \omega)} \frac{1 - g \hat{\Gamma}^{cl,q}(0, \omega \pm \Omega)}{1 - g \hat{\Gamma}^{cl,q}(0, \omega \pm 2\Omega)}.
\]

which is exactly the same with Eq. (C8).

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