



# Physics II: Overview, Processes, Production Threshold, Regions & Cuts per region

Geant4 PHENIICS & IN2P3 Tutorial,  
16 – 20 May 2022,  
Orsay

Marc Verderi  
LLR, Ecole polytechnique





# Credits...

- › Daniel Brandt, Makoto Asai, Dennis Wright (SLAC),
- › Gunter Folger (CERN), etc.
- › Ivana,
- › ...

# Introduction

- › In this presentation, we will give an overview of the physics catalog of Geant4
- › We will explain also how physics processes
  - are modeled, in term of C++ class
  - and how they are used during the tracking
- › This is for “cultural” aspects, as in general, you will not have to worry about the structure of the process class
  - Because you will be using existing physics lists
  - Which setup properly processes and take care of technical details
- › We will speak also about “cuts”, which, at the opposite are something you must care about.
  - They must be defined for some physics processes
    - › to run the simulation under time  $< \infty$
  - “Cuts” is actually a bad usage name...
  - ... and the issue is « production threshold »
  - But some care has to be taken in defining them



# Outline

## I. Physics Overview

- Overview of Geant4 physics capability

## II. Processes

- How physics processes are modeled in Geant4

## III. Production Thresholds (aka cuts)

## IV. Regions

## V. Cuts per region

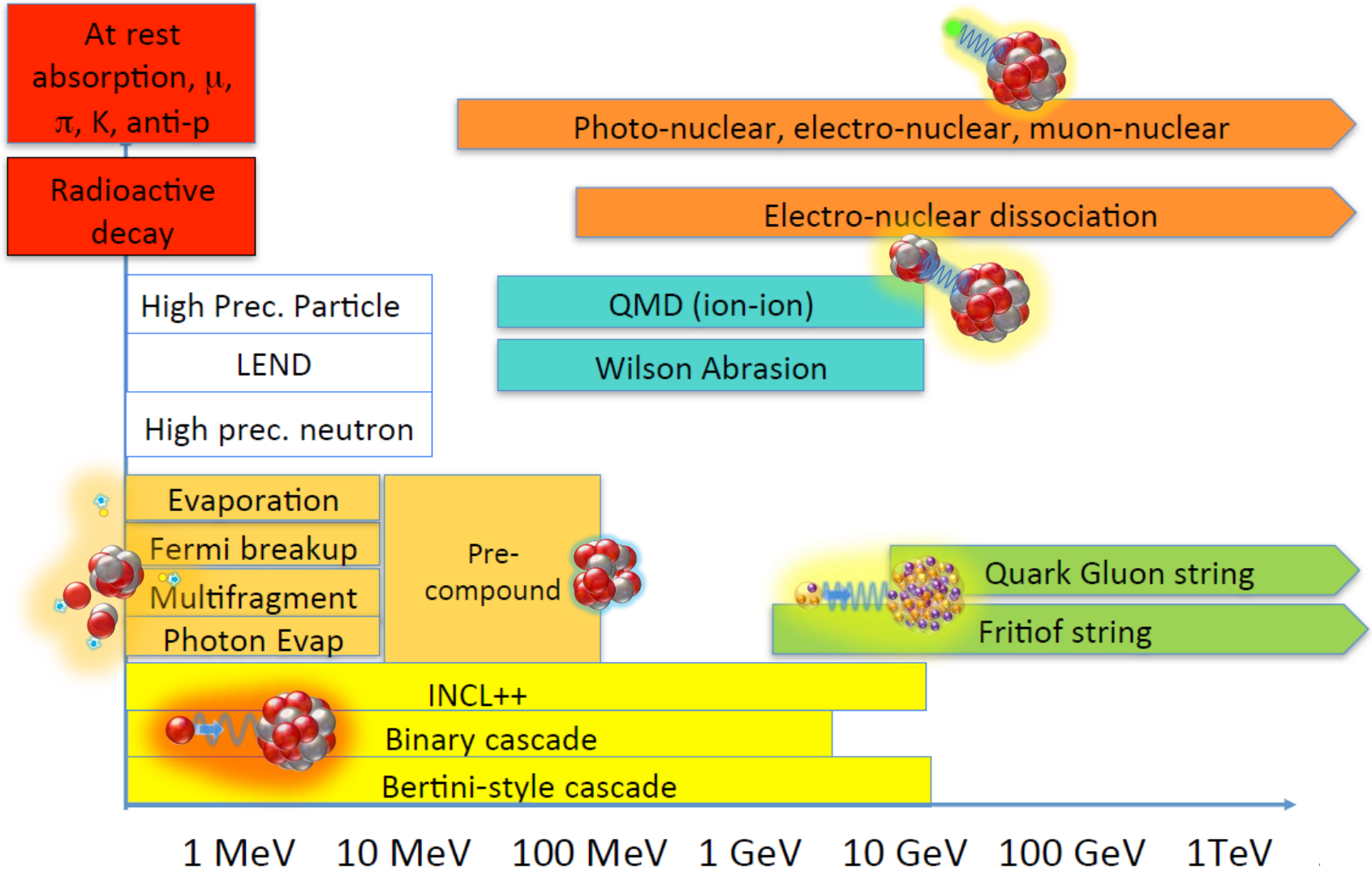


# I. Physics Overview

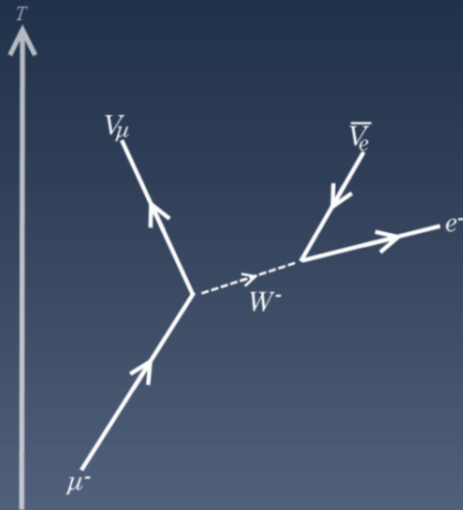
# Electromagnetic Processes

- › **Standard** : Complete set of processes covering charged particles and gammas.
  - Energy range 1 keV - ~PeV
- › **Low energy** : More precise description at low energy for  $e^+$ ,  $e^-$ ,  $\gamma$ , charged hadrons incident particle.
  - More atomic shell structure detail
  - Some processes valid down to hundreds of eV
  - Some processes not valid above 1 GeV
- › **DNA & MuElec** : for microdosimetry studies
  - Processes down to a few eV (!)
  - Plus chemistry stage for DNA
- › **Optical photon** : Long wavelength  $\gamma$  (X-ray, UV, visible)
  - Reflection, refraction, absorption, wavelength shifts, Rayleigh scattering
- › **Phonons** : under development. Acoustic phonons, for now. Suited for low-temperature (tens of mK) detectors.

# Hadronic Processes



# Decay & « Technical »

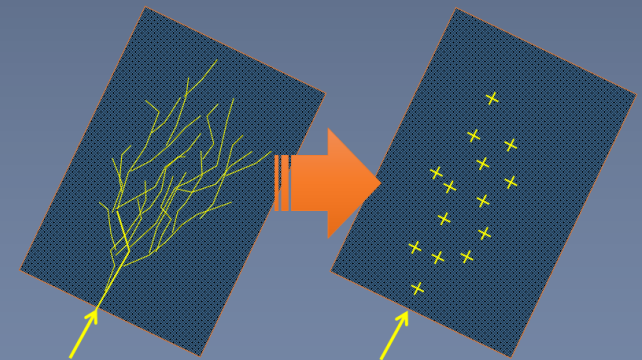


## › Decay processes

- Decay of particles of width narrow enough
  - › i.e. : exclude hadronic resonances
- weak decay (leptonic decays, semi-leptonic decays, radioactive decay of nuclei)
- electromagnetic decay (e.g.  $\pi^0$ ,  $\Sigma^0$ )

## › « Technical » processes:

- Processes without physics content but which act as interfaces for:
- Parameterization
  - › Fast Simulation fonctionnalité
  - › Hook to shortcut the detailed tracking
- Parallel geometries
  - › Limit the step on parallel geometry boundaries / switch tracking geometries
- Scoring
  - › Collect user requested information
- Biasing
  - › Modify physics behavior wrt to the reference standard one



Fast simulation : a full shower is replaced by a parametrized version of it





## II. Processes

How Geant4 models processes

# G4VProcess: 3 kind of actions (1/2)

- › Abstract class defining the common interface of **all processes** in Geant4:
  - Used by all « physics » processes
  - but is also used by the transportation, etc...
  - Defined in **source/processes/management**

- › Define **three kinds of actions**:

- **AtRest** actions:

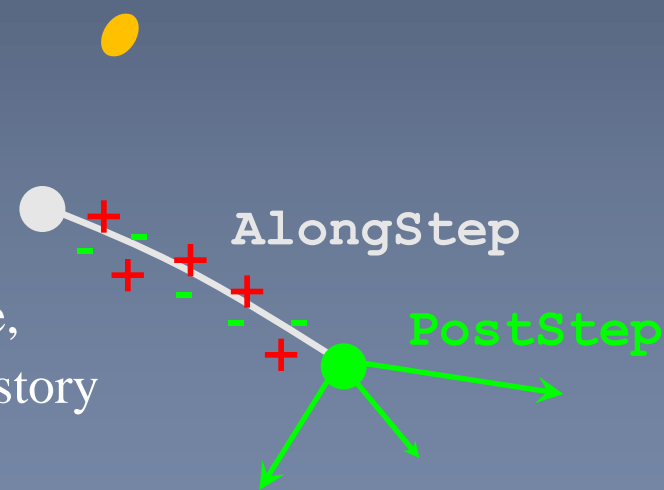
- Decay,  $e^+$  annih., absorption ...

- **AlongStep** actions:

- To describe continuous (inter)actions, occurring along the path of the particle, like ionisation; used for condensed history

- **PostStep** actions:

- For describing point-like (inter)actions, like decay in flight, hard radiation...



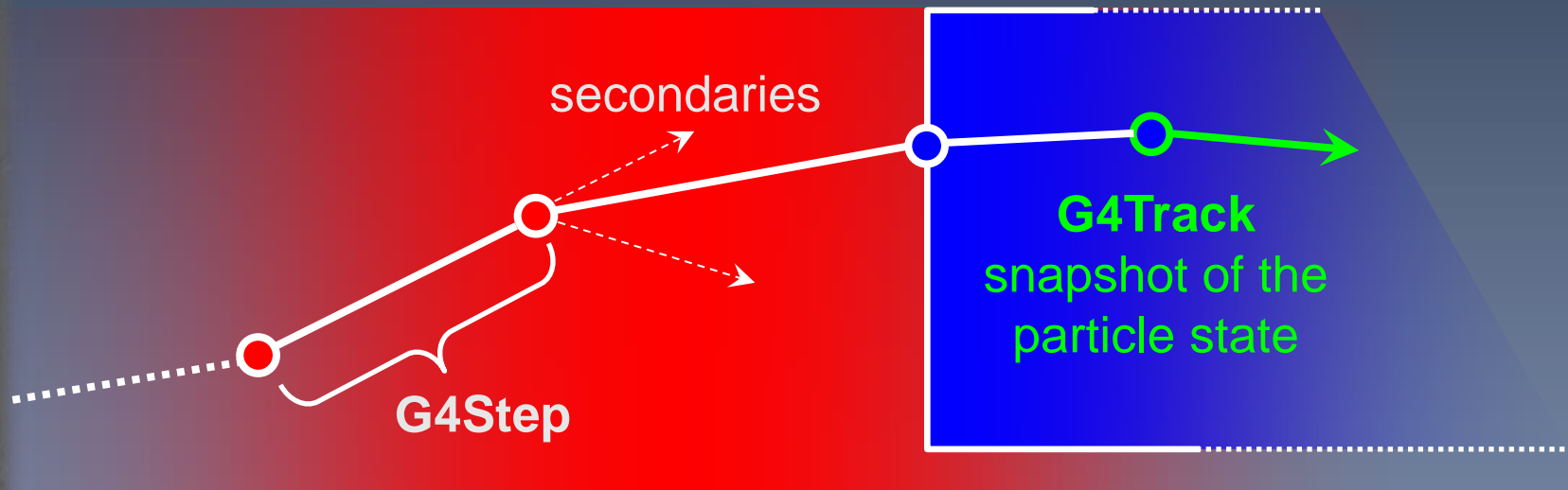
# G4VProcess: 3 kind of actions (2/2)

- › A process can implement **any combination** of the three **AtRest**, **AlongStep** and **PostStep** actions:
  - eg: decay = **AtRest** + **PostStep**
- › Each action defines **two methods**:
  - **GetPhysicalInteractionLength ()**:
    - › Used to **limit the step**:
      - either because the process « triggers » an interaction, a decay
      - or any other reasons, like fraction of energy loss, geometry boundary, user's limit ...
  - **DoIt ()**:
    - › Implements the **actual action** to be applied on the track;
    - › And the related production of secondaries.

# How Geant4 uses the processes during tracking ?

› Remember:

– Tracks are moved step by step :

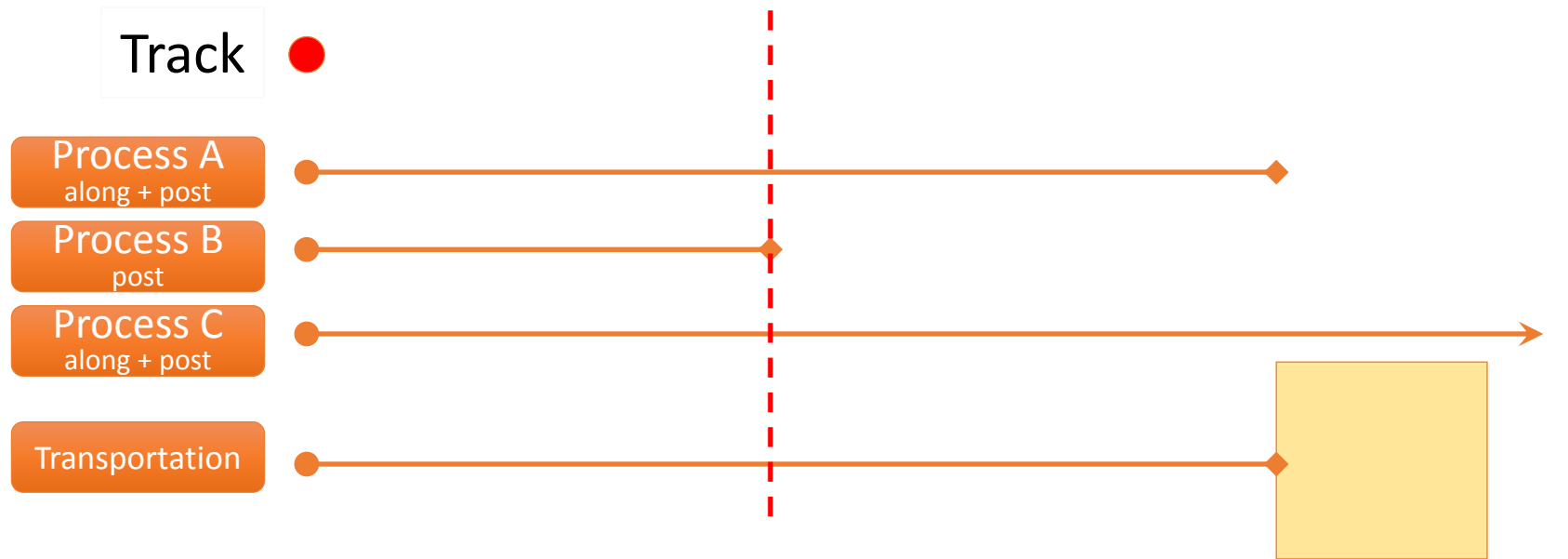


– And several processes are attached to the track

› So what happens during a step ?

# Process Handling by the Stepping

1. At the beginning of the step, the step length is determined:
  - Consider all processes attached to the current `G4Track`;
  - Define the step length as the smallest of the lengths among:
    - > All `AlongStepGetPhysicalInteractionLength()`
    - > All `PostStepGetPhysicalInteractionLength()`

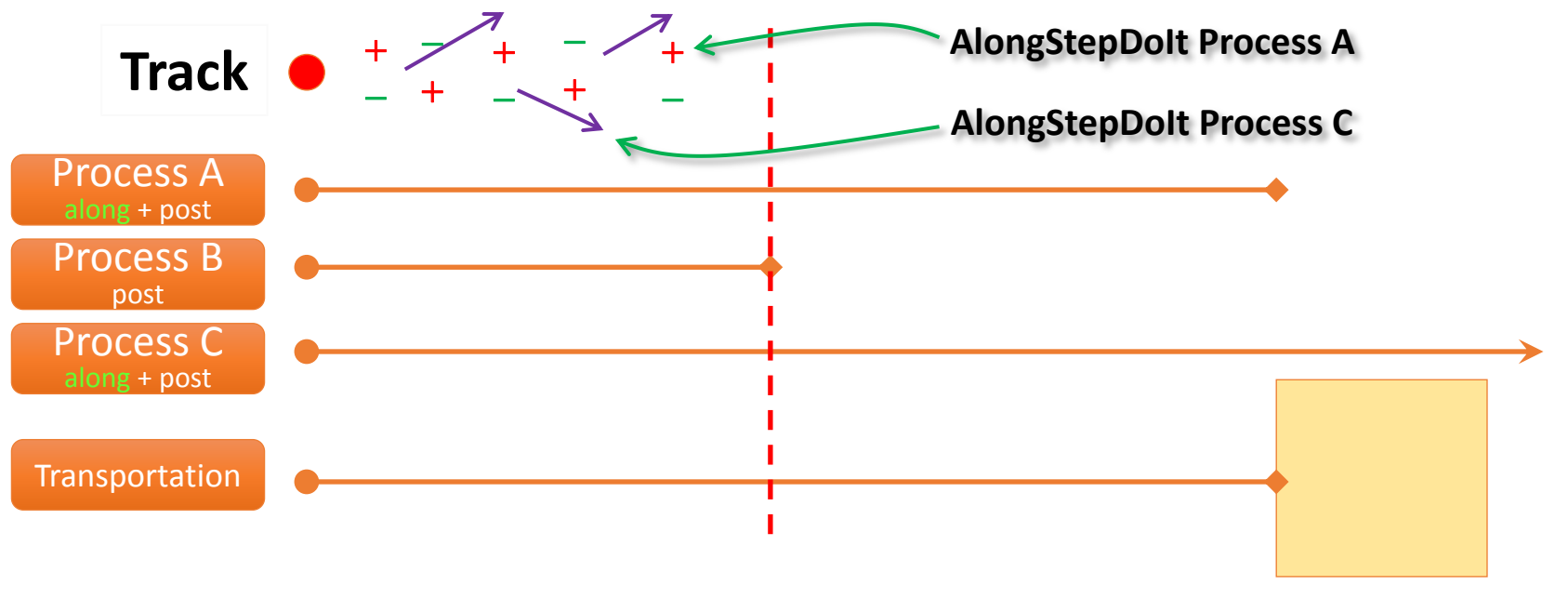


# Process Handling by the Stepping

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2. Apply all `AlongStepDoIt()` actions, « at once »:
  - Changes computed from particle state at the beginning of the step;
  - Accumulated in the `G4Step`;
  - Then applied to the `G4Track`, from the `G4Step`.

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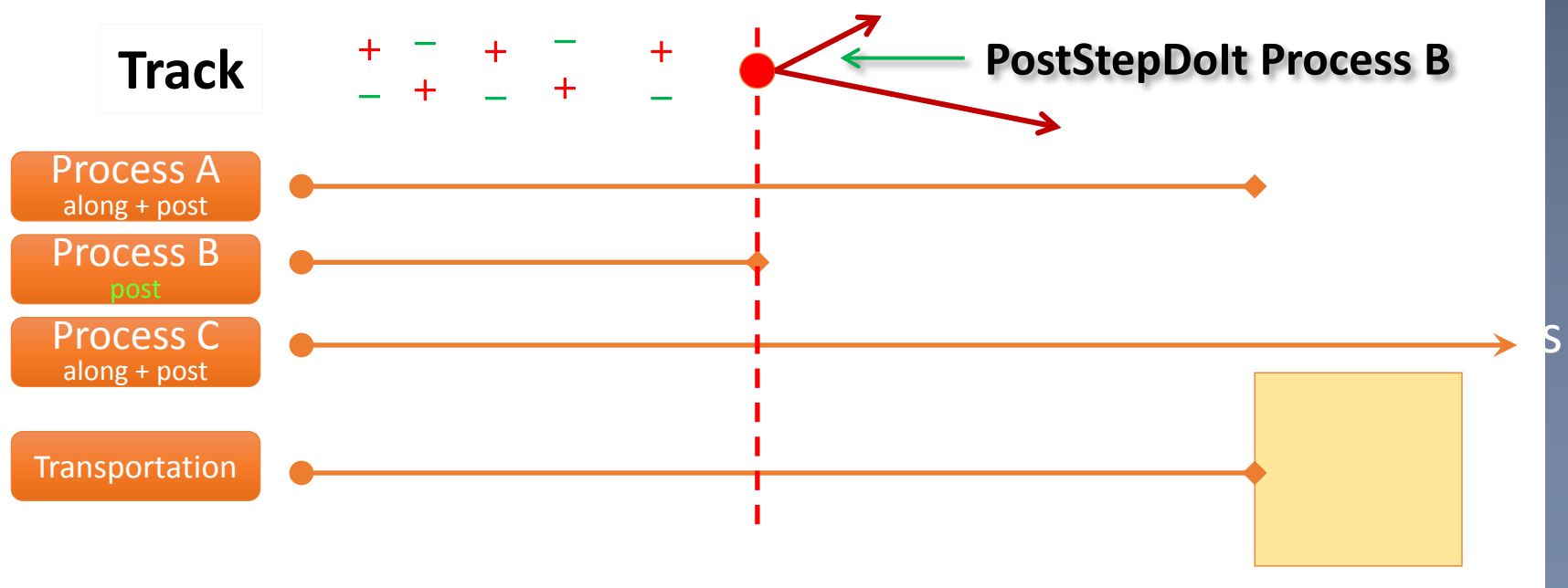
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  - Apply `PostStepDoIt()` of process which limited the step (if any);
  - And apply any other « forced » processes (not discussed here)



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### III. Production Thresholds (aka « cuts »)

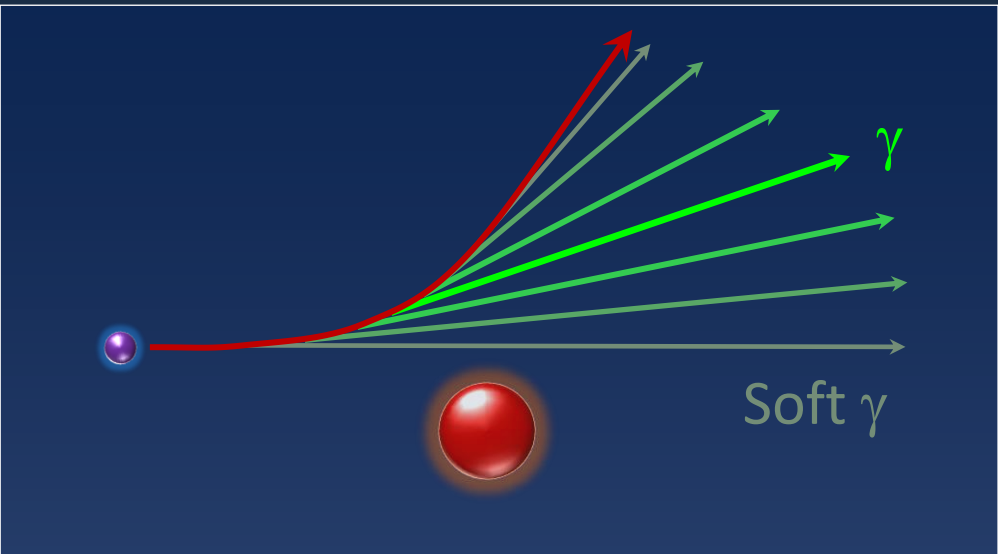
# Threshold for Secondary Production

- › Simulation accuracy limited by:
  - Intrinsic limitation of physics process modeling
    - › For example details of atomic or molecular structure, etc.
  - Finite computing power
    - › Which forces to limit the production to some # of events
    - › Which restricts the usage to models “fast enough”
      - I.e. : no way to use lattice QCD in detector simulation
    - › **Which forces to suppress the production of very low energy particles for processes having infrared divergences**
      - Infinite or very large number of produced secondaries

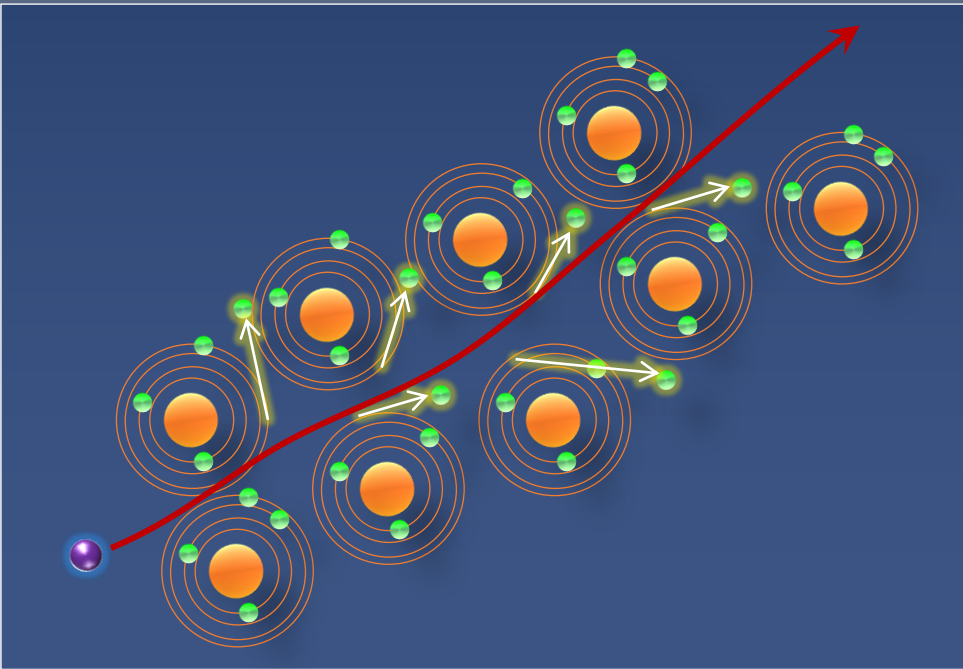
›

›

›



Bremstrahlung: actual divergence in forward production of ultra-soft gammas.



Ionisation : large production of ionisation electrons, from loosely bound atomic ones.



# Threshold for Secondary Production

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    - › **Which forces to suppress the production of very low energy particles for processes having infrared divergences**
      - Infinite or very large number of produced secondaries
- › Every simulation developer must answer the question:
  - How low can I go?
- › This is a balancing act:
  - Need to go low enough to get the physics you’re interested in
  - Can’t go too low because of CPU time consumption
- ›

# Threshold for Secondary Production

## › Geant4 solution: impose a production threshold

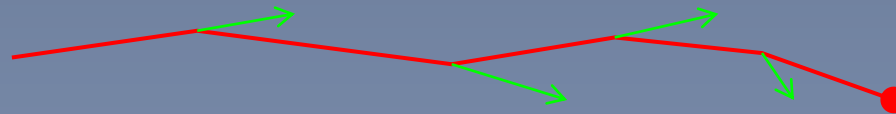
- this threshold is a distance, not an energy
  - › This a “range threshold”
- default = 0.7 mm
- What makes this “range threshold” ?
  - › In Nature, the primary particle loses energy by producing secondary electrons or gammas
    - (or by exciting atoms, molecules, etc. : not of interest here)
  - › In Geant4, the threshold makes only secondary particles able to travel  $> 0.7$  mm to be created
    - The rest is accounted for a “continuous energy loss” (condensed history)

## › Only one value (per region) of range threshold is needed for all materials

- And this distance is internally converted into the related energy thresholds by Geant4
- Conversion “range  $\rightarrow$  energy” is made to according to material

## › Near the primary particle end-point:

- When the primary becomes of too low energy to produce secondaries above threshold:
  - › discrete energy loss ceases (no more secondaries produced)
  - › the primary is tracked down to zero kinetic energy using continuous energy loss



- Note that this makes Geant4 having “no tracking cuts”

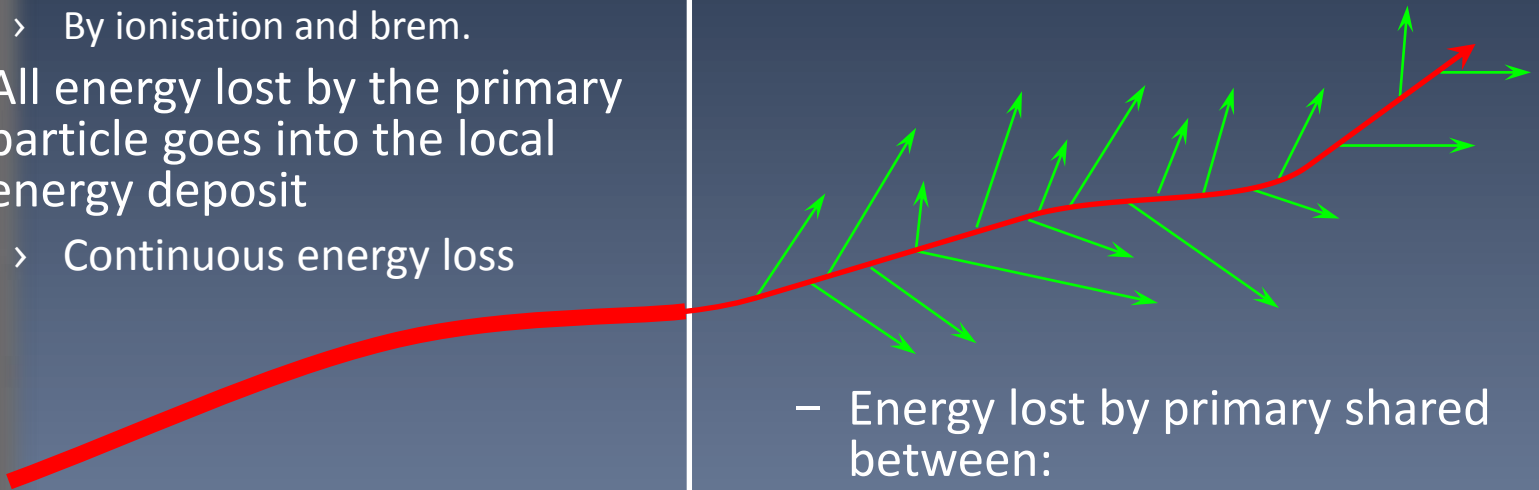
# High and low thresholds : what differences ?

## > High Threshold

- No secondary production
  - > By ionisation and brem.
- All energy lost by the primary particle goes into the local energy deposit
  - > Continuous energy loss
- You'll see as:
  - > step-> `GetTotalEnergyDeposit()` is high
  - > You don't have energy deposit elsewhere than on primary path

## > Low Threshold

- Many secondaries produced
- Energy lost by primary shared between:
  - > Local energy deposit
  - > discrete secondary production
- You'll see as:
  - > step-> `GetTotalEnergyDeposit()` is lower than before
  - > Energy deposit more scattered due to subsequent deposit of secondary particles

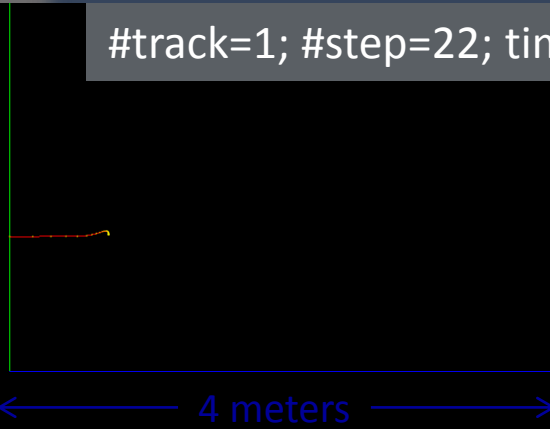




# 10 GeV $e^-$ in liquid Argon, with cuts of 1 km, 1 m, 1 mm and 1 $\mu\text{m}$

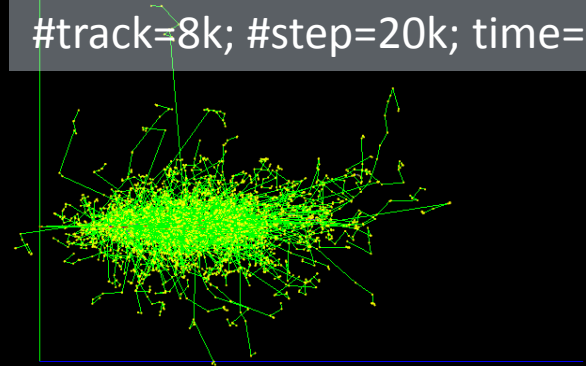
1 km

#track=1; #step=22; time~0s



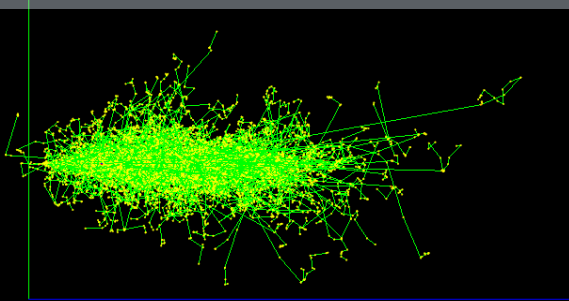
1 m

#track=8k; #step=20k; time=60ms



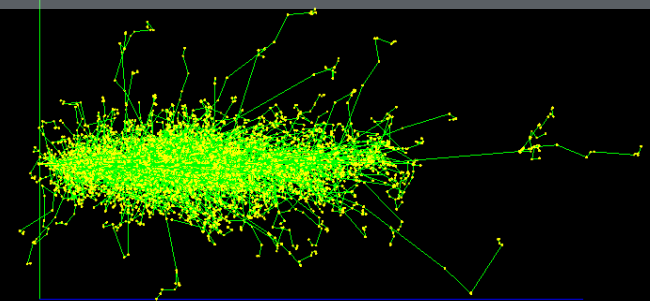
1 mm

#track=18k; #step=39k; time=90ms



1  $\mu\text{m}$

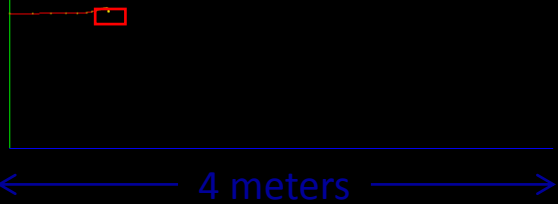
#track=724k; #step=1.5M; time=4.6s



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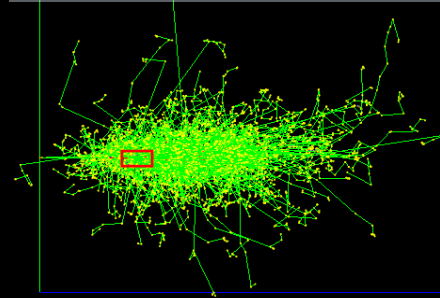
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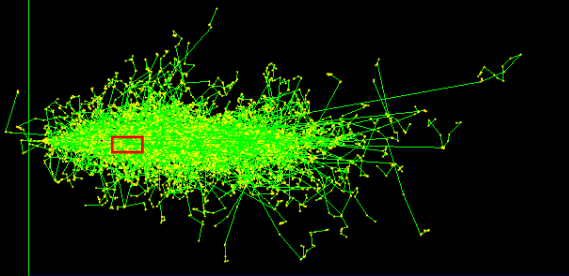
1 m

#track=8k; #step=20k; time=60ms



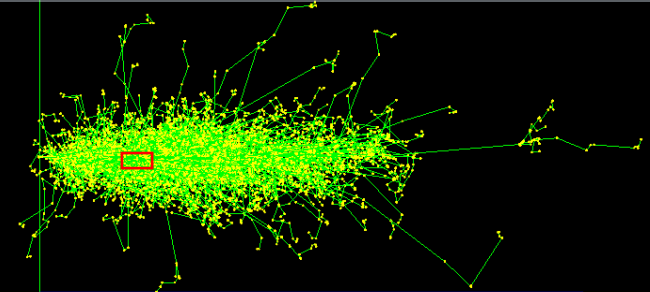
1 mm

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1  $\mu\text{m}$

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1 km

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1 mm

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1  $\mu\text{m}$

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# Assigning cuts to your simulation

- › You must assign cuts to  $\gamma$ ,  $e^-$  and  $e^+$ .
  - For  $\gamma$ 's : needed to limit production from infrared divergence of brem. process
  - For  $e^-$ 's : needed to limit high production from ionization
  - For  $e^+$ 's : bit of historical reasons (no infrared divergence process)
    - › (Plans in Geant4 to review this)
- › You may assign cuts to protons
  - To define the threshold for producing proton by **recoil in elastic collisions**
  - Threshold used for recoil ions too.
- › The easiest way to define cuts is at run time
  - On command line or with a macro
  - For  $\gamma$ ,  $e^-$  and  $e^+$  and p in one go, eg:

```
/run/setCut 2 mm
```
  - Per particle threshold, eg:

```
/run/setCutForAGivenParticle e- 0.1 mm
```
  - (later we'll add the case of "region")

# Getting information on range to energy conversion

```
/run/setCut 1 mm  
/run/beamOn 1 (to force calculations of thresholds)  
/run/dumpCouples
```

```
===== Table of registered couples =====
```

```
Index : 0      used in the geometry : Yes
```

```
Material : G4_Galactic
```

```
Range cuts      : gamma 1 mm      e- 1 mm      e+ 1 mm      proton 1 mm
```

```
Energy thresholds : gamma 990 eV      e- 990 eV      e+ 990 eV      proton 100 keV
```

```
Region(s) which use this couple :
```

```
DefaultRegionForTheWorld
```

```
Index : 1      used in the geometry : Yes
```

```
Material : G4_Pb
```

```
Range cuts      : gamma 1 mm      e- 1 mm      e+ 1 mm      proton 1 mm
```

```
Energy thresholds : gamma 101.843 keV      e- 1.36749 MeV      e+ 1.27862 MeV      proton 100 keV
```

```
Region(s) which use this couple :
```

```
DefaultRegionForTheWorld
```

```
Index : 2      used in the geometry : Yes
```

```
Material : G4_PLASTIC_SC_VINYLTOLUENE
```

```
Range cuts      : gamma 1 mm      e- 1 mm      e+ 1 mm      proton 1 mm
```

```
Energy thresholds : gamma 2.40367 keV      e- 356.639 keV      e+ 344.855 keV      proton 100 keV
```

```
Region(s) which use this couple :
```

```
DefaultRegionForTheWorld
```

```
=====
```

# Getting information on range to energy conversion

```
/run/setCut 0.01 mm  
/run/beamOn 1 (to force calculations of thresholds)  
/run/dumpCouples
```

```
===== Table of registered couples =====
```

```
Index : 0      used in the geometry : Yes
```

```
Material : G4_Galactic
```

```
Range cuts      : gamma 10 um      e- 10 um      e+ 10 um      proton 10 um  
Energy thresholds : gamma 990 ev    e- 990 ev     e+ 990 ev     proton 1 keV
```

```
Region(s) which use this couple :
```

```
DefaultRegionForTheWorld
```

```
Index : 1      used in the geometry : Yes
```

```
Material : G4_Pb
```

```
Range cuts      : gamma 10 um      e- 10 um      e+ 10 um      proton 10 um  
Energy thresholds : gamma 5.995 keV e- 58.1082 keV e+ 56.9484 keV proton 1 keV
```

```
Region(s) which use this couple :
```

```
DefaultRegionForTheWorld
```

```
Index : 2      used in the geometry : Yes
```

```
Material : G4_PLASTIC_SC_VINYLTOLUENE
```

```
Range cuts      : gamma 10 um      e- 10 um      e+ 10 um      proton 10 um  
Energy thresholds : gamma 990 ev    e- 15.1173 keV e+ 14.6763 keV proton 1 keV
```

```
Region(s) which use this couple :
```

```
DefaultRegionForTheWorld
```

```
=====
```

# Getting information on range to energy conversion

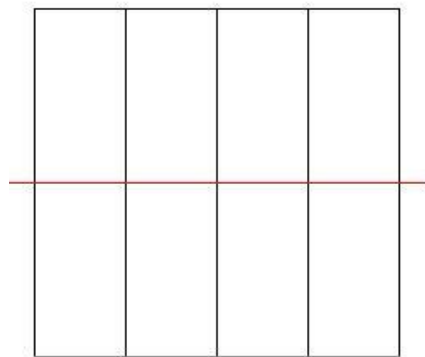
```
/run/setCut 1 nm  
/run/beamOn 1 (to force calculations of thresholds)  
/run/dumpCouples
```

```
===== Table of registered couples =====  
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Range cuts      : gamma 10 Ang      e- 10 Ang      e+ 10 Ang proton 10 Ang  
Energy thresholds : gamma 990 eV      e- 990 eV      e+ 990 eV      proton 0.1 eV  
Region(s) which use this couple :  
DefaultRegionForTheWorld  
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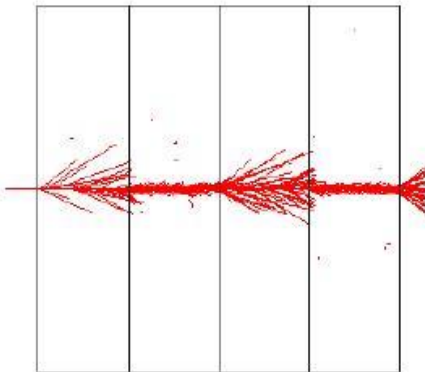
# Production Threshold vs. Energy Cut

Example: 500 MeV p in LAr-Pb Sampling Calorimeter

Energy Threshold

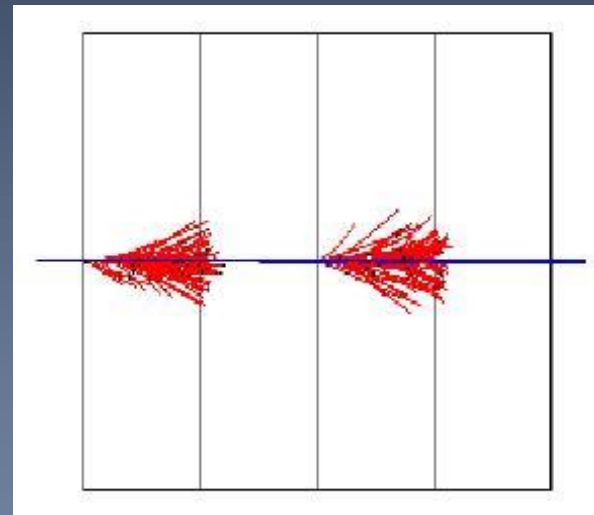


Cut = 2 MeV



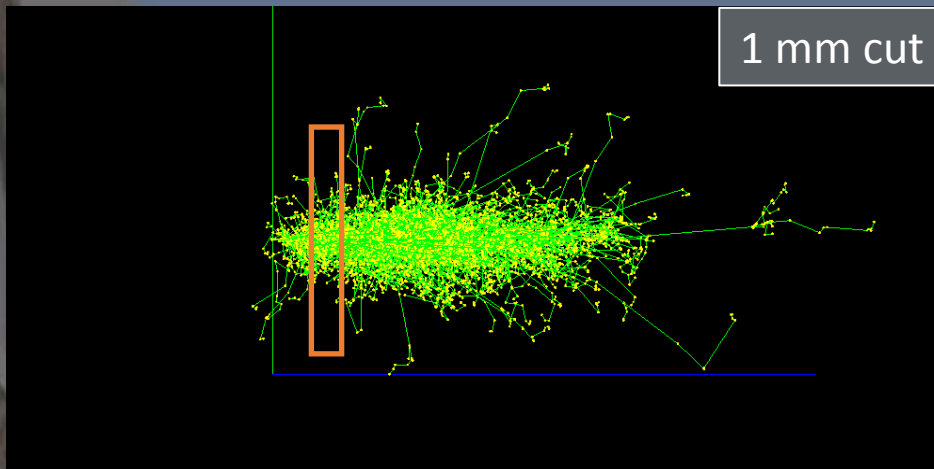
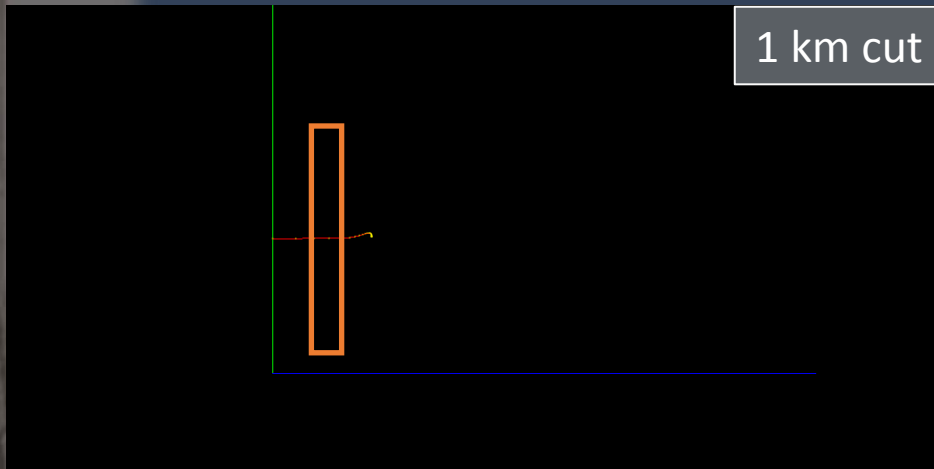
Cut = 450 keV

Geant4 Production  
Range Threshold



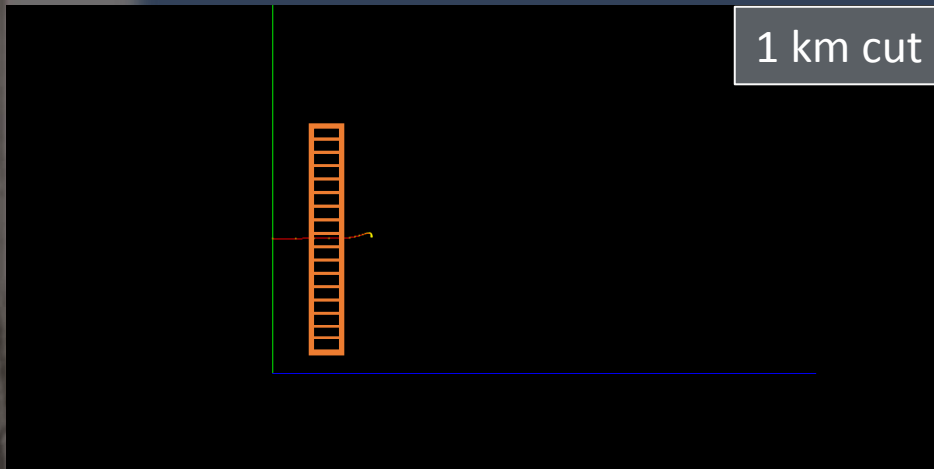


# Energy recorded

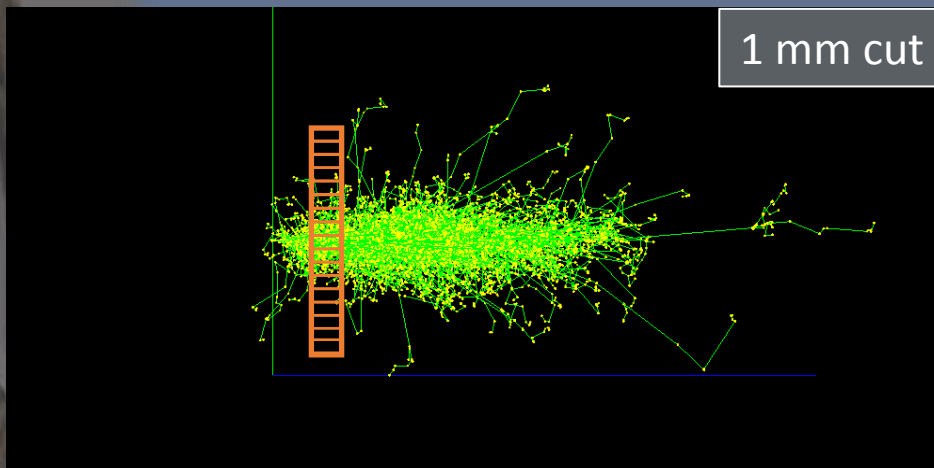


- > If recording energy deposit in a big volume
  - No difference between high and low energy thresholds

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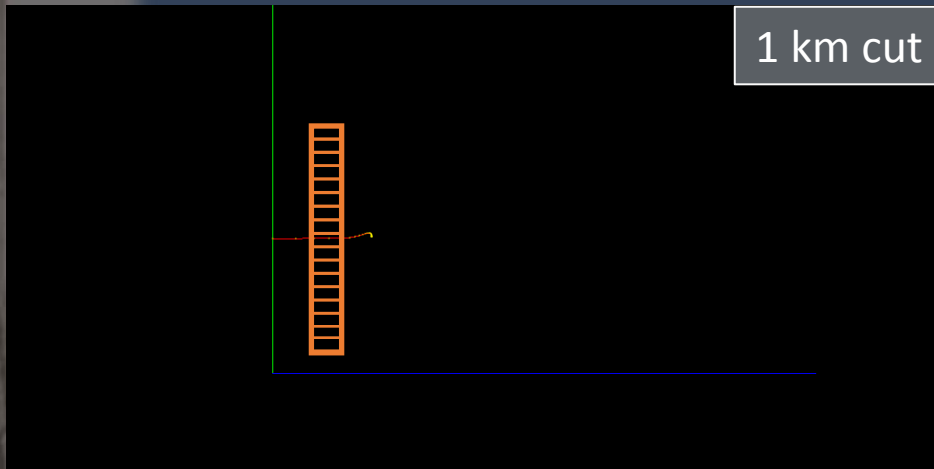


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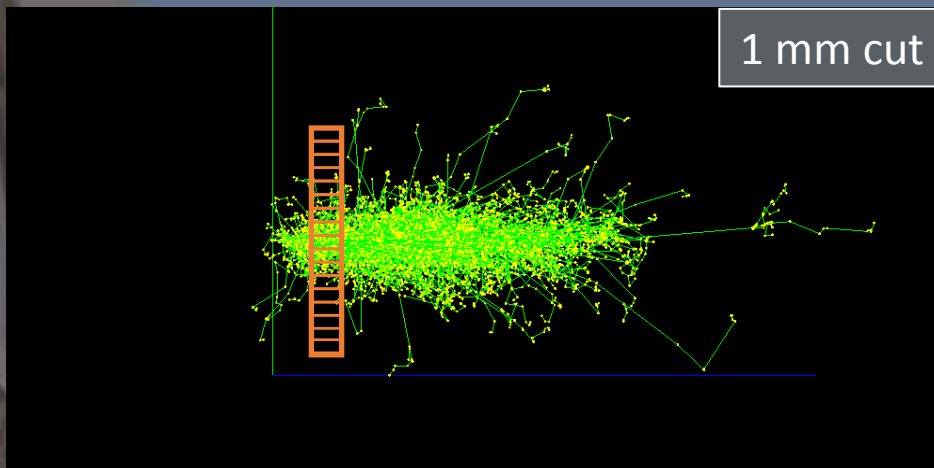


- > But if recording in small volumes
  - Big differences !

# Energy recorded



- > If recording energy deposit in a big volume
  - No difference between high and low energy thresholds



- > But if recording in small volumes
  - Big differences !

> Typically : range cut  
~ volume dimension



## IV. Regions

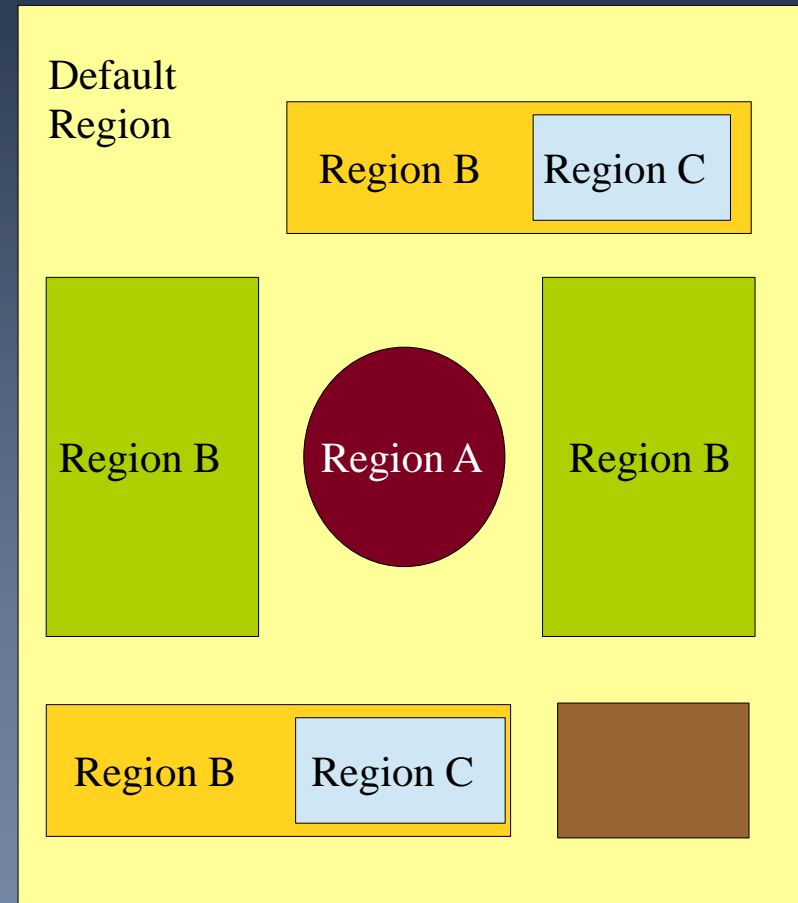
A quick geometry detour

# Concept of Regions (1/2)

- On the top of the volume hierarchy users can define regions which are selected sets of volumes, typically of sub-systems
  - E.g. barrel + end-caps of the calorimeter, “deep” areas of support structures can be a region.
- A region can be any group of volumes
- A region can hold a set of various properties:
  - Production thresholds (cuts)
  - User limits
  - User region information
  - Fast simulation manager
  - Regional user stepping action
  - Field manager

# Concept of Regions (2/2)

- A region is always associated with one or more logical volumes
- A **root logical volume** = volume associated to a region
  - All daughter volumes share the same region, unless a daughter volume itself becomes an other root.
  - A logical volume can not be shared among regions.
- World logical volume is always associated with the default region
  - Users do not need to define it.



# Creating a region, accessing it, creating a user region information object

MyDetectorConstruction.cc

```
#include "G4Region.hh"

// Create a region
G4Region* myRegion = new G4Region("MyRegion");
// Attach a logical volume to the region
myRegion->AddRootLogicalVolume(myLV);
```

MyOtherClass.cc

```
#include "G4RegionStore.hh"
#include "MyRegionInformation.hh"

// Retrieve the region by its name
G4Region* region
    = G4RegionStore::GetInstance()->GetRegion("MyRegion");

// Create some property to be assigned to a region
MyRegionInformation* myInfo = new MyRegionInformation();

// Set myInfo to the region
region->SetUserInformation(myInfo);
```

› We will see just after how to assign « cuts » to a region.



## V. Cuts per region



# Why cuts per region ?

- > Running with “as low as possible” cuts is:

- Good for physics quality
- Bad for CPU consumption

- > In large applications (ie : HEP) not all parts of detector simulation require the same level of accuracy:

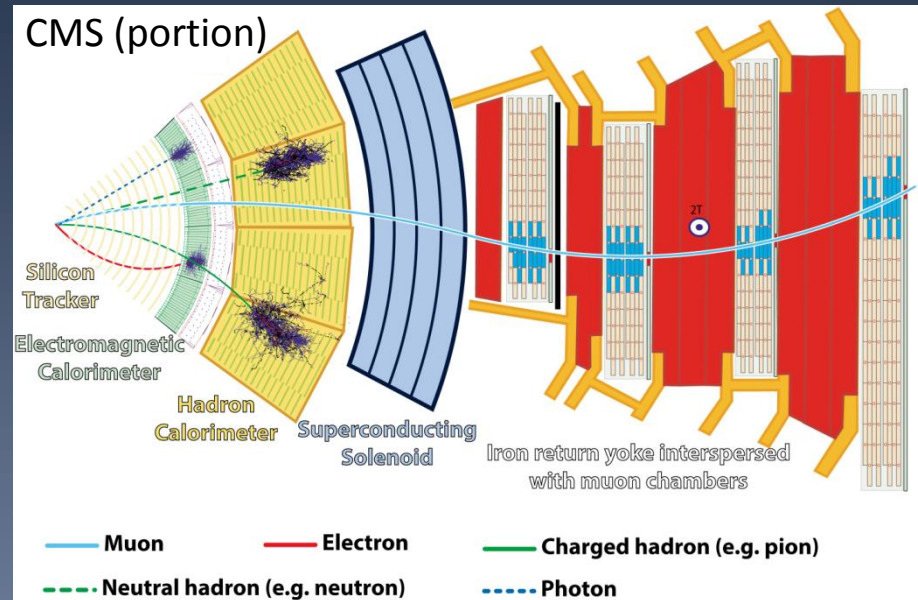
- Tracking systems:
  - > Good accuracy needed
  - > So, low cuts needed
- Hadron calorimeter:
  - > Low accuracy is enough
  - > So high cuts ok

- > Issue:

- Low cuts : Accuracy in tracking systems ✓ Processing time for hadron calorimeter ✗
- High cuts : Accuracy in tracking systems ✗ Processing time for hadron calorimeter ✓
- Medium cuts: Make everybody unhappy ;)

- > **Solution:**

- Allow “cuts per region”
- Tracking system = a region with low cuts
- Hadron calorimeter = a region with high cuts



# Assigning cuts to a region

- › Assume you define a region with name “MyRegion” in your detector construction
- › To assign cuts to it, you do:

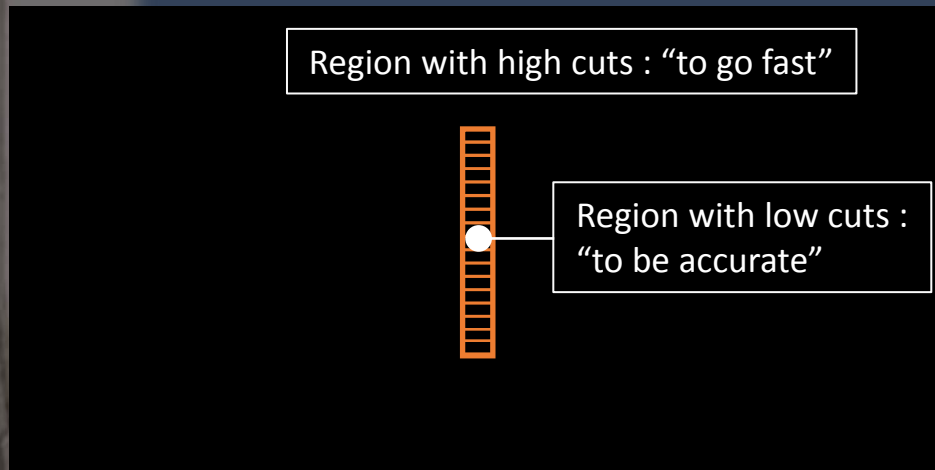
```
...  
// Create the region  
G4Region* myRegion = new G4Region("MyRegion");  
...  
...  
// Define cuts object for the new region and set values  
G4ProductionCuts* cuts = new G4ProductionCuts;  
cuts->SetProductionCut(0.01*mm); // for gamma, e+, e-, p  
// Assign cuts to region  
myRegion->SetProductionCuts(cuts);  
...
```

- › And you can change cut values with command line (or macro) as:

```
/run/setCutForRegion MyRegion 1 mm
```

- › Note that the world volume is in fact a region : it is the “default” one.

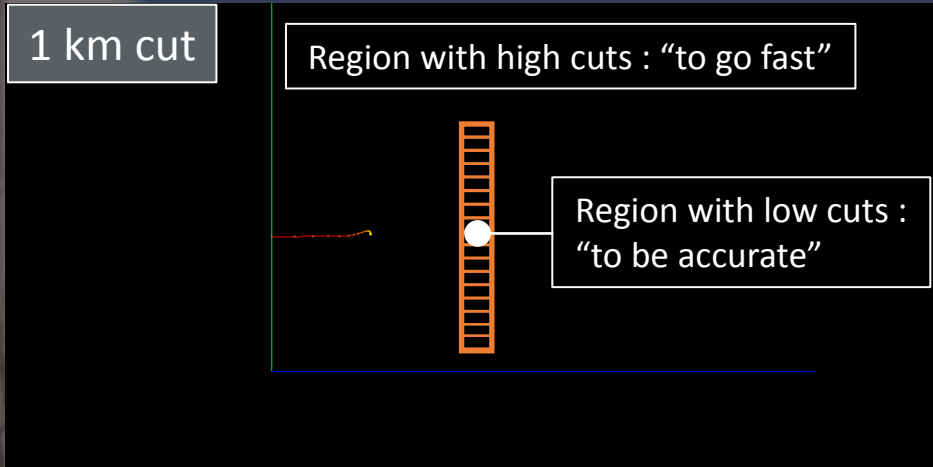
# Be critical : temptation for mistake



## > Temptation:

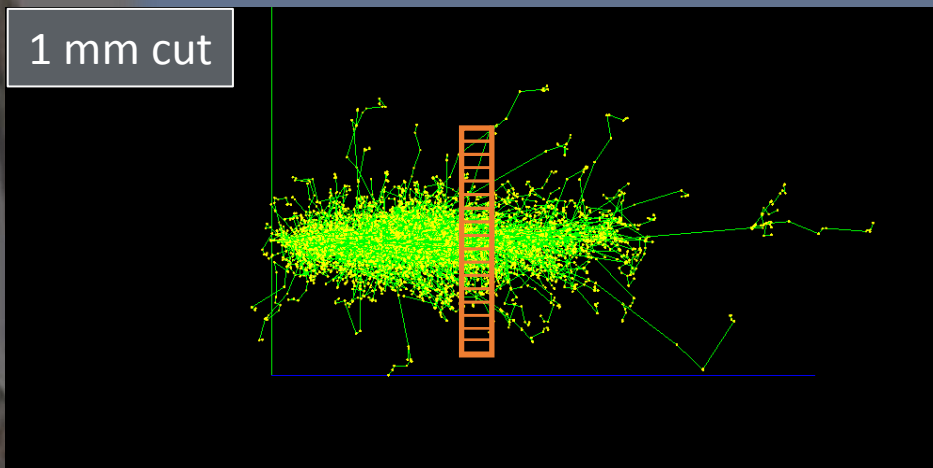
- *“Let me gain time putting a high energy threshold in the parts where I am not interested in details and putting low thresholds where I am interested in details !”*

# Be critical : temptation for mistake



## > Temptation:

- “Let me gain time putting a high energy threshold in the parts where I am not interested in details and putting low thresholds where I am interested in details !”



## > Bad idea !

- What happens in one volume is not only determined by this volume, but also by what happens ***before*** this volume.
- Our example with two extreme threshold cases makes it clear...



# Threshold for Secondary Production

- › Instead of “secondary production threshold distance” it is more convenient to simply say “cuts”
- › The cuts values are set in the SetCuts() method of your physics list
  - Either for the entire simulation (ie for the entire world volume)
  - Or per region
  - Geant4 proposes the default value of 0.7 mm
- › They can be defined with command line, eg:
  - Idle> /run/setCuts 1 mm
- › User needs to decide the best value:
  - The lower the better
    - › To be balanced with your available computing power
  - Typically range cut  $\sim$  volume dimension is fine
    - › Being careful of not having cut too severely before this volume

# Summary

- › Geant4 supplies many physics processes which cover electromagnetic, hadronic, decay physics and “technical”.
- › A unique interface, G4VProcess, allows processes to specify their nature: AtRest, Along (continuous), PostStep (discrete)
  - A process may mix several of these
- › Geant4 does not have “tracking cut”
  - Produced particles are tracked down to zero energy.
- › Geant4 makes use of a “range cut” for controlling the production of secondary particles
  - For some particles and some processes only
- › It is recommended to use a range cut  $\sim$  smallest dimension you’re interested in.

# **A QUICK ADDED INFORMATION ON PHYSICS LISTS**

# Physics List Naming Convention

- The following acronyms refer to various hadronic options
  - **FTF** -> Fritiof string model ( $>\sim 3$  GeV)
  - **QGS** -> Quark Gluon String model ( $>\sim 12$  GeV)
  - **BERT** -> Bertini-style Cascade ( $\sim < 10$  GeV)
  - **BIC** -> Binary Cascade ( $\sim < 10$  GeV)
  - **P** -> Precompound model used for nuclear de-excitation ( $\sim < 150$  MeV)
  - **HP** -> High Precision neutron model ( $< 20$  MeV)
- EM options designated by
  - No suffix : standard EM physics
  - **\_EMV** , **\_EMX** : fast options for high-energy physics
  - **\_EMY** , **\_EMZ** , **\_LIV** , **\_PEN** : more precise options, for medical and space science applications
  - **\_GS** Goudsmith-Sanderson is used for MSC of  $e^-$  and  $e^+$



# When the application starts...

- Large amount of information displayed by the physics list

```
...
FTFP_BERT : new threshold between BERT and FTFP is over the interval
  for pions :    3 to 12 GeV
  for kaons  :    3 to 12 GeV
  for proton :    3 to 12 GeV
  for neutron :  3 to 12 GeV
...
conv:   for gamma   SubType= 14  BuildTable= 1
        Lambda table from 1.022 MeV to 100 TeV, 18 bins per decade, spline: 1
        ===== EM models for the G4Region  DefaultRegionForTheWorld =====
          BetheHeitler :  Emin=          0 eV   Emax=          80 GeV
          BetheHeitlerLPM : Emin=          80 GeV  Emax=         100 TeV
...
Hadronic Processes for anti_deuteron

  Process: hadElastic
    Model:          hElasticLHEP: 0 eV /n ---> 100.1 MeV/n
    Model:          AntiAElastic: 100 MeV/n ---> 100 TeV/n
    Cr_sctns:       AntiAGlauber: 0 eV ---> 2.88022e+295 J
    Cr_sctns:       GheishaElastic: 0 eV ---> 100 TeV

  Process: anti_deuteronInelastic
    Model:          FTFP: 0 eV /n ---> 100 TeV/n
    Cr_sctns:       AntiAGlauber: 0 eV ---> 2.88022e+295 J
    Cr_sctns:       GheishaInelastic: 0 eV ---> 100 TeV

  Process: hFritiofCaptureAtRest
...

```

- The most up-to-date information you can find on a given physics list is here !

# Reference Physics Lists (1/3)

- **FTFP\_BERT**

- Recommended by Geant4 for HEP
- Contains all standard EM processes
- Uses Bertini-style cascade for hadrons  $< 5$  GeV
- Uses Fritiof model for high energies  $> 4$  GeV
- Uses Precompound + evaporation for nuclear de-excitation
- Includes neutron capture
- Includes nuclear stopping at rest of negatively charged hadrons
- Includes gamma- and electro-nuclear
- No neutron-HP, radioactive decay, optical photons

# Reference Physics Lists (2/3)

- **QGSP\_FTFP\_BERT**

- All standard EM processes
- Bertini-style cascade for hadrons  $< 8$  GeV
- Quark Gluon String model for high energies  $> 12$  GeV
- Fritiof model in between 6 – 25 GeV

- **QGSP\_BERT**

- All standard EM processes
- Bertini-style cascade for hadrons  $< 9.9$  GeV
- Quark Gluon String model for high energies  $> 12$  GeV
- Fritiof in between 9.5 – 25 GeV
- NB) We are working to extend QGS at lower energies, so that the transition with BERT can be done directly, without FTF (in this physics list)

# Reference Physics Lists (3/3)

- **QGSP\_BIC**

- Same as QGSP\_BERT, but replaces Bertini-style cascade with Binary cascade model (+ Precompound model)
- Recommended for use at energies below 200 MeV
  - Many medical applications
  - Suggested EM option: \_EMY or \_EMZ

- **FTFP\_BERT\_HP (QGSP\_BERT\_HP)**

- Same as FTFP\_BERT (QGSP\_BERT), but with the high-precision neutron model used for neutrons below 20 MeV
- Significantly slower than FTFP\_BERT (QGSP\_BERT), especially when Doppler broadening on-the-fly is used
  - There is an option to turn this off
- For radiation protection and shielding applications