

Application of Gaussian Random Field Model of Rough Surface in Gas Flow Simulation in Channels

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Abstract. Our new results on the number of crossings of gaussian random field with inclined straight line are applied to direct simulation of rarefied gas flow in a channel with rough walls by Monte-Carlo method.

Keywords: Surface roughness, rarefied gas flow, direct simulation, Monte Carlo method.

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Velocity distribution and numerical density of rarefied gas atoms in a two-dimensional (plain or axisymmetric) flow in a channel or in a nozzle at high Knudsen numbers is calculated by DSMC method. The main purpose of this paper is to study the dependence of the macro-parameters of rarefied gas molecular flow on surface roughness of the walls [1-6] and on geometrical shape of the channel. Using the expression of scattering function V on rough surface in the form $V = \hat{S} V_0$, where the roughness operator \hat{S} is fully determined by geometrical shape of roughness and by the trajectory of reflected gas atom [3-6], we expand the operator \hat{S} analytically. Expanding the local scattering function V_0 in a series in terms of the system of orthogonal functions, we account all the physical and chemical parameters of the gas and of the surface. Unlike more simple models applied by other researchers, such as polygonal-line or conical-hole based models [7-8], our approach allows us to simulate surface roughness on micro-level by a wide class of Gaussian random fields [5]. This statistical model of the roughness permits us to study the roughness operator \hat{S} applying our new results based on the random-field and random-process theory [9-10]. The complexity of the computation of continuum integrals, which must be approximated by the integrals of higher dimension [6], is tackled applying Rice series [5]. Bypassing the most principal difficulty in statistical approach, we reduce the computation time since the most cumbersome and time-consuming step of the computation can be performed prior to the simulation of gas atom trajectories. Such technique of the pre-computation of continuum integrals allows us not only minimizing computation time, but also reducing the number of the parameters. Our algorithm based on the approximation of scattering function V and of the momentum and energy exchange coefficients [1-3] eliminates the complex simulation of the geometrical-shape of the rough surface in the Monte-Carlo method. Additional advantage of considered statistical approach is that not only diffuse-specular model [6], [8], but also Cercignani-Lampis scattering kernel or phenomenological models like proposed in [10] can be applied in the local scattering function V_0 approximation. For example, fig.1 shows calculated velocity distribution $P(\theta_1)$ of gas molecules (the directions are defined by the angles θ_1) in different cross-sections of a nozzle with a gorge in the middle section for different roughness parameters σ_1 (from 0 to 0.5).

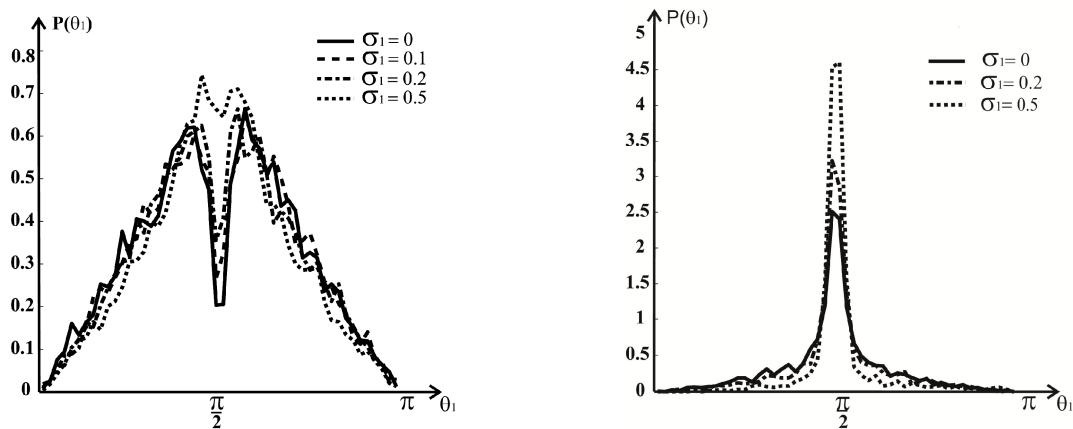


Fig.1: The distribution $P(\theta_1)$ of gas molecules velocities for different values of roughness parameter σ_1 .

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