DSMC Direct Simulation Monte-Carlo method



Orion reentry simulation (Norman, P., Valentini, P., and Schwartzentruber, T.E.,) u. Minnesota

- Gas molecules have large mean free path in comparison to size.
- Small <10-6 sec time stepping simulation.
- Simulation can divide transport phase and energy redistribution phases into two independent components.
- Transport phase \rightarrow move molecules, interact with surfaces
- Energy Phase → collide nearest neighbours
- Too many real molecules to simulate in a computer : use a statistical approach.

• Transport .

Move molecules ds = V.dt Hit surface ?

- a) reflect. Conservation of momentum.
- b) absorbed-reinjected. Conservation of energy.



- Collisions.
 - Hard Sphere
 - Variable Hard Sphere
 - Variable Soft Sphere

$$d = \left(\frac{15(mkT/\pi)^{1/2}}{2(5-2\omega)(7-2\omega)\mu}\right)^{1/2}$$



• Only simulate representative molecules.

- 1 in 10¹² for near continuum flow

- Find probability of collisions for dt for molecules in a small region.
- Select pairs for collision based on probability

(acceptance-rejection technique)

 Post collision velocities based on conservaion of energy/momentum.



- Flow macroscopic properties based on statistical sampling of simulated molecules over time.
 - ---> density (molecules per unit volume)
 - ----> Velocity (average speed of molecules)
 - ---> **Temperature...**translational,rotational, vibrational (energy of random motion, spin and internal molecule state)
 - ----> Pressure ??? (only at surfaces)

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