

# 2nd Assignment-I

- Write a program that generates random realization of a Plummer sphere (positions only).
- Write separately a routine that generates radial coordinates, and another that generates random points on the unit sphere. Combine the results to generate cartesian coordinates for each point.
- Make a plot of the theoretical radial density profile vs. a histogram of the radial distribution of particles, normalized to the volume of the spherical shell used for each radial bin.
- Determine the global *rms* deviations between both curves as a function of the size of the random realization.

# 2nd Assignment-I

Let us work with a dimensionless version of this model:

$$\delta_p(\eta) = (1 + \eta^2)^{-5/2}, \quad \text{where } \eta \equiv r / r_o, \quad \text{and } \delta_p(\eta) = \rho_p(\eta) / (3M / 4\pi r_o^3)$$

The cumulative distribution function can be obtained integrating  $\delta_p(\eta)$  in spherical shells:

$$4\pi \int_0^\eta s^2 \delta_p(s) ds = \frac{4\pi}{3} \frac{\eta^3}{(1 + \eta^2)^{3/2}} \rightarrow \frac{4\pi}{3}, \quad \text{when } \eta \rightarrow \infty$$

So we define the normalized cumulative distribution function as:

$$\mu_p(\eta) \equiv \frac{\eta^3}{(1 + \eta^2)^{3/2}}$$

The inverse function is simply,

$$\eta = \sqrt{\frac{\mu_p^{2/3}}{1 - \mu_p^{2/3}}}$$

# 2nd Assignment-I

$$\eta = \sqrt{\frac{\mu_P^{2/3}}{1 - \mu_P^{2/3}}}$$

Main program (f77)

```
*----- Plummer model positions -----  
  
open(1,file='plummer.dat',form='formatted',status='new') !open output file  
  
do i=1,Np                                !particle do-loop  
  aux = rand()                            !random number between 0 & 1  
  aux = aux**(2./3.)  
  aux = aux/(1. - aux)                    !inverse cumulative function for r  
  r = ro*Sqrt(aux)                        !random radial coordinate for Plummer  
  call randangle(x,r)                     !random direction  
  
  write(1,20) i,(x(j),j=1,3),r !dump particle info  
20  format(1x,i6,4x,3(1pe10.3,1x),4x,1pe10.3)  
end do  
  
close(1)                                  !close output file
```

# 2nd Assignment-I

## Randangle subroutine (f77)

```
Subroutine randangle(x,r)
*
* It generates the 3 coordinates x(3), of a random direction vector of user-supplied modulus r.
* Method: 2 random numbers are generated: z & theta.
* The 1st is the z-component, the 2nd, the azimuth (angle of x-y projection with x-axis)
* of a unit vector.
* The final cartesian coordinates: x(3) are obtained by simple scaling by r.
* Notice the use of z, instead of the other spherical angle to avoid a non-uniform
* Coverage of the unit sphere.
*
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*
PARAMETER (pi=3.14159265)    !pi
REAL r
REAL x(3)
REAL z,theta,aux

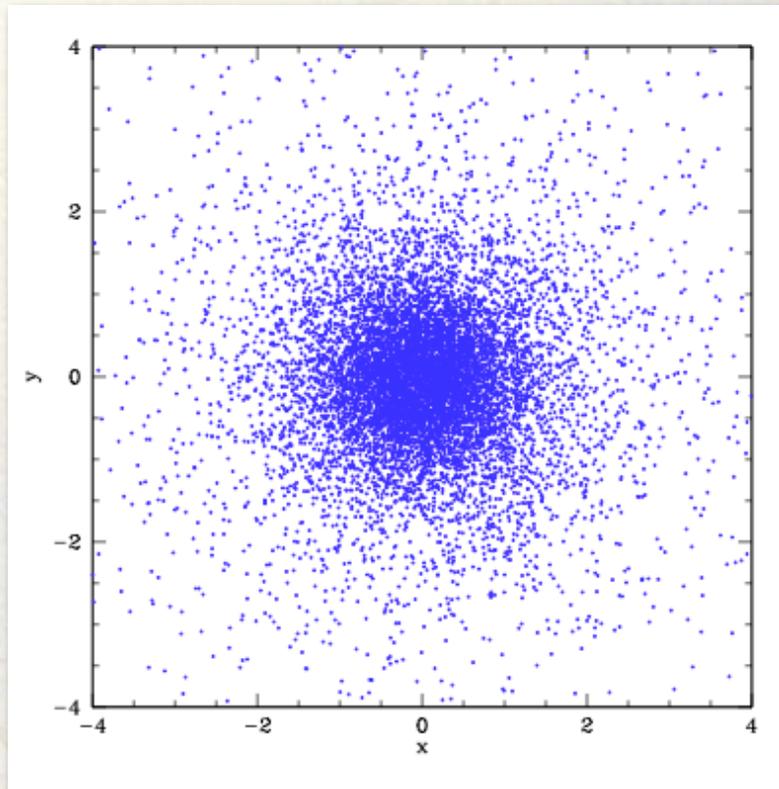
z = 2.*rand() - 1.           !random number between -1 y +1.
theta = 2.*pi*ran1(iseed)    !random number between 0 & 2*pi.
aux = Sqrt(1. - z*z)         !XY-projection of unit vector of z-coordinate
x(1) = r*cos(theta)*aux      !cartesian coordinates
x(2) = r*sin(theta)*aux
x(3) = r*z

return
end
```

# 2nd Assignment-I

Random realization of Plummer model in  $x$ - $y$  projection.

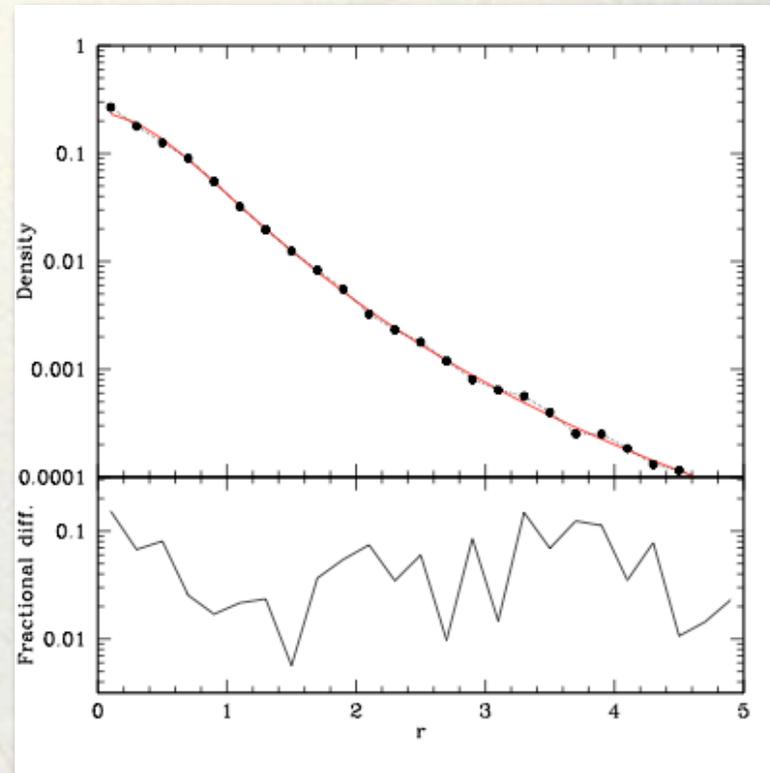
$N = 10,000$



# 2nd Assignment-I

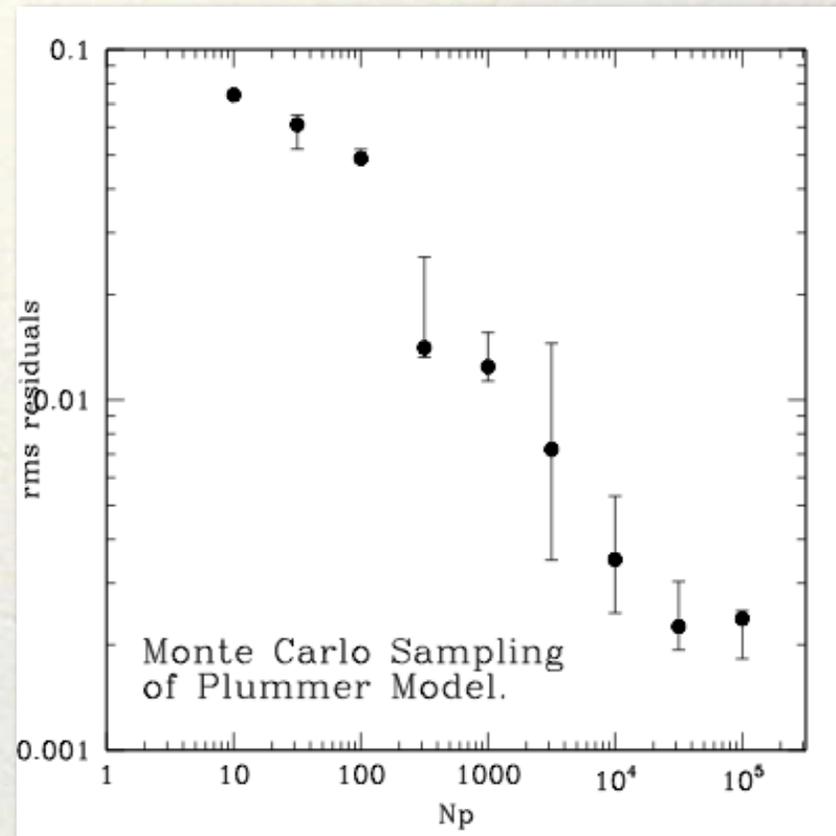
Density profile and residuals of random realization of Plummer model.

$N = 10,000$



# 2nd Assignment-I

Overall *rms* residuals of random realizations of plummer models of increasing size.



# Homework

## 2nd Assignment - II

- Write a program that approximates the value of the integral:

$$A = \frac{1}{\sqrt{2\pi}} \iint_{x^2 + y^2 \leq 1} \exp[-(x^2 + y^2) / 2] dx dy$$

using the von Neumann Rejection Technique.

- List the approximated values  $A_n$  obtained using  $n$  points, as well as the fractional errors:  $(A_n - A_o) / A_o$ ; where  $A_o$  is the exact value.
- Make a plot of the fractional error as a function of  $n$ .

# 2nd Assignment-II

## Program Aintegral

It uses the von Neumann Rejection Technique to approximate the value of the integral:

$A = \text{Int} (1/\sqrt{2\pi}) * \text{Exp}[-(x^2 + y^2)/2] \, dx dy,$   
within the domain:  $(x^2 + y^2) < 1$   
The exact value is:  $\sqrt{2\pi} * [1 - (1/\sqrt{e})] \sim 0.986281$

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```
parameter (Ao=0.986281) !exact value
real A !Approximate value
real rmax !radius of integration domain
real error !absolute fractional error
integer n !size of sample
```

```
rmax = 1.
n = 10 !initial sample size
do i=1,10
  call Integ(n,A,rmax) !von Neumann integration
  error = Abs(A-Ao)/Ao !absolute fractional error
  write(6,10) n,A,error!write result
format(1x,i9,3x,1pE11.5,2x,1pE9.4)
  n = 2*n !duplicate sample size
end do
```

```
stop
end
```

# 2nd Assignment-II

\*  
\*  
\*  
\*  
\*

```
subroutine Integ(n,A,rmax)
von Neumann integration routine:
      n:    number of points to use
      A:    returned approximate value of integral
      rmax: radius of integration domain
```

```
integer n,m
real A,rmax,fmx
real xrand,yrand,r2,frand,fraction
external func      !function to be integrated
```

```
fmx = func(0.,0.)      !maximum value of 'func' within domain
```

```
m = 0                !initialize successes counter
```

10

```
do i=1,n
  xrand = ran()      !random x within (0,1)
  yrand = ran()      !random y within (0,1)
  r2 = xrand*xrand + yrand*yrand  !respective modulus squared
  if (r2.gt.1.) go to 10  !try again if point outside domain
  frand = fmx*ran()  !random f within (0,fmx)
  if (frand.lt.func(xrand,yrand)) m = m + 1 !add to counter if good
end do
```

```
fraction = float(m)/float(n)  !ratio of sampled to under func volumes
A = 3.14159265*fraction      !multiply by sampled volume
```

```
return
```

