

Numerical Analysis of Micro-/Nanoscale Gas-Film Lubrication of Sliding Surface with Complicated Structure

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Nakamori et al. [1] found experimentally that the friction between a partly polished diamond-coated surface and a metal surface was drastically reduced to zero as relative speed increased to a few m/s, and it seems that the diamond-coated surface took off the counter surface and the sliding mechanism was the gas film lubrication. Because of that the surface roughness of the diamond coated substrates is from $0.28 \mu\text{m}$ to $0.57 \mu\text{m}$ and the distance between two sliding surface will be in the order of surface roughness or smaller, Knudsen number Kn will be larger than 0.1. Therefore, micro-/nanoscale gas flows between two sliding surfaces cannot be treated as a continuum and we use the direct simulation Monte Carlo (DSMC) method [2].

In this study, to investigate the role of the gas flow for this phenomenon, we reproduce the complicated structure of the diamond-coated surface in the computational domain by using the data measured by an atomic force microscope (AFM), and perform numerical simulations of three-dimensional micro-/nanoscale gas flows. We reproduce the diamond-coated surface by using Marching Cubes method [3]. The computational domain is shown in Fig. 1. Fig. 2 shows the pressure distribution obtained in the DSMC simulation. In the case of Fig. 2, we obtained the lift force of $2.33 \times 10^3 \text{ Pa}$, which is large enough to levitate the slider used in the experiment.

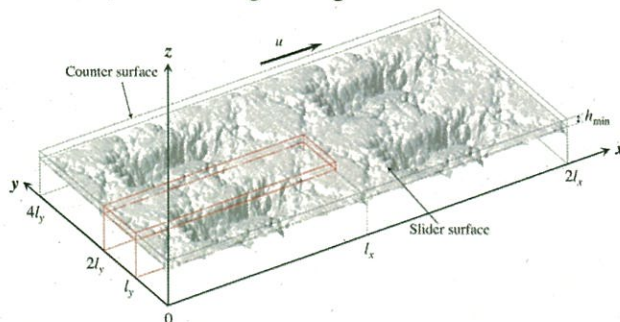


Fig. 1 Computational domain.

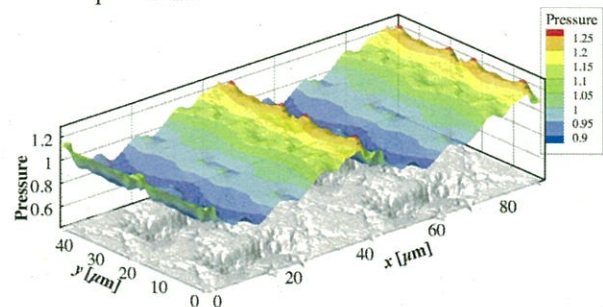


Fig.2 Pressure distribution($u=10 \text{ m/s}$, $R_a=0.21 \mu\text{m}$, $h_{min}=0.028 \mu\text{m}$).

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