

Droplets of mixtures of ^3He - ^4He at finite vorticity

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Abstract

Stimulated by recent interests in mixtures of quantum fluids, we study the properties of ^3He - ^4He mixture droplets using a phenomenological density functional. The structure of liquid drop with and without vortices is investigated using Monte Carlo simulations.

Key words: droplet, ^3He - ^4He mixture, quantum fluid

There is a burgeoning interest in droplets of quantum fluids from different directions recently. These include droplets of Bose-Einstein condensed alkali atoms confined by magnetic and optical means[1], small droplets (1000 atoms) of He in a molecular beam[2], and large droplets (mm in diameter) levitated by a magnetic field[3]. While one component drops have received much interest, multicomponent systems are beginning to be looked at. In addition to their own intrinsic interest, droplets of mixtures provide a unique opportunity to test and enhance our understanding of finite size effects and the influence of boundaries on phase segregation effects.

Current Bose condensed alkali mixtures are mostly mixtures of Bose particles. ^3He - ^4He are mixtures of Fermions and Bosons. In this work we present Monte Carlo simulations on the structure of ^3He - ^4He mixture droplets with and without vortices as a function of their relative concentration. This follows our earlier work that focuses on the simpler aspects of this system[4]. At zero vorticity, because of the quantum pressure most of the ^3He stays on the outside with a small amount of ^3He remaining inside, as is previously reported.

The total free energy for a liquid-helium mixture of ^3He and ^4He can be expressed in a density functional form.[6] We approximate the volume integral of the energy functional by a discrete sum. We sample a lat-

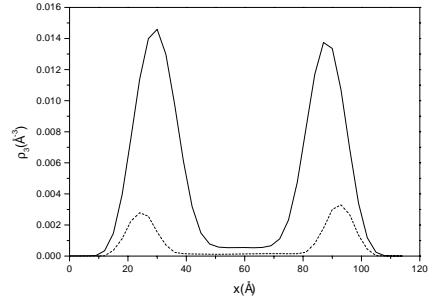


Fig. 1. Cross sections of density distributions of ^3He in the xy plane for $z=0$ with particle numbers $N_3 = 500$ and $N_4 = 3500$ (dashed line), $N_3 = 3000$ and $N_4 = 1000$ (solid line)

tice inside a sphere of radius $2R$ consisting of 40 sites along the diameter, making a total of 33398 sites. We choose the mesh size as 3 Å, which is smaller than the surface and interface thickness. We use the Metropolis algorithm, start our simulation with high temperature, then decrease it to a very low temperature step by step. This is an annealing process and the final configuration should be the lowest energy configuration. Because of the strong interactions between the helium atoms, the particle number is a very sensitive function of the chemical potential. We thus carry out the simulations with fixed particle numbers.

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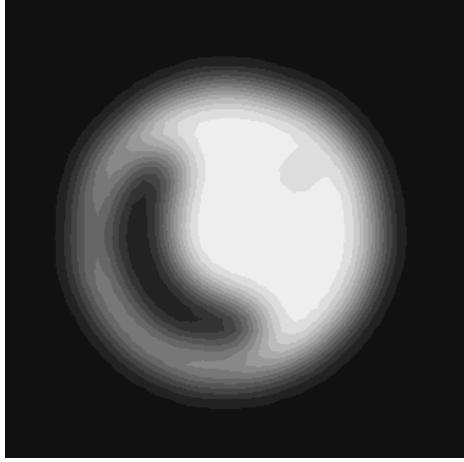


Fig. 2. Image of density distributions of ^3He with a vortex in ^4He . The vorticity is $m = 1$ and the particle numbers are: $N_3 = 3500$ and $N_4 = 500$.

Cross sections of the snap shots of the density profile for two different concentrations are shown in Fig. 1. The maximum bulk solubility of ^3He in the ^4He -rich phase is $\sim 6.6\%$ at zero temperature. When the ^3He fraction is small, there is a small amount ^3He inside the droplet, consistent with the bulk phase diagram. However, the concentration of ^3He is less than that in the bulk. We think this comes about because the ^3He inside must have the same chemical potential as that outside. Because of the free boundary condition the chemical potential of ^3He outside is lower than that in the bulk. When the ^3He number is very small, the solubility of ^3He into ^4He is much smaller than the bulk value. The effective attractive potential for ^3He atoms on the surface of ^4He provide a type of surface states, which are called Andreev states. ^3He atoms first accumulate on the free surface occupying Andreev states before start dissolving into the bulk. The recent change of the He3-4 mutual solubility in aerogels[5] is another example of the effect of boundary conditions on the phase segregation phenomena.

Taking the vortex along the z axis, the free energy functional of the vortex state is obtained with the addition of a centrifugal term $\hbar^2 \rho_4 \kappa^2 / (2m_4 r^2)$ associated with the superfluid flow where κ is the quantum circulation number and r is the distance to the z axis. At finite vorticity, ^4He moves outside because of the centrifugal force. Images of the snap shots of the density distributions at a low ^4He concentration ($N_3 = 3500$ and $N_4 = 500$) are shown in Fig. 2. The ^4He forms lumps and does seem to “wet” the ^3He completely. We emphasize that the pictures are **snap shots** and not **averages** of the density distribution. It is the density correlation between ^3He and ^4He that is being shown. Thus at low ^4He concentrations, the **density correlation** between ^3He and ^4He becomes asymmetric. Since

^4He is rotating, we expect the **average** density to be cylindrically symmetric. We think this phenomenon is related to the case of bulk ^3He - ^4He mixture, where the vortices lose their stability with increasing ^3He concentrations[?,?]. The more interesting question is how big will the damping be against the superfluid flow and how metastable is this configuration. We hope to investigate this in future publications.

This work is supported by a grant from NASA (NAG8-1427). We acknowledge helpful conversation with Ping Ao and V. Ryzhov.

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