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## SPECTRUM OF COLLECTIVE EXCITATIONS IN LIQUID ${}^4\text{He}$

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Experiments on inelastic coherent scattering of cold neutrons by simple classical liquids reveal the presence of definite elementary excitations of the density-fluctuation type in the frequency region  $\omega \sim 10^{12} \text{ sec}^{-1}$  and in the wave-number region  $Q \sim 1 - 3 \text{ \AA}$  in simple liquids [1, 2]. Theoretical calculations of the spectrum of elementary excitations for liquid argon were carried out in [3, 4]. On the other hand, it was established in [5] that the spectrum of elementary excitations in superfluid liquid helium retains a form similar to the spectrum of elementary excitations in simple classical liquids on going through the  $\lambda$  point, and only the damping of the elementary excitations increases somewhat.

It seems possible therefore to calculate the spectrum of collective excitations in liquid  ${}^4\text{He}$  in the region of high frequencies and large wave numbers, where the hydrodynamic description is not applicable [6], and an important role is played by the quasicrystalline properties of the liquid; this was the purpose of the present investigation.

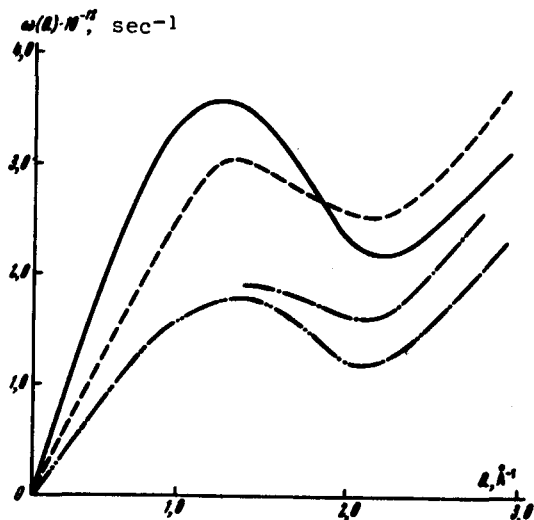
Using the method of equations of motion for two-time Green's function of the "displacement-displacement" type, we obtained the spectrum of the collective excitations in liquid  ${}^4\text{He}$  within the framework of the quasicrystalline model of the liquid [7]. The collective motions in the liquid are regarded as oscillations of the atoms in equilibrium positions, the distribution of which in space is described by the function  $g(R)$ . To take into account the strong anharmonicity due to the large zero-point oscillations, we used methods of self-consistent theory of anharmonic crystals, which make it possible to take into account all the even orders of anharmonicity [8]. The correlations of importance for solid and liquid helium, of the hard-core type, were taken into account on the basis of the approximate solution of the equation for the T-matrix described in [9]. As a result we obtained the following expressions for the longitudinal oscillation frequency:

$$\omega^2(Q) = \frac{2\pi\rho}{m} \int_0^\infty dR R^2 \int_{-1}^1 dt \Phi(R) t^2 (1 - \cos QRt), \quad (1)$$

where  $m$  is the mass of the helium atom,  $\rho = N/V$  the particle density, and  $\Phi(R)$  the longitudinal component of the pseudoharmonic force constant:

$$\Phi(R) = g(R) \left( \frac{\partial^2}{\partial R^2} + \frac{1}{R} \frac{\partial}{\partial R} \right) \int \frac{d^3r}{(2\pi/\alpha^2)^{3/2}} f(r) \exp \left[ -\frac{1}{2} \alpha^2 (r - R)^2 \right] \phi(r), \quad (2)$$

where  $g(R)$  is the distribution function of the equilibrium positions, and is assumed approximately to equal the radial distribution function obtained in



[10].  $\phi(r)$  is the (6 - 12) Lennard-Jones potential with parameters  $\epsilon = 10.22^\circ\text{K}$  and  $\sigma = 2.556 \text{ \AA}$ ,  $1/\alpha^2 = \langle (\vec{u}(0) - \vec{u}(\vec{R}))^2 \rangle$  is the rms relative displacement of the atoms from the equilibrium position, and is determined in a self-consistent manner from the frequency spectrum, and  $f(r)$  is the solution of the Bethe-Goldstone equation [9], which makes it possible to take correlations of the hard-core type into account.

In calculating the elementary-excitation spectrum,  $\alpha$  for  $T = 0^\circ\text{K}$  is determined from the condition of minimum energy of the ground state  $E(\alpha)$ . The calculations were made with a computer. The dispersion dependence for  $T = 0^\circ\text{K}$  is given in the figure (solid line), which shows for comparison the experimental curve of [5] (dotted) and the results of the calculations of [11] (dashed) and

[12] (dash-dot). We see that the results are in satisfactory agreement with experiment, thereby evidencing the applicability of the quasicrystalline model of the liquid to the description of the elementary-excitation spectrum not only in classical liquids, [3, 4], but also in the quantum liquid  $^4\text{He}$  when account is taken of appropriate corrections required by large-zero-point oscillations.

In conclusion, I consider it my pleasant duty to thank N.M. Plakida and V.B. Priezzhev for help and valuable discussions.

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#### TEMPERATURE DEPENDENCE OF THE PROBABILITY OF THE MOSSBAUER EFFECT IN YTTRIUM-IRON GARNETS

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The exchange interaction between octahedral (a) and tetrahedral (d) iron ions in ferrimagnets should exert an influence on the vibrational spectrum of the crystals [1]. We have investigated the temperature dependence of the Mossbauer-effect probability  $f'$  on  $\text{Fe}^{57}$  nuclei at temperatures 78 - 750°K in the iron garnet  $\text{Y}_3\text{Fe}_5\text{O}_{12}$  and the substituted iron garnet  $\text{Y}_3\text{Fe}_4\text{Al}_1\text{O}_{12}$ , the physical