

Abstract Submitted
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Simulation of low energy ion bombardment of alkane in low pressure RF plasmas¹ VIKTOR ZHELTUKHIN, Retired, ALBINA AZANOVA, IGOR BORODAEV, AIDOGDY SHAKHYROV², Kazan National Research Technological University — Results of molecular dynamics simulation of low energy (up to 100 eV) ion bombardment of both crystalline and amorphous alkanes in low pressure RF plasmas are described. An united-atom model of alkane chains where each site represents a CH₂ group or a CH₃ end group is used. The covalent bond between adjacent sites is modeled by a Dreiding potential. The interaction between the sites in different chains, as well as between argon and alkane sites is described by a LJ potential. A classical molecular dynamics code LAMMPS is used.

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