

PHYS:5905 Homework #5a

Please submit your solutions as a single PDF file with answers to the questions asked.

Please complete required problems before lecture on Thursday, February 21, 2019.

1. (Required) Small-Angle Collision Scattering Routine

- (a) Write a routine that performs elastic small-angle collisions of angle θ , normally distributed with a small-angle variance $\sigma(\theta)$.
- (b) This routine will take an initial velocity vector \mathbf{v}_i and scatter it away from that angle by a small θ with a random azimuthal angle ϕ about the original vector. Let us take our initial velocity vector to be $\mathbf{v}'_i = v'_0 \hat{\mathbf{x}}$, where $v'_0 = 1$.

(c) To accomplish this small-angle scattering, use the following steps:

- i. First, we rotate the velocity vector \mathbf{v}_i to the z -axis, here denoted \mathbf{v}_{inc} . This can be done by computing the direction of the velocity \mathbf{v}_i in spherical coordinates using the Cartesian to Spherical coordinate function,
`[phi1,elev1,vmag]=cart2sph(y0(4),y0(5),y0(6));`
where $\mathbf{y}_0 = [x \ y \ z \ v_x \ v_y \ v_z]$ is row vector with the initial position and velocity, the azimuthal angle `phi1` ranges over $[-\pi, \pi]$ and the elevation `elev1` ranges over $[-\pi/2, \pi/2]$. If we rotate our vector about z by `-phi1` and then about y by minus the polar angle $-\theta_1 = \text{elev1} - \pi/2$, then our vector will be aligned with z . (Verify this before moving on.)
- ii. Next, we obtain a scattering through an angle θ by rotating first about the y -axis by the angle θ using a rotation matrix

$$R_1 = \begin{pmatrix} \cos(\theta) & 0 & \sin(\theta) \\ 0 & 1 & 0 \\ -\sin(\theta) & 0 & \cos(\theta) \end{pmatrix} \quad (1)$$

We also need to allow for a rotation by ϕ about the z -axis using

$$R_2 = \begin{pmatrix} \cos(\phi) & -\sin(\phi) & 0 \\ \sin(\phi) & \cos(\phi) & 0 \\ 0 & 0 & 1 \end{pmatrix} \quad (2)$$

Using these rotation matrices, the scattered vector $\mathbf{v}_{scat} = R_2 R_1 \mathbf{v}_{inc}$

NOTE: The usual convention for rotation matrices is for a rotation of the *coordinate system* by angle ϕ . In a fixed coordinate system, the same rotation matrix will yield rotation of the vector by angle $-\phi$, which is why the negative signs in the matrices above may be opposite from where you are used to seeing them.

- iii. We can model small angle collisions by using normally distributed random numbers, using the Matlab function `randn`, to determine the angle θ with some small typical variance $\sigma(\theta)$ of the distribution. Thus, you can use
`the=randn*sigmatheta;`
to compute θ over a range with a variance $\sigma(\theta) = \pi/18$.
- iv. The azimuthal angle ϕ of the scattering is a randomly selected value with uniform probability in the range $0 \leq \phi \leq 2\pi$, using
`phi=rand*2*pi;`
- v. Finally, we perform the reverse steps of the original rotation, first $-\theta_1$ about y followed by `+phi1` about z .
- (d) (Return) Perform the scattering $n_{coll} = 1000$ times, plotting in 3D using `plot3` the track of the tip of the velocity vector (v_x, v_y, v_z) at each scattering for $\sigma(\theta) = \pi/18$. Use
`axis([-1. 1. -1. 1. -1. 1.]);`
to ensure that the entire unit sphere is included in the plot dimensions.
- (e) (Return) Produce another 3D plot using a smaller scattering angle variance, $\sigma(\theta) = \pi/180$.

2. (Required) Implementation of Monte Carlo Collisions

- (a) For electron single-particle motion and confinement in a magnetic mirror configuration, we will implement a collision model for electron-ion collisions with $m_i \gg m_e$, such that the electrons exchange no energy with the massive ions, but rather simply have a change in the direction of the velocity.
- (b) We will take a velocity-independent collision operator with a small collision frequency ν such that $\nu/\Omega \ll 1$. In this limit, collisions are a Poisson process with a probability for a collision in a timestep Δt given by

$$P_{coll} = \nu\Delta t \ll 1 \quad (3)$$

- (c) Because a collision in this limit changes the velocity of the particle discontinuously (while conserving kinetic energy), the implementation of a higher order stepping scheme that depends on derivatives at previous timesteps is problematic. (Can you discern what the problem may be?). Therefore, I recommend you use Runge-Kutta timestepping since it depends only on the current timestep values to advance the particles.
- (d) Since the probability of a collision during one timestep is low, $\nu\Delta t \ll 1$, we can achieve significant computational savings by only checking to see if a collision occurs at regular intervals T (which may be longer than a single timestep $T > \Delta t$, as long as $\nu T \ll 1$).
- (e) You are free to implement your Monte Carlo Collision operator as you see fit, but I can recommend this approach:
 - i. Integrate over a short interval of time T using the RK45 Adaptive Timestepping algorithm using a relative tolerance
`options = odeset('RelTol', 1.0e-3);`
 - ii. At the end of each interval T , check to see if the particle has escaped the magnetic mirror (with $|z| > L/2$, where L is the length between field maxima along the mirror axis). If the particle has escaped, save the time at which the particle escaped as τ and exit the time integration of the particle motion. This time τ is a measure of the confinement time in the magnetic mirror.
 - iii. At the end of each interval T , check to see if a collision has occurred.
 - iv. If the collision has occurred, rotate the velocity vector of the particle by a small angle collision with variance $\sigma(\theta) = \pi/18$ (use results of previous problem).
 - v. Integrate over the next short interval of time T using the RK45 Adaptive Timestepping algorithm, and repeat the steps above.
 - vi. Stop the time integration at T_{end} if the particle has not yet escaped.
- (f) Set up a magnetic mirror with the parameters: mirror length $L' = 10$, minimum axial field value $B'_{00} = 10$ and mirror ratio $R_m = 4$.
- (g) Choose an initial position $\mathbf{x}_0/r_L = (0, 0, 0)$, and an initial velocity $\mathbf{v}_0/v_\perp = (0, 1, 1)$ for the electron for each run of the ensemble. You may want to limit the total integration time for each run to $\Omega T_{end} = 100\pi$ and choose the collision check interval $T = T_{end}/5000$.
- (h) Run an ensemble of runs with different collisionalities over the range $0.001 \leq \nu/\Omega \leq 1$ with at least 4 different collisionalities to sample the range. Note that, for the lowest collisionality values, the particles may not escape before T_{end} , but over a sufficiently large ensemble, you should see that the mean confinement time has a dependence on the collisionality.
- (i) For each choice of collisionality, run at least 8 single-particle motion (with Monte Carlo collisions) runs to obtain good statistics for the confinement time τ .
- (j) (Return) Produce a plot of mean confinement time $\langle \tau \rangle$ with error bars giving the standard deviation $\sigma(\tau)$ vs. the normalized collision frequency ν/Ω . You will likely want to use the Matlab plotting function `errorbar` to make this plot.
- (k) (Return) Produce a plot of the position of the particle in (v_\parallel, v_\perp) space (relative to local mean magnetic field) at the point in time at which the particle is lost. The best way to do this is to write a function that inputs the vector `y0` and returns `vpar` and `vperp`.
- (l) (Optional) Plot the mean and standard deviation (`errorbar` plot) of the number of collisions as a function of normalized collision frequency ν/Ω .