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## Anisotropic Fermion Gas

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An energy term discovered by Kohn and Luttinger is shown to have a simple geometrical meaning.

KOHN and Luttinger<sup>1</sup> have pointed out that the usual linked-cluster expansion for the ground-state energy of a many-fermion system is incomplete if there are deviations from spherical symmetry, e.g., due to lattice forces. By studying a grand canonical ensemble in the limit of zero temperature, they have derived a new energy term of the second order in the interaction. At first sight it seems puzzling that this term involves only that part of the interaction which is *diagonal* in the original occupation number representation: The result should then be derivable even if all off-diagonal parts of the Hamiltonian are omitted. Such an elementary derivation will then reveal the simple meaning of the Kohn-Luttinger term.

To sketch such a derivation, we formally introduce a parameter  $\lambda$  as a factor in the interaction, viz., a "coupling strength" which may be "switched on" adiabatically. If the Fermi surface is nonspherical it will depend on  $\lambda$ . Calling its interior  $G(\lambda)$ , the ground-state energy (diagonal part) may be written

$$E(\lambda) = \int_{G(\lambda)} d^3k \epsilon_k + \frac{1}{2}\lambda \int_{G(\lambda)} d^3k \int_{G(\lambda)} d^3k' v_{\mathbf{k}\mathbf{k}'},$$

where  $v_{\mathbf{k}\mathbf{k}'} = v_{\mathbf{k}'\mathbf{k}}$ . For each value of  $\lambda$ , the Fermi surface

is defined as that surface for which  $E(\lambda)$  is a minimum:

$$\epsilon_k + \lambda \int_{G(\lambda)} d^3k' v_{\mathbf{k}\mathbf{k}'} = \mu(\lambda),$$

in conjunction with

$$\int_{G(\lambda)} d^3k = \text{const.}$$

Then, if we vary  $\lambda$ ,

$$\frac{\partial E(\lambda)}{\partial \lambda} = \frac{1}{2} \int_{G(\lambda)} d^3k \int_{G(\lambda)} d^3k' v_{\mathbf{k}\mathbf{k}'},$$

because the other terms cancel, and expressions for  $\partial^n E(\lambda)/\partial \lambda^n$  ( $n=2, 3, \dots$ ) are easily written down. They are, of course, all zero if  $G$  is a sphere and therefore independent of  $\lambda$ .  $(\partial^2 E/\partial \lambda^2)_{\lambda=0}$  corresponds precisely to the Kohn-Luttinger term [Eq. (33) in reference 1].

When off-diagonal elements of the interaction are added as perturbations, it is presumably advantageous to start from a ground-state Fermi surface which is correct to the desired order of approximation.<sup>2</sup> The possibility that the surface may not remain "sharp" (superconductor) should not be overlooked.

<sup>1</sup> W. Kohn and J. M. Luttinger, Phys. Rev. 118, 41 (1960).

<sup>2</sup> A. Klein, Phys. Rev. Letters 4, 601 (1960).