Photoinduced dynamics of excitonic order and Rabi oscillation in the two-orbital Hubbard model

Yasuhiro Tanaka$^{1,2}$ and Kenji Yonemitsu$^3$

$^1$Department of Physics, Tokyo Institute of Technology, Meguro, Tokyo 152-8550, Japan
$^2$Department of Applied Physics, Waseda University, Shinjuku, Tokyo 169-8555, Japan
$^3$Department of Physics, Chuo University, Bunkyo, Tokyo 112-8551, Japan

(Dated: July 31, 2020)

We investigate the condition for the photoinduced enhancement of an excitonic order in a two-orbital Hubbard model, which has been theoretically proposed in our previous work [Phys. Rev. B 97, 115105 (2018)], and analyze it from the viewpoint of the Rabi oscillation. Within the mean-field approximation, we simulate real-time dynamics of an excitonic insulator with a direct gap, where the pair condensation in the initial state is of BEC nature and the photoexcitation is introduced by electric dipole transitions. We first discuss that in the atomic limit our model is reduced to a two-level system that undergoes the Rabi oscillation, so that for single cycle pulses physical quantities after the photoirradiation are essentially determined by the ratio of the Rabi frequency to the pump-light frequency. Then, it is shown that this picture holds even in the case of nonzero transfer integrals where each one-particle state exhibits the Rabi oscillation leading to the enhancement of the excitonic order. We demonstrate that effects of electron-phonon interactions do not alter the results qualitatively. We also examine many-body dynamics by the exact diagonalization method on small clusters, which strongly suggests that our mechanism for the enhancement of the excitonic order survives even when quantum fluctuations are taken into account.

I. INTRODUCTION

Photoirradiation to correlated electron systems has opened a novel playground to manipulate various electronic phases. In particular, recent experimental studies have reported that electronic orders are transiently reinforced or even created by laser light $^{[1-10]}$, which indicates a clear distinction from typical photoinduced phase transitions in which they are usually suppressed. These phenomena have been observed, for instance, in materials which exhibit charge ordering $^{[1,3,10]}$, charge density wave $^{[4,8]}$, superconductivity $^{[2,4,5,7]}$, and excitonic condensation $^{[9]}$. Simultaneously, theoretical efforts to understand their mechanisms as well as to pursue a way of controlling electronic phases have been made recently, where roles of electron-electron (e-e) and/or electron-phonon (e-ph) interactions on laser-induced dynamics have been intensively studied $^{[11-22]}$. For excitonic insulators (EIs), a transient gap enhancement by photoexcitation has been observed in a candidate material $\text{Ta}_2\text{NiSe}_5$ $^{[9]}$. The EI is a state in which electrons in the conduction band and holes in the valence band form bound pairs called excitons by the Coulomb interaction, and they become a condensate. Theories of EIs have been developed in semimetals and semiconductors $^{[23-27]}$. $\text{Ta}_2\text{NiSe}_5$ is a layered semiconductor with a direct gap above $T_C = 326K$ where a second-order transition accompanied by a structural distortion occurs $^{[28,29]}$. Although the identification of an EI is a difficult task, recent experimental $^{[30,51]}$ and theoretical $^{[32-35]}$ studies have offered evidences that an EI is realized in the low temperature phase. With regards to its photoinduced phenomena, e-ph coupled systems have been investigated by mean-field theories $^{[19,22]}$ and the origin of the gap enhancement has been discussed.

In purely electronic systems without phonon degrees of freedom, we have studied $^{[21]}$ photoinduced dynamics of a direct-gap EI using a two-orbital Hubbard model in which excitonic insulators in thermal equilibrium show BEC-BCS crossover depending on the value of the interorbital Coulomb interaction $U'$. By incorporating the effects of photoexcitation through electric dipole transitions, we have shown that the enhancement of the excitonic gap occurs when the initial state is an EI in the BEC regime or a nearby band insulating state, and the pump-light frequency is close to the excitonic gap. There is an optimal value of the amplitude of the light field for inducing the gap enhancement, although its physical origin has not been clarified yet. Our study has also shown that the time evolutions of the phases of excitonic pairs in momentum space are crucially important for understanding the photoinduced behavior of the excitonic gap $^{[21]}$: They evolve basically in phase when the gap is enhanced by the laser irradiation, whereas they strongly depend on momentum when the initial EI is in the BCS regime for which the gap is suppressed.

In this paper, we elucidate the physical origin of the gap enhancement through laser-induced dipole transitions in a two-orbital Hubbard model mainly by using the time-dependent Hartree-Fock (HF) approximation. For this purpose, we consider the atomic limit in which our system is equivalent to a two-level system that exhibits the Rabi oscillation, where the dynamics of physical quantities are understood from changes in the occupation probability of the two levels. Even when we introduce nonzero transfer integrals, its photo-response is

---

$^*$tanaka.y@taoni.waseda.jp

$^1$ Present address
qualitatively unaltered as far as the initial state is near the boundary between the EI and the band insulator (BI) phases. There the EI belongs to the BEC regime and the excitonic pairs are formed locally. In momentum space, the photoinduced gap enhancement is interpreted as a consequence of a cooperative Rabi oscillation of the one-particle states. We confirm that the e-ph coupling considered in the previous theories [19, 22] has little effects on our mechanism for the gap enhancement. Moreover, we examine effects of quantum fluctuations on the dynamics by using the exact diagonalization (ED) method, which corroborates the results obtained by the HF approximation. This paper is organized as follows. In Sect. II, the two-orbital Hubbard model and the calculation method for photoinduced dynamics are introduced. The model in the atomic limit is also described. In Sect. III, the results without the phonon degrees of freedom are presented and we discuss the photoinduced dynamics in terms of the Rabi oscillation. The effects of the e-ph coupling are elucidated in Sect. IV, whereas those of quantum fluctuations are discussed in Sect. V where we give the results with the ED method. The discussion and summary are devoted to Sect. VI.

II. MODEL AND METHOD

A. Two-orbital Hubbard model

We consider a two-orbital Hubbard model in one dimension, which is defined as

\[ \hat{H} = t_c \sum_{i\sigma} (c_{i\sigma}^\dagger c_{i+1\sigma} + \text{h.c.}) + \mu_C \sum_{i\sigma} n_{i\sigma}^c + \mu_F \sum_{i\sigma} n_{i\sigma}^f + U \sum_{i} n_{i\uparrow}^c n_{i\downarrow}^c + U' \sum_{i} n_{i\uparrow}^f n_{i\downarrow}^f, \]  

(1)

where \( n_{i\sigma}^c = \alpha_{i\sigma}^c \alpha_{i\sigma}^c \) and \( n_{i\sigma}^f = n_{i\uparrow}^f + n_{i\downarrow}^f \). The intraorbital (interorbital) Coulomb interaction is denoted by \( U \) (\( U' \)). For the transfer integral \( t_f = 1 \) and \( t_c = -1 \) as in the previous study [21]. The parameter \( \mu_C(>0) \) controls the overlap between the \( c \) and \( f \) bands. When \( \mu_C > 4 \), the system with \( U = U' = 0 \) has a band structure of a direct-gap semiconductor, whereas it becomes a semimetal for \( \mu_C < 4 \). The electron density per site is fixed at \( n = 2 \).

Photoexcitation is introduced by electric dipole-allowed transitions [19, 30] that are described by the time-dependent term

\[ \hat{H}_D(\tau) = F(\tau) \sum_{i\sigma} (c_{i\sigma}^\dagger f_{i\sigma} + \text{h.c.}), \]  

(2)

which is added to Eq. (1). We define \( F(\tau) \) as

\[ F(\tau) = F_0 \sin(\omega \tau)e^{-(\tau - \tau_p)^2/\tau_w^2}, \]  

(3)

where \( \omega \) is the light frequency and we set \( \tau_p = \tau_w = \pi/\omega \). Although we mainly use the gaussian envelope for \( F(\tau) \), we also consider a rectangular envelope with which \( F(\tau) \) is defined as

\[ F(\tau) = F_0 \sin(\omega \tau)\theta(\tau)(T_{\text{IRR}} - \tau), \]  

(4)

where \( \theta(\tau) \) and \( T_{\text{IRR}} \) are the Heaviside step function and the pulse width, respectively. This form of \( F(\tau) \) enables us to interpret our results directly from the viewpoint of the Rabi oscillation. We note that the pulse shape does not qualitatively affect our results. Unless otherwise noted, we use single cycle pulses (\( T_{\text{IRR}} = 2\pi/\omega \)).

We apply the HF approximation to Eq. (1) where the excitonic order parameter and the electron density on the \( \alpha \) orbital per site are defined as \( \Delta_0 = \langle c_{i\sigma}^\dagger f_{i\sigma} \rangle \) and \( n_c = 2\langle n_{i\sigma}^c \rangle \), respectively. We have assumed that \( \langle c_{i\sigma}^\dagger f_{i\sigma} \rangle \) and \( \langle n_{i\sigma}^f \rangle \) are independent of \( i \) and \( \sigma \) [21, 57]. The total Hamiltonian in momentum representation is given as

\[ \hat{H}_{\text{tot}}^\text{HF}(\tau) = \sum_{k\sigma} \hat{H}_{k\sigma}(\tau) = \sum_{k\sigma} \Psi_{k\sigma}^\dagger h_k(\tau) \Psi_{k\sigma}, \]  

(5)

where \( \Psi_{k\sigma}^\dagger = (c_{k\sigma}^\dagger, f_{k\sigma}^\dagger) \) and \( h_k(\tau) \) is defined by

\[ h_k(\tau) = \left( \begin{array}{cc} \varepsilon_k^c & -U'\Delta_0^c + F(\tau) \\ -U'\Delta_0^f & \varepsilon_k^f \end{array} \right). \]  

(6)

In Eq. (6), \( \varepsilon_k^c = \epsilon_k^c + \frac{U}{2}n_c + U'n_f \) and \( \varepsilon_k^f = \epsilon_k^f + \frac{U}{2}n_f + U'n_c \) where \( \epsilon_k^c = 2t_c \cos k + \mu_C \) and \( \epsilon_k^f = 2t_f \cos k \) are the noninteracting energy dispersions for the \( c \) and \( f \) bands, respectively. In the ground state, \( n_c (= 2 - n_f) \) and \( \Delta_0 \) are determined self-consistently.

Photoinduced dynamics are obtained by numerically solving the time-dependent Schrödinger equation [38–41]

\[ |\psi_{k\sigma}(\tau + d\tau)| = T \exp \left[ -i \int_{\tau}^{\tau + d\tau} d\tau' \hat{H}_{k\sigma}(\tau') \right] |\psi_{k\sigma}(\tau)|, \]  

(7)

where \( |\psi_{k\sigma}(\tau)| \) denotes a one-particle state with wave vector \( k \) and spin \( \sigma \) at time \( \tau \), and \( T \) is the time-ordering operator. We use the time slice \( d\tau = 0.01 \) with \( t_f \) as the unit of energy (and \( 1/t_f \) as that of time). For a physical quantity \( X(\tau) \), its time average is denoted by \( \overline{X} \) that is calculated as

\[ \overline{X} = \frac{1}{\tau_f - \tau_i} \int_{\tau_i}^{\tau_f} X(\tau) d\tau. \]  

(8)

If \( X(\tau) \) is conserved after the photoexcitation, its value is written as \( \overline{X} \).
B. Atomic limit

In the atomic limit ($t_c = t_f = 0$), our system is reduced to a two-level system described by the Hamiltonian

$$\hat{H}^{\text{AL}}(\tau) = \sum_\sigma \Psi_\sigma^\dagger h^{\text{AL}}(\tau) \Psi_\sigma,$$

where $\Psi_\sigma = (c_\sigma, f_\sigma^\dagger)$ and $h^{\text{AL}}(\tau)$ is defined as

$$h^{\text{AL}}(\tau) = \left( \begin{array}{cc} \epsilon^c & -U'\Delta_0 + F(\tau) \\ -U'\Delta_0 + F(\tau) & \epsilon^f \end{array} \right),$$

and $\epsilon^c = \mu_C + \frac{U}{2} n_c + U' n_f$ and $\epsilon^f = \frac{U}{2} n_f + U' n_c$. In the ground state [$F(\tau) = 0$], the self-consistent equations for $n_c$ and $\Delta_0$ are written as

$$n_c = 1 - \frac{\epsilon^c - \epsilon^f}{\sqrt{(\epsilon^c - \epsilon^f)^2 + 4U'^2|\Delta_0|^2}},$$

and

$$\Delta_0 = \frac{1}{2} \left[ 1 - \frac{(\epsilon^c - \epsilon^f)^2}{(\epsilon^c - \epsilon^f)^2 + 4U'^2|\Delta_0|^2} \right]^{1/2},$$

respectively, which leads to

$$\Delta_0 = \frac{1}{2} \sqrt{n_c(2 - n_c)}.$$

With this relation, the expectation value for the energy $E^{\text{AL}}$ can be written as

$$E^{\text{AL}} = \frac{1}{2} (U - U') n_c^2 + (\mu_C - U + U') n_c + U.$$

We take $U (1/U)$ as the unit of energy (time) in the atomic limit.

III. RESULTS WITHOUT PHONONS

In this section, we show the results obtained by the HF approximation for the two-orbital Hubbard model without e-ph couplings. We first consider the case of the atomic limit and then discuss the case of nonzero transfer integrals ($t_f = -t_c = 1$).

A. Atomic limit

1. Ground state

Before the laser irradiation, we consider two phases in the ground state: an excitonic phase (EP) with $\Delta_0 \neq 0$ and a decoupled phase (DP) with $\Delta_0 = 0$ and $n_c = 0$. They correspond to EI and BI phases, respectively, when $t_c$ and $t_f$ are nonzero [21, 42]. We use $U = 1$ and $\mu_C = 0.5$, and vary $U'$ ($\geq 0$) as a parameter. Since

FIG. 1. (a) Ground-state energies for EP and DP, and (b) $\Delta_0$ as functions of $U'$ with $U = 1$ and $\mu_C = 0.5$. In (a), the energy of the magnetic phase (MP) is also shown by the dashed line.

$$\frac{\partial E^{\text{AL}}}{\partial n_c} = 0$$

gives $n_c = 1 - \mu_C/(U - U')$, we have a mean-field solution with $\Delta_0 \neq 0$ for $0 \leq U' < U'_{\text{cr}} = U - \mu_C$. In Fig. 1(a), we show the ground-state energies for the EP and the DP, the latter of which is independent of $U'$. With increasing $U'$, a transition from the EP to the DP occurs at $U' = U'_{\text{cr}}$. In fact, a magnetic phase (MP) with $\langle n_c^\sigma \rangle = 1$ and $\langle n_f^\sigma \rangle = 0$ has the ground-state energy $U' + \mu_C$, which becomes the lowest energy state for $0 \leq U' < U'_{\text{cr}}$. However, we consider only nonmagnetic initial states for the photoexcitation. The reason is that for nonzero $t_c$ and $t_f$, the photoinduced gap enhancement reported previously occurs in the vicinity of the boundary between the EI and BI phases [21] ($U' \sim U'_{\text{cr}}$) where magnetic ordered states do not appear as the ground state [43]. In this paper, we focus on the dynamics near $U' = U'_{\text{cr}}$.

In Fig. 1(b), the $U'$ dependence of $\Delta_0$ is shown, indicating that $\Delta_0$ is nonzero at $U' = 0$ and exhibits a steep decrease toward $U' = U'_{\text{cr}}$ at which it vanishes. For $U' \sim U'_{\text{cr}}$, this result is similar to that obtained with nonzero $t_c$ and $t_f$ (Fig. 10 in Sect. III B), whereas they are qualitatively different for $U' \sim 0$. The similarity near $U' = U'_{\text{cr}}$ comes from the local character of excitonic pairs: the EI is in the BEC regime of the BCS-BEC crossover [42, 44, 45]. On the other hand, for $U' \sim 0$, the EI is in the BCS regime, which cannot be described by the atomic limit.

2. Photoinduced dynamics

In Fig. 2, we show the time evolutions of $n_c$ and $|\Delta_0|$ for different values of $F_0$ with $U' = 0.45 \left( < U'_{\text{cr}} \right)$. For the

\[\Delta_0 = \frac{1}{2} \sqrt{n_c(2 - n_c)}\]
time evolutions of the real and imaginary parts of \( \Delta_0 \), see Appendix A. We use \( F(\tau) \) with the gaussian envelope [Eq. (3)] and choose \( F_0 < 0 \), although the sign of \( F_0 \) does not affect the results qualitatively [21]. The pump-light frequency is tuned to the difference between the two eigenvalues of Eq. (10) with \( \omega = 0 \). In Figs. 2(a) and 2(b), we show \( \tilde{n}_c \) and \( \tilde{|\Delta_0|} \) as functions of \( |F_0|/\omega \) for \( U' = 0.45 \) and 0.55, where the ground states are in the EP and the DP, respectively. At time \( \tau \), the wave function of the two-level system is written as

\[
|\psi(\tau)\rangle = u(\tau)c^0|0\rangle + v(\tau)f^0|0\rangle,
\]

with \(|u(\tau)|^2 + |v(\tau)|^2 = 1\) where we have omitted the spin index for brevity. By using the relations \( n_c = 2|u(\tau)|^2 \) and \( \Delta_0 = u^*(\tau)v(\tau) \), we have

\[
|\Delta_0| = \frac{1}{2}\sqrt{n_c(2-n_c)}.
\]

which holds at any \( \tau \) indicating that \( |\Delta_0| \) has its maximum value of 0.5 when \( \tilde{n}_c = 1 \). In order to examine changes in the occupation probability of the two levels, we compute the overlap between the wave function in the ground state and that after the photoexcitation. The overlap \( \alpha \) is given by

\[
\alpha = |\langle \psi(\tau = 0)|\psi(\tau)\rangle|^2
\]

\[
= 1 - \frac{1}{2}(n_c + n_G^c) + \frac{1}{2}n_c n_G^c + \Delta_0^* \Delta_0 + \Delta_G^* \Delta_0 \tag{17}
\]

where \( n_G^c \) and \( \Delta_G^c \) are \( n_c \) and \( \Delta_0 \) in the ground state, respectively. In evaluating \( \alpha \), we adjust the phase of \( \Delta_0 \) in \( |\psi(\tau = 0)\rangle \) to coincide with that of \( \Delta_0(\tau \neq 0) \). In this case, we have

\[
\alpha = 1 - \frac{1}{2}(n_c + n_G^c) + \frac{1}{2}n_c n_G^c + 2|\Delta_0| |\Delta_0| \tag{18}
\]

which is conserved after the photoexcitation. In Fig. 3(c), we show \( \tilde{\alpha} \) and the increment in the total energy \( \Delta E \) per site for \( U' = 0.45 \). For the quantities \( \tilde{n}_c \), \( \Delta E \), and \( \tilde{\alpha} \), an oscillatory behavior with respect to \( |F_0|/\omega \) is
evident, although the period of oscillation is not constant. The behavior of $|\tilde{\Delta}_0|$ appears to be more complex than that of $\tilde{n}_c$ because of the relation Eq. (16). The oscillation in $\tilde{\alpha}$ and $\Delta E$ indicates a manifestation of the Rabi oscillation in the present two-level system, which we will discuss in detail below.

3. Rabi oscillation and enhancement of excitonic order

Here we consider $F(\tau)$ with the rectangular envelope defined by Eq. (1). In Fig. 3 we show $\tilde{n}_c$, $|\tilde{\Delta}_0|$, $\Delta E$, and $\tilde{\alpha}$ as functions of $|F_0|/\omega$ where the parameters are the same as those in Fig. 3. The oscillatory behavior of these quantities is more prominent than that in Fig. 3 where the gaussian envelope is employed for $F(\tau)$. In particular, the period of the oscillation is almost constant. By using Eqs. (16) and (18), we obtain the $\alpha$ dependence of $n_c$ shown in Fig. 5. Along its curve, the position $(\tilde{\alpha}, \tilde{n}_c)$ moves depending on the value of $|F_0|/\omega$. For $U' = 0.45$, we have $(\alpha, n_c) = (1, 0.091)$ in the ground state. With increasing $|F_0|/\omega$, the position $(\tilde{\alpha}, \tilde{n}_c)$ first moves to the upper-left direction in Fig. 5 until $|F_0|/\omega \sim 0.4$ where $\tilde{n}_c$ exhibits the first peak as shown in Fig. 4(a). Reflecting the periodic behavior of $\tilde{\alpha}$, the point $(\tilde{\alpha}, \tilde{n}_c)$ goes back to the initial position at $|F_0|/\omega \sim 0.8$. The value of $\tilde{n}_c$ becomes smaller than $n_c^G$ for $0.8 < |F_0|/\omega < 1.2$ [Fig. 4(a)] where $\tilde{\alpha}$ is slightly smaller than $1$ [Fig. 4(c)]. Then, $(\tilde{\alpha}, \tilde{n}_c)$ moves to the upper-left direction again until $|F_0|/\omega = 2.0$ at which $\tilde{n}_c$ shows the second peak. For $U' = 0.55$, Eq. (13) gives $\alpha = 1 - \frac{1}{2}n_c$ since $n_c^G = \Delta_0^G = 0$. The behavior of $(\tilde{\alpha}, \tilde{n}_c)$ depending on $|F_0|/\omega$ is similar to that for $U' = 0.45$. These results show that the oscillatory behavior of physical quantities originates from that of $\tilde{\alpha}$. In order to interpret our results as the Rabi oscillation more quantitatively, we consider the case of continuous-wave (CW) lasers in the following.

In Fig. 6 we show the time profile of $n_c$ under CW excitations for $U' = 0.45$ and $\omega = 0.45$. When $|F_0|$ is small ($F_0 = -0.05$), $n_c$ shows a small oscillation around the value of $n_c^G$. As we increase $|F_0|$, a large-amplitude oscillation appears, the period of which gets shorter for larger $|F_0|$. In Fig. 7 we show the Fourier transform of $n_c$ for large $|F_0|$ ($F_0 \leq -0.1$) (see Appendix A for...
the details of the dynamics for small $|F_0|$. There is a sharp peak in each spectrum and its position denoted by $\Omega$ is nearly proportional to $|F_0|$ as shown in Fig. 8. In a two-level system driven by a CW laser, the rotating wave approximation (RWA) gives the Rabi frequency $\Omega_R$ as

$$\Omega_R = \sqrt{(\omega - E_G)^2 + F_0^2},$$

(19)

where $E_G$ is the difference between the two energy levels. At the resonance ($\omega = E_G$), we have $\Omega_R = |F_0|$. In fact, the Hamiltonian in Eq. (10) contains $n_c$ and $\Delta_0$ that are $\tau$-dependent, which is different from the conventional Rabi oscillation [46, 47]. The effects of the $\tau$-dependence of $n_c$ and $\Delta_0$ in the Hamiltonian on the dynamics are discussed in Appendix B. Considering this difference, here we will replace $\Omega_R$ at the resonance [$\omega = E_G(\tau = 0)$] by $\Omega'_R = |F_0| p$ with a coefficient $p$ [48]. This leads us to $u(\tau) \propto \sin(\Omega'_R \tau + \phi)$ so that $n_c$ is written as

$$n_c = A \sin^2(\frac{\Omega'_R}{2} \tau + \phi),$$

(20)

where $A$ and $\phi$ are constants. We fit a linear function $\Omega = p|F_0|$ to the results of $\Omega$ in Fig. 8. The fitting works well with $p$ slightly larger than 1 for both $U' = 0.45$ ($p = 1.24$) and 0.55 ($p = 1.14$). In Eq. (20), we have $\phi = 0$ ($\phi = 0$) when $U'$ is slightly smaller (larger) than $U'_c$ because of $n^G_c \sim 0$ ($n^G_c = 0$), indicating that for single cycle pulses the values of $n_c$ and $|\Delta_0|$ after the photoexcitation are governed by $\Omega'_R/\omega$. In particular, $n_c$ becomes maximum at $\Omega'_R/\omega \sim 1/2$. This relation gives $|F_0|/\omega = 0.40$ for $U' = 0.45$ and $|F_0|/\omega = 0.44$ for $U' = 0.55$, which are consistent with the results shown in Fig. 4(a). In Fig. 9, we show the $|F_0|/\omega$ dependence of $n_c$ calculated by Eq. (20) at $\tau = 2\pi/\omega$, and compare the result with $\bar{n}_c$ shown in Fig. 4(a). The quantities $A$ and $\phi$ are determined from the height of the first peak in $\bar{n}_c$ and the value of $n^G_c$. For $|F_0|/\omega \lesssim 0.8$, the results obtained by Eq. (20) reproduce those for $\bar{n}_c$ fairly well, although they deviate from each other for larger $|F_0|/\omega$, which is due to the limitation of the RWA [49].

### B. One-dimensional model

Next, we show results for the one-dimensional model with $t_f = 1$ and $t_c = -1$ in Eq. (11), for which the initial EI and BI have a direct gap [21].

#### 1. Ground state

In Fig. 10, we show $\Delta_0$ and $n_c$ as functions of $U'$ in the ground state with $U = 4$ and $\mu_C = 2.5$. As in the previous studies where the Fermi surface is perfectly nested [21, 42], an infinitesimal $U'$ produces an EI with $\Delta_0 \neq 0$. The order parameter $\Delta_0$ exhibits a maximum at $U' = 2.70$ and a transition from the EI to BI phases occurs at $U' = U'_{cr} = 3.37$ where $\Delta_0$ vanishes. Toward $U' = U'_{cr}$, $n_c$ monotonically decreases. In the BI phase, the $c$ and $f$ bands are completely decoupled so that we
2. Photoinduced dynamics

In the calculations of photoinduced dynamics, we use two sets of parameters, both of which give EIs that are located near the EI-BI phase boundary of the ground states. One is $U = 4$, $\mu_C = 2.5$ and $U' = 3.3$. The other is $U = 8$, $\mu_C = 5$ and $U' = 4.3$ where $U'_C = 4.5$. We use $F(\tau)$ with the gaussian envelope [Eq. (3)] and the pump-light frequency is tuned to the initial gap of the EI: $\omega = 1.25$ for $U = 4$ and $\omega = 1.99$ for $U = 8$. We note that as far as $U' \sim U''$, our results are qualitatively unaltered even when we choose a BI as the initial state. The system size is $N = 200$. In analogy with the case of the atomic limit, we define a $k$-dependent quantity $\alpha_k$ as follows. First, we consider the overlap between the one-particle state at time $\tau$ and that in the ground state. If we write the one-particle state as $|\psi_{k\sigma}(\tau)\rangle = u_{k\sigma}(\tau) c_{k\sigma}^\dagger (0) + v_{k\sigma}(\tau) f_{k\sigma}^\dagger (0)$ with $|u_{k\sigma}|^2 + |v_{k\sigma}|^2 = 1$, the overlap is written as

$$
|\langle \psi_{k\sigma}(\tau = 0) | \psi_{k\sigma}(\tau) \rangle|^2 = n_c^G(k) n_c(k) + [1 - n_c^G(k) ] [1 - n_c(k)] + \Delta^G(k) \Delta(k) + \Delta^G(k) \Delta^*(k),
$$

where $n_c(k)$ and $\Delta(k)$ are the momentum distribution function for $c$-electrons and the pair amplitude in $k$ space, which are written as

$$
n_c(k) = \langle c_{k\sigma}^\dagger c_{k\sigma} \rangle = |u_{k\sigma}|^2,
$$

and

$$
\Delta(k) = \langle c_{k\sigma}^\dagger f_{k\sigma} \rangle = u_{k\sigma}^* v_{k\sigma},
$$

respectively, and $n_c^G(k) |\Delta^G(k)\rangle$ is $n_c(k) |\Delta(k)\rangle$ in the ground state. Then, as in Eq. (18), we define $\alpha_k$ as

$$
\alpha_k = n_c^G(k) n_c(k) + [1 - n_c^G(k) ] [1 - n_c(k)] + 2|\Delta^G(k)| |\Delta(k)|,
$$

which is the upper limit of the overlap in Eq. (21).

After the photoexcitation, $n_c$ is conserved and it is denoted by $\bar{n}_c$, whereas the time profile of $|\Delta_0|$ exhibits an oscillation corresponding to the Higgs amplitude mode [21]. The time average of $|\Delta_0|$ is denoted by $\overline{|\Delta_0|}$, which is defined in Eq. (6). In Fig. 11 we show $\bar{n}_c$, $|\Delta_0|$, $\Delta E$, and $\tau = 0$ denoting the time average of $\alpha_k = 0$ as functions of $|F_0|/\omega$, where $k = 0$ is the location of the gap in the ground state. The time average is taken with $\tau_i = 20$ and $\tau_f = 50$. For $|F_0|/\omega \lesssim 1$, the $|F_0|/\omega$ dependence of these quantities is similar to that in the atomic limit shown in Fig. 8 indicating that the dynamics are qualitatively described by the Rabi oscillation even when the bands are formed. As shown in Fig. 11(a), $\bar{n}_c$ has a peak at $|F_0|/\omega \simeq 0.5$ which is comparable to the case of the atomic limit [Fig. 4(a)]. For large $|F_0| (|F_0|/\omega \gtrsim 1)$, a cyclic behavior of physical quantities that characterizes the Rabi oscillation becomes less clear. When we employ the rectangular envelope for $F(\tau)$, the cyclic behavior appears even in the region of large $|F_0|/\omega$ (Fig. 13).
Here we mention the choice of the values of $\tau_i$ and $\tau_f$ in Eq. (3). After the photoexcitation, physical quantities generally show oscillations in time. In the time-dependent HF method, the center of such an oscillation is almost constant because dephasing processes via electron correlations are not taken into account. In Fig. 12 we show the time profile of $|\Delta_0|$ where we use the $U = 4$, $U' = 3.3$, $\mu_C = 2.5$, $\omega = 1.25$, and $F_0 = -0.5$, which gives a large enhancement of $|\Delta_0|$. After the photoexcitation, $|\Delta_0|$ exhibits the Higgs amplitude mode with a period of about 4.0, whereas the center of its oscillation is almost constant. In order to show explicitly how the values of $\tau_i$ and $\tau_f$ affect the time average, we define the time average of a physical quantity $X(\tau)$ taken in the range from $\tau - \tau_m/2$ to $\tau + \tau_m/2$ as,

$$X_{av}(\tau) = \frac{1}{\tau_m} \int_{\tau - \tau_m/2}^{\tau + \tau_m/2} X(\tau')d\tau'.$$

The time profiles of $|\Delta_0|_{av}(\tau)$ with $\tau_m = 10$ and 30 are shown in Fig. 12 which indicates that their difference is very small for $\tau > 20$. We note that $|\Delta_0|$ with $\tau_i = 20$ and $\tau_f = 50$ presented in Fig. 11(b) corresponds to $|\Delta_0|_{av}(\tau)$ with $\tau_m = 30$ at $\tau = 35$. From these results, we confirm that, when $\tau_m = \tau_f - \tau_i$ is larger than the oscillation period of $|\Delta_0|$ and $\tau_i$ is taken sufficiently after the photoexcitation, the value of $\tau_m$ has little effects on the results. For the relevance to experiments, if we use $\tau_f = 0.4$ eV for Ta$_2$NiSe$_5$ [32], $\tau_f - \tau_i = 30$ corresponds to 50 fs, which is comparable to time resolution of recent pump-probe measurements [9]. When $|\Delta_0|$ is small after the photoexcitation, the period of the Higgs mode may become long. However, in such cases the amplitude of the Higgs mode becomes small and thus the choice of $\tau_i$ and $\tau_f$ does not largely affect the results.

![FIG. 12. Time profiles of $|\Delta_0|$ and $|\Delta_0|_{av}(\tau)$. We use $\tau_m = 10$ and 30 for $|\Delta_0|_{av}(\tau)$. The other parameters are $U = 4$, $U' = 3.3$, $\mu_C = 2.5$, $\omega = 1.25$, and $F_0 = -0.5$. The double-headed arrow indicates the range $0 < \tau < 2\tau_w$ of application of an electric field.](image)

Here we consider $F(\tau)$ with the rectangular envelope [Eq. (3)]. As in the case of the atomic limit, we discuss our results from the viewpoint of the Rabi oscillation. In Fig. 13 we show $\tilde{n}_c$, $|\Delta_0|$, $\Delta E$, and $\tau_{k=0}$ as functions of $|F_0|/\omega$. The oscillatory behavior in these quantities is more evident than that in Fig. 11 where the gaussian en-
velope is employed for $F(\tau)$. General tendencies of Figs. 13(a), 13(b), and 13(c) are similar to those of Figs. 4(a), 4(b), and 4(c), respectively. This means that if we apply Eq. (29) to the case of nonzero transfer integrals, the value of $p$ is almost unchanged from that in the atomic limit. Compared to the results with $U = 4$, the oscillatory behavior is more prominent for those with $U = 8$ where the system is closer to the atomic limit.

In Fig. 14 the time averages of the transient energy levels, $\overline{E}_{k\gamma\sigma}$, with $\gamma$ being the band index, are shown for $U = 4$. The transient energy levels $E_{k\gamma\sigma}(\tau)$ are obtained by diagonalizing Eq. (6). We use $|F_0|/\omega = 0.3$, 1.2, and 1.9, for which the values of $|\Delta_0|$ are indicated by the arrows in Fig. 13(b). For $|F_0|/\omega = 0.3$ and 1.9, the gap in $E_{k\gamma\sigma}$ is larger than that in the ground state because of the enhancement of $|\Delta_0|$, whereas it becomes smaller for $|F_0|/\omega = 1.2$ where $|\Delta_0|$ is suppressed.

Since the one-particle Hamiltonian is described by a $2 \times 2$ matrix, we can expect that the Rabi oscillation occurs for each $k$. In order to confirm this, in Fig. 13(a) we show $\overline{\tau}_k$ on the $(|F_0|/\omega, k)$ plane for $U = 4$. It is apparent that $\overline{\tau}_k$ exhibits an oscillation with respect to $|F_0|/\omega$. The oscillation amplitude depends on $k$ and is large around $k = 0$ that is the location of the initial gap, whereas the period of oscillation is nearly independent of $k$. As shown in Fig. 13 the periodic behavior of $\overline{\tau}_k$ with respect to $|F_0|/\omega$ corresponds to those in $n_c$, $|\Delta_0|$, and $\Delta E$. By using Eqs. (22) and (23), we have

$$|\Delta(k)| = \sqrt{n_c(k)(1 - n_c(k))},$$

from which the relation between $\alpha_k$ and $n_c(k)$ is obtained, as shown in Fig. 13(b) for the case of $k = 0$. A similar relation is obtained even if we choose another $k$ (not shown). In the figure, we depict $(\alpha_{k=0}, n_c(0))$ in the ground state and $(\overline{\tau}_{k=0}, n_c(0))$ for $|F_0|/\omega = 0.3$, 1.2, and 1.9. The periodic change in the position as a function of $|F_0|/\omega$ is similar to that in the atomic limit discussed in Sect. III A. These results show that the periodic behavior of $\overline{\tau}_k$ brings about that of $n_c(k)$. Thus, the $|F_0|/\omega$ dependence of physical quantities is essentially caused by the Rabi oscillation of each one-particle state.

In order to understand the $|F_0|/\omega$ dependence of $|\Delta_0|$ in Fig. 13(b) more accurately, it is necessary to discuss the phase of $\Delta(k)$ as well as the $k$ dependence of $n_c(k)$ and $\Delta(k)$. They have been shown to have an important role in determining whether the photoinduced enhancement of the excitonic gap occurs [21]. The order parameter $\Delta_0$ is related with $\Delta(k)$ by

$$\Delta_0 = \frac{1}{N} \sum_k \Delta(k),$$

and we define their phases as

$$\Delta(k) = |\Delta(k)| e^{i\theta_k},$$

and

$$\Delta_0 = |\Delta_0| e^{i\theta}.$$
is suppressed. In the former, the Rabi oscillations of a single system state with different \( k \) values work cooperatively to induce the gap enhancement.

As we have shown in the previous paper \([19]\), when \( U' \) is small and the initial EI is of BCS type, the time evolution of \( \theta_k \) induces a destructive interference to hinder the enhancement of \( |\Delta_0| \), which is also the case for \( U' = 0 \). Since the excitonic order for small \( U' \) has a long correlation length, its photoinduced dynamics cannot be understood in terms of the Rabi oscillation in the atomic limit. However, when the initial state is of BEC type, the phases \( \theta_k \) are nearly in phase and the Rabi oscillations for different \( k \) values work cooperatively to enhance \( |\Delta_0| \).

### IV. EFFECTS OF ELECTRON-PHONON COUPLING

We investigate effects of phonons on the photoinduced dynamics within the HF approximation. We consider the additional terms to Eq. (1), which are used in \([19]\),

\[
\hat{H}_{\text{eph}} = g \sum_{i\sigma} (\hat{b}_i + \hat{b}_i^\dagger) (c_{i\sigma}^\dagger f_{i\sigma} + f_{i\sigma}^\dagger c_{i\sigma}),
\]

\[
\hat{H}_p = \omega_p \sum_i \hat{b}_i^\dagger \hat{b}_i,
\]

where \( \hat{b}_i \) (\( \hat{b}_i^\dagger \)) is the annihilation (creation) operator for the phonon at the \( i \)th site. The e-ph coupling constant and the phonon frequency are denoted by \( g \) and \( \omega_p \), respectively. We define the expectation value of the lattice displacement, \( y_p = (\langle \hat{b}_i \rangle + \langle \hat{b}_i^\dagger \rangle) \), which is assumed to be independent of \( i \). The time evolution of the system is computed as follows \([19]\). For phonons, we treat them as classical variables and numerically solve the equation of motion for \( y_p \) that is written as

\[
\frac{d^2 y_p}{dt^2} = -\omega_p^2 y_p - 8g \omega_p \text{Re} \Delta_0,
\]

from which we have

\[
y_p = -\frac{8g}{\omega_p} \text{Re} \Delta_0,
\]

in the ground state. For the electronic part, we employ Eq. (7). In this section, we use \( F(\tau) \) with the gaussian envelope \([ Eq. (3) \] \). The results obtained by the rectangular envelope are given in Appendix C.

#### A. Atomic limit

First, we discuss the case of the atomic limit (\( t_f = t_c = 0 \)). In the ground state, we can show that Eq. (16) holds even in the presence of the e-ph interaction. From Eqs. (16) and (34), the ground-state energy \( E_g^{\text{AL}} \) is written as

\[
E_g^{\text{AL}} = \frac{1}{2} (U - U'_g) n_c^2 + (\mu_C + U + U'_g) n_c + U,
\]

where \( U'_g = U' - 8g^2/\omega_p \). This leads to \( n_c = 1 - \mu_C/(U - U'_g) \) and the critical value of \( U' \) for the EP-DP phase boundary is given by \( U'_c = U - \mu_C + 8g^2/\omega_p \). In Fig. 14 we show the time average of \( n_c \), which is denoted by \( \overline{n_c} \), as a function of \( |F_0|/\omega \) for \( g = 0.01 \) and \( g = 0.02 \) with \( U = 1 \), \( U' = 0.45 \), \( \mu_C = 0.5 \), and \( \omega_1 = 0.1 \). For \( g = 0.01 \) (\( g = 0.02 \)), we have \( U'_c = 0.258 \) (\( U'_c = 0.532 \)). The value of \( \omega \) is so chosen that it corresponds to the energy difference between the two levels. The time average is taken with \( \tau_f = 100 \) and \( \tau_f = 400 \) considering the long time-scale of phonons, \( 2\pi/\omega_p \). It is apparent that the Rabi oscillation appears even for nonzero \( g \). We note that this is also the case when we use the rectangular envelope for \( F(\tau) \) (Appendix C). For \( g = 0.02 \), we depict the time average of \( \overline{|\Delta_0|} \) and \( \overline{y_p} \) in Figs. 10(a) and 10(b), respectively. When \( |F_0|/\omega = 0.1 \), \( |\Delta_0| \) and \( y_p \) oscillate around the ground-state values. However, for large \( |F_0|/\omega = 0.4 \), \( |\Delta_0| \) is enhanced and \( y_p \) oscillates around zero indicating that the effect of the lattice displacement basically disappears. In Fig. 18(c), we show the trajectory of \( (S_x, S_y) \) where we define \( S_x = \text{Re} \Delta_0 \) and \( S_y = \text{Im} \Delta_0 \) in the pseudospin representation. The description of the pseudospin representation and the trajectory of \( (S_x, S_y) \) for \( g = 0 \) are given in Appendix A. For small \( |F_0|, \theta \) that is defined in Eq. (22) is confined near zero. This is because the phase mode is massive.
in the presence of the lattice displacement $|\mathbf{F}_0|$. On the other hand, $\theta$ rotates for large $|\mathbf{F}_0|$, which is qualitatively the same as that for $g = 0$ (Fig. 25 in Appendix A).

B. One-dimensional model

Next, we show results with nonzero transfer integrals ($t_f = -t_c = 1$). We compute $\bar{n}_c$ as a function of $|\mathbf{F}_0|/\omega$ for $g = 0.02$ and 0.04 where the lattice displacements in the ground state are $y_p = -0.217$ and $y_p = -0.527$, respectively. Here we use $U = 4$, $U' = 3.3$, and $\mu_C = 2.5$. The used value of $\omega$ corresponds to the initial gap. As shown in Fig. 19, the introduction of the e-ph coupling does not largely affect the $|\mathbf{F}_0|/\omega$ dependence of $\bar{n}_c$ as in the case of the atomic limit. For $g = 0.04$, we show the time profiles of $|\Delta_0|$ and $y_p$ in Figs. 20(a) and 20(b), respectively, whereas the trajectory of $(S_x, S_y)$ is shown in Fig. 20(c). We use $|\mathbf{F}_0|/\omega = 0.1$ ($|\Delta_0|/\Delta_0(\tau = 0) = 0.97$) and $|\mathbf{F}_0|/\omega = 0.4$ ($|\Delta_0|/\Delta_0(\tau = 0) = 1.53$). The results are qualitatively the same as those in the atomic limit shown in Fig. 18. These results indicate that the e-ph coupling does not have a significant role on the photoinduced gap enhancement based on the Rabi oscillation.

V. CORRELATION EFFECTS

In this section, we examine effects of the electron correlation that are ignored in the HF approximation. By using the ED method, we calculate ground-state properties and photoinduced dynamics of the two-orbital Hubbard model. We do not consider the e-ph coupling for simplicity. When we use single cycle pulses for photoexcitations, we adopt $F(\tau)$ with the gaussian envelope and the results obtained with the rectangular envelope are given in Appendix D.

A. Ground state

In the ground state, we compute the $U'$ dependence of $n_c$ where we use $U = 4$, $\mu_C = 2.5$ and the system size $N = 6$. As shown in Fig. 21, $n_c$ monotonically decreases with increasing $U'$ and it becomes zero at $U'_{cr} = 3.5$. This behavior is consistent with the HF results shown in Fig. 10 where we have $U'_{cr} = 3.37$. The qualitative difference between the HF and ED results is that for $U' < U'_{cr}$ the excitonic order parameter $\Delta_0$ is nonzero in the former.
whereas it is zero in the latter. We note that by the ED method we inevitably have a ground state with $\Delta_0 = 0$ because of the finiteness of the system. For $U' > U'_{cr}$, both methods give the BI phase with $\Delta_0 = n_c = 0$ as the ground state. In this phase, the gap $E_1 - E_0$ increases almost linearly with $U'$ as shown in Fig. 21, where $E_0$ and $E_1$ are the energies of the ground and first excited states, respectively. This behavior is also consistent with the HF results [21]. In the following, we consider the BI phase ($U' > U'_{cr}$) as the initial state before photoexcitation for comparison.

### B. Photoinduced dynamics

The time evolution of the system is obtained by numerically solving the time-dependent Schrödinger equation for the exact many-electron wave function $|\Psi(\tau)\rangle$ as

$$|\Psi(\tau + d\tau)\rangle = \exp\left[-i\tau \hat{H}_{tot}\left(\tau + \frac{d\tau}{2}\right)\right]|\Psi(\tau)\rangle,$$

where $\hat{H}_{tot}(\tau) = \hat{H} + \hat{H}_D(\tau)$ and we use $d\tau = 0.01$. We use $U = 4$, $\mu_C = 2.5$, and $U' = 3.9 > U'_{cr}$. The light frequency is set at $\omega = 0.7$ that is near the gap $E_1 - E_0 = 0.08$. In the following, we first show results with single cycle pulses and then discuss the case of CW excitations.

#### 1. Excitations with single cycle pulse

In Fig. 22 we show the time profiles of $n_c$ and $|\Delta_0|$ for different values of $F_0$ with $0 < F_0/\omega \lesssim 1$. After the photoexcitation, $n_c$ is conserved, whereas $|\Delta_0|$ keeps oscillating. The value of $n_c$ increases with increasing $F_0$, and then it decreases when we increase $F_0$ further ($F_0 = 0.6$). As shown in Fig. 22(b), there is no clear indication of a strong dephasing in the order parameter that should suppress $|\Delta_0|$ after the photoexcitation with $0 < F_0/\omega \lesssim 1$. Moreover, we do not find rapid thermalization: the oscillation in $|\Delta_0|$ persists long after the photoexcitation with $0 < F_0/\omega \lesssim 1$. Although the finite size effects may play a role, our results at this stage do not indicate that the correlation effects seriously hinder the enhancement of $|\Delta_0|$. We depict $n_c$, $|\Delta_0|$, $\Delta E$ and $\bar{\alpha}$ as functions of $F_0/\omega$ in Fig. 23 where the time average of $|\Delta_0|$ is taken with $\tau_i = 50$ and $\tau_f = 100$. Here the overlap $\alpha$ is defined by $\alpha = \langle |\Psi(\tau)|\Psi(0)\rangle^2$. After the photoexcitation, $\alpha$ is conserved and its value is denoted by $\bar{\alpha}$. Notably, our results indicate that for $F_0/\omega \lesssim 1$, the $F_0/\omega$ dependence

![FIG. 19](image1)

**FIG. 19.** $\pi_c$ as a function of $|F_0|/\omega$ for different values of $g$ with $U = 4$, $U' = 3.3$, and $\mu_C = 2.5$ for the case of nonzero transfer integrals ($t_f = -t_c = 1$). We use $\omega = 1.25$, 1.30, and 1.42 for $g = 0$, 0.02, and 0.04, respectively. The horizontal dashed lines indicate the corresponding equilibrium values.

![FIG. 20](image2)

**FIG. 20.** Time evolutions of (a) $|\Delta_0|$ and (b) $y_p$, and (c) trajectory of $(S_x, S_y)$ for $|F_0|/\omega = 0.1$ and 0.4. We use $g = 0.04$ and the other parameters are the same as those in Fig. 19. In (a) and (b), the horizontal lines indicate the corresponding equilibrium values. In (c), we depict the time domain $-20 \leq \tau \leq 300$ ($-20 \leq \tau \leq 50$) for $|F_0|/\omega = 0.1$ ($|F_0|/\omega = 0.4$), and the solid square indicates the initial position of $(S_x, S_y)$. 

![FIG. 21](image3)
that the enhancement of the HF method shown in Fig. 11. This strongly suggests that in this region the dephasing has an important role in determining the value of $|\Delta_0|$. This is in contrast to the case with $F_0/\omega \lesssim 1$ where $|\Delta_0|$ can be largely enhanced. When we use the rectangular envelope for $P(\tau)$ Eq. (13), the cyclic behavior in physical quantities becomes more evident, which we show in Appendix D.

When $U' < U'_{cr}$, our ED results do not show a clear evidence of the Rabi oscillation. Specifically, for $U' = 3.2 < U'_{cr}$, the oscillatory dependence of $a$ and $\Delta E$ on $F_0/\omega$ that appears in Fig. 23 for the case of $U' = 3.9 > U'_{cr}$ ($F_0/\omega \lesssim 1$) is less pronounced. Although the $F_0/\omega$ dependence of $|\Delta_0|$ is similar to that in Fig. 23, for $\tilde{\alpha}_c$ the finite size effect is more severe than that with $U' > U'_{cr}$ and the result with $N = 4$ is qualitatively different from that with $N = 6$ even for $F_0/\omega < 1$. We speculate that these results are due to the metallic ground state with $\Delta_0 = 0$ in the ED method. When the system is metallic ($\Delta_0 = 0$), it has basically gapless excitations and thus it is far from a two-level system.

2. Excitations with continuous-wave laser

Next, we consider the case of CW excitations and examine time evolutions of physical quantities from the viewpoint of the Rabi oscillation. In Fig. 25 we show the time profiles of $n_c$ and $\alpha$ for different values of $F_0$ with $U' = 3.9 > U'_{cr}$ for which the ground state before the photoexcitation is the BI. They exhibit an oscillation, the period of which becomes shorter with increasing $F_0$. For small $F_0$ ($F_0 \lesssim 0.1$), the time profile of $n_c$ is well described by a single sinusoidal function of the form Eq. (20) as shown in Fig. 25(a), and the minimum value in the oscillation is close to the ground-state value of $n_c$ ($= 0$). Correspondingly, a nearly sinusoidal oscillation appears in $\alpha$. It is notable that we have $\alpha \sim 1$ when $n_c \sim 0$, whereas $\alpha \sim 0$ when $n_c$ exhibits its maximum. These behaviors are consistent with the Rabi oscillation as we have discussed in Sect. III. With increasing $F_0$, the oscillatory profiles in $n_c$ and $\alpha$ gradually become more complex. For $F_0 \gtrsim 0.15$, a single sinusoidal function does not fit well to the data. Also, the minimum (maximum) in the oscillation of $n_c$ ($\alpha$) departs from its ground-state value, which is in contrast to the case with $F_0 \lesssim 0.1$.

In Fig. 26(a), we show the Fourier transform of $n_c$ that is calculated from the data for $50 \leq \tau \leq 400$. There is a sharp peak in each spectrum and its position that is denoted by $\Omega$ becomes larger for larger $F_0$. In Fig. 26(b), we plot the $F_0$ dependence of $\Omega$. For $F_0 \lesssim 0.1$, $\Omega$ is nearly proportional to $F_0$: a function $\Omega = pF_0$ with $p = 2.70$ fits well to the data. This result, in conjunction with the time profiles of $n_c$ and $\alpha$ shown in Fig. 25,
with the HF approximation where the linearity characterizing the Rabi oscillation basically appears for large $F_0/\omega$ unless the constant $\alpha$ in $\Delta_0$ dependence of physical quantities after the photoexcitation. In fact, Fig. 23 obtained with single cycle pulses indicates the signature of the Rabi oscillation for $F_0/\omega \lesssim 1$ ($F_0 \lesssim 0.5$).

Finally, we examine the correspondence between the results with CW excitations and those with single cycle pulses in the same way as we have done in Sect. III A. We apply $\Omega_0 = pF_0$ with $p = 2.70$ to Eq. (20). For $\alpha$, we use Eq. (17) with $n_0^G = \Delta_0^G = 0$. By setting $\tau = 2\pi/\omega$ in these equations, we can deduce that $\tilde{n}_c$ for single cycle pulses exhibits a minimum (maximum) at $F_0/\omega \sim 0.19$ unless the constant $A$ in Eq. (20) strongly depends on $F_0$. For $\tilde{n}_c$, this value of $F_0/\omega$ is consistent with the results shown in Fig. 24 where it exhibits a minimum at $F_0/\omega = 0.21$. For $\tilde{n}_c$, its first maximum is located at $F_0/\omega = 0.46$ which is larger than the above estimation. This discrepancy mainly comes from an increase in the amplitude of $n_\omega$ with increasing $F_0$ [Fig. 25(a)]: the $F_0$ dependence of $A$ is important in determining the maximum of $n_\omega$. This is in contrast to the time evolutions of $\alpha$ where it becomes almost zero in its first oscillation irrespective of the value of $F_0$ [Fig. 25(b)]. We note that this argument also holds for the case with the rectangular envelope where the maximum of $n_\omega$ and the minimum of $\tilde{n}_c$ are located at $F_0/\omega = 0.21$ and 0.32, respectively, as shown in Appendix D.
Finally, we discuss possible experimental observation of photoinduced gap enhancement as well as the relevance of our results to $\text{Ta}_2\text{NiSe}_5$. Recent theoretical studies\cite{32,34} have shown that various equilibrium properties of $\text{Ta}_2\text{NiSe}_5$ such as the ARPES spectra\cite{52} and the temperature dependence of magnetic susceptibility\cite{29} can be reproduced by two- or three-orbital Hubbard models. Effects of the structural distortion observed at $T_C$ have been investigated using a three-orbital Hubbard model with e-ph interactions by the HF approximation\cite{50}. It has been shown that the values of the e-ph interaction strengths needed to reproduce the experimentally observed distortion are one order of magnitude smaller than those of the transfer integrals and the e-e interaction strengths. Then, it has been argued that the EI in $\text{Ta}_2\text{NiSe}_5$ is ascribed to the BEC of electron-hole pairs which cooperatively induce the instability of the lattice distortion. These studies suggest that the photoinduced dynamics obtained in this paper based on the two-orbital Hubbard model [Eq. (1)] would be relevant to $\text{Ta}_2\text{NiSe}_5$.

In our mechanism, photoinduced gap enhancement occurs purely electronically when $\omega$ is comparable to the excitonic gap. Moreover, we have shown that e-ph couplings do not affect our results qualitatively. When $\omega$ is much larger than the excitonic gap, which is the case in recent experiments\cite{9}, a theoretical study has shown that e-ph couplings are crucially important for the appearance of the gap enhancement\cite{19}. Thus, our mechanism is considered as an alternative route to this phenomenon.

In this paper, we consider the case where the upper and lower bands have the same bandwidth ($t_c = -t_f$). However, even when the two bandwidths are different\cite{50}, we expect that the gap enhancement by the Rabi oscillation occurs as long as the initial system is a BEC-type EI or a nearby BI. This is because their dynamics should be basically understood from the real-space picture\cite{21} where the analysis in the atomic limit presented in this paper is valid.

In order to examine the relevance of our results to experiments, we estimate the number of absorbed photons per site $n_{\text{ph}}$. When $U = 4$ and $U' = 3.3$ [Fig. 11(c)], we have $\Delta E = 0.398$ for $|F_0|/\omega = 0.44$ at which $|\Delta_0|$ exhibits the first peak as a function of $|F_0|/\omega$. This corresponds to $n_{\text{ph}} = \Delta E/\omega = 0.32$. We note that a sizable gap enhancement appears with much smaller values of $n_{\text{ph}}$. For instance, 15% enhancement in $|\Delta_0|$ is obtained for $|F_0|/\omega = 0.2$ where we have $n_{\text{ph}} = 0.017$. In $\text{Ta}_2\text{NiSe}_5$, K. Okazaki et al. have reported that when the incident pump fluence is 1 mJ/cm$^2$, $n_{\text{ph}} \sim 0.1$ per Ni atom whose 3$d$ orbital hybridizes Se 4$p$ orbital and forms a hole band\cite{51}. The threshold pump fluence for the appearance of the gap enhancement reported in Ref. 9 is $F_C = 0.2$ mJ/cm$^2$, which may correspond to $n_{\text{ph}} \sim 0.02$. This suggests that the pump fluence used in the current experimental studies is enough to observe the gap enhancement based on our mechanism unless $n_{\text{ph}}$ depends largely on the value of the initial gap. However, at present a direct comparison between theoretical and experimental estimates is difficult by the following reasons. Firstly, in our model, we assume that the incident light induces the
dipole transition whereas it does not affect the intra-orbital electron motion. In order to realize this situation in real materials, the direction of light polarization as well as the crystal structure of the material are crucially important. For a material with a quasi-one dimensional structure like Ta$_2$NiSe$_5$, this indicates that the polarization of light should be perpendicular to the chain. The value of the matrix element for the dipole transition between the two bands is also important. Secondly, the pump-light frequency should be nearly tuned to the resonance condition. Note that in this case a recent theoretical study has shown that the gap enhancement does not appear when the incident light only affects the intra-orbital electron motion [22]. Thirdly, the estimation of $n_{ph}$ by the time-dependent HF method may be quantitatively inaccurate since it ignores the correlation effects [10, 52].

With regard to this point, from our ED results on small clusters with $U = 4$ and $U' = 3.9$ where the ground state is the BI, we have $n_{ph} = 0.10$ for $F_0/\omega = 0.1$ at which $|\Delta_0|$ is maximally enhanced [Fig. 23]. This value of $n_{ph}$ is comparable to the above-mentioned HF results.

We note that the gap enhancement with the help of the Rabi oscillation is irrespective of the dimensionality of the system. In fact, our results for one-dimensional systems are qualitatively unaltered even in the two-dimensional case [21]. Moreover, our ED results suggest that the Rabi-oscillation-assisted gap enhancement appears even when the effects of quantum fluctuations are considered, although how the dephasing and thermalization affect the dynamics remains as a future important problem.

In summary, we investigated dynamics of EIs induced by electric dipole transitions using the two-orbital Hubbard model. Through the HF analysis of the dynamics in the atomic limit, we have shown that the photoinduced gap enhancement in the EI for single cycle pulses reported previously [21] is explained in terms of the Rabi oscillation. The signature of the Rabi oscillation appears as a periodic behavior of physical quantities after the photoexcitation as functions of the dipole field strength $F_0$. We emphasize that although the Rabi oscillation is a one-site problem, it represents the essential feature of the photoinduced dynamics in the thermodynamic limit in the parameter range that we have considered in this paper. We have performed the ED calculations which strongly suggest the robustness of this phenomenon against the correlation effects and thus corroborate our HF results. The effects of the e-ph coupling have been examined within the HF approximation, indicating that they do not have a significant role on the gap enhancement in the present situation. Based on the present results and our previous work [21], the condition for inducing the gap enhancement is summarized as follows: (i) The initial state is an EI in the BEC regime or a BI that is located near the EI. (ii) The pump-light frequency $\omega$ is near the initial gap. (iii) There is an optimal value of $F_0$ for enhancing the excitonic gap, which satisfies the relation $\Omega_R/\omega \sim 1/2$ with the Rabi frequency $\Omega_R$.

\[ \Psi(\tau) = \psi(z, \tau) \]

\[ (A1) \]

where $\sigma_\gamma$ ($\gamma = x, y, z$) are the Pauli matrices and we omit the spin index in $\Psi_\sigma$ for brevity. With this representation, the expectation values of the pseudospin $S_\gamma = \langle \hat{S}_\gamma \rangle$ components are written as

\[ S_x(\tau) = \text{Re}\Delta_0, \]

\[ S_y(\tau) = \text{Im}\Delta_0, \]

\[ S_z(\tau) = \frac{1}{4}(n_c - n_f), \]

\[ (A2a, b, c) \]

which give $\Delta_0 = S_x + iS_y$ and $n_c = 2S_z + 1$. By using the equation of motion for the pseudospin operators, the time evolution of $S = (S_x, S_y, S_z)$ is given by

\[ \partial_\tau S = B(\tau) \times S(\tau), \]

\[ (A3) \]
for $U' = 0.45$ and $F_0' = -0.16$ for $U' = 0.55$. Such a dynamical transition has been previously reported in one-dimensional excitonic insulators within the HF theory \cite{19}.

In the following, we examine the difference between the dynamics for $|F_0| < |F_0'|$ and that for $|F_0| > |F_0'|$. First, we consider the case of $U' = 0.45$ where the initial state is in the EP. We show the trajectory of $(S_x, S_y)$ with $F_0 = -0.05$ ($|F_0| < |F_0'|$) and $-0.2$ ($|F_0| > |F_0'|$) under CW excitations with $\omega = 0.45$ in Fig. \ref{fig30}. For $F_0 = -0.05$, $(S_x, S_y)$ is bound near the ground-state position, whereas it is unbound for $F_0 = -0.2$. This corresponds to bound and unbound oscillations in $n_c$ for $F_0 = -0.05$ and $-0.2$ (Fig. \ref{fig31a}), respectively. In Fig. \ref{fig31a}, we show the Fourier transform of $n_c$, for small $|F_0| (< |F_0'|)$, indicating that $n_c$ has one slow oscillation component with frequency $\Omega_S \lesssim 0.15$ and two fast components with frequencies $\Omega^{\pm} \approx 0.15$ near $\omega$, which we can write as $\Omega^{\pm} = \omega \pm \delta\Omega$. Both $\Omega_S$ and $\delta\Omega$ increase with increasing $|F_0|$. When $|F_0|$ is small ($|F_0| < 0.06$), the peak at $\Omega^S$ is dominant, whereas
for in Fig. 29. These results indicate that the dynamics
indicated by the solid (dashed) arrows. (b) \(|F_0|\) dependence
of \(\Omega^S\). The solid line is fit to the data with \(|F_0| \leq 0.06\). The
fitting result in Fig. 31 for \(U' = 0.45\) is also shown by the
dashed line.

those at \(\Omega^f\pm\) become dominant for \(0.06 < |F_0| < |F_0^c|\).
As we increase \(|F_0|\) further (\(|F_0| > |F_0^c|\)), the spectra
change drastically as we have shown in Fig. 31. In Fig. 31(b), we show the \(|F_0|\) dependence of \(\Omega^S\). When \(|F_0| \leq 0.06\), \(\Omega^S\) is proportional to \(|F_0|\) and we have \(\Omega^S = p|F_0|\)
with \(p = 0.63\) that is different from the value \((p = 1.24)\)
obtained in Sect. III A for \(|F_0| > |F_0^c|\). The region of \(|F_0| \leq 0.06 < |F_0| < |F_0^c|\) where the value of \(p\) largely deviates coincides with that where \(\Omega^S\) exhibits the abrupt increase in Fig. 31. These results indicate that the dynamics for \(|F_0| < |F_0^c|\) have a character different from that for \(|F_0| > |F_0^c|\).

Next, we show results with \(U' = 0.55\) where the initial
state is in the DP. The trajectory of \((S_x, S_y)\) for \(F_0 = -0.05\) \((|F_0| < |F_0^c|)\) and \(F_0 = -0.25\) \((|F_0| > |F_0^c|)\) with \(\omega = 0.6\) is depicted in Fig. 32. Similar to the case of \(U' = 0.45\), \((S_x, S_y)\) is bound near its initial position for \(|F_0| < |F_0^c|\), whereas it is unbound for \(|F_0| > |F_0^c|\). In Fig. 31(a), we show the Fourier transform of \(n_c\) for \(|F_0| < |F_0^c|\).

The dominant oscillation components in \(n_c\) have frequencies \(\Omega^f = \omega \pm \delta\omega\), and there is a slow oscillation component with \(\Omega^S = 2\delta\omega\) whose amplitude is higher order in \(|F_0|\). When \(|F_0|\) is small, we can solve Eq. 3 in the lowest order of \(F_0\) with the initial condition \(S(\tau = 0) = (0, 0, -1/2)\) as

\[
\begin{align*}
S^x(\tau) &= \frac{F_0}{\omega^2 - a^2} (a \sin \omega \tau - \omega \sin a \tau), \quad (A5a) \\
S^y(\tau) &= \frac{\omega F_0}{\omega^2 - a^2} (\cos a \tau - \cos \omega \tau), \quad (A5b) \\
S^z(\tau) &= \frac{\omega F_0}{\omega^2 - a^2} \left[ \frac{1 - \cos(\omega + a) \tau}{\omega + a} + \frac{1 - \cos(\omega - a) \tau}{\omega - a} + \frac{1 - \cos 2\omega \tau}{2\omega} \right], \quad (A5c)
\end{align*}
\]

where \(a = \mu_C - U + U'\). From Eq. (A5c), we find that for \(|F_0| \rightarrow 0\), \(\delta \Omega \rightarrow a = 0.05\) and there is an oscillation component with frequency \(2\omega = 1.2\), which are consistent with the numerical results shown in Fig. 33(a). As we increase \(|F_0|\), \(\Omega^S\) increases. \(\Omega^S = 0.1 + 10|F_0|^2\) is well fit to the data as shown in Fig. 33(b). These results indicate that the dynamics for \(|F_0| < |F_0^c|\) is essentially different from that for \(|F_0| > |F_0^c|\) as in the case of \(U' = 0.45\). In fact, the spectra of \(n_c\) for \(|F_0| > |F_0^c|\) (not shown) are largely different from those for \(|F_0| < |F_0^c|\).

Appendix B: Effects of \(\tau\)-dependence of \(n_c\) and \(\Delta_0\) in Eq. (10) on the dynamics

As we have mentioned in III A 3, Eq. (10) possesses \(\tau\)-dependent mean-field order parameters from which the time evolution operator is constructed. In order to examine how their \(\tau\)-dependence affects the dynamics in the atomic limit, we artificially replace \(n_c\) and \(\Delta_0\) in the time evolution operator by \(n_c^G\) and \(\Delta_0^G\), respectively, and compute the time profile of \(n_c\) under CW excitations. The parameters we used are the same as those in Fig. 6. The results are shown in Fig. 34(a). Compared with Fig. 4, a large oscillation in \(n_c\) appears even when \(|F_0|\) is small. From the Fourier spectra shown in Fig. 34(b), we obtain \(\Omega \sim p|F_0|\) with \(p = 0.86\). This result indicates that the \(\tau\) dependence of the order parameters is important.
in determining the dynamics for small $|F_0| (< 0.1)$. However, it does not alter the dynamics qualitatively for larger $|F_0|$ where the Rabi oscillation appears in Fig. 6. These facts give a reason why the quantitative difference between the value of $\tilde{n}_c$ for single cycle pulses and that of $n_c$ at $\tau = 2\pi/\omega$ computed from Eq. (20) becomes large for small $|F_0| (|F_0|/\omega \lesssim 0.3)$, which can be seen in Fig. 9.

Appendix C: HF results in the presence of phonons for the case of rectangular-envelope pulse

We show the $|F_0|/\omega$ dependence of $\tilde{n}_c$ in the presence of the e-ph coupling when we use $F(\tau)$ with the rectangular envelope. In Fig. 35 the results in the atomic limit are depicted. The parameters are the same as those in Fig. 17. For the one-dimensional model with $t_f = 1$ and $t_c = -1$, the results are shown in Fig. 36 where the parameters are the same as those in Fig. 19. From Figs. 35 and 36 we confirm that the e-ph coupling has little effects on the $|F_0|/\omega$ dependence of $\tilde{n}_c$ as in the case of the gaussian-envelope pulse shown in Figs. 17 and 19.

Appendix D: ED results for the case of rectangular-envelope pulse

In Fig. 37 we show $\tilde{n}_c$, $|\Delta_0|$, $\Delta E$, and $\tilde{\alpha}$ as functions of $F_0/\omega$ obtained by the ED method when we use $F(\tau)$ with the rectangular envelope. The parameters are the same as those in Fig. 23. For $F_0/\omega < 1$, the $F_0/\omega$ dependence of these quantities is similar to those in Fig. 23 indicating that the pulse shape does not significantly affect our results as in the case of the HF method. The cyclic behavior is evident even for $F_0/\omega > 1$, although in
FIG. 36. Same plot as Fig. 19 except that we use $F(\tau)$ with this region the increase (decrease) in $\tilde{n}_c$ and $\Delta E$ ($\tilde{\alpha}$) does not correspond to the large enhancement in $|\Delta_0|$, which is in contrast to the results with the HF method shown in Fig. 13. This is caused by the dephasing discussed in Sect. V, which suppresses $|\Delta_0|$. In fact, for $F_0/\omega \sim 1.6$ and 3.2 of the results with $N = 6$, where $\tilde{n}_c$ and $\Delta E$ exhibit a peak and $\tilde{\alpha} \sim 0$, we have confirmed that the time profile of $|\Delta_0|$ is similar to that in Fig. 21.

**ACKNOWLEDGMENTS**

This work was supported by JSPS KAKENHI Grant Nos. JP15H02100, JP16K05459, JP19K3247, and JP20K03841, MEXT Q-LEAP Grant No. JP-MXS0118067426, JST CREST Grant No. JPMJCR1901, and Waseda University Grant for Special Research Projects (Project No. 2020C-280).

[1] K. Onda, S. Ogihara, K. Yonemitsu, N. Maeshima, T. Ishikawa, Y. Okimoto, X. Shao, Y. Nakano, H. Yamochi, G. Saito, and S. Koshihara, Phys. Rev. Lett. 101, 067403 (2008).
[2] D. Fausti, R. L. Tobey, N. Dean, S. Kaiser, A. Dienst, M. C. Hoffmann, S. Pyon, T. Takayama, H. Takagi, and A. Cavalleri, Science 331, 189 (2011).
[3] T. Ishikawa, Y. Sagae, Y. Naitoh, Y. Kawakami, H. Itoh, K. Yamamoto, K. Yakushi, H. Kishida, T. Sasaki, S. Ishihara, Y. Tanaka, K. Yonemitsu, and S. Iwai, Nat. Commun. 5, 5528 (2014).
[4] W. Hu, S. Kaiser, D. Nicoletti, C. R. Hunt, I. Gierz, M. C. Hoffmann, M. Le Tacon, T. Loew, B. Keimer, A. Cavalleri, Nat. Mater. 13, 705 (2014).
[5] S. Kaiser, C. R. Hunt, D. Nicoletti, W. Hu, I. Gierz, H. Y. Liu, M. Le Tacon, T. Loew, D. Haug, B. Keimer, A. Cavalleri, Phys. Rev. B 89, 184516 (2014).
[6] L. Stojechevska, I. Vaskivskyi, T. Mertelj, P. Kusar, D. Svetin, S. Brazovskii, and D. Mihailovic, Science 344, 177 (2014).
[7] M. Miranov, A. Cantaluppi, D. Nicoletti, S. Kaiser, A. Perucchi, S. Lupi, P. Di Pietro, D. Pontiroli, M. Riccio, S. R. Clark, D. Jakob, A. Cavalleri, Nature 530, 461 (2016).
[8] A. Singer, S. K. K. Patel, R. Vukajlov, V. Uhlir, J. Wingert, S. Festsens, D. Zhu, J. M. Glownia, H. T. Lemke, S. Nelson, M. Kozina, K. Rossnagel, M. Bauer, B. M. Murphy, O. M. Magnussen, E. E. Fullerton, and O. G. Shyryko, Phys. Rev. Lett. 117, 056401 (2016).
[9] S. Mor, M. Herzog, D. Golez, P. Werner, M. Eckstein, N. Katayama, M. Nohara, H. Takagi, T. Mizokawa, C. Monney, and J. Stahler, Phys. Rev. Lett. 119, 086401 (2017).
[10] Y. Kawakami, Y. Yoneyama, T. Amano, H. Itoh, K. Yamamoto, Y. Nakamura, H. Kishida, T. Sasaki, S. Ishihara, Y. Tanaka, K. Yonemitsu, and S. Iwai, Phys. Rev. B 95, 201105(R) (2017).
[11] H. T. Lu, S. Sota, H. Matsueda, J. Bonca, and T. Tohyama, Phys. Rev. Lett. 109, 197401 (2012).
[12] N. Tsuji, T. Oka, H. Aoki, and P. Werner, Phys. Rev. B 85, 155124 (2012).
[13] H. Hashimoto, H. Matsueda, H. Seo, and S. Ishihara, J. Phys. Soc. Jpn. 83, 123703 (2014).
[14] H. Hashimoto, H. Matsueda, H. Seo, and S. Ishihara, J. Phys. Soc. Jpn. 84, 113702 (2015).
[15] H. Yanagiya, Y. Tanaka, and K. Yonemitsu, J. Phys. Soc. Jpn. 84, 094705 (2015).
[16] M. Nakagawa and N. Kawakami, Phys. Rev. Lett. 115, 165303 (2015).
[17] K. Yonemitsu, J. Phys. Soc. Jpn. 86, 024701 (2017).
[18] K. Ido, T. Ohgroe, and M. Imada, Sci. Adv. 3, e1700718 (2017).
[19] Y. Murakami, D. Golez, M. Eckstein, P. Werner, Phys. Rev. Lett. 119, 247601 (2017).
[20] K. Oya and A. Takahashi, Phys. Rev. B 97, 115147 (2018).
[21] Y. Tanaka, M. Daira, and K. Yonemitsu, Phys. Rev. B 97, 115105 (2018).
[22] T. Tanabe, K. Sugimoto, and Y. Ohta, Phys. Rev. B 98, 235127 (2018).
[23] N. F. Mott, Phil. Mag. 6, 287 (1961).
[24] R. S. Knox, Solid State Phys. Suppl. 5, 100 (1963).
[25] D. Jerome, T. M. Rice, and W. Kohn, Phys. Rev. 158, 462 (1967).
[26] B. I. Halperin and T. M. Rice, Rev. Mod. Phys. 40, 755 (1968).
[27] J. Kunes, J. Phys.: Cond. Mat. 27, 335201 (2015).
[28] Y. Wakisaka, T. Sudayama, K. Takubo, T. Mizokawa, M. Arita, H. Namatame, M. Taniguchi, N. Katayama, M. Nohara, and H. Takagi, Phys. Rev. Lett. 103, 286402 (2009).
[29] F. J. Di Salvo, C. H. Chen, R. M. Fleming, J. V. Waszczak, R. G. Dunn, S. A. Sunshine, James A. Ibers, J. Less-Common Met. 116, 51 (1986).
[30] Y. F. Lu, H. Kono, T. I. Larkin, A. W. Rost, T. Takayama, A. V. Boris, B. Keimer, and H. Takagi, Nat. Commun. 8, 14408 (2017).
[31] S. Li, S. Kawai, Y. Kobayashi, and M. Itoh, Phys. Rev. B 97, 165127 (2018).
FIG. 37. Same plots as Fig. 23 by the ED method except that we use $F(\tau)$ with the rectangular envelope.

[32] K. Seki, Y. Wakisaka, T. Kaneko, T. Toriyama, T. Konishi, T. Sudayama, N. L. Saini, M. Arita, H. Namatame, M. Taniguchi, N. Katayama, M. Nohara, H. Takagi, T. Mizokawa, and Y. Ohta, Phys. Rev. B 90, 155116 (2014).