

# Time-dependent adiabatic $GW$

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May 24, 2023

## 1 From Keldysh formalism to quantum master equation

Existing researches [1] utilizing TD-aGW carry out the calculation within the framework of quantum master equation (QME)

$$\dot{\rho} + i[\rho, H] = \int_{-\infty}^t dt' F[\rho(t')] \quad (1)$$

which is the equation of motion of the reduced single-electron density matrix

$$\rho(\mathbf{r}, \mathbf{r}', t) = \langle \psi^\dagger(\mathbf{r}') \psi(\mathbf{r}) \rangle = -iG^<(\mathbf{r}, t; \mathbf{r}', t). \quad (2)$$

QME can be derived heuristically when the system contains well-defined quasiparticles. The generality of QME and the flow from the most generalized Keldysh formalism to quantum kinetic theories is discussed in [2, 3, 4, 5]. In this section we give a brief summary of the derivation and coverage as well as possible simplifications of QME.

In Keldysh formalism we still have the Dyson equation

$$G = G_0 + G_0 \Sigma G = G_0 + G \Sigma G_0, \quad (3)$$

although now  $G$ ,  $G_0$  and  $\Sigma$  are  $2 \times 2$  matrices due to the need of anti-time-ordered, lesser and greater Green functions besides the time-ordered Green function in deriving a diagrammatic perturbation theory for non-equilibrium Green functions [6]; alternatively they can be defined on the Keldysh contour [7]. Note that the choice of  $H_0$  has some arbitrariness at this point: we may split the full many-body self-energy correction  $\Sigma^{\text{full}}$  into an “effective field” which has no retardation and is inserted into  $H_0$  and a “real” interaction part  $\Sigma$ . By Lengreth’s rules [7], we reduce (3) on the Keldysh contour into the following equation on the physical time axis:

$$G^< = G_0^< + G_0^R \Sigma^R G_0^< + G^R \Sigma^< G_0^A + G^< \Sigma^A G^A, \quad G^{A,R} = G_0^{A,R} + G_0^{A,R} \Sigma^{A,R} G^{A,R}. \quad (4)$$

By applying

$$\overleftarrow{G}_0^{-1} = -i\overleftarrow{\partial}_t - H_0, \quad \overrightarrow{G}_0^{-1} = i\overrightarrow{\partial}_t - H_0 \quad (5)$$

to the left and right hand sides of the equation about  $G^<$  and utilizing the fact that  $G_0^< \overleftarrow{G}_0^{-1} = \overrightarrow{G}_0^{-1} G_0^< = 0$ , we obtain

$$i\partial_T G^< - [H_0, G^<] = \Sigma^R G^< + \Sigma^< G^A - G^R \Sigma^< - G^< \Sigma^A, \quad (6)$$

where we have redefined the time variables as

$$T = \frac{t_1 + t_2}{2}, \quad t = t_1 - t_2. \quad (7)$$

Roughly speaking,  $T$  is related to the time variance of the external driving field or the relaxation of the system, while  $t$  corresponds to the time variable in equilibrium Green function, and often Fourier transform is done in  $t$ , and thus  $G^< = G^<(T, \omega)$ .

The “precursor” equation (6) is already very close to QME, its left hand side being the left hand side of QME after integrating over  $\omega$ . Three approximations are needed however to close the equation: the explicit form of  $\Sigma[G]$ , representation of  $G^{A,R}$  with  $G^<$  and  $\Sigma^{A,R}$ , and reconstruction of  $G^<$  using  $\rho(T)$  i.e.  $G^<(t_1 = t_2 = T)$ . It can be proved that the last is always possible [3], with the lowest order approximation being the Generalized Kadanoff–Baym Ansatz, and a closed QME can therefore be obtained [4, 7]. Note that in principle this procedure does not require the system to have well-defined electron-like quasiparticle.

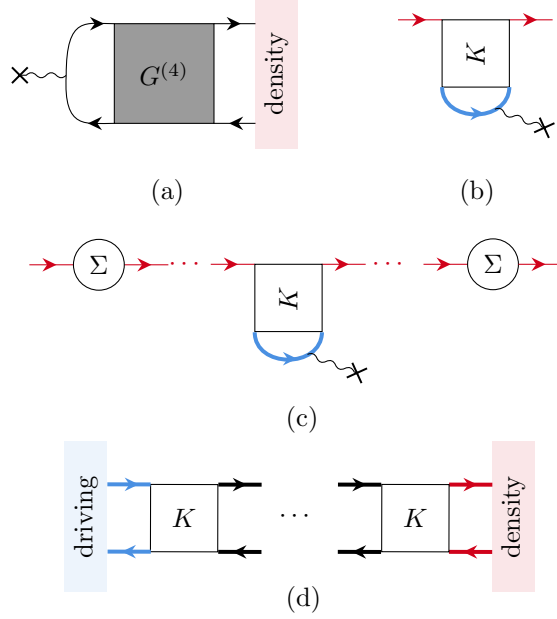


Figure 1: Feynman diagrams representing linear response. (a) Linear response of electron density (red rectangle) is given by plating a line representing the external field into one bare propagator. (b) Single-electron self-energy, rewritten as a convolution between the four-point kernel function (i.e. self-energy of renormalized electron-hole pair) and the renormalized Green function (thick blue line) with the influence of the driving field. The electron lines that are connected to the external legs in (a) are colored red. Note that the blue line has its own self-energy correction, and the external field line is connected to one bare propagator within the thick blue line. (c) Inserting (b) into (a). (d) The diagram we get after evaluating the derivative with respect to the external field i.e. after removing the external field line. Since the thick blue line in (b) and (c) has its own self-energy correction which is also represented by (b), we get accurately diagrams representing renormalized electron-hole pair; note that  $K$  can be obtained by removing a renormalized Green function from  $\Sigma$ .

QME can be used to study higher-order correlations in a system without explicitly constructing the relevant Green functions. The linear response, for example, reveals  $G^{(4)}$ , since the first order response can be diagrammatically obtained by choosing an arbitrary propagator and plant an external field line into it (Figure 1 (a)). Since the single-electron self-energy can always be written into the convolution of the two-electron kernel function and the renormalized Green function (Figure 1 (b)), the latter containing its own self-energy correction and hence a two-electron kernel function, by taking the linear response i.e. breaking a propagator in a single-electron self-energy, we get a two-electron kernel function with the same level of accuracy (Figure 1 (c, d)).

Two further approximations reduce QME to quantum Boltzmann equation (QBE) [2, 7]. Under the assumption that well-defined quasiparticles always exist and hence we have

$$G^<(\mathbf{X}, \mathbf{p}, T, \omega) = 2\pi\delta(\omega - E_{\mathbf{p}}(\mathbf{X}))f(\mathbf{X}, \mathbf{p}, T), \quad (8)$$

where Wigner transform has been applied. Using the condition of smooth external driving force, by gradient expansion, the left hand side of QME becomes the usual form of collisionless Boltzmann equation. With both the quasiparticle assumption the right hand side of QME becomes the Fermi golden rule, which is prevalently used as the collision integral [8]. Deviations from the quasiparticle approximation, e.g. strong electron-phonon interaction-induced renormalization, may still lead to QBE, although with modified collision integral [2, 9].

## 2 The COHSEX approximation

### References

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