Molecular Dynamics Simulation Report

Optimized Lattice Parameter: 5.2823e-10 m

Melting Point: 135 K

Boiling Point: 150 K

Total Runtime: 29.60 seconds

Best Verlet Algorithm: Leapfrog Verlet (Adaptive)

Best Time Step: 1.00e-14 s

Reason: Lowest average potential energy (-1.32e-20 J) and minimal energy drift (1.34e-25 J) during preliminary simulation.