

Correlation effects on the third-frequency-moment sum rule of electron liquids

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The exact form of the third-frequency-moment sum rule is used to obtain the long-wavelength limit of the dynamic local-field correction of electron liquids within the formalism of Gross and Kohn. The effects of correlation modify the high-frequency behavior of the local-field correction typically by a factor 2–3 at metallic densities.

Recently, Gross and Kohn¹ extended the density-functional formalism to treat the dynamic response of electron liquids within the local-density approximation. They constructed a linear density-density response function in such a way that it satisfied some exact conditions, which include the compressibility sum rule and the third-frequency-moment ($\langle\omega^3\rangle$) sum rule. In the application of the latter sum rule, however, they used an approximate form in which correlation effects were partially neglected.² The purpose of the present Brief Report is to obtain the frequency-dependent local-field correction in the long-wavelength limit by using the exact form of the $\langle\omega^3\rangle$ sum rule.³

The $\langle\omega^3\rangle$ sum rule⁴ gives a constraint on the asymptotic form of the local-field correction $G(q, \omega)$ in the long-wavelength limit,⁵

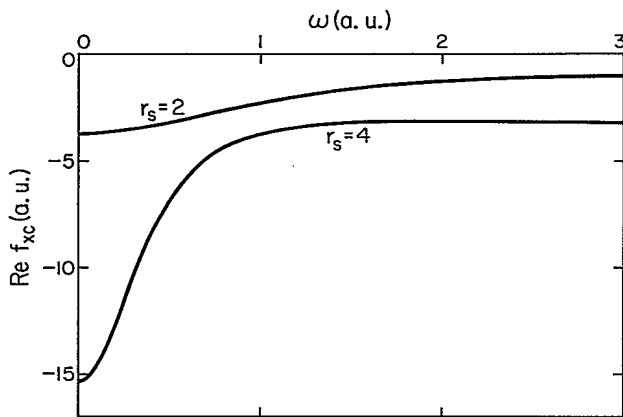
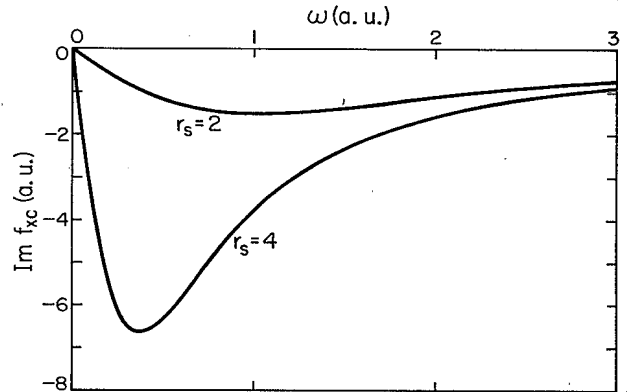
$$\lim_{q \rightarrow 0} G(q, \infty) = -\frac{q^2}{2\pi n e^2} \left(\frac{2}{15} \langle V \rangle + \langle E_{\text{kin}} \rangle - \langle E_{\text{kin}} \rangle_0 \right), \quad (1)$$

where $\langle V \rangle$ and $\langle E_{\text{kin}} \rangle$ are the average potential and kinetic energies per particle, respectively, and $\langle E_{\text{kin}} \rangle_0$ is the

noninteracting Fermi-gas value of the latter. Using the virial forms for $\langle V \rangle$ and $\langle E_{\text{kin}} \rangle$ one obtains

$$\lim_{q \rightarrow 0} f_{\text{xc}}^h(q, \infty) = -\frac{4}{5} n^{2/3} \frac{d}{dn} \left[\frac{\epsilon_{\text{xc}}}{n^{2/3}} \right] + 6n^{1/3} \frac{d}{dn} \left[\frac{\epsilon_{\text{xc}}}{n^{1/3}} \right], \quad (2)$$

where $f_{\text{xc}}^h(q, \omega) = -(4\pi/q^2)G(q, \omega)$, and ϵ_{xc} is the exchange-correlation energy per particle. Neglecting the second term on the right-hand side of Eq. (2), which corresponds to neglecting the difference between $\langle E_{\text{kin}} \rangle$ and $\langle E_{\text{kin}} \rangle_0$ in Eq. (1), one recovers the expression $f_{\infty}(n)$ of Ref. 1. It should be noted that the second term of Eq. (2) is entirely due to correlation; the exchange part, $\epsilon_{\text{x}} = -\frac{3}{4}(3/\pi)^{1/3}n^{1/3}$, does not contribute. If the second term of Eq. (2) is included in the parametrization (14) of Ref. 1, the coefficients $a(n)$ and $b(n)$ are considerably changed. In place of Figs. 1 and 2 of Ref. 1, one now obtains Figs. 1 and 2 below. At $\omega=0$, there is no difference. On the other hand, the high-frequency limits of $\text{Re } f_{\text{xc}}^h$ are reduced by factors of 2.6 and 3.2 for $r_s=2$ and 4, respectively, leading to a much stronger frequency dependence of f_{xc}^h than in the approximation of Ref. 1. The

FIG. 1. Real part of the parametrization for $f_{\text{xc}}^h(q=0, \omega)$.FIG. 2. Imaginary part of the parametrization for $f_{\text{xc}}^h(q=0, \omega)$.

difference between $f_{xc}(r, \omega=0)$, as used in the original paper by Zangwill and Soven,⁶ and $f_{xc}(r, \bar{\omega})$ at an appropriate frequency $\bar{\omega}$, is now estimated to be about 2% to 6%.

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