Theory of Single Susceptibility for Near-field Optics
Equally Associated with Scalar and Vector Potentials

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Abstract

A nonlocal response theory was developed to describe a many-electron system within the neighborhood of a nanostructure radiating the longitudinal and transverse electric fields, which are fundamentally reduced to the scalar and vector potentials (SP and VP). The coexistence of the SP and VP incidences distinguishes such a near-field optical system from the ordinary optical system, in which only the VP (under the Coulomb gauge) incidence survives far from the light source. This fact is the motivation for equal treatment of the SP and VP as the cause of the response in the near-field optical system. In the semiclassical treatment, the linear and nonlinear single susceptibilities are derived in the form of Heisenberg operators by the functional derivatives of the action integral of the matter with respect to the SP and VP. These single susceptibilities relate the SP and VP (as the cause) to the induced charge and current densities (as the result), and guarantee charge conservation and gauge invariance; this theory is free from gauge-fixing. It is necessary to consider the quantum many-electron effect (exchange-correlation effect) to make the ground state bounded in the non-perturbed system. This is done by employing the fundamental idea of density functional theory, instead of the ordinary unequal treatment of the SP and VP, that is, remaking the SP into a Coulomb interaction between electron charges. Applying the present linear response theory to the non-metallic material in a limited near-field optical system reveals that the electric field with the associated permittivity is not suitable quantity to describe the response, instead, the SP and VP with associate single susceptibility are essential.

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I. INTRODUCTION

This paper develops a nonlocal response theory adequate for near-field optics (NFO) in the semiclassical treatment. The linear and nonlinear single susceptibilities are derived systematically by the functional derivatives of the action integral of the matter with respect to the scalar and vector potentials (SP and VP). These linear and nonlinear single susceptibilities relate the SP and VP (as the cause) to the induced charge and current densities (as the result), and guarantee charge conservation and gauge invariance. The present single susceptibilities and associated induced charge and current densities are given in the form of Heisenberg operators.

In Ref.[1], the present author discussed the linear single susceptibility, its application to an one-electron optical system, and a naive idea of employing the density functional theory. This paper is its generalization including systematic derivation of linear and nonlinear single susceptibilities in the form of Heisenberg operator, a simple proof of charge conservation and gauge invariance guaranteed by such the susceptibilities, and application to a many-electron system with detailed discussion on the density functional theory.

The introduction below contains the followings: §I A reveals the necessity of the single susceptibility, instead of the electric permittivity and magnetic permeability. §I B points out the preference to equal treatment of the SP and VP as the cause of response in NFO, instead of the unequal treatment in ordinary optics under the Coulomb gauge. §I C explains the difficulty of constructing the response theory in NFO, which inevitably connected to a many-electron problem via the SP. §I D represents the purpose of this paper.

A. The necessity of the single susceptibility

As the cause of response, it is natural and essential to use the SP and VP, which represent for the electromagnetic (EM) field in the Hamiltonian for quantum electrodynamics. Three reasons are given below for the inapplicability of the electric and magnetic fields as the cause of response. First, there exist such systems that cannot be described in terms of the electric and/or magnetic fields, namely, the superconductor system with the Meissner effect[2] and the coherent electron system with the Aharanov-Bohm effect[3]. A limited NF optical system is another example, as shown in the one-electron system in Ref.[1] (and will be shown in a
many-electron system in §VI of this paper).

Second, the constitutive equations with the electric permittivity and magnetic permeability give relationships between redundant degrees of freedom. Actually, the essential source of the EM field is the three components of charge density and the transverse current density. The longitudinal current density is excluded because it can be determined through charge conservation law, once the charge density is known. However, the polarization and magnetization as the source of the EM field have totally six components, which include the redundancy. So that the associated constitutive equations using the two susceptibilities include the constraint condition for the redundancy, of which the physical meaning is not declared. This situation is physically unreasonable and should be fixed by the constitutive equation using a single susceptibility associated with the proper degrees of freedom.

Third, as first claimed by Cho[4, 5] for the low-symmetry systems with chirality (such as the NF optical system with a skewed nanostructure), the ordinary two constitutive equations are not available because the electric and magnetic responses become indistinguishable. He also revealed that this error cannot be fixed by the Drude-Born-Fedorov formulas[6], which extends the two constitutive equations adding the cross terms of the electric-field-induced magnetization and the magnetic-field-induced polarization.

Therefore, from a general viewpoint, it is essential to employ a single susceptibility with the SP and VP, instead of the electric permittivity and magnetic permeability with the electric and magnetic fields.

**B. The preference to equal treatment of the SP and VP in a NF optical system**

Suppose a small-scale material is placed in the vicinity of a nanostructure, which functions as a light source (FIG.1). In such a system, under the NF incidence condition, the target material is exposed to the longitudinal and transverse electric fields simultaneously, whereas in a system under the far-field incidence condition, the target material is exposed only to the transverse field, which survives far from the light source. Therefore, the coexistence of longitudinal and transverse electric fields distinguishes such a system under the NF incidence condition from that under the far-field incidence condition.

Here, the longitudinal electric field originates from the charge density on the nanostructure, obeys Coulomb’s law, and has a non-radiative nature to localize around the nanos-
structure. On the other hand, the transverse electric field originates from the transverse current density on the nanostructure, obeys Ampere-Maxwell law and Faraday’s law, and has a radiative nature allowing it to propagate far from the light source, accompanied by the magnetic field. (The longitudinal current density is determined via the charge conservation law, once the charge density is known, and is not an independent source.) Therefore, the two incidences coexisting in an NF optical system have distinct properties.

Furthermore, owing to the non-relativistic nature of the system, the SP and VP appear in a different manner in the Hamiltonian, (for example, Eq. (30) in §IV,) which governs the electron response. Considering that the SP and VP under the Coulomb gauge represent the longitudinal and transverse electric fields, respectively, one may confirm that the two types of incidences in NFO cause different responses; see §VI for an explicit demonstration.

Therefore, it is reasonable to treat SP and VP equally as the cause of response in the NF optical system. Up to now, there has been no such theoretical framework for equally treating the SP and VP. The reason for this lies in the many-electron problem inevitably related to NFO via the SP (the longitudinal electric field), as is mentioned in the next subsection.

FIG. 1: Optical systems under near- and far-field incidences (the left and right side figures, respectively). The former system is exposed simultaneously to the incident longitudinal and transverse electric fields (fundamentally represented by the scalar and vector potentials, respectively, under the Coulomb gauge), whereas the latter system is exposed to only the transverse field (the vector potential under the Coulomb gauge).
C. A many-electron problem inevitably related with NFO

The relationship between NFO and many-electron problem has not been well recognized, although the problem of how to consider the Coulomb interaction in response function has remained for a long time[7]. In the usual Hamiltonian for a many-electron system under the Coulomb gauge, the SP is rewritten as the interaction between the electron charge density operators, and only the VP is considered as the cause of the response. This unequal treatment of the SP and VP is needed to consider the quantum many-electron effect (the so-called exchange-correlation effect) to construct the ground and excited states as the proper bound states in a many-electron system. This usual procedure to treat the non-relativistic many-electron system is compatible with ordinary optics, where the electron system of interest is far from the light source, and the SP incidence is negligible. By contrast, in an NF optical system, the usual approach results in a difficulty of understanding the response to the SP incidence, because both the SP incidence (radiated by the nanostructure) and the inherent SP (originating from the particle charge) are built into the two-body Coulomb interaction, and the two contributions are indistinguishable. To make matters worse, the Coulomb interaction in itself is so difficult to treat that it is often ignored, without considering it includes the effect of the SP incidence.

For the NF optical system, there are two existing approaches based on certain single susceptibilities (the nonlocal response functions). Cho formulated a single susceptibility that relates the transverse VP (as the cause) to the current density (as the result), and applied it to various optical systems[8]. Additionally, a modification that considers the longitudinal electric field incidence in NF optical systems has been proposed in Chap. 5 of Ref.[5]. Keller formulated the linear single susceptibility, which relates the transverse electric field and the incident part of the longitudinal electric field (as the cause) to the current density (as the result) [9].

In the above two existing formulations, the SP under the Coulomb gauge (or the longitudinal electric field), except the linear-dependence of the incidence, is rewritten as the two-body Coulomb interaction in the usual manner. Therefore, the response to the SP, in principle, can be rigorously considered via the Coulomb interaction if the many-electron problem is properly solved, whereas the response to the VP incidence under the Coulomb gauge (or the transverse electric field incidence) is treated in the perturbative manner. In
this type of approach, it is essential to solve the many-electron problem, in particular, for
the nonlinear process related with the SP. Even if the Coulomb interaction is properly con-
sidered, the unequal treatment may make it difficult to regulate the perturbation order of
the responses and to understand the role of the SP incidence.

As a result, the response theory in NFO is inevitably relates to the many-electron problem,
which causes difficulty.

D. The purpose of this paper

§I A-§IC lead to the logical fallacy to use the ordinary two susceptibilities with the electric
and magnetic fields, and the preference to use the single susceptibility equally associated
with the SP and VP, considering properly the many-electron effect in NF optical systems,
although the ordinary two susceptibilities have been widely used both in ordinary optics
and NFO. To best understand the fundamental physics in NFO, it is essential to develop
an adequate response theory. For this purpose, the present paper defines and characterizes
a single susceptibility equally associated with the SP and VP based on the action integral
from scratch.

The contents of this paper are as follows: §II defines the linear and nonlinear single
susceptibilities equally associated with the SP and VP, starting from the action integral. §III
shows that the present susceptibility respects both charge conservation and gauge invariance
in a general manner. §IV derives the linear and nonlinear single susceptibilities in the form
of the Heisenberg operators. §V shows that the many-electron effect in the present response
theory may be supported by the density functional theory to prepare the non-perturbed
ground state as well as a complete set of many-electron states. §VI applies the present linear
response theory to a simplified many-electron system, and show that the electric field with
the associated permittivity is not suitable to describe the response of a limited NF optical
system with a non-metallic material, so that the SP and VP with the single susceptibility
is essential. §VII provides a summary of this work. Two appendices are included: §A and
§B provides calculation details of §II and §VI, respectively.
II. DEFINITION OF NEW SINGLE SUSCEPTIBILITY

Based on the Lagrangian formulation of non-relativistic quantum electrodynamics, we define the single susceptibility, which relates the SP and VP (as the cause) to the induced charge and current densities (as the result). Furthermore, it is shown that this susceptibility guarantees that charge conservation and gauge invariance hold; see the next section. The action integral for non-relativistic quantum electrodynamics is:

\[
\mathcal{I} = \mathcal{I}_{\text{mat}} + \mathcal{I}_{\text{EM}},
\]

where

\[
\begin{align*}
\mathcal{I}_{\text{mat}} &\equiv \int \frac{d^4x}{c} \left\{ \hat{\psi}_{\alpha}^\dagger(x)(i\hbar \partial_t - q\phi(x))\hat{\psi}_{\alpha}(x) - \frac{1}{2m} \left( \frac{\hbar}{i} \partial_t - qA_i(x) \right) \hat{\psi}_{\alpha}(x) \right. \\
&\quad \left. - \phi(x)\rho^{(\text{EXT})}(x) + A_i(x)j_i^{(\text{EXT})}(x) - \hat{\psi}_{\alpha}^\dagger(x)\psi^{(\text{AUX})}(x)\hat{\psi}_{\alpha}(x) \right\}, \\
\mathcal{I}_{\text{EM}} &\equiv \int \frac{d^4x}{c} \left\{ \frac{\epsilon_0}{2} (\partial_t A_i(x) + \partial_i \phi(x)) (\partial_t A_i(x) + \partial_i \phi(x)) \right. \\
&\quad \left. - \frac{\epsilon_0 c^2}{2} \epsilon_{ijk} \partial_j A_k(x) \epsilon_{ilm} \partial_l A_m(x) \right\},
\end{align*}
\]

where \(m\) and \(q(= -e)\) are the electron mass and charge, \(c\) is the speed of light, \(\phi, A\) are the SP and VP, which are assumed to be classical field in the semiclassical treatment, \(\hat{\psi}_{\alpha}, \hat{\psi}_{\alpha}^\dagger\) are the electron field operators with the spin state \(\alpha\) (one of the two spin states; so called "up" and "down" states), and \(\rho^{(\text{EXT})}, j^{(\text{EXT})}\) are the nuclear charge and the current densities, respectively, which possibly generate inherent EM field. A static auxiliary potential \(\psi^{(\text{AUX})}(x)\) is null for now, but is introduced here for the discussion in §V concerning the density functional theory to consider the quantum many-electron effect (the exchange-correlation effect), \(\epsilon_{ijk}\) is an antisymmetric tensor, and the Einstein rule is used for indices of vector and Grassmann fields, that is, summation should be executed over repeated indices. At this first stage of investigation, the interaction between spin polarization and the EM field is ignored. The soundness of the above action integral is confirmed by its Euler equations, which will soon be derived.

The electron field operators are considered as quantized Grassmann fields. The Grassmann field satisfies \([\hat{\psi}_{\alpha}(r, t), \hat{\psi}_{\beta}^\dagger(r', t')]_+ = 0 [10]\), and corresponds to the "classical" field of the electron. These operators become the creation and annihilation operators of the electron in quantum theory (the quantized Grassmann fields), introducing the anti-commutation
The charge conservation law below holds, and is checked through explicit calculation:

\[ 0 = \partial_t \hat{\rho}(x) + \partial_i \hat{j}_i(x) = 0. \]  

In the four-element representation, Eqs.(6) and (7) become:

\[ (\delta^\mu \Box - \partial^\mu \partial_\nu) A^\nu(x) = \frac{1}{\epsilon_0 c} (\hat{j}^\mu(x) + j^{\text{(EXT)}}(x)), \]  

where

\[ \hat{j}^\mu = (c\hat{\rho}, \hat{j}), \quad \hat{j}^\mu = (c\hat{\rho}, -\hat{j}), \]

\[ A^\mu = (\phi, cA), \quad A_\mu = (\phi, -cA), \]

\[ \partial^\mu = (1/c \partial_\tau, -\nabla), \quad \partial_\mu = (1/c \partial_\tau, \nabla), \]

\[ \Box = \partial^\mu \partial_\mu = 1/c^2 \partial^2_\tau - \Delta, \quad \text{etc.} \]
Although Lorentz invariance is not maintained in the non-relativistic theory, we use the four-element notation to simply represent charge conservation and gauge invariance. For example, Eqs.(8)-(10) become:

\[ \hat{j}^\mu(x) = -c^2 \frac{\delta}{\delta A_\mu(x)} \mathcal{I}_{\text{mat}}, \]  

\[ \partial_\mu \hat{j}^\mu(x) = 0. \]  

The action integral, Eq.(1) is invariant under the following gauge transformation:

\[ A^\mu(x) \rightarrow A^\mu(x) - c \partial^\mu \eta(x), \]

\[ \hat{\psi}_\alpha(x) \rightarrow e^{\frac{i}{\hbar} q \eta(x)} \hat{\psi}_\alpha(x), \quad \hat{\psi}^\dagger_\alpha(x) \rightarrow \hat{\psi}^\dagger_\alpha(x) e^{-\frac{i}{\hbar} q \eta(x)}, \]  

where \( \eta(x) \) is the gauge function. From the point of view of Noether’s theorem[11], the gauge invariance of the action integral is the cause of the charge conservation law, Eq.(10) or Eq.(14).

Let us separate the EM field into two parts:

\[ A^\mu(x) = A^{(0)\mu}(x) + \Delta A^\mu(x), \]  

where \( A^{(0)\mu} \) is the static, non-perturbative EM potential satisfying Eqs.(6) and (7), and \( \Delta A^\mu(x) \) is the perturbative EM potential. Under this variation of the EM field, let us re-optimize the action integral of the matter, \( \mathcal{I}_{\text{mat}}[\hat{\psi}^\dagger_\alpha, \hat{\psi}_\alpha, A^\mu] \). That is, we re-optimize the electron field operator satisfying Eqs.(4) and (5) under \( A^{(0)\mu} + \Delta A^\mu(x) \). In the above procedure, the variation of the action integral of the matter is expressed by the total functional derivative with respect to \( A^\mu(x) \):

\[ \frac{\delta}{\delta A^\mu(x)} \mathcal{I}_{\text{mat}}[\hat{\psi}^\dagger_\alpha[A^\nu], \hat{\psi}_\alpha[A^\nu], A^\nu] \bigg|_{A^\nu = A^{(0)\nu}} = \]

\[ = \left[ \frac{\delta}{\delta A_\mu(x)} \right]_{\text{explicit}} \mathcal{I}_{\text{mat}} + \int d^4x' \delta \hat{\psi}^\dagger_\alpha(x') \delta \hat{\psi}_\alpha(x') \delta \mathcal{I}_{\text{mat}} 
+ \int d^4x' \delta \mathcal{I}_{\text{mat}} / \delta \hat{\psi}_\alpha(x') \delta \hat{\psi}^\dagger_\alpha(x') \bigg|_{A^\nu = A^{(0)\nu}} \]

\[ = -\frac{1}{c^2} \hat{j}^\mu(x; [A^{(0)\mu}]), \]  

where the first term in the second expression is the variation explicitly caused by the perturbative EM field, and the second and third terms are the implicit variations, created through
owing to Eqs.(A1) and (A2) in Appendix A, the derivatives of the action integral of the matter, the following extension of Eq.(18) holds, because of Eq.(A2) in Appendix A. In the same manner as for higher order total functional derivative of the action integral of the matter is simply the current density in the non-perturbed system. Furthermore, the second order total functional derivative is calculated as follows:

\[
\frac{\delta}{\delta A^{\mu_1}(x_1)} \frac{\delta}{\delta A^{\mu}(x)} I_{\text{mat}}[\hat{\psi}_\alpha^\dagger[A^\nu], \hat{\psi}_\alpha[A^\nu], A^\nu]_{A^\nu = A^{(0)\nu}} = \left[ \frac{\delta}{\delta A^{\mu_1}(x_1)} \left( \frac{\delta}{\delta A^{\mu}(x)} I_{\text{mat}} \right) \bigg|_{\text{explicit}} \right] + \int d^4x' \frac{\delta}{\delta A^{\mu_1}(x_1)} \left( \frac{\delta \hat{\psi}_\alpha^\dagger(x') \delta \hat{\psi}_\alpha(x')}{\delta A^{\mu}(x)} \right)_{A^\nu = A^{(0)\nu}}
\]

\[
= - \frac{1}{c^2} \frac{\delta \hat{j}^\mu(x; [A^\nu])}{\delta A^{\mu_1}(x_1)} \bigg|_{A^\nu = A^{(0)\nu}},
\]

where the second and third terms in the second expression are null. Actually, the integrand of the second term is:

\[
\left[ \left( \frac{\delta}{\delta A^{\mu_1}(x_1)} \right) \frac{\delta \hat{\psi}_\alpha^\dagger(x')}{\delta A^{\mu}(x)} \right] \delta \hat{\psi}_\alpha^\dagger(x') \delta I_{\text{mat}} + \frac{\delta \hat{\psi}_\alpha^\dagger(x')}{\delta A^{\mu}(x)} \left( \frac{\delta}{\delta A^{\mu_1}(x_1)} \delta \hat{\psi}_\alpha(x') \right)_{A^\nu = A^{(0)\nu}}
\]

The first term in this equation is null because of Eq.(4), and the second term is also null because of Eq.(A2) in Appendix A. In the same manner as for higher order total functional derivatives of the action integral of the matter, the following extension of Eq.(18) holds, owing to Eqs.(A1) and (A2) in Appendix A,

\[
\frac{\delta^{n+1} I_{\text{mat}}[\hat{\psi}_\alpha^\dagger[A^\nu], \hat{\psi}_\alpha[A^\nu], A^\nu]}{\delta A^{\mu_1}(x_1) \cdots \delta A^{\mu_n}(x_n)} \bigg|_{A^\nu = A^{(0)\nu}} = - \frac{1}{c^2} \frac{\delta^n \hat{j}^\mu(x; [A^\nu])}{\delta A^{\mu_1}(x_1) \cdots \delta A^{\mu_n}(x_n)} \bigg|_{A^\nu = A^{(0)\nu}}.
\]

To define the single susceptibility, suppose the system under the non-perturbative EM field \( A^{(0)\mu}(x) \) is exposed to the perturbative EM field \( \Delta A^\mu(x) \). The non-perturbative EM field \( A^{(0)\mu} \) is a solution of the coupled equations Eqs.(4)-(7), namely, Heisenberg’s equation and Maxwell’s wave equations, and is assumed to be a static solution existing in the ground state. On the other hand, the total EM field \( A^{(0)\mu} + \Delta A^\mu \) is not necessarily a solution of Maxwell’s wave equations, Eqs.(6) and (7), that is, \( \Delta A^\mu \) is introduced as a virtual variation. The induced current density is the variation from the current density in the non-perturbative
From Eqs.(19) and (20), the linear and nonlinear single susceptibility operators are defined as:

\[
\hat{j}_{\mu}(x; [A^{(0)}] + \Delta A^{\nu}) - \hat{j}_{\mu}(x; [A^{(0)}]) = \int d^4 x_1 \frac{\delta \hat{j}_{\mu}(x; [A^{\nu}])}{\delta A^{\mu_1}(x_1)} \bigg|_{\nu = A^{(0)}} \Delta A^{\mu_1}(x_1) \\
+ \frac{1}{2!} \int d^4 x_1 \int d^4 x_2 \frac{\delta^2 \hat{j}_{\mu}(x; [A^{\nu}])}{\delta A^{\mu_1}(x_1) \delta A^{\mu_2}(x_2)} \bigg|_{\nu = A^{(0)}} \Delta A^{\mu_1}(x_1) \Delta A^{\mu_2}(x_2) \\
+ \frac{1}{3!} \int d^4 x_1 \int d^4 x_2 \int d^4 x_3 \frac{\delta^3 \hat{j}_{\mu}(x; [A^{\nu}])}{\delta A^{\mu_1}(x_1) \delta A^{\mu_2}(x_2) \delta A^{\mu_3}(x_3)} \bigg|_{\nu = A^{(0)}} \Delta A^{\mu_1}(x_1) \Delta A^{\mu_2}(x_2) \Delta A^{\mu_3}(x_3) \\
+ \cdots .
\]

From Eqs.(19) and (20), the linear and nonlinear single susceptibility operators are defined as:

\[
\hat{\chi}^{\mu}_{\mu_1}(x, x_1, x_1) \equiv \frac{\delta \hat{j}_{\mu}(x; [A^{\nu}])}{\delta A^{\mu_1}(x_1)} \bigg|_{\nu = A^{(0)}} \\
= -e^2 \left. \frac{\delta^2 \mathcal{T}_{\text{mat}}}{\delta A_{\mu}(x) \delta A^{\mu_1}(x_1)} \right|_{\nu = A^{(0)}} \Delta A^{\mu_1}(x_1),
\]

\[
\hat{\chi}^{\mu}_{\mu_1 \mu_2}(x, x_1, x_2) \equiv \frac{1}{2!} \frac{\delta^2 \hat{j}_{\mu}(x; [A^{\nu}])}{\delta A^{\mu_1}(x_1) \delta A^{\mu_2}(x_2)} \bigg|_{\nu = A^{(0)}} \\
= -e^2 \left. \frac{\delta^3 \mathcal{T}_{\text{mat}}}{\delta A_{\mu}(x) \delta A^{\mu_1}(x_1) \delta A^{\mu_2}(x_2)} \right|_{\nu = A^{(0)}} \Delta A^{\mu_1}(x_1) \Delta A^{\mu_2}(x_2),
\]

\[
\hat{\chi}^{\mu}_{\mu_1 \cdots \mu_n}(x, x_1, \cdots, x_n) \equiv \frac{1}{n!} \frac{\delta^n \hat{j}_{\mu}(x; [A^{\nu}])}{\delta A^{\mu_1}(x_1) \cdots \delta A^{\mu_n}(x_n)} \bigg|_{\nu = A^{(0)}} \\
= -e^2 \left. \frac{\delta^{n+1} \mathcal{T}_{\text{mat}}}{\delta A_{\mu}(x) \delta A^{\mu_1}(x_1) \cdots \delta A^{\mu_n}(x_n)} \right|_{\nu = A^{(0)}} \Delta A^{\mu_1}(x_1) \cdots \Delta A^{\mu_n}(x_n),
\]

The susceptibility is defined using a small amount of variation, \(\Delta A^{\mu}\). That is, the EM field does not in general satisfy its Euler equation, Eq.(11), while the electron field operators satisfy Eqs.(4) and (5). To evaluate the real EM field, \(\Delta A^{\mu}\) must be determined and a further procedure is required to solve the coupled equations, with the constitutive equations in terms of the susceptibility and Maxwell’s wave equations Eqs.(6) and (7). This procedure is provided in a self-consistent manner, as performed by K.Cho[8] using his single susceptibility.
III. CHARGE CONSERVATION LAW AND GAUGE INVARIANCE OF THE SINGLE SUSCEPTIBILITY

In the last expressions in Eqs.(21)-(23) the coordinates \(x_1, x_2, \cdots\) for the cause (the perturbative EM field) and the coordinates \(x\) for the result (the induced current density) are symmetric. Charge conservation for the induced charge density holds to each order of the perturbation because of Eq.(10) or Eq.(14) and Eqs.(20)-(23); this is described by the derivative of the coordinate for the result, \(x\):

\[
\partial_\mu \hat{\chi}^\mu_{\mu_1 \cdots}(x, x_1, \cdots) = 0. \tag{24}
\]

The symmetry of the coordinates between the result and the cause leads to the following equation concerning the derivative of any coordinate for the cause, e.g., \(x_1\):

\[
\partial^{\mu_1} \hat{\chi}^\mu_{\mu_1 \cdots}(x, x_1, \cdots) = 0. \tag{25}
\]

Equation (25) means that the susceptibility guarantees that gauge invariance is respected. That is, the resultant charge and current densities are independent of the chosen gauge. To confirm this fact, consider the convolution integral of the single susceptibility with the perturbative EM field, in a certain gauge, e.g.,

\[
\int d^4x_1 \hat{\chi}^\mu_{\mu_1 \cdots}(x, x_1, \cdots) \Delta A^{\mu_1}(x_1). \tag{26}
\]

A gauge transformation of \(\Delta A\) to \(\Delta A'\) in another gauge is expressed as:

\[
\Delta A^{\mu_1}(x_1) = \Delta A'^{\mu_1}(x_1) - c \partial^{\mu_1} \eta(x_1), \tag{27}
\]

where \(\eta\) is the gauge function. Equation (26) leads to:

\[
\int d^4x_1 \hat{\chi}^\mu_{\mu_1 \cdots}(x, x_1, \cdots) \Delta A^{\mu_1}(x_1)
\]

\[
= \int d^4x_1 \hat{\chi}^\mu_{\mu_1 \cdots}(x, x_1, \cdots) \Delta A'^{\mu_1}(x_1) + c \int d^4x_1 \partial^{\mu_1} \hat{\chi}^\mu_{\mu_1 \cdots}(x, x_1, \cdots) \eta(x_1)
\]

\[
= \int d^4x_1 \hat{\chi}^\mu_{\mu_1 \cdots}(x, x_1, \cdots) \Delta A'^{\mu_1}(x_1). \tag{28}
\]

The contribution of the gauge function vanishes in the convolution integral. Thus, the gauge of the perturbative EM field may be freely selected. This means that the susceptibility is independent of the chosen gauge and, in practice, one may select a gauge that is most convenient for calculation.
In Eq. (2), if the factor \( i\hbar \hat{\psi}^\dagger_\alpha(x) \) of the first term is regarded as the canonical momentum of \( \hat{\psi}_\alpha(x) \), then the Hamiltonian density may be determined as the Legendre transformation from the Lagrangian density, that is:

\[
\hat{H} = \int \! d^3x \left( \frac{\hbar}{i} \partial_t - qA_t(x) \right) \hat{\psi}^\dagger_\alpha(x) \left( \frac{\hbar}{i} \partial_t - qA_t(x) \right) \hat{\psi}_\alpha(x) + q\phi(x) \hat{\psi}^\dagger_\alpha(x) \hat{\psi}_\alpha(x) + h.c.
\]

(30)

This Hamiltonian governs the motion of electron field operators. Assuming that the non-perturbative EM field \( \phi^{(0)}, A^{(0)} \) is the static EM field existing in the ground state of a many-electron system, the Hamiltonian, \( \hat{H} \) may be separated into a non-perturbative part, \( \hat{H}^{(0)} \) and a perturbative part, \( \hat{V} \) as follows:

\[
\hat{H}^{(0)} = \int \! d^3x \left( \frac{\hbar}{i} \partial_t - qA_t^{(0)}(x) \right) \hat{\psi}^\dagger_\alpha(x) \left( \frac{\hbar}{i} \partial_t - qA_t^{(0)}(x) \right) \hat{\psi}_\alpha(x) + q\phi^{(0)}(x) \hat{\psi}^\dagger_\alpha(x) \hat{\psi}_\alpha(x) + h.c.
\]

(31)

\[
\hat{V}(t) \equiv \hat{H} - \hat{H}^{(0)} = \int \! d^3x \hat{\nu}(x),
\]

\[
= \int \! d^3x \left\{ \Delta \phi(x) q\hat{\psi}^\dagger_\alpha(x) \hat{\psi}_\alpha(x) - \Delta A_t(x) \frac{q}{2m} \left( \hat{\psi}^\dagger_\alpha(x) \left( \frac{\hbar}{i} \partial_t - qA_t^{(0)}(x) \right) \hat{\psi}_\alpha(x) + h.c. \right) \right. \\
+ \left. \frac{q}{2m} \Delta A_t(x) \Delta A_t(x) q\hat{\psi}^\dagger_\alpha(x) \hat{\psi}_\alpha(x) \right\}
\]

(32)
\( \hat{j}_\mu (x) \bigg|_{A=A^{(0)}} \) in Eq.(32) is the current density Eq.(29), with the explicitly-appeared VP being replaced by that in the non-perturbed system. The tensor Eq.(33) represents the non-relativistic effect. Actually, this tensor is the analogue of the four-element Kronecker delta, but brings inequality of the temporal and spatial coordinates.

Here, the field operators in the interaction picture (the asymptotic field operators) \( \hat{\psi}^{(in)}_\alpha, \hat{\psi}^{(in)} \) are governed by the non-perturbative Hamiltonian \( \hat{H}^{(0)} \) and coincide with the field operators in the Heisenberg picture, \( \hat{\psi}_\alpha, \hat{\psi} \) at the infinite past time, \( t \to -\infty \), assuming the adiabatic switch-on. The unitary operator \( \hat{U}(t, -\infty) \) is the time-evolution operator of the states in the interaction picture, and relates the operators between the Heisenberg and interaction pictures as follows:

\[
\begin{align*}
\hat{\psi}_\alpha (x) &= \hat{U}^{-1}(t, -\infty) \hat{\psi}^{(in)}_\alpha (x) \hat{U}(t, -\infty), \\
\hat{\psi}_\alpha^\dagger (x) &= \hat{U}^{-1}(t, -\infty) \hat{\psi}^{(in)t}_\alpha (x) \hat{U}(t, -\infty),
\end{align*}
\]

where

\[
\hat{U}(t, -\infty) = \lim_{t_0 \to -\infty} \hat{U}(t, t_0) = \lim_{t_0 \to -\infty} \hat{U} e^{\frac{i}{\hbar} \int_{t_0}^{t} dt' \hat{V}^{(in)}(t')},
\]

\( \hat{V}^{(in)}(t') \equiv \hat{V}(\hat{\psi}^{(in)}_\alpha, \hat{\psi}^{(in)}; t') \)

Combining Eq.(34) and Eq.(29), the four-element current density operator in the interaction picture may be defined as: \( \hat{j}^{(in)\mu}(x) = (c \hat{\rho}^{(in)}(x), \hat{j}^{(in)}(x)) \). These charge and current densities do not satisfy the charge conservation law, except for \( A = A^{(0)} \), and are merely convenient tools used for obtaining the expansion of the retarded product of the Heisenberg operators.

\[
\begin{align*}
\hat{j}_\mu (x) &= \hat{U}^{-1}(t, -\infty) \hat{j}^{(in)\mu}(x) \hat{U}(t, -\infty), \\
\hat{j}^{(in)\mu}(x) &= \begin{cases} 
\frac{c q \hat{\psi}^{(in)}_\alpha(x)}{2m} \left( \frac{\hbar}{i} (-\partial^\mu) - \frac{q}{c} A^\mu(x) \right) \hat{\psi}^{(in)}_\alpha(x) + \text{h.c.} & \text{for } \mu = 0, \\
\hat{\psi}^{(in)t}_\alpha(x) & \text{for } \mu = 1, 2, 3. 
\end{cases}
\end{align*}
\]

To obtain the perturbative expansion (the retarded product series) of the Heisenberg operators, let us introduce an operator in the intermediate picture, where \( \hat{U}(t, t_0) \) will be used instead of \( \hat{U}(t, -\infty) \):

\[
\begin{align*}
\hat{\rho}^\bullet (x; t_0) &= \hat{U}^{-1}(t, t_0) q \hat{\psi}^{(in)t}_\alpha(x) \hat{\psi}^{(in)}_\alpha(x) \hat{U}(t, t_0), \\
\hat{j}^\bullet_t (x; t_0) &= \hat{U}^{-1}(t, t_0) \frac{q}{2m} \hat{\psi}^{(in)t}_\alpha(x) \left( \frac{\hbar}{i} \partial_t - q A_t(x) \right) \hat{\psi}^{(in)}_\alpha(x) \hat{U}(t, t_0) + \text{h.c.}
\end{align*}
\]

The corresponding four-element current density is

\[
\hat{j}^\bullet^\mu(x; t_0) = (c\hat{\rho}^\bullet(x; t_0), \hat{j}^\bullet_t(x; t_0))
\]
As \( t_0 \rightarrow -\infty \), these operators coincide with those of the Heisenberg picture, while at \( t_0 = t \), they coincide with those of the interaction picture:

\[
\hat{j}^\bullet \mu (x; -\infty) = \hat{j}^\mu (x),
\]
\[
\hat{j}^\bullet \mu (x; t) = \hat{j}^{(in)} \mu (x).
\]  (37)  (38)

Next, let’s investigate the time evolution of \( \hat{j}^\bullet \mu \) as a function of \( t_0 \).

\[
\partial_{t_0} \hat{j}^\bullet \mu (x; t_0) = \{ \partial_{t_0} \hat{U}^{-1}(t, t_0) \} \hat{j}^{(in)} \mu (x) \hat{U}(t, t_0) + \hat{U}^{-1}(t, t_0) \hat{j}^{(in)} \mu (x) \{ \partial_{t_0} \hat{U}(t, t_0) \}
\]
\[
= \frac{1}{i \hbar} \hat{V}^{(in)}(t_0) \hat{U}^{-1}(t, t_0) \hat{j}^{(in)} \mu (x) \hat{U}(t, t_0) + \hat{U}^{-1}(t, t_0) \hat{j}^{(in)} \mu (x) \hat{U}(t, t_0) - \frac{1}{i \hbar} \hat{V}^{(in)}(t_0)
\]
\[
= - \frac{1}{i \hbar} \left[ \hat{j}^\bullet \mu (x; t_0), \hat{V}^{(in)}(t_0) \right]
\]

Integrating over \([t_0, t]\), approximating iteratively using Eq.(38), and changing the region of multi-integration, we obtain:

\[
\hat{j}^\bullet \mu (x; t_0) = \hat{j}^{(in)} \mu (x) + \frac{1}{i \hbar} \int_{t_0}^{t} dt_1 \left[ \hat{j}^\bullet \mu (x; t_1), \hat{V}^{(in)}(t_1) \right]
\]
\[
= \hat{j}^{(in)} \mu (x) + \frac{1}{i \hbar} \int_{t_0}^{t} dt_1 \left[ \hat{j}^{(in)} \mu (x), \hat{V}^{(in)}(t_1) \right]
\]
\[
+ \left( \frac{1}{i \hbar} \right)^2 \int_{t_0}^{t} dt_1 \int_{t_1}^{t} dt_2 \left[ \left[ \hat{j}^{(in)} \mu (x), \hat{V}^{(in)}(t_2) \right], \hat{V}^{(in)}(t_1) \right]
\]
\[
+ \left( \frac{1}{i \hbar} \right)^3 \int_{t_0}^{t} dt_1 \int_{t_1}^{t} dt_2 \int_{t_2}^{t} dt_3 \left[ \left[ \left[ \hat{j}^{(in)} \mu (x), \hat{V}^{(in)}(t_3) \right], \hat{V}^{(in)}(t_2) \right], \hat{V}^{(in)}(t_1) \right] + \ldots
\]
\[
= \hat{j}^{(in)} \mu (x) + \frac{1}{i \hbar} \int_{t_0}^{t} dt_1 \left[ \hat{j}^{(in)} \mu (x), \hat{V}^{(in)}(t_1) \right]
\]
\[
+ \left( \frac{1}{i \hbar} \right)^2 \int_{t_0}^{t} dt_1 \int_{t_0}^{t} dt_2 \left[ \left[ \hat{j}^{(in)} \mu (x), \hat{V}^{(in)}(t_2) \right], \hat{V}^{(in)}(t_1) \right]
\]
\[
+ \left( \frac{1}{i \hbar} \right)^3 \int_{t_0}^{t} dt_1 \int_{t_0}^{t} dt_2 \int_{t_0}^{t} dt_3 \left[ \left[ \left[ \hat{j}^{(in)} \mu (x), \hat{V}^{(in)}(t_3) \right], \hat{V}^{(in)}(t_2) \right], \hat{V}^{(in)}(t_1) \right] + \ldots
\]

Then, taking the limit \( t_0 \rightarrow -\infty \), the above equation yields the retarded product of the Heisenberg operators, as follows:

\[
\hat{j}^\mu (x) = \hat{j}^{(in)} \mu (x) + \frac{1}{i \hbar \epsilon} \int_{ct_1 \in (-\infty, ct]} d^4 x_1 \left[ \hat{j}^{(in)} \mu (x), \hat{v}^{(in)}(x_1) \right]
\]
\[
+ \left( \frac{1}{i \hbar \epsilon} \right)^2 \int_{ct_1 \in (-\infty, ct]} d^4 x_1 \int_{ct_2 \in (-\infty, ct_1]} d^4 x_2 \left[ \left[ \hat{j}^{(in)} \mu (x), \hat{v}^{(in)}(x_1) \right], \hat{v}^{(in)}(x_2) \right]
\]
\[
+ \left( \frac{1}{i \hbar \epsilon} \right)^3 \int_{ct_1 \in (-\infty, ct]} d^4 x_1 \int_{ct_2 \in (-\infty, ct_1]} d^4 x_2 \int_{ct_3 \in (-\infty, ct_2]} d^4 x_3 \left[ \left[ \left[ \hat{j}^{(in)} \mu (x), \hat{v}^{(in)}(x_1) \right], \hat{v}^{(in)}(x_2) \right], \hat{v}^{(in)}(x_3) \right]
\]
in Eq. (36) is of zeroth and first order for $\mu$ and $\hat{\psi}$ dependence, which exists in the first and second terms of Eq. (39): replacing $\hat{\psi}$ is of first and second order. The linear single susceptibility operator comes from the EM potential. In Equation (39), the dependence of the EM potential through $\hat{j}^{(in)}(x)$ in the form of Heisenberg operator by the functional derivative of Eq. (39) with respect to the EM potential through $\hat{j}^{(in)}(x)$ and $\hat{\psi}^{\dagger}$ by $\hat{\psi}^{(in)}$, $\hat{\psi}^{(in)\dagger}$, respectively. Next, let us derive the single susceptibility in the form of Heisenberg operator by the functional derivative of Eq. (39) with respect to the EM potential. In Equation (39), the dependence of the EM potential through $\hat{j}^{(in)}(x)$ in Eq. (36) is of zeroth and first order for $\mu \in \{1, 2, 3\}$, and dependence through $\hat{v}^{(in)}(x_1)$ is of first and second order. The linear single susceptibility operator comes from the $A^1$-dependence, which exists in the first and second terms of Eq. (39):

$$\hat{\chi}^{\mu}_{\mu_1}(x, x_1) = \left. \frac{\delta \hat{j}^{(in)}(x)}{\delta A^{\mu_1}(x_1)} \right|_{A = A^{(0)}}$$

$$= \left. \frac{-q}{mc^2} \hat{\delta}^{\mu}_{\mu_1} \hat{\delta}^4(x - x_1) \hat{j}^{(in)0}(x) + \frac{1}{i\hbar c^2} \theta(ct - ct_1) \left[ \hat{j}^{(in)\mu}(x), \hat{j}^{(in)\mu_1}(x_1) \right] \right|_{A = A^{(0)}}$$

where $\hat{j}^{(in)\mu}(x) = \hat{j}^{(in)\mu}(x) \big|_{A = A^{(0)}}$.

The Heisenberg operators of the nonlinear single susceptibilities, to second and higher order, are as follows. To avoid any confusion in the case of two times coinciding, the long and explicit expressions are given, without using the time ordering operator.

$$2! \hat{\chi}^{\mu}_{\mu_1\mu_2}(x, x_1, x_2) = \left. \frac{\delta^2 \hat{j}^{(in)}(x)}{\delta A^{\mu_1}(x_1) \delta A^{\mu_2}(x_2)} \right|_{A = A^{(0)}}$$

$$= \frac{1}{i\hbar c^2} \frac{-q}{mc^2} \left\{ \delta(ct - ct_1)\theta(ct - ct_2) \frac{\hat{\delta}^{\mu}_{\mu_2}}{\hat{\delta}^4(x - x_1)} \left[ \hat{j}^{(in)0}(x), \hat{j}^{(in)\mu_1}(x_2) \right] \right.$$ 

$$+ \delta(ct - ct_2)\theta(ct - ct_1) \frac{\hat{\delta}^{\mu}_{\mu_1}}{\hat{\delta}^4(x - x_2)} \left[ \hat{j}^{(in)0}(x), \hat{j}^{(in)\mu_1}(x_1) \right]$$

$$+ \theta(ct - ct_1)\delta(ct_1 - ct_2) \frac{\hat{\delta}^{\mu_1}_{\mu_2}}{\hat{\delta}^4(x_1 - x_2)} \left[ \hat{j}^{(in)0}(x), \hat{j}^{(in)\mu_1}(0) \right] \right\}$$

$$+ \left( \frac{1}{i\hbar c^2} \right)^2 \left\{ \theta(ct - ct_1)\theta(ct_1 - ct_2) \left[ \left[ \hat{j}^{(in)\mu}(x), \hat{j}^{(in)\mu_1}(x_1) \right], \hat{j}^{(in)\mu_2}(x_2) \right] \right.$$ 

$$+ \theta(ct - ct_2)\theta(ct_2 - ct_1) \left[ \left[ \hat{j}^{(in)\mu}(x), \hat{j}^{(in)\mu_2}(x_2) \right], \hat{j}^{(in)\mu_1}(x_1) \right] \right\}. \quad (43)$$
\[
3! \chi^\mu_{\mu_1\mu_2\mu_3}(x, x_1, x_2, x_3) = \frac{\delta^3 j^\mu(x)}{\delta A^{\mu_1}(x_1) \delta A^{\mu_2}(x_2) \delta A^{\mu_3}(x_3)} \bigg|_{A=A(0)}
= \frac{1}{i\hbar c^2} \left\{ \begin{array}{c}
\theta(ct - ct_2)\delta(ct - ct_1)\delta(ct_2 - ct_3)\tilde{\delta}_{\mu_1}^\mu \delta^3(x - x_1)\tilde{\delta}_{\mu_2\mu_3}^\mu \delta^3(x_2 - x_3) \left[ j^{(in0)}(x), j^{(in0)}_{\mu_2}(x_2) \right] \\
+ \theta(ct - ct_3)\delta(ct - ct_2)\delta(ct_3 - ct_1)\tilde{\delta}_{\mu_2}^\mu \delta^3(x - x_2)\tilde{\delta}_{\mu_3\mu_1}^\mu \delta^3(x_3 - x_1) \left[ j^{(in0)}(x), j^{(in0)}_{\mu_3}(x_3) \right] \\
+ \theta(ct - ct_1)\delta(ct - ct_3)\delta(ct_1 - ct_2)\tilde{\delta}_{\mu_3}^\mu \delta^3(x - x_3)\tilde{\delta}_{\mu_1\mu_2}^\mu \delta^3(x_1 - x_2) \left[ j^{(in0)}(x), j^{(in0)}_{\mu_1}(x_1) \right]
\end{array} \right. \\
+ \left( \frac{1}{i\hbar c^2} \right)^2 \frac{-q}{mc^2} \left\{ \begin{array}{c}
\delta(ct - ct_1)\theta(ct_1 - ct_2)\theta(ct_2 - ct_3)\tilde{\delta}_{\mu_1}^\mu \delta^3(x - x_1) \left[ j^{(in0)}(x), j^{(in0)}_{\mu_2}(x_2) \right, j^{(in0)}_{\mu_3}(x_3) \right] \\
+ \delta(ct - ct_1)\theta(ct_1 - ct_3)\theta(ct_3 - ct_2)\tilde{\delta}_{\mu_1}^\mu \delta^3(x - x_1) \left[ j^{(in0)}(x), j^{(in0)}_{\mu_3}(x_3) \right, j^{(in0)}_{\mu_2}(x_2) \right] \\
+ \delta(ct - ct_2)\theta(ct_2 - ct_3)\theta(ct_3 - ct_1)\tilde{\delta}_{\mu_2}^\mu \delta^3(x - x_2) \left[ j^{(in0)}(x), j^{(in0)}_{\mu_3}(x_3) \right, j^{(in0)}_{\mu_1}(x_1) \right] \\
+ \delta(ct - ct_2)\theta(ct_2 - ct_1)\theta(ct_1 - ct_3)\tilde{\delta}_{\mu_2}^\mu \delta^3(x - x_2) \left[ j^{(in0)}(x), j^{(in0)}_{\mu_1}(x_1) \right, j^{(in0)}_{\mu_3}(x_3) \right] \\
+ \delta(ct - ct_3)\theta(ct_3 - ct_1)\theta(ct_1 - ct_2)\tilde{\delta}_{\mu_3}^\mu \delta^3(x - x_3) \left[ j^{(in0)}(x), j^{(in0)}_{\mu_2}(x_2) \right, j^{(in0)}_{\mu_1}(x_1) \right] \\
+ \delta(ct - ct_3)\theta(ct_3 - ct_2)\theta(ct_2 - ct_1)\tilde{\delta}_{\mu_3}^\mu \delta^3(x - x_3) \left[ j^{(in0)}(x), j^{(in0)}_{\mu_1}(x_1) \right, j^{(in0)}_{\mu_2}(x_2) \right]
\end{array} \right.
\]
\[ +\theta(ct - ct_3)\theta(ct_3 - ct_2)\theta(ct_2 - ct_1) \left\{ \left[ j^{(\text{in})}_{\mu_1}(x), j^{(\text{in})}_{\mu_2}(x), j^{(\text{in})}_{\mu_3}(x) \right] \right\} \]

The charge conservation, Eq.(24) and gauge invariance, Eq.(25) are respected in Eqs.(42)-(44). This fact is successfully checked after long and tedious calculations; a supplementary document is provided for details.

V. THE GROUND STATE IN DENSITY FUNCTIONAL THEORY AND SINGLE SUSCEPTIBILITY

The linear and nonlinear single susceptibilities are the expectation values of the corresponding operators, Eqs.(42)-(44), using the ground state in the non-perturbed electron system, which is specified by the simplified conditions in this paper:

\[ A(x) = A^{(0)}(x) = 0, \quad j^{(\text{EXT})}(x) = 0, \quad \phi(x) = \phi^{(0)}(x) \quad \text{and} \quad \rho^{(\text{EXT})}(x) \quad \text{are static}. \quad (45) \]

Let us explain how density functional theory[13, 14] may allow us to prepare the ground state and the complete set of the states in a many-electron system, refining the naive idea in Ref. [1]. For that purpose, we need the electron field operators together with the SP and VP satisfying the coupled equations, Eqs.(4)-(9). However, in the semiclassical treatment of the present theory, Eqs.(8) and (9) are replaced with their expectation values using the ground state, which we seek now on. Due to this procedure, the quantum many-electron effect, the so-called exchange-correlation effect is ignored. Therefore, the solution of Eqs.(4)-(9) as it is may not reproduce the electron charge density of the proper ground state, \( \rho_{\text{GS}}(r) \), which is obtained using the ordinary Hamiltonian including the two-body Coulomb interaction, converted from the SP under the Coulomb gauge. Such the electron density \( \rho_{\text{GS}}(r) \), in turn, brings about the proper SP \( \phi^{(0)}(x) \) under the Coulomb gauge. Suppose that the proper electron charge density \( \rho_{\text{GS}}(r) \) is already known under the ordinary Hamiltonian.

Now, we like to seek for the ground state \( |0\rangle \) in need, adjusting the auxiliary potential \( v^{(\text{AUX})}(r) \) to make the electron charge density fit the proper one:

\[ \langle 0|\hat{\rho}(x)|0 \rangle = \rho_{\text{GS}}(r). \quad (46) \]

Such a situation in Eq.(46) is assumed by Kohn and Sham in the density functional theory[14]. That is, Eqs.(4) and (5) are equivalent to Eq.(2.8) in Ref.[14] [the Kohn-Sham equation], if \( v^{(\text{AUX})}(r) \) is regarded as the so-called exchange-correlation potential.
For details, one may prepare the spin-orbital function $\varphi_k(r)$ ($k, \alpha$ stands for the orbital and spin states) as the eigenstate of the Kohn-Sham equation with the eigenenergy $\hbar \omega_k$. Under the conditions of Eq.(45), the Kohn-Sham equation is,

$$0 = \left( \hbar \omega_k - q\phi(0)(r) - \frac{\hbar}{2m}i\partial_i - v^{(\text{AUX})}(r) \right) \varphi_k(r),$$

(47)

where $v^{(\text{AUX})}(r)$ is set to the exchange-correlation potential, that guarantees Eq.(46). Then, $\hat{\psi}_\alpha(x) = \sum_k \varphi_k(r) \hat{a}_{k\alpha}(t)$ satisfies Eq.(4) under the condition Eq.(45), where $\hat{a}_{k\alpha}$ is the operator to annihilate the electron associated with the spin-orbital $\varphi_k(r)$ in the non-perturbative system. Considering $\{ \varphi_k(r) \}$ as a complete set of the one-electron functional space, the ground state with the electron number $n$ in the present theory is constructed as the single Slater determinant,

$$|0\rangle = \lim_{t_0 \to -\infty} \frac{1}{\sqrt{n!}} \prod_{k\alpha} \hat{a}_{k\alpha}^{(\text{in})}\dagger(t_0) |\text{vac}\rangle,$$

(48)

where $|\text{vac}\rangle$ is the vacuum state, and the indices $k\alpha$ scan over the $n$ spin-orbitals from the lowest eigenenergies. Furthermore, under the fixed $v^{(\text{AUX})}(r)$ and $\phi(0)(r)$, one may consider all the possible combination of $n$ spin-orbitals and obtain the normalized orthogonal complete set $\{|m\rangle | m = 0, 1, 2, \cdots \}$ in terms of all the possible single Slater determinants.

On the above logic, one should know the proper electron charge density $\rho_{GS}(r)$ beforehand to determine $v^{(\text{AUX})}(r)$, which is the universal functional of the electron density[13, 14]. In practice, however, one may solve the Kohn-Sham equation, possibly under the local density approximation for $v^{(\text{AUX})}(r)$, and reconsider the resulting charge density as $\rho_{GS}(r)$.

The expectation value of the single susceptibility operator is, $\langle 0 | \hat{\chi}_{\mu_1}(x, x_1) |0\rangle$, and, for example, the linear susceptibility becomes:

$$\langle 0 | \hat{\chi}_{\mu_1}^\mu(x, x_1) |0\rangle = -\frac{q}{mc^2} \delta_{\mu\mu_1} \delta^4(x - x_1) \langle 0 | \hat{j}^{(\text{in})\mu}(x) |0\rangle$$

$$+ \frac{1}{i\hbar c^2} \theta(ct - ct_1) \langle 0 | \left[ \hat{j}^{(\text{in})\mu}(x), \hat{j}^{(\text{in})\mu_1}(x_1) \right] |0\rangle.$$  

(49)

Next, to evaluate the products of two (or more) current density operators, e.g., the second term in Eq.(49), we may use the projection operator $\hat{1} = \sum_m |m\rangle\langle m|$. Now, the expectation
value in the second term of Eq. (49) becomes,

\[
\langle 0 | \left[ j^{(in0)}(x) , \tilde{j}^{(in0)}_{\mu_1}(x_1) \right] | 0 \rangle \\
= \sum_m \left\{ \langle 0 | j^{(in0)}(x) | m \rangle \langle m | \tilde{j}^{(in0)}_{\mu_1}(x_1) | 0 \rangle - \langle 0 | \tilde{j}^{(in0)}_{\mu_1}(x_1) | m \rangle \langle m | j^{(in0)}(x) | 0 \rangle \right\} \\
= \sum_m \lim_{t_0 \to \infty} \left\{ \langle 0 | e^{\frac{-i}{\hbar} \tilde{H}(t-t_0)} \tilde{j}^{(in0)}_{\mu_1}(x_1) | t=t_0 \rangle e^{\frac{i}{\hbar} \tilde{H}(t-t_0)} | 0 \rangle \langle m | e^{\frac{-i}{\hbar} \tilde{H}(t-t_0)} \tilde{j}^{(in0)}_{\mu_1}(x_1) | t=t_0 \rangle e^{\frac{i}{\hbar} \tilde{H}(t-t_0)} | 0 \rangle \right\} \\
\quad \quad - \langle 0 | e^{\frac{-i}{\hbar} \tilde{H}(t_1-t_0)} \tilde{j}^{(in0)}_{\mu_1}(x_1) | t_1=t_0 \rangle e^{\frac{i}{\hbar} \tilde{H}(t_1-t_0)} | m \rangle \langle m | e^{\frac{-i}{\hbar} \tilde{H}(t_1-t_0)} \tilde{j}^{(in0)}_{\mu_1}(x_1) | t_1=t_0 \rangle e^{\frac{i}{\hbar} \tilde{H}(t_1-t_0)} | 0 \rangle \right\} \\
= \sum_m \left\{ e^{\frac{i}{\hbar}(E_m-E_0)(t-t_1)} \langle 0 | \tilde{j}^{(in0)}_{\mu_1}(x_1) | t_1=-\infty \rangle | m \rangle \langle m | \tilde{j}^{(in0)}_{\mu_1}(x_1) | t_1=-\infty \rangle | 0 \rangle \right\} \\
\quad \quad - e^{\frac{i}{\hbar}(E_m-E_0)(t-t_1)} \langle 0 | \tilde{j}^{(in0)}_{\mu_1}(x_1) | t_1=-\infty \rangle | m \rangle \langle m | \tilde{j}^{(in0)}_{\mu_1}(x_1) | t_1=-\infty \rangle | 0 \rangle \right\}. \tag{50}
\]

In the induced charge and current densities obtained from the convolution integral of Eq. (50) with the perturbative EM field, the energy denominator will appear as shown in §VI.

In the above theoretical framework, |m\rangle’s are simply the members of the complete set, and, in principle, do not carry physical meaning of excited states of a many-electron system. Considering that the density functional theory concerns only the ground state of the many-electron system, the above treatment is a sound application of density functional theory to the response theory adequate for NFO. Remark that the variational principle based on Eq. (1) cannot determine the auxiliary potential, v^{(AUX)}(x) but is determined with the help of another theory, namely, the density functional theory.

As a summary, the quantum many-electron effect is temporally ignored in the present semiclassical theory, but is compensated with the support of the density functional theory. In other words, the SP inherently existing in the electron system is separated as \phi^{(0)}(x) and v^{(AUX)}(r), and the SP incidence may be treated equally with the VP incidence. Note that, \phi^{(0)}(x) is under the Coulomb gauge but the SP and VP incidences may be gauge-free, that is, the present response theory is still free from gauge-fixing.

VI. APPLICATION: A LOGICAL FALLACY TO USE THE ELECTRIC FIELD IN NEAR-FIELD OPTICS

Under non-resonant conditions in the optical near field, non-metallic materials cause various phenomena not observed in conventional optics, such as highly efficient light emission from indirect-transition-type semiconductors (LED[17, 18] and Laser[18, 19]), chem-
ical reaction with insufficient photon energy (chemical vapor deposition[20], optical NF lithography[21], optical NF etching[22]), frequency up-conversion[23, 24], non-adiabatic effect beyond forbidden transition (local energy concentration[25], nano-photonic gate device[26]), and gigantic magneto-optical rotation of the LED[18, 27].

These experimental results draw attention to a fundamental role of the non-resonant condition in NFO. We have no complete answer at this stage but the application of the present response theory to a many-electron system in NFO shows a logical fallacy to use the electric field and the electric permittivity, and the necessity to use the EM potential and the associated single susceptibility. The discussion of the one-electron system appeared in Ref.[1], but is concisely reviewed below in §VI B-§VI E because the many-electron version in §VIF may be simply a recast of the one-electron version, owing to the density functional theory. For the readability, calculation details are given in Appendix B.

A. Classification of optical systems

First, let us classify the optical systems. The two systems under near- and far-field incidence conditions in FIG.1 are subdivided into two classes depending on the near- or far-field observation condition. These four classes are listed in TABLE I, together with a summary of the results mentioned below. In particular, the systems of (I′) and (II′) are the limiting cases of null longitudinal incidence of the systems (I) and (II), respectively. Thus, in the systems (I′) and (II′), the longitudinal response vanishes and the difference in response may not be observed. In the following, therefore, we focus mainly on systems (I) and (II), in which longitudinal incidence exists.

B. Susceptibilities associated with longitudinal and transverse electric fields

Applying the present linear response theory and the long wave approximation (LWA) to the spinless one-electron system with two levels on a small scale, the induced charge and current densities (as a result of the response), \( \Delta \rho(\mathbf{r}, t) \) and \( \Delta j(\mathbf{r}, t) \), are described as the total derivative with respect to the longitudinal and transverse electric fields (as the cause of the response), \( \Delta E^{(l)}(\mathbf{0}, t) \) and \( \Delta E^{(t)}(\mathbf{0}, t) \), where \( \mathbf{0} \) is the representative position in the
TABLE I: Classification of optical systems by distance from the target material to the light source and distance from that to the observation point, together with a summary of the results; the validity of the electric field as the cause of the response.

| Source: Δρ and Δj | Near-field observation | Far-field observation | Source: Δj |
|-------------------|------------------------|----------------------|-------------|
| Near-field incidence: ΔE^{(l)} + ΔE^{(t)} | (I) NF optical system | (II) NF optical system |          |
| Validity of the electric field | non-resonant / resonant | non-resonant / resonant | OK / OK |
| Far-field incidence: ΔE^{(t)} | (I') NF optical system | (II') conventional optical system | |
| Validity of the electric field | non-resonant / resonant | non-resonant / resonant | OK / OK |

where the partial derivative coefficients, $\chi^{\cdots\cdots}(r, \omega)$'s are susceptibilities associated with the longitudinal and transversal electric fields. In Eq.(52), the time derivatives of the two types of electric fields, namely, $\Delta \dot{E}_{j}^{(l)}(0, t)$ and $\Delta \dot{E}_{j}^{(t)}(0, t)$, are regarded as the causes, instead of the two types of electric fields themselves. The magnetic response will appear in the higher order of the LWA and vanishes in Eqs.(51) and (52) representing the leading order; see Refs.[4, 5] and §VI G. For the present spinless electron system, the electron field operators, $\hat{\psi}^\dagger(x)$, $\hat{\psi}(x)$ is reconsidered as $\hat{\psi}^\dagger(x)$, $\hat{\psi}(x)$, respectively, eliminating the index of the spin state, $\alpha$.

To evaluate the susceptibilities in Eqs.(51) and (52), the two levels are assumed to be the ground and excited states in the non-perturbed system with eigenenergies, $\hbar \omega_0$ and $\hbar \omega_1$, and orbitals, $\varphi_0(r)$ and $\varphi_1(r)$, respectively. Those orbitals are assumed to be bound states expressed by real functions, carry well-defined and distinct spatial parities (even and odd parities), and form the normalized orthogonal complete set. The excitation energy is $\hbar \Delta \omega_1 \equiv \hbar \omega_1 - \hbar \omega_0 > 0$; this finite excitation energy means that the target is a non-metallic material, such as a molecule, nano-structured semiconductor and insulator.
The susceptibilities in Eqs. (51) and (52) are derived in §B.1, and those leading to the induced charge density result in the following:

\[ \chi_j^{\rho-\ell}(\mathbf{r},\omega) = \chi_j^{\rho-t}(\mathbf{r},\omega) = 2q^2 \frac{\eta}{\eta^2 - 1} \frac{1}{\hbar \omega} D_j \varphi_0(\mathbf{r}) \varphi_1(\mathbf{r}), \quad (53) \]

where \( \eta \equiv \frac{\hbar \Delta \omega_1}{\hbar \omega} = \frac{\text{excitation energy}}{\text{photon energy}} \), and

\[ D_i \equiv \int d^3 \mathbf{r} \varphi_1(\mathbf{r}) r_i \varphi_0(\mathbf{r}). \quad (55) \]

This means that the responses to the longitudinal and transverse electric fields are common, such that the induced charge density has a linear relationship with the total electric field, namely, \( \Delta \rho(\mathbf{r},t) = \chi_j^{\rho-\ell}(\mathbf{r},\omega) \left( \Delta E_j^{\ell}(0,t) + \Delta E_j^{t}(0,t) \right) \).

The susceptibilities leading to the induced current density are not so simple and result in the following:

\[ \chi_j^{j-\ell}(\mathbf{r},\omega) = q^2 \frac{\hbar^2}{m} \frac{1}{\eta^2 - 1} \frac{1}{(\hbar \omega)^2} D_j \left( \partial_i \varphi_1(\mathbf{r}) \varphi_0(\mathbf{r}) - \varphi_1(\mathbf{r}) \partial_i \varphi_0(\mathbf{r}) \right), \quad (56) \]

\[ \chi_j^{j-t}(\mathbf{r},\omega) = \eta^2 \chi_j^{j-\ell}(\mathbf{r},\omega) - q^2 \frac{\hbar^2}{m} \frac{1}{(\hbar \omega)^2} \varphi_0(\mathbf{r}) \varphi_0(\mathbf{r}). \quad (57) \]

The susceptibility to the transverse electric field, Eq. (57), is composed of two terms. The first term, namely, the resonant term, includes the energy denominator enhanced under the resonant condition, \( \eta \approx 1 \), as in the susceptibility to the longitudinal electric field, Eq. (56). The second term, namely, the non-resonant term, does not include such a resonance factor.

### C. Equal responses under the resonant condition

Under the condition \( \eta \approx 1 \) in all cases in TABLE I, Eq. (57) is dominated by the resonant term (the first term) over the non-resonant term (the second term) and asymptotically equals Eq. (56).

\[ \chi_{ij}^{j-\ell}(\mathbf{r},\omega) \approx \chi_{ij}^{j-\ell}(\mathbf{r},\omega). \quad (58) \]

Equation (58) together with Eq. (53) reveal the equivalency of the responses to the longitudinal and transverse electric fields, so that the total electric field is regarded as the cause of the response in all the optical systems under the resonant condition listed in TABLE I.
D. Equal responses under the far-field observation condition

In the system (II) and (II') in TABLE I, the far field to be observed is insensitive to the details of the source but is determined by the spatial average of the source. Under the LWA, such an average can be achieved by the spatial average of the susceptibilities. Detailed calculations are shown in §B2 and the results are as follows:

\[
\bar{\chi}_{j}(r, \omega) = \chi_{j}(r, \omega) = 0, \quad (59)
\]

\[
\bar{\chi}^{(t)}_{ij}(r, \omega) = \frac{1}{mV} \left( \frac{q^2\hbar^2}{\Delta \omega_1} - (\hbar \omega)^2 \right), \quad (60)
\]

where the overline represents the spatial average and \( V \) is the volume of the target material. From Eqs.(59) and (60), one may not observe different responses to the two types of incidences under the far-field observation condition. The null response represented in Eq.(59) is reasonable because the induced charge density yields the longitudinal electric field, which has a non-radiative nature and vanishes in the far-field regime.

E. Unequal responses under the non-resonant, NF incidence, and NF observation conditions

The different responses to the longitudinal and transverse electric fields claimed in §1B may be detected only in the system (I) in TABLE I under the non-resonant condition, which is just the compliment to the popular optical systems under the resonant condition or the far-field incidence condition or the far-field observation condition. In the NF optical system (I) with a non-metallic material under the non-resonant condition, the total electric field is not the cause of the response; therefore, the response may not be described by the ordinary constitutive equation, namely, the linear relationship between the polarization and "electric field" via the electric permittivity, so that the single susceptibility is essential to treat separately the longitudinal and transverse incidences.

F. Extension to the many-electron system

The above one-electron model is very simple and the responses may be modified in a many-electron system or a low-symmetry system. However, the difference in the responses
to the two types of electric fields originates in the non-relativistic nature of the system (as stated in §1B), and should survive in actual NF optical systems with non-metallic materials (the materials with finite excitation energy). Actually, the results revealed in §VI B-§VIE are applicable to the corresponding many-electron system, considering the auxiliary potential \( v^{(\text{AUX})}(x) \) to construct the orbitals using the Kohn-Sham equation (46), and replacing the complete orthogonal set composed of the one-electron ground and excited states, 
\[
\{ \hat{a}_0^{(\text{in}0)}(-\infty)|\text{vac}\rangle, \hat{a}_1^{(\text{in}0)}(-\infty)|\text{vac}\rangle \} \quad (-\infty \text{ means the time of the infinite past})
\]
to the corresponding set, composed of two single Slater determinants, 
\[
\{ |0\rangle, \hat{a}_1^{(\text{in}0)}(-\infty)\hat{a}_0^{(\text{in}0)}(-\infty)|0\rangle \}
\]
where \(|0\rangle\) is the ground state in the density functional theory as defined in §V, and \(\hat{a}_0^{(\text{in}0)}\) [\(\hat{a}_0^{(\text{in}0)}\dagger\)] and \(\hat{a}_1^{(\text{in}0)}\) [\(\hat{a}_1^{(\text{in}0)}\dagger\)] are the annihilation [and creation] operators associated by the the highest occupied molecular orbital (HOMO) and the lowest unoccupied molecular orbital (LUMO), respectively, determined by the Kohn-Sham equation (46). Owing to the density functional theory, recasting the formulation in the one-electron system brings that in the many-electron system, if the HOMO and LUMO dominate the excitation.

As a result, the many-electron system of a non-metallic material under the non-resonant, NF incidence, and NF observation conditions may not be described in terms of the electric field and the associated permittivity. Instead, the EM potential and the single susceptibility are essential.

G. Comparison with the existing theories

In NFO, the response to the longitudinal electric field is discussed in Chap. 5 in Ref.[5] and Chap. 9 in Ref.[9], as mentioned in §1C. The present work is a further comparison of the responses to the two-types of electric field, considering the non-resonant condition.

Another logical fallacy to use the electric and magnetic fields is pointed out by Cho, as briefly mentioned in §1B. In Refs.[4, 5], Cho derived a Taylor series of the nonlocal response function[8] under the LWA, and assigned the electric permittivity and magnetic permeability in the macroscopic constitutive equation as the term of order \(\mathcal{O}(ka)^0\) (the leading order) and \(\mathcal{O}(ka)^2\), respectively, where \(ka \ll 1, 2\pi/k\) is the light wavelength, and \(a\) is the representative size of the material. Furthermore, he pointed out that the ordinary two susceptibilities are irrational because the separability of the electric and magnetic responses not applicable and the term of order \(\mathcal{O}(ka)^1\) appears in a chiral symmetric system, including
a NF optical system with a low-symmetric nanostructure. The present demonstration is concerned with the logical fallacy, which appears in the electric response (the leading order from the viewpoint of Cho) in NFO under a non-resonant condition.

H. A remark on applying the finite differential time domain (FDTD) method to NF optical systems

The macroscopic constitutive equations in terms of the electron permittivity and magnetic permeability have been widely employed to calculate the optical near field in the FDTD method[16]. One may notice that the permittivity in the FDTD method carries a simple spatial dependence and leads to some quantitative error. Actually, the microscopic susceptibilities, for example, Eq.(53), Eq.(56), and Eq.(57), have rippling spatial distributions originating from the orbitals.

In the case of the NF optical system (I) in TABLE I with a non-metallic material under the non-resonant condition, the situation is more serious because the concept electric field is not available, such that it is a logical fallacy to use the macroscopic constitutive equation. Thus, a novel simulation method is necessary, in particular, for the NF optical system with a non-metallic material.

I. Why this fallacy has been missed for a long time?

Why has the comparison of responses to the two types of electric fields not been addressed in NF optical theory? First, in the long history of optics, the NF optical system (I) in TABLE I under a non-resonant condition has been out of focus. Such a system could not be resolved until the technical difficulty of NF observation was overcome. Additionally, resonance phenomena continue to attract attention. Furthermore, even in NFO, there has been less emphasis on non-metallic materials, as opposed to metallic materials, which are essential for plasmonics.

The second reason is that the ordinary Hamiltonian for a many-electron system does not include the longitudinal electric field, which is rewritten to the two-body Coulomb interaction, as stated in §1C. With this Hamiltonian, the non-linear response to the longitudinal electric field (the SP under the Coulomb gauge) incidence accompanies the Coulomb inter-
action, and is ignored or unequally treated compared with the response to the transverse electric field (the VP under the Coulomb gauge).

J. Summary of this section

In the NF optical system (I) in TABLE I, the responses to the longitudinal and transverse electric fields should be separately treated, and in a more general view point beyond the LWA and linear response theory, it is essential to employ the linear and nonlinear single susceptibilities, considering both of the SP and VP equally as the cause of response.

To the best of our knowledge, the NF optical system with non-metallic material under the non-resonant condition, namely, the system (I) in TABLE I, is the third example that cannot be described in terms of the electric field and/or magnetic field, after the superconductor system with the Meissner effect[2] and the electron system with the Aharonov-Bohm effect[3], as mentioned in §I A.

VII. SUMMARY

1. Aiming to investigate electron response in NFO, we define the linear and nonlinear single susceptibilities, equally considering the SP and VP as the cause of the response.

2. It is shown that the present single linear and nonlinear susceptibilities guarantee charge conservation and gauge invariance.

3. The linear and nonlinear susceptibilities in the form of Heisenberg operators are derived systematically by means of the functional derivatives of the action integral of the matter with respect to the SP and VP.

4. It is shown that the density functional theory may be used in the non-perturbed system and support to prepare the ground state and a complete set of states, which in turn are used to evaluate the expectation values of the operators of the linear and nonlinear susceptibilities.

5. Applying the present response theory to a simplified model system, it is shown that the single susceptibility is essential to describe the response of the optical system
with non-metallic material under the non-resonant, NF incidence, and NF observation conditions.

Some remaining problems meriting further investigation include:

1. Applying the present response theory to actual non-resonant NF optical systems with a non-metallic material in Refs. [17]-[27] to explore the mechanism leading to the outstanding experimental results such as the high-efficient light emission and gigantic magneto-optical effect, etc.

2. Developing a constitutive equation based on the single susceptibility which can aid experimentalists in NFO as a substitute for the electric permittivity and magnetic permeability of ordinary optics.

3. Developing a practical simulator for the many-electron system in NFO, using the present response theory with the support of the density functional theory, as the replacement of the FDTD simulation method,

4. Extending the response theory to treat the spin-polarization system in NFO, based on the Pauli or Dirac equation.

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Appendix A: Optimization of Electron Field Operators Under Arbitrary EM Potential

Under a given EM potential, \( A^\nu \), the electron field operator optimized to satisfy Eq.(4) is considered as the functional of \( A^\nu \), i.e., \( \hat{\psi}_\alpha(x; [A^\nu]), \hat{\psi}_\alpha^\dagger(x; [A^\nu]) \). Then, the next equation holds for \( n = 0, 1, 2, \ldots \): \( \delta \hat{\psi}_\alpha^\dagger(x') \delta I_{\text{mat}} \bigg|_{\hat{\psi}_\alpha, \hat{\psi}_\alpha^\dagger, A^\nu} = 0 \), \( \delta A^\mu_1(x_1) \cdots \delta A^\mu_n(x_n) / \delta A^\nu = 0 \), \( \delta I_{\text{mat}} \bigg|_{\hat{\psi}_\alpha, \hat{\psi}_\alpha^\dagger, A^\nu} = 0 \).

Proof: Equation (4) should be hold both under \( A^{(0)\nu} \) (non-perturbative EM potential) and under \( A^{(0)\nu} + \Delta A^\nu \), therefore,

\[
\delta \hat{\psi}_\alpha^\dagger(x') \delta I_{\text{mat}} \bigg|_{\hat{\psi}_\alpha, \hat{\psi}_\alpha^\dagger, A^\nu} = 0, \\
\delta A^\mu_1(x_1) \cdots \delta A^\mu_n(x_n) / \delta A^\nu = 0.
\]

Taylor expansion leads to:

\[
\sum_{n=0}^{\infty} \frac{1}{n!} \int d^4 x_n \cdots \int d^4 x_1 \frac{\delta^n \left( \delta \hat{\psi}_\alpha^\dagger(x') \delta I_{\text{mat}} \right)_{\hat{\psi}_\alpha, \hat{\psi}_\alpha^\dagger, A^\nu}}{\delta A^\mu_n(x_n) \cdots \delta A^\mu_1(x_1)_{\hat{\psi}_\alpha, \hat{\psi}_\alpha^\dagger, A^\nu}} = 0,
\]

Considering this equation as the identity with respect to \( \Delta A^\mu(x) \) results in Eq.(A1). Equation (A2) is proved in the same manner, starting from Eq.(5).

Appendix B: Calculation details in §VII

Here we provide the calculation details in §VII, including the derivation of the unfamiliar relationship Eq.(B14) between two types of dipole transition matrix elements.

1. Derivation of the constitutive equations, Eqs.(51) and (52), and the susceptibilities, Eq.(53), Eqs.(56) and (57)

The incident SP and VP, \( \Delta \phi(r, t) \) and \( \Delta A_i(r, t) \), are assumed to be monochromatic with the angular momentum \( \omega \), and are expressed using the Coulomb gauge and LWA as follows:

\[
\Delta \phi(r, t) = \Delta \phi(r) \cos \omega t = \left( \Delta \phi(0) - \Delta E^{(t)}(0) \cdot r \right) \cos \omega t, \\
\Delta A(r, t) = \Delta A(r) \sin(\omega t + \xi) = -\frac{1}{\omega} \Delta E^{(t)}(0) \sin(\omega t + \xi),
\]

30
where \( \xi \) is the phase difference between the two incident potentials. In the spinless one-electron system, the linear response theory with Eqs.(20) and (42) leads to the Heisenberg operators of the induced charge and current densities, as follows in the three-element representation:

\[
\begin{align*}
\Delta \hat{\rho}(\mathbf{r}, t) &= \int_{-\infty}^{t} dt_1 \int d^3 r_1 \left\{ \frac{1}{i\hbar} \left[ \hat{\rho}^{(in0)}(\mathbf{r}, t), \hat{\rho}^{(in0)}(\mathbf{r}_1, t_1) \right] \Delta \phi(\mathbf{r}_1, t_1) \right. \\
&\quad \left. - \frac{1}{i\hbar} \left[ \hat{\rho}^{(in0)}(\mathbf{r}, t), \hat{j}^{(in0)}_{i1}(\mathbf{r}_1, t_1) \right] \Delta A_{i1}(\mathbf{r}_1, t_1) \right\}, \\
\Delta \hat{j}_i(\mathbf{r}, t) &= \int_{-\infty}^{t} dt_1 \int d^3 r_1 \left\{ \frac{1}{i\hbar} \left[ \hat{j}^{(in0)}_i(\mathbf{r}, t), \hat{\rho}^{(in0)}(\mathbf{r}_1, t_1) \right] \Delta \phi(\mathbf{r}_1, t_1) \right. \\
&\quad \left. - \frac{1}{i\hbar} \left[ \hat{j}^{(in0)}_i(\mathbf{r}, t), \hat{j}^{(in0)}_{i1}(\mathbf{r}_1, t_1) \right] \Delta A_{i1}(\mathbf{r}_1, t_1) \right\} - \frac{q}{m} \hat{\rho}^{(in0)}(\mathbf{r}, t) \Delta A_i(\mathbf{r}, t).
\end{align*}
\]

(B3, B4)

The last term in Eq.(B4) originates from the non-relativistic nature of the system and is needed to maintain charge conservation law.

Evaluating the expectation value of Eqs.(B3) and (B4) using the ground state \([\phi_0(\mathbf{r}) \text{ in Eq.}(B8)]\) and substituting Eqs.(B1) and (B2) leads to Eqs.(51) and (52), in which the causes of the responses are the two types of electric fields and their temporal derivatives, defined as

\[
\begin{align*}
\Delta E_j^{(t)}(0, t) &\equiv \Delta E_j^{(t)}(0) \cos \omega t, \quad \Delta E_j^{(t)}(0, t) \equiv \Delta E_j^{(t)}(0) \cos(\omega t + \xi), \\
\Delta \dot{E}_j^{(t)}(0, t) &\equiv \partial \Delta E_j^{(t)}(0, t), \quad \Delta \dot{E}_j^{(t)}(0, t) \equiv \partial \Delta E_j^{(t)}(0, t).
\end{align*}
\]

(B5, B6)

In the above, no magnetic response appears because it is the higher order in the LWA as revealed by Cho [4, 5]. To obtain susceptibilities, Eq.(53),Eqs.(56) and (57) using the two-level model, we take the expectation values of Eqs.(B3) and (B4) using the ground state \(\varphi_0(\mathbf{r})\), insert the projection operator [the left side of the second equation in Eq.(B7)], between the two operators in the commutators, and integrate over the domains of \(t_1\) and \(\mathbf{r}_1\). We assume that the two orbitals are real functions, and form the normalized orthogonal complete set:

\[
\int d^3 r \varphi_m(\mathbf{r}) \varphi_n(\mathbf{r}) = \delta_{mn}, \quad \sum_m \varphi_m(\mathbf{r}) \varphi_m(\mathbf{r}') = \delta^3(\mathbf{r} - \mathbf{r}'),
\]

(B7)

where \(\varphi_m(\mathbf{r})\) satisfies,

\[
\hat{H}^{(0)} \varphi_m(\mathbf{r}) = \hbar \omega_m \varphi_m(\mathbf{r}), \quad (m = 0, 1).
\]

(B8)
Having real orbitals infers even temporal parity, such that there is a null VP (or magnetic field) in the non-perturbed system. To derive the susceptibilities associated with the transversal electric field in Eqs.(53) and (57), we use the well-known linear relationship between the two types of dipole transition matrix elements,

\[ C_i \equiv \int d^3r \left( \partial_i \varphi_1(r) \varphi_0(r) - \varphi_1(r) \partial_i \varphi_0(r) \right) = \frac{2m}{\hbar^2} \hbar \Delta \omega_1 D_i. \]  

Equation (B9) is derived from the matrix element of Heisenberg equation for dipole charge density:

\[ \frac{\partial}{\partial t} r_j \hat{\rho}^{(in0)}(r, t) = \frac{1}{i\hbar} \left[ r_j \hat{\rho}^{(in0)}(r, t), \hat{H}^{(0)} \right], \]  

using \( \hat{\rho}^{(in0)}(r, t) = e^{-\frac{\hat{H}^{(0)}}{\hbar}} \rho^{(in0)}(r, 0) e^{+\frac{\hat{H}^{(0)}}{\hbar}} \) and the projection operator, i.e., the second equation in Eq.(B7) satisfying Eq.(B8).

2. Derivation of the spatial average of the susceptibilities, Eqs.(59) and (60)

The following replacements in Eq.(53), Eqs.(56) and (57) lead to Eqs.(59) and (60):

\[ \varphi_0(r) \varphi_1(r) \rightarrow \frac{1}{V} \int d^3r \varphi_0(r) \varphi_1(r) = 0, \]  

\[ \partial_i \varphi_1(r) \varphi_0(r) - \varphi_1(r) \partial_i \varphi_0(r) \rightarrow \frac{1}{V} \int d^3r \partial_i \varphi_1(r) \varphi_0(r) - \varphi_1(r) \partial_i \varphi_0(r) = \frac{1}{V} C_i, \]  

\[ \varphi_0(r) \varphi_0(r) \rightarrow \frac{1}{V} \int d^3r \varphi_0(r) \varphi_0(r) = \frac{1}{V}. \]  

To derive Eq.(60), we additionally use the trade-off relationship between the two types of dipole transition matrix elements,

\[ D_i C_j = \delta_{ij}. \]  

This is effective in the two-level system with well-defined parity and derived from the quantum-mechanical commutation relationship:

\[ [r_i, \frac{\hbar}{i} \partial_j] = i\hbar \delta_{ij}, \quad \text{i.e.,} \quad r_i \left( \frac{\hbar}{i} \partial_j \cdots \right) + \frac{\hbar}{i} \partial_j (r_i \cdots) = i\hbar \delta_{ij} \cdots. \]  

Inserting the projection operator between \( r_i \) and \( \frac{\hbar}{i} \partial_j \), and eliminating the null integrals caused by mismatched parity result in Eq.(B14). From Eq.(B9) and Eq.(B14), \( D_i \) and \( C_i \) are specified as

\[ D_i = \frac{1}{C_i} = \frac{\hbar}{\sqrt{2m \hbar \Delta \omega_1}}. \]
(We do not use Eq.(B16) in this paper.)

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Supplementary Document for
"Theory of Single Susceptibility for Near-field Optics
Equally Associated with Scalar and Vector Potentials"

Check for Charge Conservation Law and Gauge Invariance

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Abstract

Charge conservation law and gauge invariance are explicitly checked for the linear and nonlinear single susceptibilities derived in the main text.

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S. CHARGE CONSERVATION LAW AND GAUGE INVARIANCE OF LINEAR AND NONLINEAR FOUR-ELEMENT SINGLE SUSCEPTIBILITY

S 1. Linear single susceptibility: Eq.(40) in the main text

To show that four-element linear single susceptibility guarantees the charge conservation law, Eq.(24) in the main text, suppose the four-element divergence of Eq.(40) in the main text, considering

\[
\partial_\mu \hat{\chi}^{\mu\mu_1}(x, x_1) = 0,
\]

\[
\frac{-q}{mc^2} \delta(c t - c t_1) \bar{\delta}_{\mu_1} \partial_\mu \left( \delta^3(x - x_1) \hat{j}^{(in)0}(x) \right)
\]

\[
+ \frac{1}{i\hbar c^2} \delta(c t - c t_1) \left[ \hat{j}^{(in)0}(x), \hat{j}^{(in)0}_{\mu_1}(x_1) \right]
\]

\[
= 0.
\]

In the second term of the second hand, we use the following commutation relationship at the same time :

\[
\delta(c t - c t_1) \left[ \hat{j}^{(in)0}(x), \hat{j}^{(in)0}_{\mu_1}(x_1) \right] = -i\hbar c^2 \frac{-q}{mc^2} \delta(c t - c t_1) \bar{\delta}_{\mu_1} \partial_\mu \left( \delta^3(x - x_1) \hat{j}^{(in)0}(x) \right)
\]

\[
= -i\hbar c^2 \frac{-q}{mc^2} \delta(c t - c t_1) \bar{\delta}_{\mu_1} \left( \partial_\mu \delta^3(x - x_1) \right) \hat{j}^{(in)0}(x_1)
\]

The proof of Eq.(S2) is as follows: If \( \mu_1 = 0 \) in the left hand side of Equation (S2), it is the commutator between charge density operator at the same time, and is zero.

\[
\left[ \hat{j}^{(in)0}(x), \hat{j}^{(in)0}_{\mu_1}(x_1) \right]_{t_1 = t} = c^2 q^2 \left[ \hat{\psi}^\dagger_\alpha(x) \hat{\psi}_\alpha(x), \hat{\psi}^\dagger_\alpha(x_1) \hat{\psi}_\alpha(x_1) \right]_{t_1 = t} = 0.
\]

One may check Eq.(S3) by a straightforward calculation using the anti-commutation relation of electron field operators at the same time,

\[
\left[ \hat{\psi}_\alpha(x), \hat{\psi}^\dagger_\alpha(x_1) \right]_{+, t_1 = t} = \delta^3(x - x_1).
\]

Next, if \( \mu_1 = i \in \{1, 2, 3\} \) in the left hand side of Equation (S2), the commutator in three-element representation becomes as follows:

\[
\left[ \hat{j}^{(in)0}(x), \hat{j}^{(in)0}_{\mu_1=i}(x_1) \right]_{t_1 = t}
\]

\[
= -cq \frac{q}{2m} \left[ \hat{\psi}^\dagger_\alpha(x) \hat{\psi}_\alpha(x), \hat{\psi}^\dagger_\alpha(x_1) \left( \frac{\hbar}{i} \partial^1_1 - q A^0_1(x_1) \right) \hat{\psi}_\alpha(x_1)
\]

\[
+ \left( \frac{\hbar}{-i} \partial^1_1 - q A^0_1(x_1) \right) \hat{\psi}^\dagger_\alpha(x_1) \right) \hat{\psi}_\alpha(x_1) \right]_{t_1 = t}
\]
As the term includes $qA_i^{(0)}(x_1)$ is zero following Eq.(S3), let us treat the term including the derivative.

\[
\left[\hat{\gamma}^{(in0)}_{\mu}(x), \hat{j}^{(in0)}_{\mu_1=x_1}(x_1)\right]_{t_1=t} = -\frac{cq}{2}i\hbar c^2\frac{-q}{mc^2}\lim_{x_1^*\to x_1} \partial_{x_1}^* \left[ \hat{\psi}_\alpha^+(x) \hat{\psi}_\alpha(x) + \hat{\psi}_\alpha^+(x_1) \hat{\psi}_\alpha(x) \right]_{t_1=t} \\
= -\frac{cq}{2}i\hbar c^2\frac{-q}{mc^2} \lim_{x_1^*\to x_1} \partial_{x_1}^* \left[ \left( \delta^3(x-x_1) \left( \hat{\psi}_\alpha^+(x) \hat{\psi}_\alpha(x_1) + \hat{\psi}_\alpha^+(x_1) \hat{\psi}_\alpha(x) \right) \right.ight.

\[
\left. \left. \left. + (\partial_3 \delta^3(x-x_1)) \left( \hat{\psi}_\alpha^+(x) \hat{\psi}_\alpha(x_1) + \hat{\psi}_\alpha^+(x_1) \hat{\psi}_\alpha(x) \right) \right)_{t_1=t} \right] \\
= -\frac{cq}{2}i\hbar c^2\frac{-q}{mc^2} \partial_{x_1} \left( \delta^3(x-x_1) \left( \hat{\psi}_\alpha^+(x) \hat{\psi}_\alpha(x_1) + \hat{\psi}_\alpha^+(x_1) \hat{\psi}_\alpha(x) \right) \right)_{t_1=t} \\
= -cq\hbar c^2\frac{-q}{mc^2} \partial_{x_1} \left( \delta^3(x-x_1) \hat{\psi}_\alpha^+(x_1) \hat{\psi}_\alpha(x_1) \right) = -i\hbar c^2\frac{-q}{mc^2} \delta^3(x-x_1) \hat{j}^{(in0)0}(x) \right) \\
= -cq\hbar c^2\frac{-q}{mc^2} \partial_{x_1} \left( \delta^3(x-x_1) \hat{\psi}_\alpha^+(x_1) \hat{\psi}_\alpha(x_1) \right)_{t_1=t} = -i\hbar c^2\frac{-q}{mc^2} \delta^3(x-x_1) \hat{j}^{(in0)0}(x_1)_{t_1=t},
\]

where the last two-way expressions Eqs.(S5) and (S6) are in four-element representation instead of three-element representation. Summarizing Eqs.(S3) and (S6) result in Eq.(S2). As a result, the present four-element linear susceptibility, Eq.(40) in the main text maintains the charge conservation law, Eq.(24) in the main text.

For the proof for the gauge invariance, Eq.(25) in the main text, of the linear susceptibility, suppose the four-element divergence with respect to $x_1$. Then, using Eq.(S2) with the replacement, $x \leftrightarrow x_1$ and the relation, $\tilde{\delta}^\mu_1 \partial_\mu = -\tilde{\delta}_\mu \partial^\mu$, one may obtain:

\[
\partial_\mu \chi^\mu_{\mu_1}(x, x_1) = -\frac{q}{mc^2} \delta(ct - ct_1) \left( \tilde{\delta}^\mu_1 \partial_\mu \delta^3(x-x_1) \right) \hat{j}^{(in0)0}(x) \\
= -\frac{1}{i\hbar c^2} \delta(ct - ct_1) \left[ \hat{j}^{(in0)0}(x), \hat{j}^{(in0)0}_0(x_1) \right]
\]

As shown above, the linear susceptibility Eq.(40) in the main text maintains the gauge invariance Eq.(25) in the main text.
S 2.  Second order nonlinear single susceptibility: Eq.(41) in the main text

Next, let us show that the charge conservation law is satisfied by the second order nonlinear single susceptibility, Eq.(41) in the main text. Operating $\partial_\mu = \delta^{\mu}_0 \partial_0 + \delta^{\mu}_1 \partial_1 + \delta^{\mu}_2 \partial_2 + \delta^{\mu}_3 \partial_3$ to Eq.(41) in the main text and considering $\partial_\mu \hat{\gamma}^{(in0)}\mu(x) = 0$ (the charge conservation law for the current density operator in the non-interacting system), the surviving terms are those the operator $\partial_\mu$ operates on the step function or delta function in front of the commutator, and operates on $\hat{\gamma}^{(in0)}\mu(x)$ in the commutator.

$$\partial_\mu 2! \hat{\chi}^{\mu}_{\mu_1\mu_2}(x, x_1, x_2) =$$

$$\frac{1}{\hbar c^2} \frac{-q}{mc^2} \left\{ \delta(c - c_1)\theta(c - c_2) \left[ \tilde{\delta}^{\mu}_{\mu_1} \partial_\mu \left( \delta^3(x - x_1) \left[ \hat{j}^{(in0)}\mu_0(x), \hat{j}^{(in0)}\mu_2(x_2) \right] \right) \right] \right\} (S8a)

+ \delta(c - c_2)\theta(c - c_1) \left[ \tilde{\delta}^{\mu}_{\mu_2} \partial_\mu \left( \delta^3(x - x_2) \left[ \hat{j}^{(in0)}\mu_0(x), \hat{j}^{(in0)}\mu_1(x_1) \right] \right) \right] (S8b)

+ \delta(c - c_1)\delta(c_1 - c_2) \left[ \delta^3(x_1 - x_2) \left[ \hat{j}^{(in0)}\mu_0(x), \hat{j}^{(in0)}\mu_1(x_1) \right] \right] (S8c)

+ \left( \frac{1}{\hbar c^2} \right)^2 \left\{ \delta(c - c_1)\theta(c_1 - c_2) \left[ \left[ \hat{j}^{(in0)}\mu_0(x), \hat{j}^{(in0)}\mu_2(x_2) \right], \hat{j}^{(in0)}\mu_1(x_1) \right] \right\}. (S8e)$$

Applying Eq.(S2), the third term (S8c) vanishes, and the fourth and fifth terms (S8d) and (S8e) cancel the first and second terms (S8a) and (S8b), respectively.

As a result, the second order nonlinear single susceptibility operator Eq.(41) in the main text maintains the charge conservation law, Eq.(24) in the main text.

To check the gauge invariance of the second order nonlinear single susceptibility operator, let us operate $\partial^{\mu_1}$ to Eq.(41) in the main text.

$$\partial^{\mu_1} 2! \hat{\chi}^{\mu}_{\mu_1\mu_2}(x, x_1, x_2) =$$

$$\frac{1}{\hbar c^2} \frac{-q}{mc^2} \left\{ \delta(c - c_1)\theta(c - c_2) \left[ \tilde{\delta}^{\mu}_{\mu_1} \partial^{\mu_1} \delta^3(x - x_1) \left[ \hat{j}^{(in0)}\mu_0(x), \hat{j}^{(in0)}\mu_2(x_2) \right] \right] \right\} (S9a)

- \delta(c - c_2)\delta(c - c_1) \left[ \tilde{\delta}^{\mu}_{\mu_2} \delta^3(x - x_2) \left[ \hat{j}^{(in0)}\mu_0(x), \hat{j}^{(in0)}\mu_1(x_1) \right] \right] (S9b)

+ \theta(c - c_1)\delta(c_1 - c_2) \left[ \delta^3(x_1 - x_2) \left[ \hat{j}^{(in0)}\mu_0(x), \hat{j}^{(in0)}\mu_1(x_1) \right] \right] \right\} (S9c)

+ \left( \frac{1}{\hbar c^2} \right)^2 \left\{ (-\delta(c - c_1)\theta(c_1 - c_2) + \theta(c - c_1)\delta(c_1 - c_2)) \left[ \left[ \hat{j}^{(in0)}\mu_0(x), \hat{j}^{(in0)}\mu_0(x_1) \right], \hat{j}^{(in0)}\mu_2(x_2) \right] \right\} (S9d)

- \theta(c - c_2)\delta(c_2 - c_1) \left[ \left[ \hat{j}^{(in0)}\mu_0(x), \hat{j}^{(in0)}\mu_2(x_2) \right], \hat{j}^{(in0)}\mu_0(x_1) \right] \right\}. (S9e)$$
Replacing the fifth term (S9e) using the next Jacobi identity:

\[
\begin{bmatrix}
\dot{j}^{(in0)} (x), \dot{j}^{(in0)} \mu_0 (x_2)
\end{bmatrix}, \dot{j}^{(in0)} \mu_0 (x_1)
\]

\[
= - \left[ \left[ \dot{j}^{(in0)} \mu_0 (x_2), \dot{j}^{(in0)} 0 (x_1) \right], \dot{j}^{(in0)} \mu_0 (x) \right] - \left[ \left[ \dot{j}^{(in0)} 0 (x_1), \dot{j}^{(in0)} \mu_0 (x) \right], \dot{j}^{(in0)} \mu_2 (x_2) \right]
\]

then, the first term in the right hand side of Eq.(S10) with Eq.(S2) offsets the term (S9c), and the second term in the right hand side of Eq.(S10) offsets the second term in (S9d).

The first term in (S9d) offsets the first term, (S9a), considering the commutation relation at the simultaneous time, \( \delta(ct - ct) \left[ \dot{j}^{(in0)} \mu_0 (x), \dot{j}^{(in0)} 0 (x_1) \right] \) and Eq.(S2) (remark the change of upper or lower subscript). The second term (S9b) vanishes by means of Eq.(S2).

As a result, the second order nonlinear single susceptibility operator Eq.(41) in the main text maintains the gauge invariance, Eq.(25) in the main text.

S 3. Third order nonlinear single susceptibility: Eq.(42) in the main text

With respect to the third order nonlinear single susceptibility, let us check the charge conservation law. Operating \( \partial_\mu \) to Eq.(42) in the main text,

\[
\partial_\mu 3! \hat{\chi}_{\mu_1 \mu_2 \mu_3}(x, x_1, x_2, x_3) = \frac{1}{i\hbar c^2} \left( \frac{-q}{mc^2} \right)^2 
\]

\[
\left\{ \theta(ct - ct) \delta(ct - ct) \delta(ct_1 - ct_2) \delta(ct_2 - ct_3) \delta_{\mu_2 \mu_3} \delta^3(x_2 - x_3) \delta_{\mu_1 \mu_2} \delta^3(x - x_1) \right\}
\]

\[
\left[ \dot{j}^{(in0)} 0 (x), \dot{j}^{(in0)} 0 (x_2) \right] \] 

(S11a)

\[
+ \theta(ct - ct_1) \theta(ct_1 - ct) \delta(ct_2 - ct_3) \delta_{\mu_2 \mu_3} \delta^3(x_1 - x_2) \delta_{\mu_1 \mu_2} \delta^3(x - x_3) \right\}
\]

(S11b)

\[
\left[ \dot{j}^{(in0)} 0 (x), \dot{j}^{(in0)} 0 (x) \right] \] 

(S11c)

\[
\left\{ \delta(ct - ct_1) \theta(ct_1 - ct) \theta(ct_2 - ct_3) \delta_{\mu_1 \mu_2} \partial_\mu \right] \left[ \dot{j}^{(in0)} 0 (x), \dot{j}^{(in0)} 0 (x_2) \right] \] 

(S11d)

\[
\left\{ \delta(ct - ct_1) \theta(ct_1 - ct) \theta(ct_2 - ct_3) \delta_{\mu_1 \mu_2} \partial_\mu \right] \left[ \dot{j}^{(in0)} 0 (x), \dot{j}^{(in0)} 0 (x_2) \right] \] 

(S11e)

\[
\left\{ \delta(ct - ct_1) \theta(ct_1 - ct) \theta(ct_2 - ct_3) \delta_{\mu_1 \mu_2} \partial_\mu \right] \left[ \dot{j}^{(in0)} 0 (x), \dot{j}^{(in0)} 0 (x) \right] \] 

(S11f)

\[
\left\{ \delta(ct - ct_1) \theta(ct_1 - ct) \theta(ct_2 - ct_3) \delta_{\mu_1 \mu_2} \partial_\mu \right] \left[ \dot{j}^{(in0)} 0 (x), \dot{j}^{(in0)} 0 (x_2) \right] \] 

(S11g)

\[
\left\{ \delta(ct - ct_1) \theta(ct_1 - ct) \theta(ct_2 - ct_3) \delta_{\mu_1 \mu_2} \partial_\mu \right] \left[ \dot{j}^{(in0)} 0 (x), \dot{j}^{(in0)} 0 (x) \right] \] 

(S11h)

\[
\left\{ \delta(ct - ct_1) \theta(ct_1 - ct) \theta(ct_2 - ct_3) \delta_{\mu_1 \mu_2} \partial_\mu \right] \left[ \dot{j}^{(in0)} 0 (x), \dot{j}^{(in0)} 0 (x_2) \right] \] 

(S11i)

\[
\left\{ \delta(ct - ct_1) \theta(ct_1 - ct) \theta(ct_2 - ct_3) \delta_{\mu_1 \mu_2} \partial_\mu \right] \left[ \dot{j}^{(in0)} 0 (x), \dot{j}^{(in0)} 0 (x) \right] \] 

(S11j)
The term (S11p) offsets the term (S11d), applying Eq.(S2) to the most inner commutator in (S11p). In the same manner, the terms (S11q)-(S11u), respectively, offsets the terms (S11e)-(S11i), applying Eq.(S2). The terms (S11j)-(S11l) vanishes, applying Eq.(S2) to the inner commutator at the simultaneous time. The term (S11m) offsets the term (S11a), applying Eq.(S2) to the most inner commutator. In the same manner, the terms (S11n)-(S11o), respectively, offsets (S11b)-(S11c), using Eq.(S2).

As a result, the third order nonlinear single susceptibility operator Eq.(42) in the main text maintains the charge conservation law, Eq.(24) in the main text.

To check the gauge invariance of the third order nonlinear single susceptibility operator, let us operate $\partial^{\mu_1}$ to Eq.(42) in the main text.

$$\partial^{\mu_1} 3! \hat{\chi}^\mu_{\mu_1 \mu_2 \mu_3}(x, x_1, x_2, x_3) =$$

$$\frac{1}{\hbar c^2} \left( \frac{-q}{mc^2} \right)^2 \left\{ \theta (ct-c_2) \delta (ct-c_1) \delta (ct_2 - ct_3) \left[ \delta^{\mu_1} \delta^{\mu_2} \delta^{\mu_3} (x - x_1) \right] \delta_{\mu_1 \mu_2 \mu_3} \delta^3 (x_2 - x_3) \left[ \hat{j}^{(in0)} \delta^0 (x_1), \hat{j}^{(in0)} \delta^0 (x_2) \right] \right\} . \quad (S12a)$$
\begin{align}
\delta(t - ct_1)\theta(t_{c1} - ct_2)\theta(t_{c2} - ct_3) \left( \delta^{\mu}_{\nu_1} \partial^{\mu_1} \delta^3(x - x_1) \right) \left[ \begin{array}{c}
\hat{j}^{(in) 0}_0(x), \hat{j}^{(in) 0}_{\mu_2}(x_2), \hat{j}^{(in) 0}_{\mu_3}(x_3)
\end{array} \right] \\
+ \delta(t - ct_1)\theta(t_{c1} - ct_3)\theta(t_{c3} - ct_2) \left( \delta^{\mu}_{\nu_1} \partial^{\mu_1} \delta^3(x - x_1) \right) \left[ \begin{array}{c}
\hat{j}^{(in) 0}_0(x), \hat{j}^{(in) 0}_{\mu_3}(x_3), \hat{j}^{(in) 0}_{\mu_2}(x_2)
\end{array} \right] \\
- \delta(t - ct_2)\theta(t_{c2} - ct_3)\delta(t_{c3} - ct_1) \delta^{\mu_2}_{\mu_3} \delta^3(x - x_2) \left[ \begin{array}{c}
\hat{j}^{(in) 0}_0(x), \hat{j}^{(in) 0}_{\mu_3}(x_3), \hat{j}^{(in) 0}_0(x_1)
\end{array} \right] \\
+ \left( -\delta(t - ct_2)\delta(t_{c2} - ct_3)\theta(t_{c3} - ct_1) + \delta(t - ct_2)\theta(t_{c2} - ct_3)\delta(t_{c3} - ct_1) \right) \\
+ \left( -\delta(t - ct_3)\delta(t_{c3} - ct_1)\theta(t_{c1} - ct_2) + \delta(t - ct_3)\theta(t_{c3} - ct_1)\delta(t_{c1} - ct_2) \right) \\
+ \left( -\delta(t - ct_3)\delta(t_{c3} - ct_2)\theta(t_{c2} - ct_3) + \delta(t - ct_3)\theta(t_{c3} - ct_2)\delta(t_{c2} - ct_3) \right) \\
+ \left( -\delta(t - ct_1)\theta(t_{c1} - ct_2)\delta(t_{c2} - ct_3) + \theta(t - ct_1)\delta(t_{c1} - ct_2)\theta(t_{c2} - ct_3) \right) \\
\left( \delta^{\mu}_{\nu_3} \delta^3(x - x_3) \left[ \begin{array}{c}
\hat{j}^{(in) 0}_0(x), \hat{j}^{(in) 0}_{\mu_2}(x_2), \hat{j}^{(in) 0}_0(x_1)
\end{array} \right] \right) \\
+ \theta(t - ct_1)\delta(t_{c1} - ct_2)\theta(t_{c2} - ct_3)\delta^{\mu}_{\nu_2} \partial^{\mu_1} \delta^3(x - x_2) \left[ \begin{array}{c}
\hat{j}^{(in) 0}_0(x), \hat{j}^{(in) 0}_{\mu_3}(x_3), \hat{j}^{(in) 0}_0(x_1)
\end{array} \right] \\
- \theta(t - ct_2)\delta(t_{c2} - ct_3)\delta(t_{c3} - ct_1) \delta^{\mu_2}_{\mu_3} \partial^{\mu_1} \delta^3(x - x_2) \left[ \begin{array}{c}
\hat{j}^{(in) 0}_0(x), \hat{j}^{(in) 0}_{\mu_3}(x_3), \hat{j}^{(in) 0}_0(x_1)
\end{array} \right] \\
+ \theta(t - ct_3)\delta(t_{c3} - ct_1)\theta(t_{c1} - ct_2) \delta^{\mu_2}_{\mu_3} \partial^{\mu_1} \delta^3(x - x_1) \left[ \begin{array}{c}
\hat{j}^{(in) 0}_0(x), \hat{j}^{(in) 0}_{\mu_3}(x_3), \hat{j}^{(in) 0}_0(x_1)
\end{array} \right] \\
+ \theta(t - ct_3)\delta(t_{c3} - ct_2)\delta(t_{c2} - ct_3) \delta^{\mu_2}_{\mu_3} \partial^{\mu_1} \delta^3(x - x_2) \left[ \begin{array}{c}
\hat{j}^{(in) 0}_0(x), \hat{j}^{(in) 0}_{\mu_3}(x_3), \hat{j}^{(in) 0}_0(x_1)
\end{array} \right] \\
+ \theta(t - ct_2)\theta(t_{c2} - ct_3)\delta^{\mu}_{\nu_2} \partial^{\mu_1} \delta^3(x - x_1) \left[ \begin{array}{c}
\hat{j}^{(in) 0}_0(x), \hat{j}^{(in) 0}_{\mu_3}(x_3), \hat{j}^{(in) 0}_0(x_1)
\end{array} \right] \\
+ \theta(t - ct_3)\delta(t_{c3} - ct_1)\theta(t_{c1} - ct_2) \delta^{\mu}_{\nu_3} \partial^{\mu_1} \delta^3(x - x_1) \left[ \begin{array}{c}
\hat{j}^{(in) 0}_0(x), \hat{j}^{(in) 0}_{\mu_3}(x_3), \hat{j}^{(in) 0}_0(x_1)
\end{array} \right] \\
+ \left( \frac{1}{\hbar \omega_0} \right)^3
\end{align}
In the following, we prove the next equation, which leads to the gauge invariance.

\[
(S12p) + (S12d) + (S12j) + (S12s) + (S12r) + (S12n) = 0, \quad (S13)
\]

\[
(S12q) + (S12e) + (S12l) + (S12t) + (S12u) + (S12o) = 0, \quad (S14)
\]

\[
(S12m) + (S12a) + (S12k) = 0, \quad (S15)
\]

\[
(S12i) + (S12c) + (S12h) = 0, \quad (S16)
\]

\[
(S12f) + (S12b) + (S12g) = 0. \quad (S17)
\]

1. Eq.(S13): The first term of \((S12p)\) offsets \((S12d)\), using Eq.(S2). To the inner double commutator in the second term of \((S12p)\), the nest Jacobi identity is applied:

\[
\left[ \left[ \hat{j}^{(in0)} \mu (x), \hat{j}^{(in0)}_0 (x_1) \right], \hat{j}^{(in0)} \mu_2 (x_2) \right], \hat{j}^{(in0)}_3 (x_3) \quad (S18)
\]

\[
= - \left[ \left[ \hat{j}^{(in0)}_0 (x_1), \hat{j}^{(in0)} \mu_2 (x_2) \right], \hat{j}^{(in0)} \mu (x) \right], \hat{j}^{(in0)}_3 (x_3) \]

\[
- \left[ \left[ \hat{j}^{(in0)} \mu_2 (x_2), \hat{j}^{(in0)} \mu (x) \right], \hat{j}^{(in0)}_0 (x_1) \right], \hat{j}^{(in0)}_3 (x_3) \right].
\]

Furthermore, the inner commutator (, assuming at the simultaneous time) in the first term of Eq.(S18), one may apply Eq.(S2). The part including this factor in the second term of \((S12p)\) offsets the term \((S12j)\). In the second term of \((S12p)\), the part including the second term of Eq.(S18) offsets the first term of \((S12s)\). Up to now, \((S12p) + (S12d) + (S12j) + \text{the first term of } (S12s) = 0\) has been shown.

Next, to the outer double commutator in the second term of \((S12r)\), let us apply the next Jacobi identity,

\[
\left[ \left[ \hat{j}^{(in0)} \mu (x), \hat{j}^{(in0)} \mu_2 (x_2) \right], \hat{j}^{(in0)} \mu_3 (x_3) \right], \hat{j}^{(in0)}_0 (x_1) \quad (S19)
\]

\[
= - \left[ \left[ \hat{j}^{(in0)} \mu_3 (x_3), \hat{j}^{(in0)}_0 (x_1) \right], \left[ \hat{j}^{(in0)} \mu (x), \hat{j}^{(in0)} \mu_2 (x_2) \right] \right]
\]

\[
+ \left[ \left[ \hat{j}^{(in0)} \mu (x), \hat{j}^{(in0)} \mu_2 (x_2) \right], \hat{j}^{(in0)}_0 (x_1) \right], \hat{j}^{(in0)}_3 (x_3) \right].
\]

The commutator in the first term of Eq.(S19): \(\left[ \hat{j}^{(in0)} \mu_3 (x_3), \hat{j}^{(in0)}_0 (x_1) \right]\) is the commutator at the simultaneous time, therefore, we may use Eq.(S2). The part including this factor in \((S12r)\) offsets the term \((S12n)\). In \((S12r)\), the part including the second term of Eq.(S19) offsets the second term of \((S12s)\). Up to now, \((S12r) + (S12n) + \text{the second term of } (S12s) = 0\) is shown.

Together with the previous result, Eq.(S13) holds.
2. **Eq.(S14):** This equation is Eq.(S13) with the replacement \( x_2 \leftrightarrow x_3 \) and \( \mu_2 \leftrightarrow \mu_3 \), therefore, Eq.(S14) holds.

3. **Eq.(S15):** The first term of (S12m) offsets (S12a), using Eq.(S2).

   To the double commutator in the second term of (S12m), the nest Jacobi identity is applied:
   \[
   \left[ \left[ \hat{j}^{(in0)} \mu(x), \hat{j}^{(in0)}_0(x_1) \right], \hat{j}^{(in0)}_0(x_2) \right] = - \left[ \left[ \hat{j}^{(in0)}_0(x_1), \hat{j}^{(in0)}_0(x_2) \right], \hat{j}^{(in0)}(x) \right] - \left[ \left[ \hat{j}^{(in0)}_0(x_2), \hat{j}^{(in0)}(x) \right], \hat{j}^{(in0)}_0(x_1) \right] \tag{S20}
   \]

   In the above, we use the inner commutator (at the simultaneous time) in the first term of the second hand side becomes zero, using Eq.(S2). The second term of (S12m) includes the factor of Eq.(S20) and offsets (S12k).

   As a result, Eq.(S15) holds.

4. **Eq.(S16):** To the double commutator in (S12i), we apply the next Jacobi identity:
   \[
   \left[ \left[ \hat{j}^{(in0)}_0(x), \hat{j}^{(in0)}_0(x_1) \right], \hat{j}^{(in0)}_0(x_2) \right] = - \left[ \left[ \hat{j}^{(in0)}_0(x_1), \hat{j}^{(in0)}_0(x_2) \right], \hat{j}^{(in0)}(x) \right] - \left[ \left[ \hat{j}^{(in0)}_0(x_2), \hat{j}^{(in0)}(x) \right], \hat{j}^{(in0)}_0(x_1) \right] \tag{S21}
   \]

   The inner commutator in the first term of Eq.(S21) is the commutator at the simultaneous time, therefore, we may use Eq.(S2). The part including this factor in (S12i) offsets the term (S12c). In (S12i), the part including the second term of Eq.(S21) offsets the second term of (S12h). The first term of (S12h) is zero, because the inner commutator included in this term is commutator at the simultaneous time and leads to zero, using Eq.(S2).

   Therefore, Eq.(S16) holds.

5. **Eq.(S17):** This equation is Eq.(S16) with the replacement \( x_2 \leftrightarrow x_3 \) and \( \mu_2 \leftrightarrow \mu_3 \), therefore, Eq.(S17) holds.

As the summery, Eqs.(S13)-(S17) hold and the third order nonlinear single susceptibility Eq.(42) in the main text maintains the gauge invariance Eq.(25) in the main text.