



ROOT : A super quick recap

1-Dimensional Histograms : TH1I,TH1F,TH1D

```
TH1F* name = new TH1F("name","Title", Bins, lowest bin, highest bin);
```

• Example:

```
TH1F* h1 = new TH1F("h1","x distribution",100,-4,4);
```

```
Float_t x = 0;
```

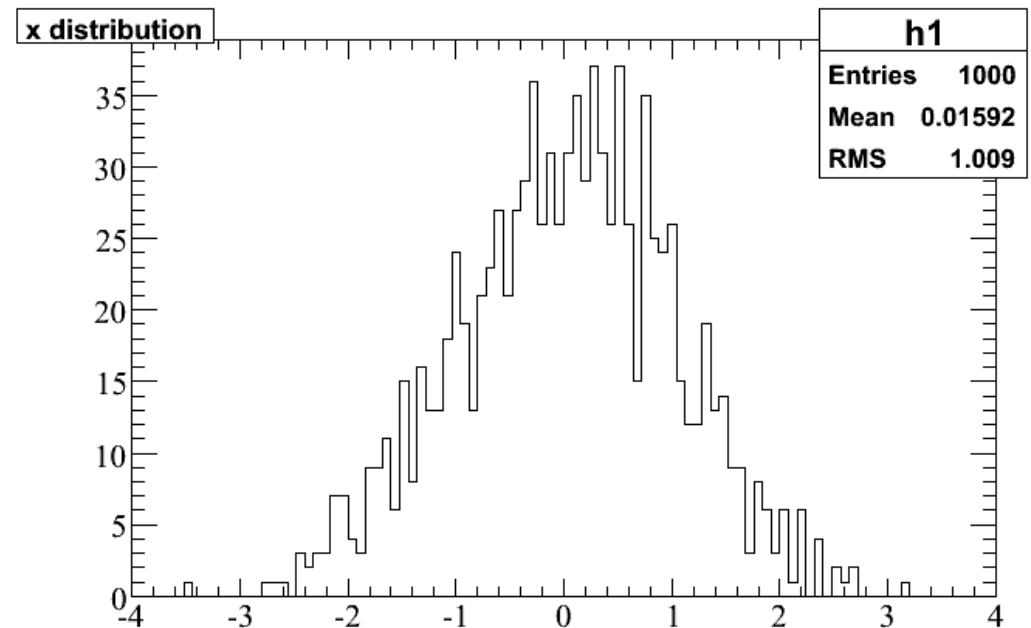
```
for(Int_t i=0; i<1000; i++){
```

```
    x=gRandom->Gaus(0,1);
```

```
    h1->Fill(x);
```

```
}
```

```
h1->Draw();
```



The complete list of Methods can be found in:
<http://root.cern.ch/root/html/TH1F.html>

Histograms

```
// using different constructors
```

```
TH1I* h1 = new TH1I("h1", "h1 title", 100, 0.0, 4.0);
```

```
TH2F* h2 = new TH2F("h2", "h2 title", 40, 0.0, 2.0,30, -1.5, 3.5);
```

```
TH3D* h3 = new TH3D("h3", "h3 title", 80, 0.0, 1.0,100, -2.0, 2.0,50, 0.0, 3.0);
```

```
// cloning a histogram
```

```
TH1F* hc = (TH1F*)h1->Clone("title of the cloned histogram");
```

```
// projecting histograms
```

```
// the projections always contain double values !
```

```
TH1D* hx = h2->ProjectionX(); // ! TH1D, not TH1F
```

```
TH1D* hy = h2->ProjectionY(); // ! TH1D, not TH1F
```

Histograms : A complex example

```
// histograms filled and drawn in a loop
void hsum() {

TCanvas *c1 = new TCanvas("c1","The HSUM example",200,10,600,400);
c1->SetGrid();

// Create some histograms.
TH1F *total  = new TH1F("total","This is the total distribution",100,4,4);
TH1F *main   = new TH1F("main","Main contributor",100,-4,4);
TH1F *s1     = new TH1F("s1","This is the first signal",100,-4,4);
TH1F *s2     = new TH1F("s2","This is the second signal",100,-4,4);

total->Sumw2(); // store the sum of squares of weights
total->SetMarkerStyle(21);
total->SetMarkerSize(0.7);
main->SetFillColor(16);
s1->SetFillColor(42);
s2->SetFillColor(46);
```

Macro: hSum.C

Histograms : A complex example

```
// Fill histograms randomly

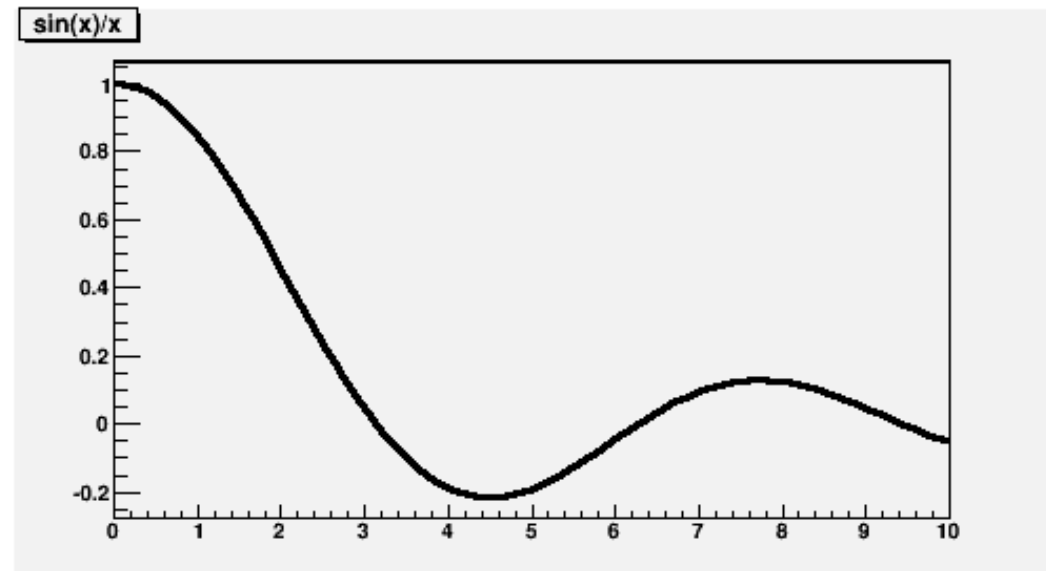
gRandom->SetSeed();

Float_t xs1, xs2, xmain;
for (Int_t i=0; i<10000; i++) {
    xmain = gRandom->Gaus(-1,1.5);
    xs1   = gRandom->Gaus(-0.5,0.5);
    xs2   = gRandom->Landau(1,0.15);
    main->Fill(xmain);
    s1->Fill(xs1,0.3);
    s2->Fill(xs2,0.2);
    total->Fill(xmain);
    total->Fill(xs1,0.3);
    total->Fill(xs2,0.2);
}
total->Draw("sameaxis"); // to redraw axis hidden by the fill area
```

Macro: hSum.C

Formulas and Functions

```
void formula2() {
    TCanvas *c1 = new TCanvas("c1","Example with Formula",500,500);
    // Create a one dimensional function and draw it //
    TF1 * fun1 = new TF1("fun1","sin(x)/x",0,10);
    c1->cd(1);
    fun1->Draw();
    cout << " Deriv " << fun1->Derivative(2.) << endl;
    cout << " Integral " << fun1->Integral(0.,3.) << endl;
    cout << " Func val " << fun1->Eval(1.2456789) << endl;
}
```



Macro: formula2.C

Creating a TF1 with Parameters

The second way to construct a **TF1** is to add parameters to the expression. Here we use two parameters:

```
root[] TF1 *f1 = new TF1("f1","[0]*x*sin([1]*x)",-3,3);
```

The parameter index is enclosed in square brackets. To set the initial parameters explicitly you can use:

```
root[] f1->SetParameter(0,10);
```

This sets parameter 0 to 10.

You can also use `SetParameters` to set multiple parameters at once.

```
root[] f1->SetParameters(10,5);
```

This sets parameter 0 to 10 and parameter 1 to 5.

We can now draw the **TF1**:

```
root[] f1->Draw();
```

Creating a TF1 with a User Function

```
// define a function with 3 parameters
Double_t fitf(Double_t *x,Double_t *par){
    Double_t arg = 0;
    if (par[2] != 0) arg = (x[0] - par[1])/par[2];
    Double_t fitval = par[0]*TMath::Exp(-0.5*arg*arg);
    return fitval;
}

void userfunctexample() {
    // Create a TF1 object using the function defined above. The last
    // parameter specify the number of parameters for the function.
    TF1 * func = new TF1("fit",fitf,-3,3,3);
    // set the parameters to the mean and RMS of the histogram
    func->SetParameters(500,0.,0.5);
    // give the parameters meaningful names
    func->SetParNames ("Constant","Mean_value","Sigma");
    // call TH1::Fit with the name of the TF1 object
    func->Draw();
}
```


Write and Read a ROOT file

- **Write (in a macro):**

```
TFile *myfile = new TFile("fillrandom.root", "RECREATE");  
myfile->cd();  
form1->Write();  
sqroot->Write();  
h1f->Write();  
myfile->Close();
```

- **Read (in a macro):**

```
TFile *pfile = new TFile("fillrandom.root");  
TH1F * ph1 = (TH1F *) pfile->Get("h1f"); // take the object from the TFile
```

- **Read (in the terminal):**

```
root [0] TFile *pfile = new TFile("fillrandom.root");  
root [1] pfile->ls(); // it is useful only to know what is stored inside your  
TFile if you have no idea of how it has been created.  
root [2] TH1F * ph1 = (TH1F *) pfile->Get("h1f"); // take the object from the  
TFile
```

TBrowser : A GUI to analyze the TFile

```
root [0] new TBrowser()
```

The screenshot displays the ROOT Object Browser interface. On the left, a file tree shows the directory structure, with the 'h1f' object selected under the path 'student/root/2013'. The main canvas, titled 'Test random numbers', shows a histogram of data points. A statistics box for the 'h1f' object is visible in the top right corner of the canvas area.

| h1f | |
|---------|-------|
| Entries | 10000 |
| Mean | 3.646 |
| RMS | 1.84 |

Command
Command (local):

```
stefano@venere:~/student/root/2013$ root -l
root [0] TBrowser pippo
root [1] (class TFile*)0x989c8f0
```

Passing arguments to a macro

```
// define a function with 3 parameters
Double_t fitf(Double_t *x, Double_t *par) {
    Double_t arg = 0;
    if (par[2] != 0) arg = (x[0] - par[1])/par[2];
    Double_t fitval = par[0]*TMath::Exp(-0.5*arg*arg);
    return fitval;
}

void userfunctexample(Int_t min = -3, Int_t max = 3) {
    // Create a TF1 object using the function defined above. The last
    // parameter specify the number of parameters for the function.
    TF1 * func = new TF1("fit", fitf, min, max, 3);
    // set the parameters to the mean and RMS of the histogram
    func->SetParameters(500, 0., 0.5);
    // give the parameters meaningful names
    func->SetParNames ("Constant", "Mean_value", "Sigma");
    // call TH1::Fit with the name of the TF1 object
    func->Draw();
    TString nameout = "Funz.";
    nameout += min;
    nameout += ".";
    nameout += ".root";
    TFile *fo = new Tfile(nameout, "RECREATE");
    fo->cd();
    func->Write();
    fo->Close();
}
```

Documentation

- <https://wwwusers.ts.infn.it/~lea/lacd2016.html>
- root.cern.ch
- In any case

Documentation

- <https://wwwusers.ts.in>
- root.cern.ch
- In any case



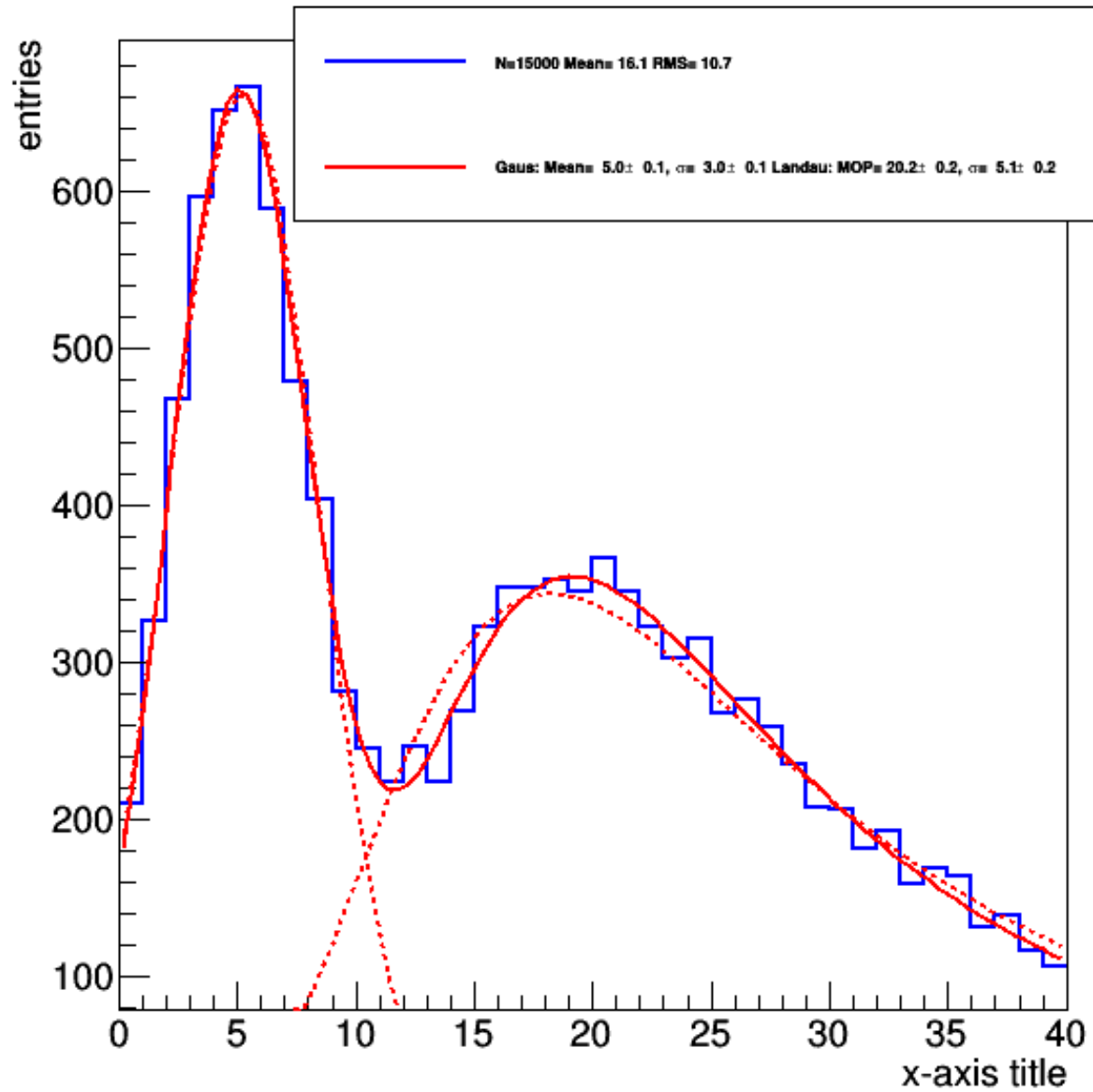
**KEEP
CALM
AND
ASK
GOOGLE**

Esercitazione 10

Esercizio 1

- Fill a histogram randomly ($n=10,000$) with a Landau distribution with a most probable value at 20 and a “width” of 5 (use the ROOT website to find out about the Landau function)
- Fill the same histogram randomly ($n=5,000$) with a Gaussian distribution centered at 5 with a “width” of 3
- Write a compiled script with a fit function that describes the total histogram nicely (it might be a good idea to fit both peaks individually first and use the fit parameters for a combined fit)
- Add titles to x- and y-axis
- Include a legend of the histogram with number of entries, mean, and RMS values
- Add text to the canvas with the fitted function parameters
- Draw everything on a square-size canvas (histogram in blue, fit in red)
- Save as png, eps, and root file

Esercitazione 10





Simulations with ROOT

Thanks to Prof. Massimo Maserà for the next slides

Monte Carlo technique

- Monte Carlo refers to any procedure that makes use of random numbers.
- Monte Carlo methods are used in:
 - Simulation of natural phenomena
 - Simulation of experimental apparatus
 - Numerical analysis (e.g. Integral evaluation)
- What is a Random Number?
 - Is e.g. 3 a random number?
 - A sequence of random numbers is a set of numbers that have nothing to do with the other numbers in the sequence

Monte Carlo technique

- General procedure:
 - A sequence of m random numbers with uniform distribution in the $[0,1]$ interval is extracted.
 - This sequence is used to produce a second sequence distributed according to a generic $f(x)$, which is the pool of simulated data

Pseudo-Random numbers

- The sequence is reproducible, because the algorithm for the generation is deterministic.
- General characteristic of a random generator:
 - Statistical independence
 - “Long” repetition period
 - The sequence looks random, when indeed it is not

Random generators in ROOT

Are implemented in the **TRandom** class: fast generator with a short (10^9) period, based of the linear congruential method.

- TRandom1: inherits from TRandom and implements RANLUX
- TRandom2: inherits from TRandom has a period of 1026, but only 3 32 bits word
- TRandom3: inherits from TRandom and implements the Mersenne-Twister generator which has a period of $2^{19937}-1$ ($\approx 10^{6002}$).

Here are the CPU times obtained using the four random classes on an Ixplus machine with an Intel 64 bit architecture and compiled using gcc 3.4:

| | TRandom (ns/call) | TRandom1 (ns/call) | TRandom2 (ns/call) | TRandom3 (ns/call) |
|----------------------|-----------------------------|------------------------------|------------------------------|------------------------------|
| Rndm() | - | - | 6 | 9 |
| Gaus() | 31 | 161 | 35 | 42 |
| Rannor() | 116 | 216 | 126 | 130 |
| Poisson(m=10) | 147 | 1161 | 162 | 239 |
| Poisson(m=10) UNURAN | 80 | 294 | 89 | 99 |

BAD

SLOW

FAST

GOOD
(default)

Simulation of a radioactive decay of a single nucleus

- Radioactive decay is an intrinsic random process: the probability of decay is constant (independent of the age of nuclei)
- The probability that a nucleus decays in the time Δt is p :

$$p = \alpha \Delta t \quad (\text{per } \alpha \Delta t \ll 1)$$

- Let's consider a system initially having N_0 unstable nuclei: how does the number of nuclei vary with time?

Simulation of a radioactive decay of a single nucleus

- The algorithm which has to be implemented is the following

LOOP from $t=0$ to T , step Δt

Reset the number of decayed nuclei (N_{dec}) in the current time bin

LOOP over each remaining parent nucleus (from 0 to N_{tot})

Decide if the nucleus decays:

if($\text{gRandom} \rightarrow \text{Rndm}() < \alpha \Delta t$)

increment the number of decayed nuclei in this time bin

endif

END loop over nuclei

$N_{\text{tot}} = N_{\text{tot}} - N_{\text{dec}}$

Plot N_{tot} vs t

END loop over time

END

Esercitazione 10 – Esercizio 2 (Decay.C)

- Write a macro to implement the algorithm shown in the previous slide. Show the number of remaining nuclei as a function of time for the two following cases:
 - $N_0=1000$, $\alpha=0.01 \text{ s}^{-1}$, $\Delta t=1 \text{ s}$
 - $N_0=50000$, $\alpha=0.03 \text{ s}^{-1}$, $\Delta t=1 \text{ s}$
- Show the results in linear and logarithmic scale, for t between 0 and 300 seconds
- Compare the results with the expected result

$$dN = -N\alpha dt$$



$$N = N_0 e^{-\alpha t}$$

Possible Solution

- In the next slide a possible solution implemented using a macro in ROOT is shown
- The ROOT classes are used to generate random numbers, to store informations in the histograms and for input/output operations
- The macro (decay.C) is composed by two functions
- It can be interpreted or executed by CLANG (CINT)


```
#if !defined(_CINT_) || defined(_MAKECINT_)
#include <TF1.h>
#include <TFile.h>
#include <TH1D.h>
#include <TMath.h>
#include <TRandom3.h>
#include <Riostream.h>
#endif
```

```
// Declare function
Double_t exponential(Double_t *x, Double_t *par);
//
void Decay(Int_t n0 = 50000, Double_t alpha = 0.03, Double_t Delt = 1.0, Double_t ttot = 300, Int_t seed = 95689){

gRandom->SetSeed(seed);
Int_t Nbins = static_cast<Int_t>(ttot/Delt); // number of time intervals
cout << "Numbersof bins: "<<Nbins<<" di ampiezza "<<Delt<<" s";
Double_t timetot = Delt*Nbins; // totale time = ttot
cout<<" Tempo totale "<<timetot<<endl;
// histogram booking
TH1D *h1 = new TH1D("h1","Remaining nuclei",Nbins+1,-Delt/2.,timetot+Delt/2.);
h1->SetFillColor(kOrange-4);
//Theoretical function
TF1 *fteo = new TF1("fteo",exponential,0.,timetot,2);
fteo->SetLineColor(kRed);
Double_t N0 = n0;
Double_t ALPHA = alpha;
fteo->SetParameters(N0,ALPHA);
fteo->SetParNames("normalizzazione","coefficiente");

Double_t prob = alpha*Delt; //probability
h1->Fill(0.,static_cast<double>(n0));
for(Double_t time=Delt; time<timetot+Delt/2.; time+=Delt){
    Int_t ndec = 0;
    for(Int_t nuclei=0; nuclei<n0;nuclei++)if(gRandom->Rndm()<prob)ndec++;
    n0-=ndec;
    h1->Fill(time,static_cast<double>(n0));
}

TFile *file = new TFile("decay.root","recreate");
h1->Write();
fteo->Write();
h1->Draw("histo");
fteo->Draw("same");
file->Close();
}

//
Double_t exponential(Double_t *x, Double_t *par){
    Double_t xx = x[0];
    return par[0]*exp(-par[1]*xx);
}
```

Header files, need to compile the macro

```

#if defined(__CINT__) || defined(__MAKECINT__)
#include <TF1.h>
#include <TFile.h>
#include <TH1D.h>
#include <TMath.h>
#include <TRandom3.h>
#include <Riostream.h>
#endif

// Declare function
Double_t exponential(Double_t *x, Double_t *par);
//
void Decay(Int_t n0 = 50000, Double_t alpha = 0.03, Double_t Delt = 1.0, Double_t ttot = 300, Int_t seed = 95689){

  gRandom->SetSeed(seed);
  Int_t Nbins = static_cast<Int_t>(ttot/Delt); // number of time intervals
  cout << "Numbersof bins: "<<Nbins<<" di ampiezza "<<Delt<<" s";
  Double_t timetot = Delt*Nbins; // totale time = ttot
  cout<<" Tempo totale "<<timetot<<endl;
  // histogram booking
  TH1D *h1 = new TH1D("h1","Remaining nuclei",Nbins+1,-Delt/2.,timetot+Delt/2.);
  h1->SetFillColor(kOrange-4);
  //Theoretical function
  TF1 *fteo = new TF1("fteo",exponential,0.,timetot,2);
  fteo->SetLineColor(kRed);
  Double_t N0 = n0;
  Double_t ALPHA = alpha;
  fteo->SetParameters(N0,ALPHA);
  fteo->SetParNames("normalizzazione","coefficiente");

  Double_t prob = alpha*Delt; //probability
  h1->Fill(0.,static_cast<double>(n0));
  for(Double_t time=Delt; time<timetot+Delt/2.; time+=Delt){
    Int_t ndec = 0;
    for(Int_t nuclei=0; nuclei<n0;nuclei++)if(gRandom->Rndm()<prob)ndec++;
    n0-=ndec;
    h1->Fill(time,static_cast<double>(n0));
  }

  TFile *file = new TFile("decay.root","recreate");
  h1->Write();
  fteo->Write();
  h1->Draw("histo");
  fteo->Draw("same");
  file->Close();
}

//
Double_t exponential(Double_t *x, Double_t *par){
  Double_t xx = x[0];
  return par[0]*exp(-par[1]*xx);
}

```

Functions are declared before their implementation. Default values can be Passed as "default" argument of the function

```

#if !defined(_CINT_) || defined(_MAKECINT_)
#include <TF1.h>
#include <TFile.h>
#include <TH1D.h>
#include <TMath.h>
#include <TRandom3.h>
#include <Riostream.h>
#endif

// Declare function
Double_t exponential(Double_t *x, Double_t *par);
//
void Decay(Int_t n0 = 50000, Double_t alpha = 0.03, Double_t Delt = 1.0, Double_t ttot = 300, Int_t seed = 95689){

gRandom->SetSeed(seed);
Int_t Nbins = static_cast<Int_t>(ttot/Delt); // number of time intervals
cout << "Numbersof bins: "<<Nbins<<" di ampiezza "<<Delt<<" s";
Double_t timetot = Delt*Nbins; // totale time = ttot
cout<<" Tempo totale "<<timetot<<endl;
// histogram booking
TH1D *h1 = new TH1D("h1","Remaining nuclei",Nbins+1,-Delt/2.,timetot+Delt/2.);
h1->SetFillColor(kOrange-4);
//Theoretical function
TF1 *fteo = new TF1("fteo",exponential,0.,timetot,2);
fteo->SetLineColor(kRed);
Double_t N0 = n0;
Double_t ALPHA = alpha;
fteo->SetParameters(N0,ALPHA);
fteo->SetParNames("normalizzazione","coefficiente");

Double_t prob = alpha*Delt; //probability
h1->Fill(0.,static_cast<double>(n0));
for(Double_t time=Delt; time<timetot+Delt/2.; time+=Delt){
    Int_t ndec = 0;
    for(Int_t nuclei=0; nuclei<n0;nuclei++)if(gRandom->Rndm()<prob)ndec++;
    n0-=ndec;
    h1->Fill(time,static_cast<double>(n0));
}

TFile *file = new TFile("decay.root","recreate");
h1->Write();
fteo->Write();
h1->Draw("histo");
fteo->Draw("same");
file->Close();
}

```

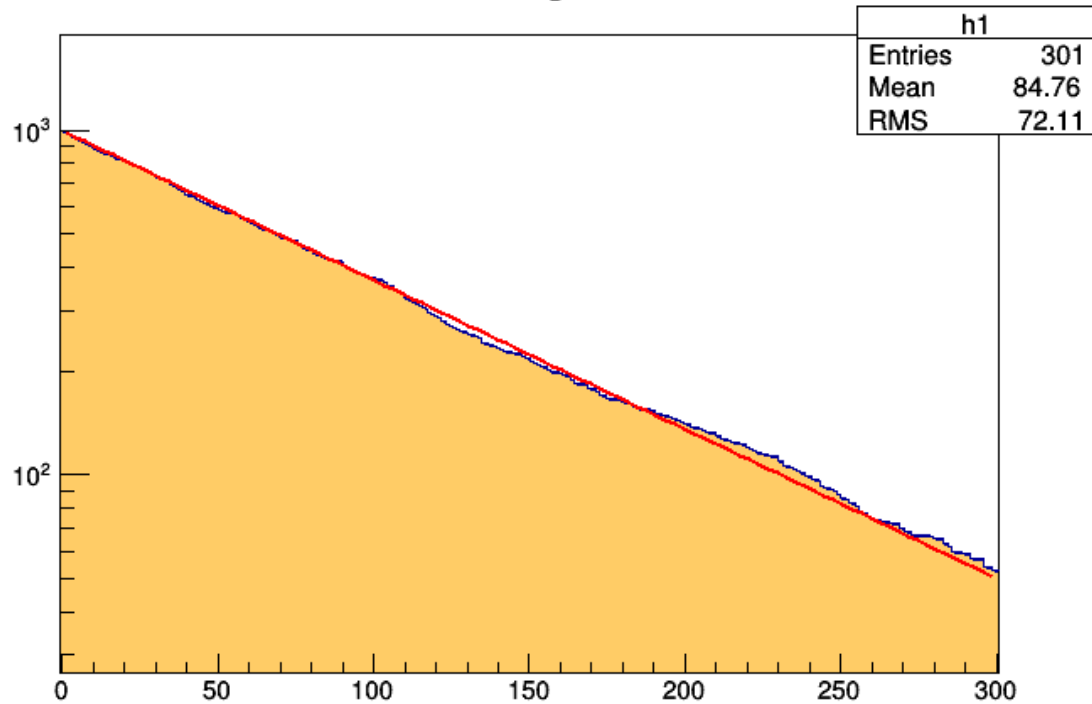
Definition of the "exponential function"
That can be used in the TF1 definition

```

//
Double_t exponential(Double_t *x, Double_t *par){
    Double_t xx = x[0];
    return par[0]*exp(-par[1]*xx);
}

```

Remaining nuclei

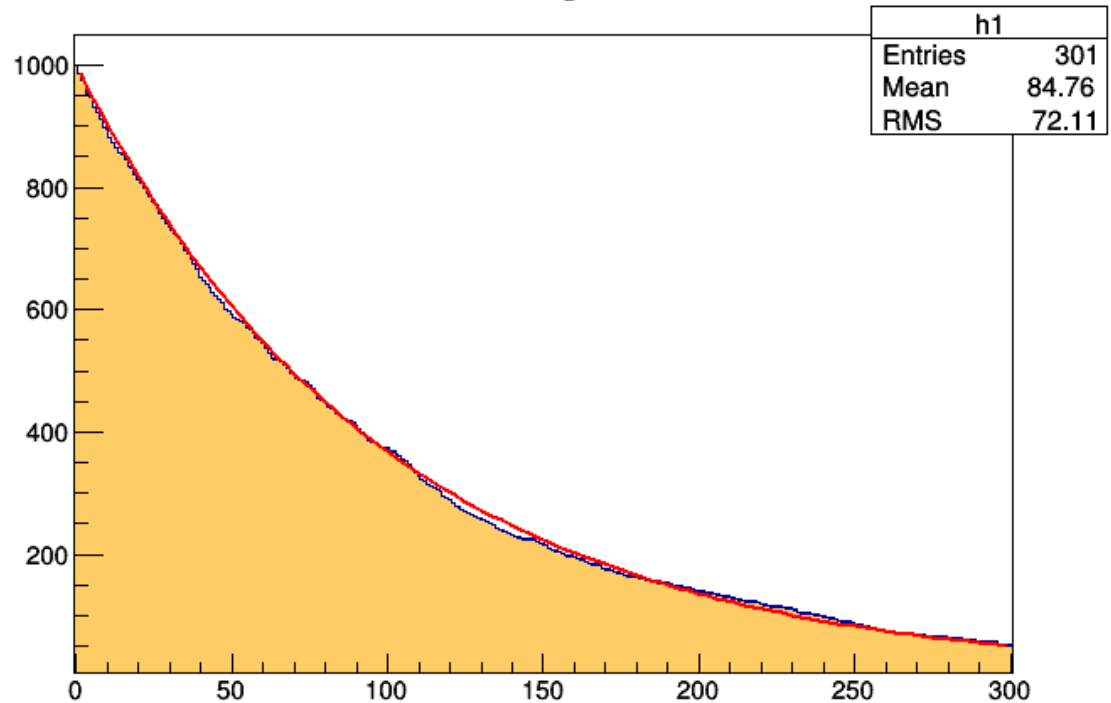


Continuous line: expected exponential.
Histogram: simulation result.

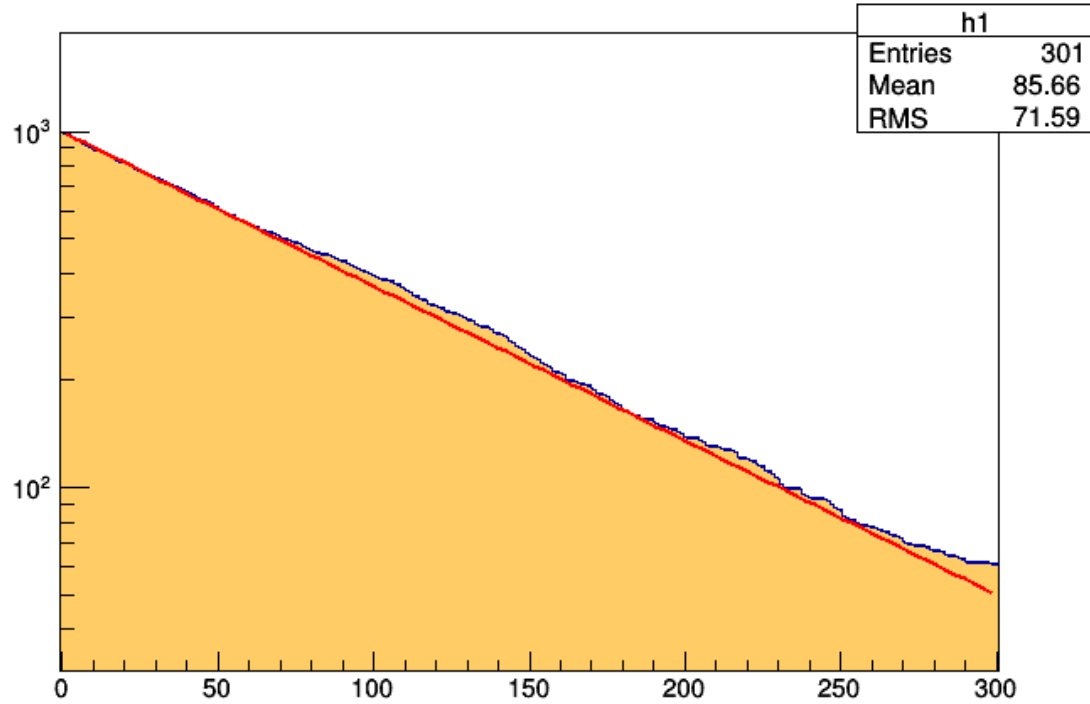
Result for:
 $N_0=100$,
 $\alpha=1 \times 10^{-2} \text{ s}^{-1}$
 $\Delta t=1 \text{ s}$

Statistical fluctuation are very important

Remaining nuclei



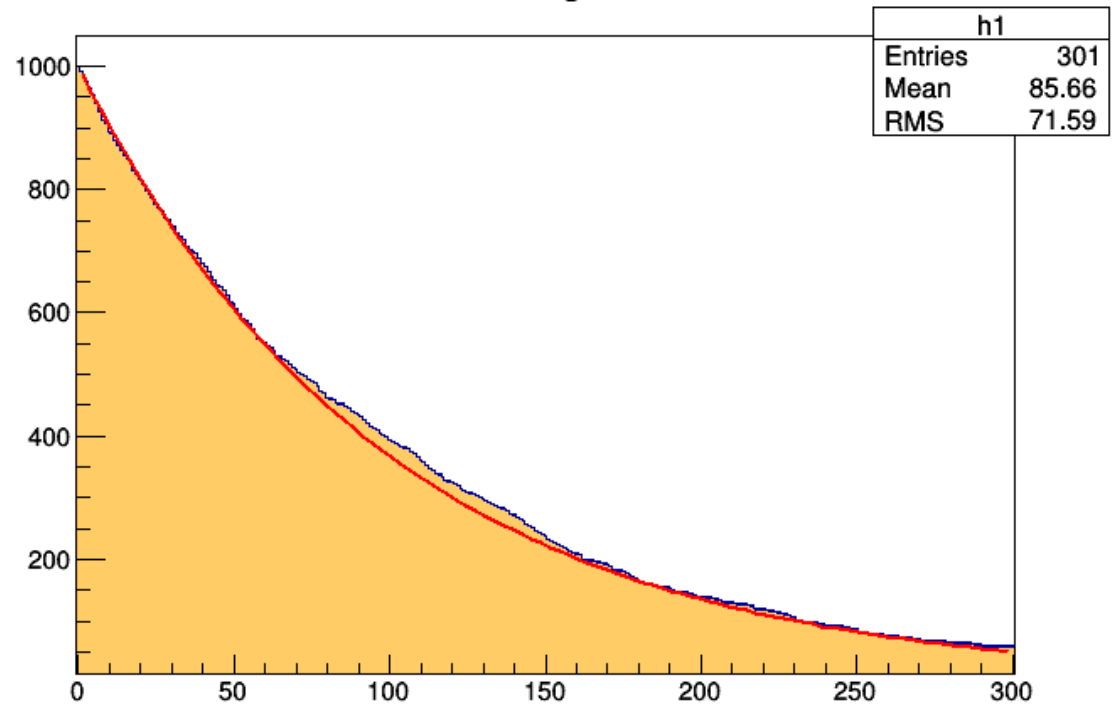
Remaining nuclei



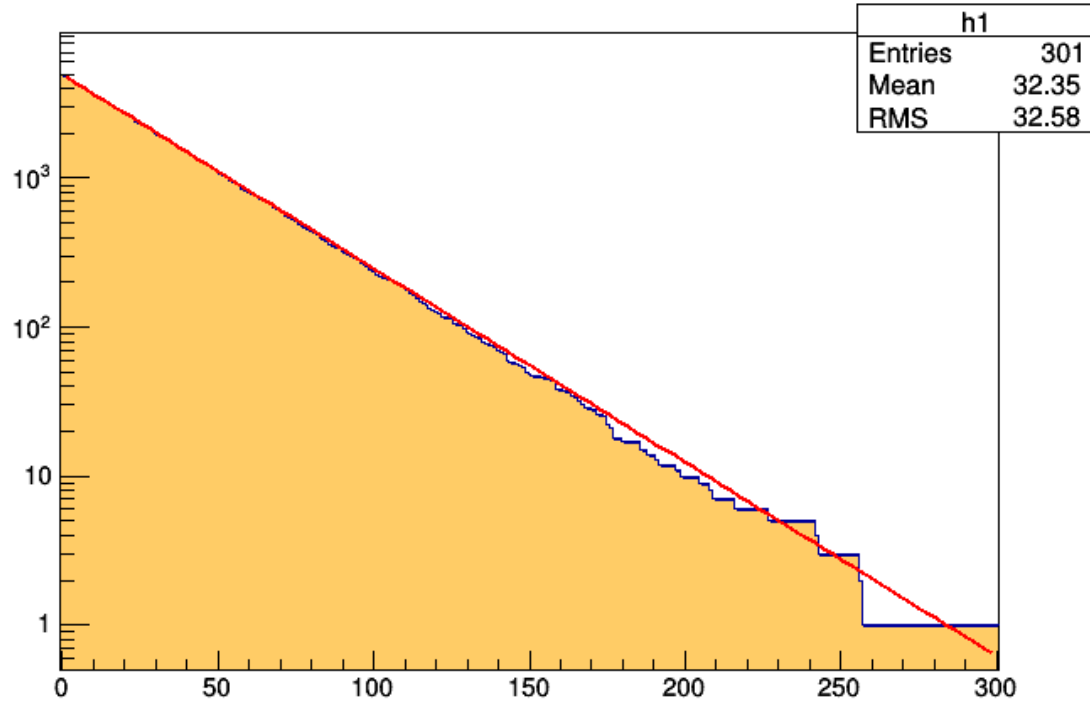
Same parameters as before,
but different seed: the results
are different!

Risultato per:
 $N_0=100$,
 $\alpha=1 \times 10^{-2} \text{ s}^{-1}$
 $\Delta t=1 \text{ s}$

Remaining nuclei



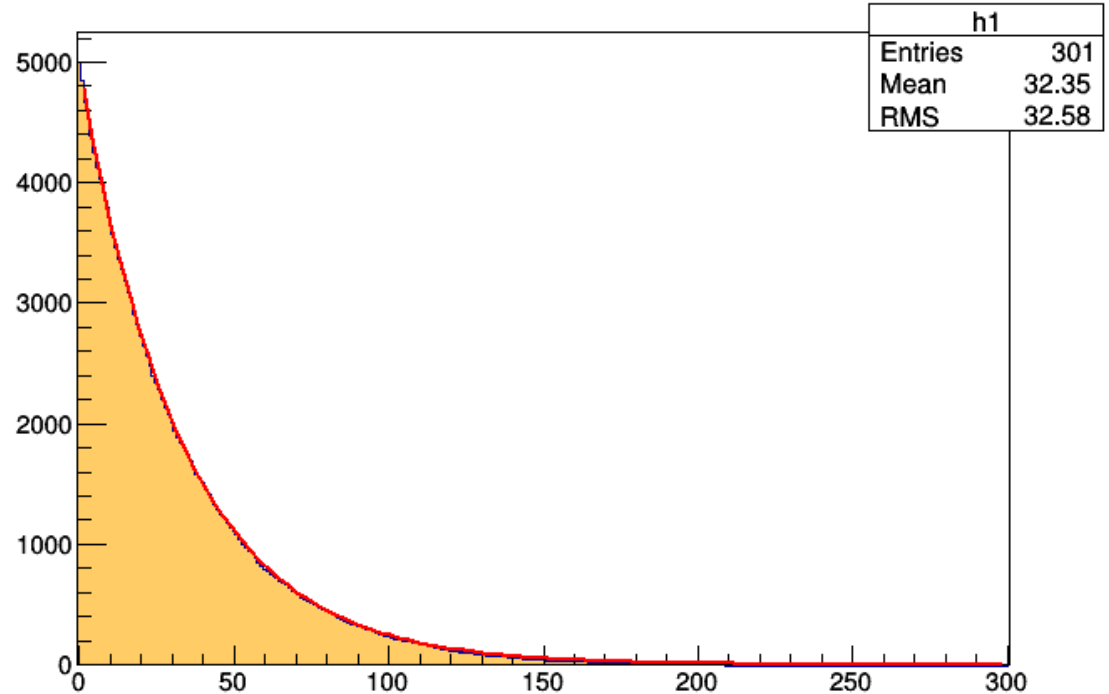
Remaining nuclei



The importance of fluctuations depends on the number N of residual nuclei

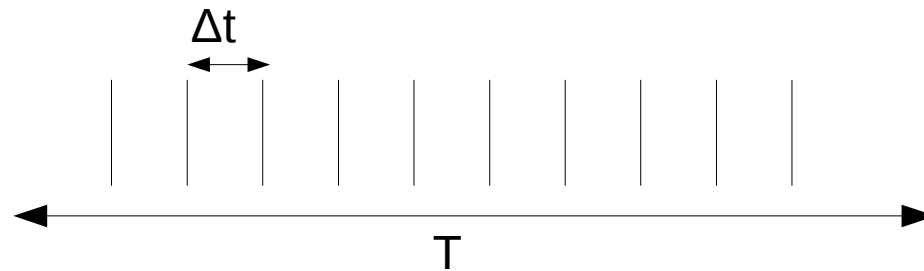
Result for:
 $N_0=5000$,
 $\alpha=3\times 10^{-2} \text{ s}^{-1}$
 $\Delta t=1 \text{ s}$

Remaining nuclei



How many decays in a fixed time interval T?

- Let's assume the number of decays in time T is much less than the number of parent nuclei. (i.e. let's assume a constant probability to observe a decay)
- The time T can be broken into m shorter intervals of duration Δt ($\Delta t/m$):



- The probability to observe 1 decay in time Δt is :

$$p = \beta \Delta t$$

How many decays in a fixed time interval T?

$$p = \beta \Delta t$$

- $\beta = \alpha N$ (Note that α is the decay probability)
- Δt must be small enough so that $\beta \Delta t \ll 1$
- The probability of observing n decays in time T is given by the binomial distribution:

$$P(\text{n decays in } m \text{ intervals}) = \binom{m}{n} p^n (1-p)^{(m-n)}$$

$$P = \frac{m!}{n!(m-n)!} \left(\frac{\beta T}{m}\right)^n \left(1 - \frac{\beta T}{m}\right)^{(m-n)}$$

How many decays in a fixed time interval T?

- In the limit of large m ($m \rightarrow \infty$ and $\Delta t \rightarrow 0$)

$$\left. \begin{aligned} \left(1 - \frac{\beta T}{m}\right)^m &\rightarrow e^{-\beta T} \\ \left(1 - \frac{\beta T}{m}\right)^{-n} &\rightarrow 1 \\ \frac{m!}{(m-n)!} &\rightarrow m^n \end{aligned} \right\}$$

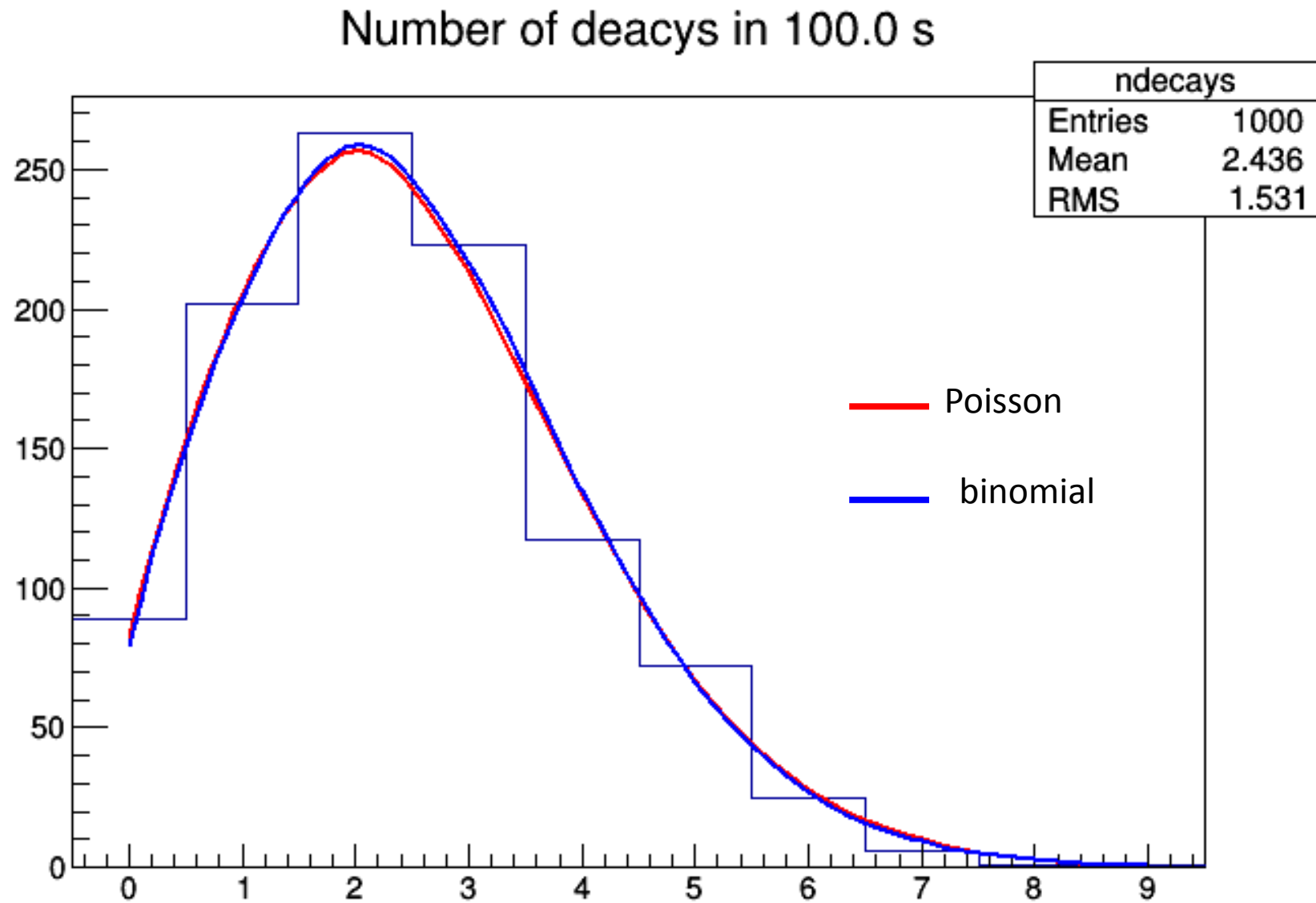
$$P(n; \beta T) = \frac{(\beta T)^n}{n!} e^{-\beta T} = \frac{\mu^n e^{-\mu}}{n!}$$

Known as Poisson distribution

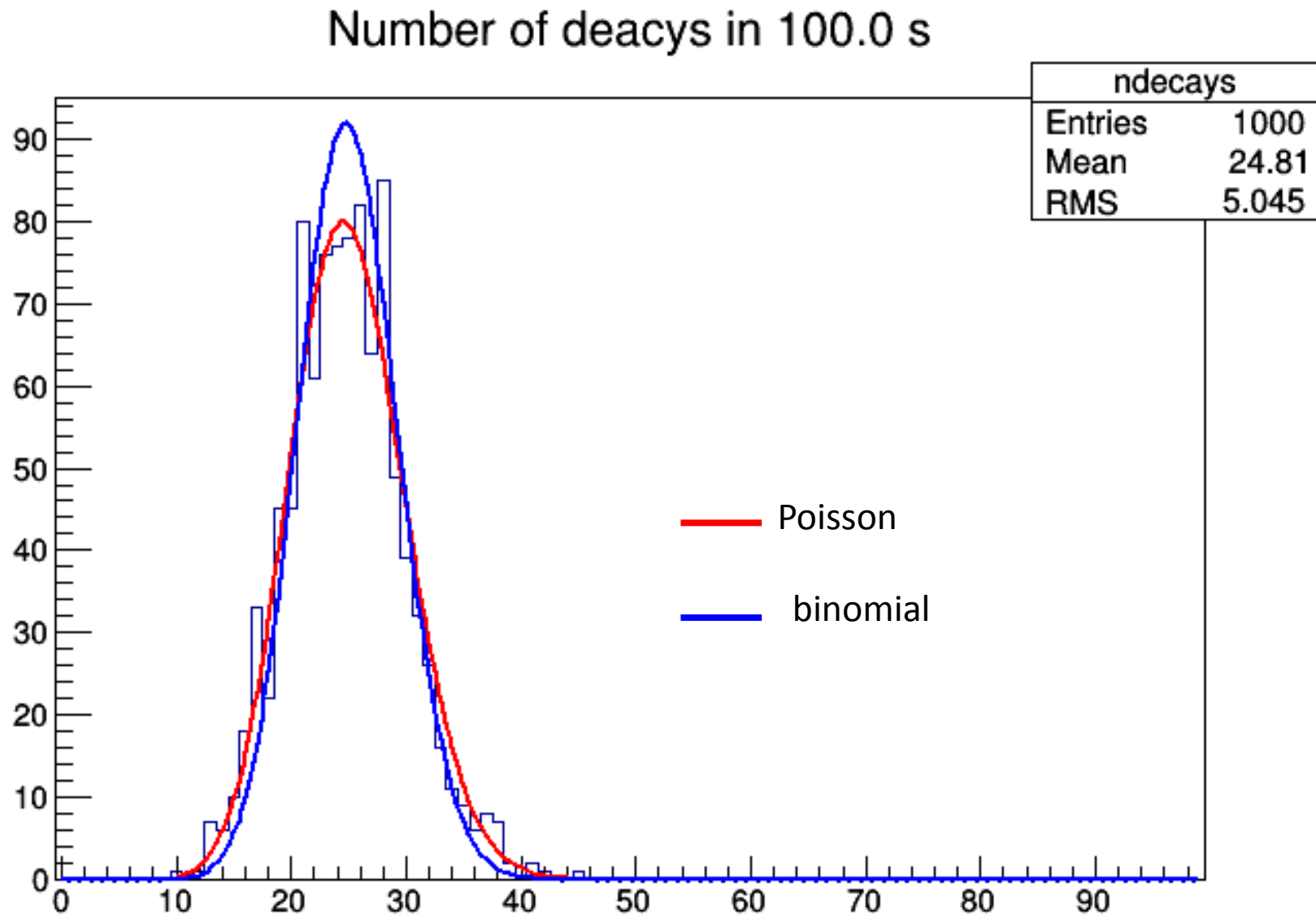
Esercitazione 10 – Esercizio 3 (poisson.C)

- Modify the program written for exercise 2 to simulate an experiment that counts the number of decays observed in a time interval, T .
- Allow the experiment to be repeated and histogram the distribution of the number of decays for the following two cases:
 $N_0=1000$, $\alpha=2.5 \times 10^{-5} \text{ s}^{-1}$, $\Delta t=1 \text{ s}$, $T = 100 \text{ s}$
 $N_0=1000$, $\alpha=2.0 \times 10^{-4} \text{ s}^{-1}$, $\Delta t=1 \text{ s}$, $T = 100 \text{ s}$
- Compare the distribution of n for 1000 experiments with Binomial and Poisson distributions.

Result for: $N_0=1000$, $\alpha=2.5 \times 10^{-5}$, $\Delta t=1$ s, $T=100$ s



Result for: $N_0=1000$, $\alpha=2.5 \times 10^{-4}$, $\Delta t=1$ s, $T=100$ s



```

// Function declaration
double Decay(int n0,double alpha, double Delt,double timetot); // Decay simulation

double poissonian(double xx, double norm, double p); // POISSONIAN density
double binomial(double xx, double norm, double ntrials, double prob); // BINOMIAL density
void poisson(int n0 = 1000,double alpha = 2.5e-5, double Delt = 1.,double time=100.,unsigned int seed = 1234); //main function

////////////////////////////////////
void poisson(int n0,double alpha, double Delt,double time,unsigned int seed){

    gRandom->SetSeed(seed);

    TString title = Form("Number of deacys in %4.1f s",time);
    int nbins = static_cast<int>(n0*alpha*time*4);
    cout<<"numer of bins"<<nbins<<endl;
    double high = nbins-0.5;
    TH1F *ndecays = new TH1F("ndecays",title,nbins,-0.5,high);

    //theoretical function (poissonian)
    double mu = alpha*n0*time;
    title =Form("Poissonian distribution with parameter %10.4g ",mu);

    TH1F *fteo = new TH1F("fteo",title,nbins,-0.5,high);

    //theoretical function 2(binomial)
    double ntrials = time/Delt;
    double prob = alpha*n0*Delt;
    title = Form("Binomial with parametris p= %10.4g e m = %8.2g",prob,ntrials);

    TH1F *fteo2 = new TH1F("fteo2",title,nbins,-0.5,high);

    cout<<"Numer of nuclei: " <<n0<<endl;
    cout<<"Parameter of poissonian distribution " <<mu<<endl;
    cout<<"Parameters of binomial distribution p= " <<prob<<" m =" << ntrials<<endl;

    // Fill histo with theoretical functions
    for (double x=0.;x<high;x+=1.)
        fteo->Fill(x,poissonian(x,knorm,mu));
    for (double x=0.;x<high;x+=1.)
        fteo2->Fill(x,binomial(x,knorm,ntrials,prob));

    // Simulation of the decay
    for(int exper = 0; exper<knorm; exper++){
        if((exper%100)==0)cout<<"Processing event n. " <<exper<<endl;
        double nodecay=Decay(n0,alpha,Delt,time);
        ndecays->Fill(nodecay);
    }
}

```

```

////////////////////////////////////
double Decay(int n0,double alpha, double Delt,double timetot){
    double prob = alpha*Delt; //probabilita
    int n0init = n0;
    for(double time=0.; time<timetot; time+=Delt){
        int ndec = 0;
        for(int nuclei=0; nuclei<n0;nuclei++)if(gRandom->Rndm()<prob)ndec++;
        n0-=ndec;
    }
    return static_cast<double>(n0init-n0);
}
////////////////////////////////////
double poissonian(double xx, double norm, double p){
    return norm*(TMath::PoissonI(xx,p));
    // return norm*(TMath::Power(param,xx)*TMath::Exp(-param-Gamma(xx+1)));
}
////////////////////////////////////
double binomial(double xx, double norm, double ntrials, double prob) {

    Double_t n = TMath::Binomial(ntrials,xx);
    n *= TMath::Power(prob,xx)*TMath::Power((1-prob),(ntrials-xx));
    n *= norm;
    return n;
}

```