# Introduction to Monte Carlo methods and its applications to Particle Physics

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#### Overview

- The Monte Carlo method
- How to install Geant4
- Some basic features of C++
- A short introduction to Geant4

For these slides I took inspiration from:

- M. Asai (SLAC, Stanford)
- A. Dotti (SLAC, Stanford)
- S. Incerti (CNRS, Bordeaux)
- L. Pandola (INFN-LNS, Catania)
- C. Pistillo (LHEP, Bern)
- S. Rahatlou (Sapienza, Roma)
- <u>www.cplusplus.com</u>

 You can download examples and slides from: <u>http://www.roma1.infn.it/~mancinit/</u> <u>Teaching/LabFisNucl-2019/</u>

The Monte Carlo trolley, or FERMIAC, was invented by Enrico Fermi and constructed by Percy King. The drums on the trolley were set according to the material being traversed and a random choice between fast and slow neutrons. Another random digit was used to determine the direction of motion, and a third was selected to give the distance to the next collision. The trolley was then operated by moving it across a two-dimensional scale drawing of the nuclear device or reactor assembly being studied.

The trolley drew a path as it rolled, stopping for changes in drum settings whenever a material boundary was crossed. This infant computer was used for about two years to determine, among other things, the change in neutron population with time in numerous types of nuclear systems.

> Image from "The Beginning of Monte Carlo Method, N. Metropolis 1987

# Introduction to the Monte Carlo method

The "FERMIAC"

## Monte Carlo methods

- It is a mathematical approach using a sequence of random numbers to solve a problem
- Stochastic quantities (e.g.: average value of *mm* of rain)
- Deterministic problems (definite integral)
- Generate N random "points"  $\vec{x_i}$  in the problem space

• Calculate 
$$\langle f \rangle = \frac{1}{N} \sum_{i=1}^{N} f(\vec{x_i})$$
 and  $\langle f^2 \rangle = \frac{1}{N} \sum_{i=1}^{N} f^2(\vec{x_i})$ 

## Monte Carlo methods

- Comte de Buffon (1777): needle tossing experiment to calculate the  $\pi$ ;
- Laplace (1886): random points in a rectangle to calculate  $\pi$ ;
- Fermi (1930): random approach to calculate the properties of the newly discovered neutron;
- Manhattan project (40's): simulations during the initial developments of thermonuclear weapons;
- Von Neumann and Ulam coined the term 'Monte Carlo' (1949);
- Exponential growth of the electronic computers (40's-60's);
- Berger (1963): first complete coupled electron-photon transportation code 'ETRAN'.

#### How to calculate an integral

- Randomly choose couple of numbers  $(x_i, y_i)$  from the range and the domain, respectively, of the function f
- The fraction of points where  $y_i \leq f(x_i)$  is equal to the fraction of the area below the function
- Technique proposed by Von Neumann, known as the "acceptance-rejection method"
- It is used to generate random numbers for an arbitrary Probability Density Function (PDF)

#### Example of an integral

• What if you have to calculate the integral of a function as:

$$f(x) = 1.4 \left[ \sin^4(x) \cos^2\left(\frac{x}{3}\right) + \sin^6\left(\frac{x}{4}\right) \right] \exp\left(-\frac{x}{8}\right)$$



# Example of an integral

- Using the acceptance-rejection method
- The orange points are the accepted (*n*)
- The blue are the rejected
   (N = accepted + rejected)



#### Let's calculate $\pi$

$$f(x) = \sqrt{(1 - x^2)}$$

$$A_{circ} = \int_{-1}^{1} f(x) dx = \pi \frac{r^2}{2}$$

$$A_{rect} = \Delta y \cdot \Delta x$$

$$\frac{n}{N} \propto \frac{A_{circ}}{A_{rect}} = \frac{\pi r^2/2}{\Delta x \cdot \Delta y}$$

$$\pi \approx 2\frac{n}{N}\frac{\Delta x \cdot \Delta y}{r^2} = 4\frac{n}{N}$$



## Let's calculate $\pi$

- Is there a way to speed up the convergence of the computation?
- Use the symmetry!

 This is the method for calculating π was proposed by Laplace in "Théorie Analytique des Probabilitiés" (1825)!





#### Use the simmetry!



#### Inverse transform sampling

- If a PDF f is integrable (called **cumulative**, F)
- and the cumulative is invertible  $F^{-1}$
- It is possible to sample x accordingly to f:

$$x = F^{-1}(u)$$

where *u* is uniformly distributed

## Random Numbers Generators

- At the core of all Monte Carlo calculations is some mechanism to produce a long sequence of random numbers r<sub>i</sub> that are uniformly distributed over the open interval [0, 1)
- Digital computers, by design, are incapable of producing random results
- A true random sequence could, in principle, be obtained by coupling to our computer some external device that would produce a truly random signal
- However, use of such a random number generator would not be practical!
- Imposible to debug!



### Pseudo-random Number Generators

 Such a generator is a deterministic algorithm that, given the previous numbers (usually just the last number) in the sequence, the next number can be efficiently calculated

$$x_{n+1} = f(x_n, x_{n-1}, ..., x_0)$$

- *x*<sub>0</sub> is called "**seed**"
- A unique seed returns a unique random number sequence
- It is important to use a **new seed every time** that a random selection is initiated
- A typical error is the use of the same seed for multiple generation, which leads to the generation of the same sample of random numbers

## Pseudo-random Number Generators

- Because the set of numbers directly representable in the computer is finite, the sequence will necessarily repeat
- The length of the sequence prior to beginning to repeat is called **period**
- Many pseudo-random number generators have been proposed and used over the years in a wide variety of Monte Carlo work.
- Designing better random number generators (and test them) is still an active area of research

# Bad PRNG: RANDU

- Linear congruential PRNG used since '60
- Was the most widely used random number generator in the world
- Developed by IBM
- Three-dimensional plot of 100,000 values generated with RANDU
- Each point represents 3 consecutive pseudorandom values



not only the period is important!

#### Simple case: decay in flight

- Suppose a  $\pi^+$  with momentum p
- The life time is a random value with a pdf

$$f(t) = \frac{1}{\tau} \exp\left(-\frac{1}{\tau}\right)$$

• Therefore, *t* can be sampled from the inverse of the cumulative:

$$t = F^{-1}(r) = -\tau \ln(1 - r)$$
$$r \in [0, 1)$$

## Simple case: decay in flight

• Select the decay channel:

Mode		Fraction ( $\Gamma_i / \Gamma$ )		
$\Gamma_1$	$\mu^+ u_\mu$	[1]	(99.98770 ± 0.00004)%	
$\Gamma_2$	$\mu^+ u_\mu\gamma$	[2]	$(2.00 \pm 0.25) \times 10^{-4}$	
$\Gamma_3$	$e^+\nu_e$	[1]	$(1.230 \pm 0.004) \times 10^{-4}$	[table from PDG]
$\Gamma_4$	$e^+ \nu_e \gamma$	[2]	$(7.39 \pm 0.05) \times 10^{-7}$	
$\Gamma_5$	$e^+  u_e \pi^0$		$(1.036 \pm 0.006) \times 10^{-8}$	
$\Gamma_6$	$e^+\nu_e e^+e^-$		$(3.2 \pm 0.5) \times 10^{-9}$	
$\Gamma_7$	$e^+\nu_e\nu\overline{\nu}$		$< 5 \times 10^{-6}$	

- In the CM frame the decay is isotropic  $\theta \in [0,\pi); \quad \phi \in [0,2\pi)$
- Finally, Lorentz-boost in the Lab. frame
- 4 random numbers for one decay!

#### Problem

- Why did I sampled  $\theta$  and  $\phi$  in the CM frame?
- What if I sample uniformly  $\theta \in [0, 2\pi)$ ;  $r \in [0, 1)$ ?
- The extracted points gather in the centre

 A uniform distribution in polar coordinates is not uniform in orthogonal coordinate system



- It is the most common application of MC in Particle Physic
- Assume that all the possible interactions are known
- The distance *s* between two subsequent interactions is distributed as  $p(s) = \mu \exp(-\mu s)$

• Being  $\mu$  a property of the medium

- $\mu$  is proportional to the probability of an interaction per unit length, therefore:
  - is proportional to the **total cross section**  $\mu = N\sigma = N\sum_{i} \sigma_{i} = \sum_{i} \mu_{i}$
  - μ<sub>i</sub> are the partial cross section of all the competing processes
  - depends on the **density** of the material
     (N is the number of scattering centres in the medium)

- Divide the particle trajectory in "steps"
  - Straight free-flight tracks along the step
  - Could be limited by geometry boundaries
- Sampling the step length accordingly to  $\,p(s)\,$
- Sampling the interaction at the end of the step
- Sampling the interaction accordingly to  $\,\mu_i/\mu$
- Sampling the final state using the physics model of the interaction i
  - Update the properties of the primary particle
  - Add the possible secondaries produced (to be tracked later)

- Follow all secondaries, until absorbed (or leave the geometry)
- $\mu$  depends on the energy (cross sections do!)



#### Tracking, not so easy...

- This basic recipe doesn't work well for charged particles
- The cross sections of some processes (ionisation and bremsstrahlung) is very high, so the steps would be very small
- In each interaction only a small fraction of energy is lost and the effect on the particle are small
- A lot of CPU time used to simulate many interactions having small effects

# The solution: approximate

- Simulate explicitly interactions only if the energy loss is above a threshold  $E_0$  (**hard** interactions)
  - Detailed simulation
- The effects of all sub-threshold interactions is described cumulatively (**soft** interactions)
- Hard interactions occur much less frequently than soft interactions

#### Flowchart of an event



...luckily enough, somebody else already implemented the tracking algorithms for us (and much more)



# A short introduction to Geant4

A sketch of the ATLAS MC simulation

# Geant4 (GEometry ANd Traking)

- Developed by an International Collaboration
  - Established in 1998
  - Approximately 100 members, from Europe, US and Japan
  - <u>http://geant4.org</u>
- Open source
- Written in C++ language
  - Takes advantage from the Object Oriented software technology

[Geant4, a simulation toolkit Nucl. Inst. and Methods Phys. Res. A, 506 250-303

Geant4 developments and applications Transaction on Nuclear Science 53, 270-278]



## Geant4 applications

- Physics experiments
- but also:
  - Hadrontherapy
  - Radiobiology
  - and many others...

atomistic view of a dinucleosome irradiated by a single 100 keV proton Image from M. A. Bernal et al Physica Medica, vol. 31, no. 8, pp. 861–874, Dec. 2015.





# Geant4, further applications

- Radio-protection in space mission
- Shielding for satellites
- Single event upset and radiation damages to electronics
- Simulations for nuclear spallation sources
- Radioactive waste



First slide of the talk "ESA Geant4 R&D Activities from the Geant4 Space User Workshop Hiroshima, 26 August 2015





Figure from M. Sawant, COTS Journal Jan. 2012



#### How to install Geant4

some tips

#### Check on website

• You can find the most recent tips on:

http://www.roma1.infn.it/~mancinit/?action=3Teaching/ Suggestions on Geant4 installation



#### Prerequisites on CentOS7



- sudo yum update
- sudo yum groupinstall "Development Tools"
- sudo yum install xerces-c xerces-c-devel
- sudo yum install centos-release-scl-rh
- sudo yum install qt qt-x11 qt-devel
- optional:
  - sudo yum install devtoolset-7 scl enable devtoolset-7 bash
  - sudo yum install python27 python27-pyhton-pip python27-pyhton-tools python27-numpy

# CMake3 for CentOS7



- cmake is needed to compile and use Geant4
- the version available on the CentOS7 repository is too old, we have to compile it
- wget https://github.com/Kitware/CMake/releases/ download/v3.13.2/cmake-3.13.2.tar.gz
- tar -xvzf cmake-3.13.2.tar.gz
- cd cmake-3.13.2
- ./bootstrap && gmake
- Add to ~/.bashrc the line: export PATH=/home/soft/cmake-3.13.2/bin:\$PATH

### Prerequisites on RedHat



- sudo yum update
- sudo yum groupinstall "Development Tools"
- sudo yum install xerces-c xerces-c-devel libXmu libXmu
- sudo yum install cmake
- sudo yum install qt qt-x11 qt-devel
- optional:
  - sudo yum install devtoolset-7 scl enable devtoolset-7 bash
  - sudo yum install python27 python27-pyhton-pip python27-pyhton-tools python27-numpy


- sudo apt update
- sudo apt install build-essential git
- sudo apt install cmake
- sudo apt install qt4-default qt4-devtools libxerces-c-dev

# Prerequisites on Mac Os X



- install XCode and type in a terminal: xcode-select -install
- install XQuartz (<u>https://www.xquartz.org/</u>)
- install brew: /usr/bin/ruby -e "\$(curl -fsSL https:// raw.githubusercontent.com/Homebrew/install/ master/install)"
- brew doctor; brew update
- brew install cmake xerces-c wget
- brew install carlomt/repo/qt04



- wget <a href="http://cern.ch/geant4-data/releases/geant4.10.05.tar.gz">http://cern.ch/geant4-data/releases/</a>
- tar -xvzf geant4.10.05.tar.gz
   this will create geant4.10.05/
- mkdir geant4.10.05-build
- mkdir geant4.10.05-install
- mkdir geant4-data
- cd geant4.10.05-build

# Compile and install



- cmake -DCMAKE\_BUILD\_TYPE=RelWithDebInfo

   DGEANT4\_INSTALL\_DATA=ON
   DGEANT4\_INSTALL\_DATADIR=../geant4-data
   DGEANT4\_BUILD\_MULTITHREADED=ON
   DGEANT4\_USE\_GDML=ON -DGEANT4\_USE\_QT=ON
   DGEANT4\_BUILD\_CXXSTD=c++11
   DCMAKE\_INSTALL\_PREFIX=../geant4.10.05-install
   ../geant4.10.05
- make -j `nproc`
- make install
- Add to your .bashrc (~/.bash\_profile on a Mac): export G4DIR=~/geant4.10.05-install
   \$G4DIR/bin/geant4.sh
   alias g4make='cmake -DGeant4\_DIR=\$G4DIR'

\$("#word-list-out").e(" "); meru-stu(s, par h(); Var c = l(), a = " ", d = parseInt(\$("#limit\_val").a()), f = parseInt(\$("slid").a()) function("rand:" + f); d < f && (f = d, function("check rand\u00f3\u00f3rand: \* f + "tops: \* d))</pre> var n = [], d = d - f, e;**if (0 < c.length)** { for (var g = 0;g < c.length;g++) {</pre> e = m(b, c[g]), -1 < e && b.splice(e, 1);</pre> b.unshift({use\_wystepuje:"parameter", word:c[g]}); for (g = 0;g < c.length;g++) {</pre>

C++

Some basic features of

70

17

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127

228

229

[slides made getting inspiration from http://www.cplusplus.com]

### Just an introduction

- This is not a C++ course
- Just few information useful to understand the Geant4 examples
- For a complete course: <u>http://www.roma1.infn.it/people/rahatlou/programmazione++/</u>

### Few things about C++

- A general-purpose programming language
- Has imperative, **object-oriented** and generic programming features
- Provides facilities for low-level memory manipulation
- In 1983, "C with Classes" was renamed to "C++" (++ being the increment operator in C)
- Initially standardised in 1998 (current standard is C++17 but the most used is C++11)

### Classes

 Classes are an expanded concept of data structures: like data structures, they can contain data members, but they can also contain functions as members

Like Plato's ideas (the idea of apple), classes have generic attributes (e.g. color). Each instance (this Golden Delicious apple) of the class have a specific attribute (e.g. yellow)

class Apple {
 public:
 void setColor(color);
 color getColor();

private: color fColor; double fWeight;

### Example of class usage



### Example of class usage



### Example of class usage

```
#include <iostream>
using std::cout;
```

```
class Rectangle {
    int width, height;
    public:
        void set_values (int,int);
        int area() {return width*height;}
};
void Rectangle::set_values (int x, int y)
{
    width = x;
    height = y;
}
```

```
int main () {
   Rectangle rect;
   rect.set_values (3,4);
   cout << "area: " << rect.area();</pre>
```

Hyperuranion (ὑπερουράνιος τόπος) literally: "place beyond heaven"





# What if I want to protect the rectangle properties (the dimensions), once instantiated?



### Constructors

```
#include <iostream>
using std::cout;
class Rectangle {
    int width, height;
  public:
    Rectangle(int x, int y);
    int area() {return width*height;}
};
Rectangle::Rectangle(int x, int y)
{
  width = x;
  height = y;
}
int main () {
  Rectangle rect(3,4);
  cout << "area: " << rect.area();</pre>
  return 0;
```

Using the constructor and removing the setting method

### Constructors

```
#include <iostream>
using std::cout;
class Rectangle {
    int width, height;
  public:
    Rectangle(int x, int y);
    int area() {return width*height;}
};
Rectangle::Rectangle (int x, int y) :
width(x), height(y) { }
int main () {
  Rectangle rect(3,4);
  cout << "area: " << rect.area();</pre>
  return 0;
}
```

Better implementation!

### Inheritance

- Classes in C++ can be extended, creating new classes which retain characteristics of the base class
- This process, known as inheritance, involves a base class and a derived class
- The derived class inherits the members of the base class, on top of which it can add its own members



### Inheritance, an example

```
class Polygon {
  protected:
    int width, height;
  public:
    void set_values (int a, int b)
    { width=a; height=b; }
};
```

```
class Rectangle: public Polygon
{
   public:
    int area ()
    {
      return width*height;
   }
};
```

```
class Triangle: public Polygon
{
   public:
    int area()
    {
      return width*height/2;
   }
};
```

### Protected and not private!

- The protected access specifier used in class Polygon is similar to private. Its only difference occurs in fact with inheritance:
- When a class inherits another one, the members of the derived class can access the protected members inherited from the base class, but not its private member
- By declaring width and height as protected instead of private, these members are also accessible from the derived classes
   Rectangle and Triangle, instead of just from members of
   Polygon
- If they were public, they could be accessed just from anywhere

### Let's use the classes...

```
#include <iostream>
using std::cout;
using std::endl;
int main () {
   Rectangle rect;
   Triangle trgl;
   rect.set_values (4,5);
   trgl.set_values (4,5);
   cout << rect.area() << endl;
   cout << trgl.area() << endl;
   return 0;
}</pre>
```

have a look at the example <u>https://github.com/carlomt/inheritance\_example</u> for more details

# CMake



- a cross-platform free and open-source software application for managing the build process of software using a compiler-independent method
- supports directory hierarchies and multiple libraries
- can locate executables, files, and libraries
- <u>https://cliutils.gitlab.io/modern-cmake/</u>
- use a version of CMake that came out after your compiler
- since CMake will dumb itself down to the minimum required version in your CMake file, installing a new CMake, even system wide, is pretty safe



Lets get back to Geant4

# Geant4 is a toolkit

- Geant4 is a toolkit (= a collection of tools)
  - i.e. you **cannot** run it out of the box
  - You must write an application, which uses Geant4
- Consequences:
  - There are **no** such concepts as "Geant4 **defaults**"
  - You must provide the necessary information to configure your simulation
  - · You must deliberately choose which Geant4 tools to use
- Guidance: many examples are provided
  - **Basic Examples**: overview of Geant4 tools
  - Advanced Examples: Geant4 tools in real-life applications

#### Reminder...



### Geant4 way of tracking

- Force step ending at geometry boundaries
- All AlongStep processes co-occur
- The PostStep compete, i.e.: only one is selected



### Geant4 way of tracking

- If particle is at rest chose one of the AtRest processes
- The secondaries are saved in the stack
- To be further tracked with a last in first out approach



# You MUST:

- Describe your experimental set-up
- Provide the primary particles input to your simulation
- Decide which particles and physics models you want to use out of those available in Geant4 and the precision of your simulation (cuts to produce and track secondary particles)



### You may also want to:

- Interact with Geant4 kernel to control your simulation
- Visualise your simulation configuration or results
- Produce histograms, tuples etc. to be further analysed





### Jargons

- Run, event, track, step, step point
- Process
  - At rest, along step, post step
- Cut = production threshold
- Sensitive detector, score, hit, hits collection,



# A run in Geant4

- As an analogy of the real experiment, a run of Geant4 starts with "Beam On"
- A run is a collection of events which share the same detector and physics conditions
- G4RunManager class manages
   processing a run



G4UserRunAction is the optional user hook



 $\begin{aligned} \mathcal{I} &= -\frac{1}{4} F_{\mu\nu} F^{\mu\nu} \\ &+ i F \mathcal{D} \mathcal{V} + h.c. \\ &+ \mathcal{Y}_i \mathcal{Y}_{ij} \mathcal{Y}_j \mathcal{P} + h.c. \\ &+ |\mathcal{D}_{\mu} \mathcal{P}|^2 - V(\mathcal{P}) \end{aligned}$ 

### Physics processes

an overview...

### The G4VProcess

- All physics processes derive from G4VProcess
- G4VProcess is an abstract class
- It defines the common interface of all processes in Geant4
- Three kind of "actions":
  - AlongStep
     all the soft interactions
  - PostStep
     all the hard interactions
  - · AtRest

decays, e+ annihilation



# $\gamma$ model inventory

- Many models available for each process
- Differ for energy range, precision and CPU speed
- Final state generators

Model	Emin	Emax
G4LivermoreRayleighModel	100 eV	10 PeV
G4PenelopeRayleighModel	100 eV	10 GeV
G4KleinNishinaCompton	100 eV	10 TeV
G4KleinNishinaModel	100 eV	10 TeV
G4LivermoreComptonModel	100 eV	10 TeV
G4PenelopeComptonModel	10 keV	10 GeV
G4LowEPComptonModel	100 eV	20 MeV
G4BetheHeitlerModel	1.02 MeV	100 GeV
G4PairProductionRelModel	10 MeV	10 PeV
G4LivermoreGammaConversionModel	1.02 MeV	100 GeV
G4PenelopeGammaConversionModel	1.02 MeV	10 GeV
G4PEEFluoModel	l keV	10 PeV
G4LivermorePhotoElectricModel	10 eV	10 PeV
G4PenelopePhotoElectricModel	10 eV	10 GeV

# ElectroMagnetic models

- The same physics processes can be described by different models
- For instance: Compton scattering can be described by
  - G4KleinNishinaCompton
  - G4LivermoreComptonModel (low-energy, based on the Livermore database)
  - G4PenelopeComptonModel (low-energy, based on the Penelope analytical model)
  - G4LivermorePolarizedComptonModel (low-energy, Livermore database with polarization)
  - G4PolarizedComptonModel (Klein-Nishina with polarization)
  - G4LowEPComptonModel (full relativistic 3D simulation)
- Different models can be combined, so that the appropriate one is used in each given energy range (à performance optimization)

# EM Physics constructors

. . .

G4EmStandardPhysics – default G4EmStandardPhysics\_option1 – HEP fast but not precise G4EmStandardPhysics option2 – Experimental G4EmStandardPhysics option3 – medical, space G4EmStandardPhysics\_option4 – optimal mixture for precision G4EmLivermorePhysics G4EmLivermorePolarizedPhysics **Combined** Physics Standard > 1 GeV G4EmPenelopePhysics **LowEnergy** < 1 GeV G4EmLowEPPhysics G4EmDNAPhysics\_option...

 Advantage of using of these classes – they are tested on regular basis and are used for regular validation

# Hadronic processes

- At rest
  - Stopped muon, pion, kaon, anti-proton
  - Radioactive decay
  - Particle decay (decay-in-flight is PostStep)
- Elastic
  - Same process to handle all long-lived hadrons (multiple models available)

Inelastic

- Different processes for each hadron (possibly with multiple models vs. energy)
- Photo-nuclear, electro-nuclear, mu-nuclear
- Capture
  - Pion- and kaon- in flight, neutron
- Fission

### Hadronic physics challenge

- Three energy regimes
  - < 100 MeV</pre>
  - resonance and cascade region (100 MeV 10 GeV)
  - > 20 GeV (QCD strings)
- Within each regime there are several models
- Many of these are phenomenological

### Hadronic models

- Two families of builders for the high-energy part
  - **QGS**, or list based on a model that use the Quark Gluon String model for high energy hadronic interactions of protons, neutrons, pions and kaons
  - **FTF**, based on the FTF (FRITIOF like string model) for protons, neutrons, pions and kaons
- Three families for the cascade energy range
  - **BIC**, binary cascade
  - **BERT**, Bertini cascade
  - INCLXX, Liege Intranuclear cascade model
#### ParticleHP

- Data-driven approach for inelastic reactions for n (in place since many years, named NeutronHP) p, d, t, <sup>3</sup>He and  $\alpha$
- Data based on TENDL-2014 (charged particles)
   and ENDFVII.r1 (neutrons).
- For neutrons, includes information for elastic and inelastic scattering, capture, fission and isotope production
- Range of applicability: from thermal energies up to 20 MeV
- Very precise tracking, but also very slow
- Use it with care: thermal neutron tracking is very CPU-demanding

#### Haronic model inventory



http://geant4.cern.ch/support/proc\_mod\_catalog/models

### Physics List

- A class which collects all the particles, physics processes and production thresholds needed for your application
- It tells the run manager how and when to invoke physics
- It is a very flexible way to build a physics environment
  - user can pick the particles he wants
  - user can pick the physics to assign to each particle
- But, user must have a good understanding of the physics required
- Omission of particles or physics could cause errors or **poor simulation**

#### There is not default, but...

- Geant4 provides several "production physics lists" which are routinely validated and updated with each release these should be considered only as starting points which you may need to validate or modify for your application
- There are currently 19 packaged physics lists available
- 6 reference physics lists:
  - FTFP\_BERT, FTFP\_BERT\_HP QGSP\_BERT, QGSP\_BERT\_HP, QGSP\_BIC QGSP\_FTFP\_BERT

# Naming...

- The following acronyms refer to various hadronic options
  - QGS -> Quark Gluon String model (>~20 GeV) FTF -> Fritiof string model (>~5 GeV) BIC -> Binary Cascade (<~ 10 GeV) BERT -> Bertini-style cascade (<~ 10 GeV)</li>
- HP -> High Precision neutron model ( < 20 MeV) P -> G4Precompund model used for de-excitation
- EM options designated by
  - no suffix: standard EM physics
  - EMV suffix: older but faster EM processes
     other suffixes for other EM options

#### **Production Physics Lists**

- FTFP\_BERT
  - recommended by Geant4 for HEP
  - contains all standard EM processes
  - uses Bertini-style cascade for hadrons < 5 GeV</li>
  - uses FTF (Fritiof) model for high energies ( > 4 GeV)
- QGSP\_BERT
  - all standard EM processes
  - Bertini-style cascade up to 9.9 GeV
  - QGS model for high energies (> ~18 GeV) FTF in between

#### **Production Physics Lists**

- QGSP\_BIC
  - same as QGSP\_BERT, but replaces Bertini cascade with Binary cascade and G4Precompound model
  - recommended for use at energies below 200 MeV (many medical applications)
- FTFP\_BERT\_HP
  - same as FTFP\_BERT, but with high precision neutron model used for neutrons below 20 MeV
  - significantly slower than FTFP\_BERT when full thermal cross sections used there's an option to turn this of
  - for radiation protection and shielding applications

#### Other Physics Lists

- If primary particle energy in your application is < 5 GeV (for example, clinical proton beam of 150 MeV)
  - start with a physics list which includes BIC or BERT
  - e.g. QGSP\_BIC, QGSP\_BERT, FTFP\_BERT, etc.
- If neutron transport is important
  - start with physics list containing "HP"
  - e.g. QGSP\_BIC\_HP, FTFP\_BERT\_HP, etc.
- If you're interested in Bragg curve physics
  - use a physics list ending in "EMV" or "EMX" or "EMY"
  - e.g. QGSP\_BIC\_EMY

#### An event in Geant4

- An event is the basic unit of simulation in Geant4
- **G4Event** class represents an event. It has following objects at the end of its (successful) processing
  - List of primary vertices and particles (as input)
  - Hits and Trajectory collections (as output)
- G4EventManager class manages
   processing an event
- G4UserEventAction is the optional user hook



#### A track in Geant4

- Track is a snapshot of a particle
- It has physical quantities of current instance only. It does not record previous quantities
- It's not a collection of steps. Instead, a track is being updated by steps
- G4TrackingManager manages processing a track, a track is represented by G4Track class
- G4UserTrackingAction is the optional user hook

#### A step in Geant4

- A step is a variation of a track
- Has two points (pre and post step points)
- In case a step is limited by a boundary, the end point stands on the boundary, and it logically belongs to the next volume
- Boundary processes such as transition
   radiation or refraction could be simulated



- G4SteppingManager class manages processing a step, a step is represented by G4Step class
- G4UserSteppingAction is the optional user hook

#### Particle in Geant4

- A particle in Geant4 is represented by three layers of classes:
  - G4Track
    - Geometrical information (position)
  - · G4DynamicalParticle

- 000
- Dynamic physical properties (momentum, energy, spin...)
- G4ParticleDefinition
  - · Static properties (charge, mass, life time)



#### Sampling the step size

- In Geant4, particle transportation is a process as well, by which a particle interacts with geometrical volume boundaries and field of any kind
- Each particle has its own list of applicable processes. At each step, all processes listed are invoked to get proposed physical interaction lengths
- The process which requires the shortest interaction length limits the step

# Let's cut it out... (cuts in MC)

- The traditional Monte Carlo solution is to set a tracking cut-off in energy:
  - a particle is stopped when its energy goes below it



- its residual energy is deposited at that point
- Imprecise stopping and energy deposition location
- Particle and material dependence

# Let's cut it out... (cuts in Geant4)

- Geant4 does not have tracking cuts i.e.: all tracks are tracked down to 0 energy
- A Cut in Geant4 is a production threshold
- It is applied only for physics processes that have infrared divergence
  - Bremsstrahlung
  - Ionisation  $e^{-}$  ( $\delta$  rays)
  - Protons from hadronic elastic scattering



#### A range cut

- The threshold is a **distance**!
- Default = 1 mm
- Particles unable to travel at least the range cut value are not produced



- Sets the "spatial accuracy" of the simulation
- Production threshold is internally converted to an energy threshold for each material

# Cut in energy

- 460 keV
- good for LAr
- not for Pb

- 2 MeV
- good for Pb
- not for Lar



#### Cut in range

- 1.5 mm
- ~460 KeV in
- ~2 MeV in P

run with the hares and hunt with the hounds... (good for both!)

01 cm



# Setting the cuts

Optional method in G4VPhysicsList

```
void MyPhysicsList::SetCuts()
{
    //G4VUserPhysicsList::SetCuts();
    defaultCutValue = 0.5 * mm;
    SetCutsWithDefault();
    SetCutValue(0.1 * mm, "gamma");
    SetCutValue(0.01 * mm, "e+");
    G4ProductionCutsTable::GetProductionCutsTable()
        ->SetEnergyRange(100*eV, 100.*GeV);
}
```

- not all models are able to work with very low production thresholds
- an energy threshold limit is used,
- its default value is set to 990 eV.
- You can change this value

#### Cuts UI command

# Universal cut (whole world, all particles)
/run/setCut 10 mm

# Override low-energy limit
/cuts/setLowEdge 100 eV

# Set cut for a specific particle (whole world)
/run/setCutForAGivenParticle gamma 0.1 mm

# Set cut for a region (all particles)
/run/setCutForARegion myRegion 0.01 mm

# Print a summary of particles/regions/cuts
/run/dumpCouples

#### Cuts per region

- Complex detector may contain many different sub-detectors involving:
  - finely segmented volumes
  - position-sensitive materials (e.g. Si trackers)
  - · large, undivided volumes (e.g. calorimeters)
- The same cut may not be appropriate for all of these
- User can define regions (independent of geometry hierarchy tree) and
   assign different cuts for each region
- A region can contain a subset of the logical volumes

#### To limit the step

- To have more precise energy deposition
- To increase precision in magnetic field
- Include G4StepLimiter in your physics list
  - as a Physics process
  - compete with the others

# System of Units

- Internal unit system used in Geant4 is completely hidden not only from user's code but also from Geant4 source code implementation
- Each hard-coded number must be multiplied by its proper unit:



- radius = 10.0 \* cm; E = 1.\*GeV;
- To get a number, it must be divided by a proper unit:
  - G4cout<< "E dep: "<< eDep/MeV <<" [MeV]"<<G4endl;

#### User classes

- You have to write the main()
- Initialisation classes:
  - G4VUserDetectorConstruction
  - G4VUserPhysicsList
  - G4VUserActionInitialization
- Action classes
  - G4VUserPrimaryGeneratorAction
  - G4UserRunAction
  - G4UserEventAction
  - G4UserStackingAction
  - G4UserTrackinAction
  - G4UserSteppingAction



#### classes written in red are mandatory!

#### Your program

- Geant4 does not provide a main()
- In your main(), you have to
- Construct G4RunManager (sequential mode)
- Set user mandatory initialisation classes to RunManager
  - · G4VUserDetectorConstruction
  - · G4VUserPhysicsList
  - G4VUserActionInitialization
- You can define VisManager, (G)UI session,
- You can initialise optional user action classes

## Multi-threading in Giant



#### Parallel

# Multithread



- Speed-up of simulations even on a laptop
- More efficient usage of memory: all cores only read geometry and physics of the Geant4 simulation (no duplication)



# How to survive in the real life

Some tips, you will find your way...

#### The Geant4 website

- www.geant4.org
- Future and past events (e.g.: this course)
- User support

   (one of the best
   friends for Geant4
   users...)



#### The Geant4 website

- <u>www.geant4.org/</u> <u>geant4/support.2</u>
- Documentation
- Application
   Developers (you!)
   guide



#### Application Developers user guide



#### Application Developers user guide

- <u>http://geant4-userdoc.web.cern.ch/geant4-userdoc/</u> <u>UsersGuides/ForApplicationDeveloper/html/index.html</u>
- Introduces new Users to the Geant4 toolkit
- Describes the most useful tools
- Describes hoot set-up and run a simulation application
- Intended as an overview of the toolkit, not an exhaustive treatment

#### The Geant4 website

- www.geant4.org
- Documentation
- Application
   Developers (you!)
   guide
- Physics Reference Manual



#### Physics Reference Manual



#### Physics Reference Manual

- <u>http://geant4-userdoc.web.cern.ch/geant4-userdoc/</u> <u>UsersGuides/PhysicsReferenceManual/html/index.html</u>
- A reference for toolkit Users and developers who wish to consult and study the physics of an interaction/model
- Present the theoretical formulation, model or parameterisation of the physics interactions provided by Geant4

#### The Geant4 website

- www.geant4.org
- Documentation
- Application
   Developers (you!)
   guide
- Physics Reference Manual
- Physics Lists Guide



# Physics List Guide

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🖬 PhysicsL	istGuide	Contents:	
G	4	<ul> <li>Physics List Guide</li> <li>Bibliography</li> <li>Reference Physics Lists</li> <li>FTFP_BERT Physics List</li> <li>QBBC Physics List</li> <li>QGSP_BERT Physics List</li> <li>Shielding Physics List</li> </ul>	
10. Search docs		<ul> <li>EM physics constructors</li> <li>EM Opt0</li> <li>EM Opt1</li> <li>EM Opt2</li> </ul>	
CONTENTS: Physics List Guide Reference Physics Lists	ls	<ul> <li>EM Opt3</li> <li>EM Opt4</li> <li>EM Liv</li> <li>EM Pen</li> <li>EM GS</li> </ul>	
Electromagnetic phys	ics constructors	<ul> <li>EM SS</li> <li>EM DNA</li> <li>Tables by constructor</li> <li>Tables by particle</li> </ul>	
# Doxygen

- <u>http://geant4-</u> <u>userdoc.web.cern.ch/</u> <u>geant4-userdoc/</u> <u>Doxygen/</u> <u>examples\_doc/html/</u> <u>hierarchy.html</u>
- · All the class interfaces





#### Examples

an overview...

### Examples omnia divisa est in partes trees...

- Basic set of examples is oriented to novice users and covering the most typical use-cases of a Geant4 application with keeping simplicity and ease of use
- **Extended** set of examples may require some additional libraries besides of Geant4. This set covers many specific use cases for actual detector simulation
- Advanced set of examples covers the use-cases typical of a "toolkit"- oriented kind of development, where real complete applications for different simulation studies are provided; may require additional third party products to be built

## Where?

- Where to find the examples:
  - \$G4DIR/examples/basic
  - \$G4DIR/examples/extended
  - \$G4DIR/examples/advanced



## Basic examples

Code name	Few Characteristics
Example B1	<ul> <li>Simple geometry with a few solids</li> <li>Geometry with simple placements (G4PVPlacement)</li> <li>Scoring total dose in a selected volume user action classes</li> <li>Geant4 physics list (QBBC)</li> </ul>
Example B2	<ul> <li>Simplified tracker geometry with global constant magnetic field</li> <li>Geometry with simple placements (G4PVPlacement) and parameterisation (G4PVParameterisation)</li> <li>Scoring within tracker via G4 sensitive detector and hits</li> <li>Geant4 physics list (FTFP_BERT) with step limiter</li> <li>Started from novice/N02 example</li> </ul>
Example B3	<ul> <li>Schematic Positron Emitted Tomography system</li> <li>Geometry with simple placements with rotation (G4PVPlacement)</li> <li>Radioactive source</li> <li>Scoring within Crystals via G4 scorers</li> <li>Modular physics list built via builders provided in Geant4</li> </ul>
Example B4	<ul> <li>Simplified calorimeter with layers of two materials</li> <li>Geometry with replica (G4PVReplica)</li> <li>Scoring within layers in four ways: via user actions, via user own objects via G4 sensitive detector and hits and via scorers</li> <li>Geant4 physics list (FTFP_BERT)</li> <li>Histograms (ID) and ntuple saved in the output file</li> <li>Started from novice/N03 example</li> </ul>
Example B5	<ul> <li>A double-arm spectrometer with wire chambers, hodoscopes and calorimeters with a local constant magnetic field</li> <li>Geometry with placements with rotation, replicas and parameterisation</li> <li>Scoring within wire chambers, hodoscopes and calorimeters via G4 sensitive detector and hits</li> <li>Geant4 physics list (FTFP_BERT) with step limiter</li> <li>UI commans defined using G4GenericMessenger</li> <li>Histograms (ID, 2D) and ntuple saved in the output file</li> <li>Started from extended/analysis/A01</li> </ul>



## Basic examples

	Code name	Few Characteristics
Basic!	Example BI	<ul> <li>Simple geometry with a few solids</li> <li>Geometry with simple placements (G4PVPlacement)</li> <li>Scoring total dose in a selected volume user action classes</li> <li>Geant4 physics list (QBBC)</li> </ul>
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A bit complex	Example B3	<ul> <li>Schematic Positron Emitted Tomography system</li> <li>Geometry with simple placements with rotation (G4PVPlacement)</li> <li>Radioactive source</li> <li>Scoring within Crystals via G4 scorers</li> <li>Modular physics list built via builders provided in Geant4</li> </ul>
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# My example

- <u>https://github.com/</u> <u>carlomt/AnaEx01</u>
- A modified version of extended/AnaEx01
- output in root files
- input in GPS
- example of cosmic muons...



#### Courses

- Geant4 Advanced Course CERN (Geneva)
   26-27 March 2019
   <u>https://indico.cern.ch/event/789510/</u>
- Geant4 Course at the 16th Seminar on Software for Nuclear, Sub-nuclear and Applied Physics
   Porto Conte (Alghero)
   26-31 May 2019
   <a href="https://agenda.infn.it/event/17240/">https://agenda.infn.it/event/17240/</a>

thank you for your attention!