A diagrammer’s note on DC conductivity of anisotropic Fermi liquids for beginners: Maebashi-Fukuyama formula and Taylor formula

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Abstract

Formal but exact DC conductivity formulae for anisotropic Fermi liquids are reviewed. One is the Maebashi-Fukuyama formula based on the Fermi-surface harmonics. The other is the Taylor formula based on the scattering eigenfunction. In comparison to these two formulae the current-vertex-correction in the fluctuation-exchange approximation is shown to be a bad vision caused by an inconsistent approximation.

1 Introduction

The DC conductivity formula obtained by the Fermi-liquid theory is shown to be equivalent to the one obtained by the linearized Boltzmann equation in the textbook [1]. This proof is done for the case of spherical Fermi surface. Its extension to the case of anisotropic Fermi surface is done by Maebashi and Fukuyama [2]. The Maebashi-Fukuyama formula is formal but exact.

On the other hand, Taylor [3] derived a formal but exact solution of the linearized Boltzmann equation.

These two works employ opposite strategies.

The Maebashi and Fukuyama formula is on the basis of the Fermi-surface harmonics which is the polynomial of the velocity. Then the interaction is expanded in terms of the velocity.

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The Taylor formula is on the basis of the scattering eigenfunction. In this formula the interaction is diagonalized but the velocity is expanded in terms of the scattering eigenfunction.

By these two formulae we know the exact forms of the DC conductivity of anisotropic Fermi liquids. Consequently it becomes evident that there is no room for the current-vertex-correction discussed in the fluctuation-exchange (FLEX) approximation [4].

2 Maebashi-Fukuyama formula

2-1 Fermi-surface harmonics expansion

We start from the Fermi-liquid formula [2] for the DC conductivity

\[
\sigma_{xx} = 2e^2 \sum_p \sum_{p'} v_x(p)(K''_{pp'})^{-1} v_x(p') \left( -\frac{\partial f_p}{\partial \epsilon_p} \right),
\]

which is (2.34) of [2] and equivalent to the result of the linearized Boltzmann equation, (58) of [5].

Using the Fermi-surface harmonics [6, 7], \(\psi_L(p)\), (1) is written as

\[
\sigma_{xx} = 2e^2 \sum_p \sum_{p'} \psi_1(p) \sum_L \sum_{L'} (K''_{LL'})^{-1} \psi_L(p) \psi_{L'}(p') \psi_1(p') \langle v_x^2 \rangle \left( -\frac{\partial f_p}{\partial \epsilon_p} \right),
\]

where \(\psi_1(p) = v_x(p)/\sqrt{\langle v_x^2 \rangle}\) and

\[
(K'')^{-1}_{pp'} = \sum_L \sum_{L'} (K''_{LL'})^{-1} \psi_L(p) \psi_{L'}(p').
\]

The normalization factor is defined as

\[
\langle v_x^2 \rangle = \frac{1}{N(\epsilon)} \sum_p v_x(p)^2 \delta(\epsilon - \epsilon_p),
\]

with

\[
N(\epsilon) = \sum_p \delta(\epsilon - \epsilon_p).
\]

By the orthonormal condition

\[
\frac{1}{N(\epsilon)} \sum_p \psi_L(p) \psi_{L'}(p) \delta(\epsilon - \epsilon_p) = \delta_{LL'},
\]
(2) is written as

\[ \sigma_{xx} = 2e^2 \int d\epsilon N(\epsilon) \int d\epsilon' N(\epsilon') (K''_{11})^{-1} \langle v_x^2 \rangle \left(-\frac{\partial f}{\partial \epsilon}\right). \]  

Here we have used the relation

\[ \sum_p = \int d\epsilon \sum_p \delta(\epsilon - \epsilon_p). \]  

Introducing the memory-function matrix

\[ (M'')^{-1}_{11} = \int d\epsilon' N(\epsilon') (K''_{11})^{-1}, \]  

we obtain

\[ \sigma_{xx} = 2e^2 \int d\epsilon N(\epsilon) (M'')^{-1}_{11} \langle v_x^2 \rangle \left(-\frac{\partial f}{\partial \epsilon}\right). \]  

Finally (11) is written as

\[ \sigma_{xx} = 2e^2 \sum_p v_x(p) (M'')^{-1}_{11} v_x(p) \left(-\frac{\partial f_p}{\partial \epsilon_p}\right). \]  

In conclusion the DC conductivity is determined by the memory-function matrix [2].

The microscopic calculation of the memory-function matrix based on the Fermi-liquid theory is carried out by Maebashi and Fukuyama [2, 8]. To obtain a finite DC conductivity we have to take the contribution of the Umklapp process into the calculation of \( M'' \). Next we have to perform the matrix inversion from \( M'' \) to \( (M'')^{-1} \).

2-2 Collision term

In the previous subsection it is clarified that the DC conductivity is determined by the memory-function matrix. In the Boltzmann equation it appears in the collision term. Thus we investigate the collision term after Allen [6] to obtain the feeling of the matrix structure.

The collision term \( I_p \) for \( n_p \) is given by

\[ I_p = \sum_{p'} K''_{pp'} \left(n_{p'} - n_p\right), \]  

as (47) of [5]. Using the Fermi-surface harmonics (12) is written as

\[ I_p = \sum_{p'} \sum_L \sum_{L'} K''_{LL'} \psi_L(p) \psi_{L'}(p') \sum_{L''} n_{L''} \left[\psi_{L''}^*(p') - \psi_{L''}^*(p)\right]. \]  

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By the orthonormal condition (6) and \( \psi_0(p) = 1 \), (13) is written as

\[
I_p = \int d\epsilon' N(\epsilon') \sum_L \sum_{L'} \sum_{L''} K_{LL'}^{''} n_{L''} \psi_L(p) \left\{ \delta_{L'L''} - \delta_{L'0} \psi_{L''}(p) \right\}.
\]

Introducing the memory-function matrix

\[
M_{LL'}^{''} = \int d\epsilon' N(\epsilon') K_{LL'}^{''},
\]

(14) is written as

\[
I_p = \sum_L \psi_L(p) \left\{ \sum_{L'} M_{LL'}^{''} n_{L'} - \sum_{L''} \sum_{L'} C_{L''L'L} M_{L''0}^{''} n_{L'} \right\},
\]

where \( C_{L''L'L} \) is the Clebsch-Gordan coefficient defined by

\[
\psi_L(p) \psi_{L'}(p) = \sum_{L''} C_{L''L'L} \psi_{L''}(p).
\]

Thus we obtain the collision term \( I_L \) for \( n_L \) as

\[
I_L = \sum_{L'} M_{LL'}^{''} n_{L'} - \sum_{L''} \sum_{L'} C_{L''L'L} M_{L''0}^{''} n_{L'}.
\]

For anisotropic Fermi surfaces \( n_L \) couples to the other \( n_{L'} \)s via the memory-function matrix so that we have to perform the matrix inversion, \( M'' \rightarrow (M'')^{-1} \), to obtain the conductivity.

The case of the spherical Fermi surface is special and we do not have to perform the matrix inversion. By the symmetry \( K_{pp'}^{''} \) only depends on the angle between \( p \) and \( p' \) so that the memory-function matrix becomes a diagonal matrix. Since \( C_{0L'L'}^{''} = \delta_{L'L'} \) which is obtained from (17) by using \( \psi_0(p) = 1 \), we obtain

\[
I_L = 4\pi \left\{ w_L - w_0 \right\} n_L,
\]

where we put \( M_{LL'}^{''} = 4\pi w_L \delta_{LL'} \). This is the result shown in [9]. The DC conductivity is determined by \( n_1 \) and the transport life-time \( \tau \) is given by

\[
\frac{1}{\tau} = 4\pi \left\{ w_0 - w_1 \right\}.
\]

2-3 Vector mean free path
The use of the vector mean free path $\Lambda$ is convenient \[3, 10\] for the discussion of the symmetry property of the conductivity tensor. In the following all the vectors are defined on the Fermi surface. The conductivity tensor $\sigma$ is expressed in the dyadic form as \[2, 8, 3, 10\]

$$\sigma = 2e^2 \sum_p v_p \Lambda_p \left( -\frac{\partial f_p}{\partial \epsilon_p} \right),$$

(21)

where

$$\Lambda_p = \sum_{p'} (K'')^{-1}_{pp'} v_{p'},$$

(22)

with $v_p = \left( v_x(p), v_y(p), v_z(p) \right)$.

For systems with cubic symmetry the conductivity tensor reduces to a scalar $\sigma$ \[3\]. This means the integrand in (21) is proportional to the scalar product $v_p \cdot v_p$. Thus introducing the anisotropic life time $\tau_p$, the vector mean free path is written as

$$\Lambda_p = \tau_p v_p.$$  

(23)

Namely, the direction of $\Lambda_p$ is the same as that of $v_p$. Substituting (23) into (22) we obtain

$$\tau_p = \sum_{p'} (K'')^{-1}_{pp'} \frac{v_p \cdot v_{p'}}{v_p \cdot v_p},$$

(24)

This relation (24) is a mere rewriting and has no practical use. Using this anisotropic life time the conductivity is expressed as

$$\sigma_{xx} = 2e^2 \sum_p v_x(p) \tau_p v_x(p) \left( -\frac{\partial f_p}{\partial \epsilon_p} \right).$$

(25)

The Fermi-liquid formula \[11\] by Yamada and Yosida corresponds to this case. The appearance of (25) is simple but it is none other than (11). Actually

$$\sigma_{xx} = 2e^2 \sum_p \frac{v_p \cdot v_p}{3} \tau_p \left( -\frac{\partial f_p}{\partial \epsilon_p} \right) = 2e^2 \sum_p \sum_{p'} (K'')^{-1}_{pp'} \frac{v_p \cdot v_{p'}}{3} \left( -\frac{\partial f_p}{\partial \epsilon_p} \right),$$

(26)

in the case of cubic symmetry.

For systems without cubic symmetry the proportional relation (23) does not hold \[3, 10\] so that we have to use the Maebashi-Fukuyama formula discussed above.
For systems with spherical symmetry $\tau_p$ reduces to a constant $\tau$. In this case the inverse relation of (22) \( v_p = \sum_{p'} K''_{pp'} \Lambda_{p'} \), (27) reduces to \( v_p = \tau \sum_{p'} K''_{pp'} v_{p'} \). (28) Thus we obtain \[
\frac{1}{\tau} = \sum_{p'} K''_{pp'} (\hat{p} \cdot \hat{p}') ,
\] (29) with $v_p = p/m$ and $\hat{p} = p/|p|$. Since \[
K''_{pp'} = \frac{1}{\tau_0} \delta_{pp'} - W_{pp'} ,
\] (30) and \[
\frac{1}{\tau_0} = \sum_{p'} W_{pp'} ,
\] (31) (29) is written as \[
\frac{1}{\tau} = \sum_{p'} W_{pp'} (1 - \cos \theta) ,
\] (32) with $\cos \theta = \hat{p} \cdot \hat{p}'$. This expression (32) is none other than (20).

3 Taylor formula

3-1 Setup

Taylor [3] introduced the vector mean free path $\Lambda$ and wrote down the DC conductivity tensor $\sigma$ in the dyadic form as
\[
\sigma = 2e^2 \sum_p v(p) \Lambda(p) \left( -\frac{\partial f_p}{\partial \epsilon_p} \right) .
\] (33)
In the following all the vectors are defined on the Fermi surface. The vector mean free path is determined by
\[
v(p) = \sum_{p'} Q(p, p') [\Lambda(p) - \Lambda(p')] .
\] (34)
The Fermi-liquid theory works [2] on the calculation of the collision operator $Q$. Hereafter we only consider the $x$-component of \( (34) \)

$$v_x(p) = \sum_{p'} Q(p, p')[\Lambda_x(p) - \Lambda_x(p')]$$ \hspace{1cm} (35)

Introducing the inner product

$$\langle a, b \rangle \equiv \sum_p \left( - \frac{\partial f_p}{\partial \epsilon_p} \right) a(p)b(p),$$ \hspace{1cm} (36)

the DC conductivity is compactly written as \[1\] \( (37) \)

$$\sigma_{xx} = 2e^2 \langle v_x, \Lambda_x \rangle.$$ \hspace{1cm} (37)

### 3-2 Loose explanation

Loosely \( (35) \) is written as

$$v_x = \left( \frac{1}{\tau} - \hat{Q} \right) \Lambda_x,$$ \hspace{1cm} (38)

where

$$\hat{Q} \Lambda_x \equiv \sum_{p'} Q(p, p') \Lambda_x(p'),$$ \hspace{1cm} (39)

and

$$\frac{1}{\tau} \Lambda_x \equiv \sum_{p'} \frac{1}{\tau(p)} \delta_{pp'} \Lambda_x(p'),$$ \hspace{1cm} (40)

with

$$\frac{1}{\tau(p)} \equiv \sum_{p'} Q(p, p').$$ \hspace{1cm} (41)

The inverse of \( (38) \) is

$$\Lambda_x = \left( \frac{1}{\tau} - \hat{Q} \right)^{-1} v_x = \tau v_x + \tau \hat{Q} \tau v_x + \tau \hat{Q} \tau \hat{Q} \tau v_x + \cdots.$$ \hspace{1cm} (42)

The perturbative expansion for the DC conductivity is shown in Fig. 1.

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1. See, for example, Ashcroft and Mermin: Solid State Physics (1976) Chap. 16 - PROBLEM 4.
If we can find the scattering eigenfunction $a_s(p)$ which satisfies
\[
\tau(p) \sum_{p'} Q(p, p') a_s(p') = \alpha_s a_s(p), \tag{43}
\]
with the eigenvalue $\alpha_s$, the inversion (42) is readily accomplished. To proceed we expand the velocity in terms of the scattering eigenfunction as
\[
\tau(p)v_x(p) = \sum_s \beta_s a_s(p). \tag{44}
\]
Employing (43) and (44) in (42) we obtain
\[
2 \Lambda_x(p) = \sum_s \left[ 1 + \alpha_s + \alpha_s^2 + \cdots \right] \beta_s a_s(p), \tag{45}
\]
To proceed further we have to establish the orthonormality of the scattering eigenfunction.

3-3 Tight explanation
First we rewrite (43) as
\[
\sum_{p'} \sqrt{\tau(p)} Q(p, p') \sqrt{\tau(p')} a_s(p') \sqrt{\tau(p')} = \alpha_s \frac{a_s(p)}{\sqrt{\tau(p)}}, \tag{46}
\]
Since $\sqrt{\tau(p)} Q(p, p') \sqrt{\tau(p')}$ is a Hermitian matrix, the function $a_s(p)/\sqrt{\tau(p)}$ constitutes an orthogonal basis. Thus the orthonormal condition is written as
\[
\left( \frac{a_s}{\sqrt{\tau}}, \frac{a_{s'}}{\sqrt{\tau}} \right) = \delta_{ss'}. \tag{47}
\]

$\tau \hat{Q} v_x = \tau(p) \sum_{p'} Q(p, p') \sum_s \beta_s a_s(p') = \sum_s \beta_s a_s a_s(p)$, and so on.

$^3$ $Q(p, p')$ itself is a Hermitian matrix.
Employing (44) and (45) in (37) we obtain

$$\sigma_{xx} = 2e^2 \sum_{s'} \sum_s \beta_{s'} \frac{1}{1 - \alpha_s} \beta_s \left( \frac{a_{s'}}{\tau}, a_s \right).$$

(48)

Finally making use of the orthonormal condition (47) we obtain the Taylor formula for the DC conductivity

$$\sigma_{xx} = 2e^2 \sum_s \beta_s \frac{1}{1 - \alpha_s} \beta_s.$$  

(49)

The repeated scatterings shown in Fig. 1 result in the renormalization of the life time $\tau$.

4 Current-vertex-correction in FLEX approximation

The purpose of this note is to clarify the exact expression for the DC conductivity of anisotropic Fermi liquids.

Consequently it becomes evident that there is no room for the current-vertex-correction discussed in the FLEX approximation [4].

Other deficiencies of the FLEX approximation have been also discussed previously [12, 13].

References

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