



Physics processes in general

CERN Geant4 User's Workshop

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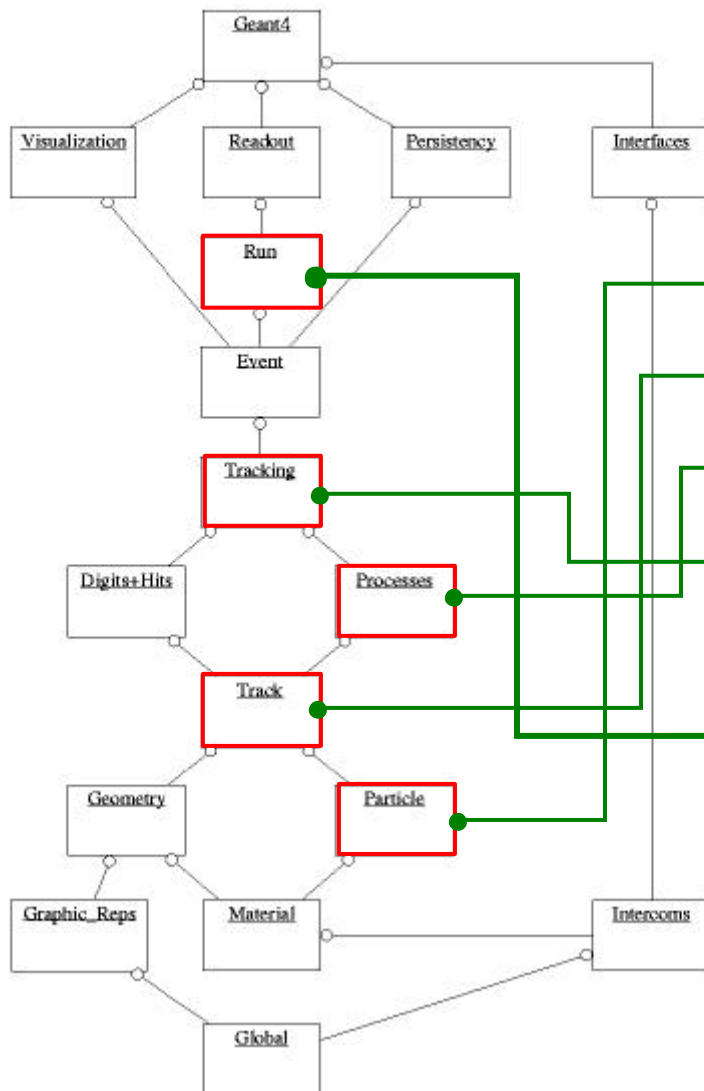
Derived from a talk by Marc Verderi
Ecole Polytechnique - LLR



Introduction

- Present the ingredients needed to understand how to build a « physics list », which is the physics setup:
 - ◆ It is the place where the user tells what particles, processes and production cuts will be used in his/her application;
 - ◆ This is a **mandatory** and **critical** user's task;
- We will go through several aspects regarding the « heart » of GEANT4;
- Present only the 'theoretical' aspect of processes:
 - ◆ Presentations on the 'concrete' processes follow;

involved



👉 The ingredients presented sit in these categories:

- Particles
- Track
- Processes
- We will show how they are handled by the **Tracking**;

👉 The « physics list » interface sits in the Run category.

Layout

- I. What is tracked;
 - Definition of particles;
 - G4Track;
- II. The process interface;
 - G4VProcess;
 - How processes are used by the stepping;
- III. The production cuts;
- IV. Building the « physics lists ».
- V. User-defined limits



I. What is tracked in GEANT4;

Speak about:

G4ParticleDefinition;

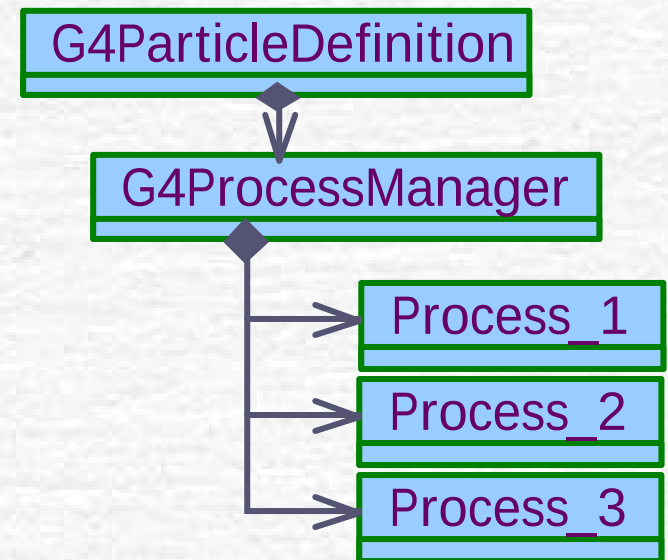
G4DynamicParticle;

G4Track;



G4ParticleDefinition

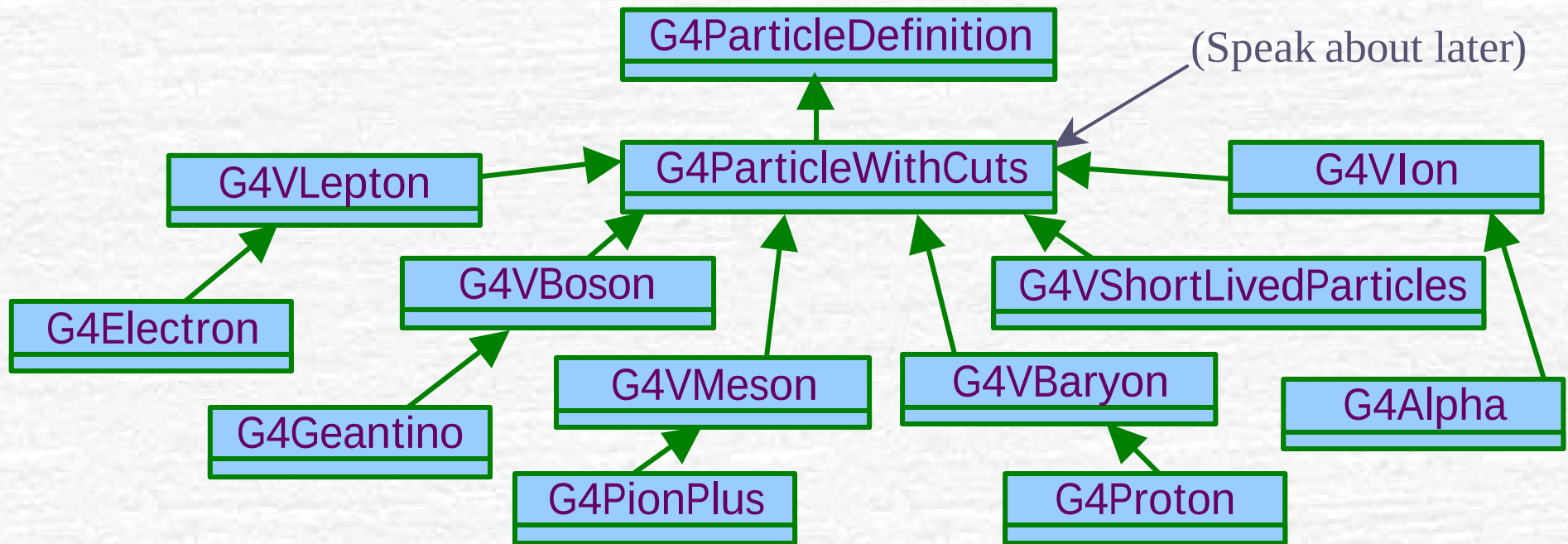
- ☛ The particle types in GEANT4 are described by the **G4ParticleDefinition** class;
 - Class defined in **source/particles/management**;
- ☛ Describes the « **intrinsic** » particle properties:
 - Mass, width, spin, lifetime...
- ☛ Describes its « **sensitivity** » to physics:
 - This is realized by a **G4ProcessManager**;
 - Attached to the **G4ParticleDefinition**;
 - The **G4ProcessManager** manages the **list of processes** the **user wants** the particle to be **sensitive to**;
 - Note that **G4ParticleDefinition** doesn't know by itself its sensitivity to physics.



Concrete G4ParticleDefinition (1)

- G4ParticleDefinition is the base class for defining concrete particles:

- Several layers are defined:



Concrete G4ParticleDefinition (2)

- Most common particles, with lifetime large enough, are implemented as static classes:
 - Like `G4Electron`, K^0_S , gamma, pions, but also α ...
 - To allow –say- electrons in the simulation, the following call should be made in the « physics list »:
 - `G4Electron::ElectronDefinition();`
- Heavy ions are created **on the fly** by processes;
 - Too many ions to have a class per ion !
 - An ion is tracked, and then its « ion type » disappears;
 - Ions are all created from the static class `G4GenericIon`. To allow heavy ions in the simulation, you should call:
 - `G4GenericIon::GenericIonDefinition();`
- Resonances (`G4VShortLivedParticles`) are also created on the fly. Similar calls to the definitions (excited baryons, mesons ...) should be made.

Example: G4Electron class (1)

- Extract from `source/particles/leptons/include/G4Electron.hh`

```
class G4Electron : public G4VLepton
{
public:
    static G4Electron* ElectronDefinition();
private: //hide constructor as private
    G4Electron(
        const G4String&    aName,          G4double    mass,
        G4double           width,          G4double    charge,
        G4int              iSpin,          G4int       iParity,
        G4int              iConjugation,    G4int       ilsospin,
        G4int              ilsospin3,      G4int       gParity,
        const G4String&     pType,          G4int       lepton,
        G4int              baryon,         G4int       encoding,
        G4bool             stable,         G4double    lifetime,
        G4DecayTable       *decaytable);
private:
    static G4Electron theElectron;
    ...
};
```

Example: G4Electron class (2)

- Extract from [source/particles/leptons/src/G4Electron.cc](#)

```
G4Electron::G4Electron(
    const G4String&    aName,      G4double    mass,
    G4double          width,      G4double    charge,
    G4int             iSpin,      G4int        iParity,
    G4int             iConjugation, G4int        iIsospin,
    G4int             iIsospin3,   G4int        gParity,
    const G4String&    pType,      G4int        lepton,
    G4int             baryon,      G4int        encoding,
    G4bool            stable,      G4double    lifetime,
    G4DecayTable      *decaytable )
: G4VLepton(    aName,    mass,    width,    charge,    iSpin,    iParity,
                iConjugation, iIsospin, iIsospin3, gParity, pType,
                lepton,    baryon, encoding, stable, lifetime, decaytable )
{ SetParticleSubType("e"); }

G4Electron G4Electron::theElectron(
    "e-", 0.51099906*MeV, 0.0*MeV, -1.*eplus,
    1,    0,    0,
    0,    0,    0,
    "lepton", 1,    0,    11,
    true,    -1.0,    NULL);

G4Electron* G4Electron::ElectronDefinition(){return &theElectron;}
```

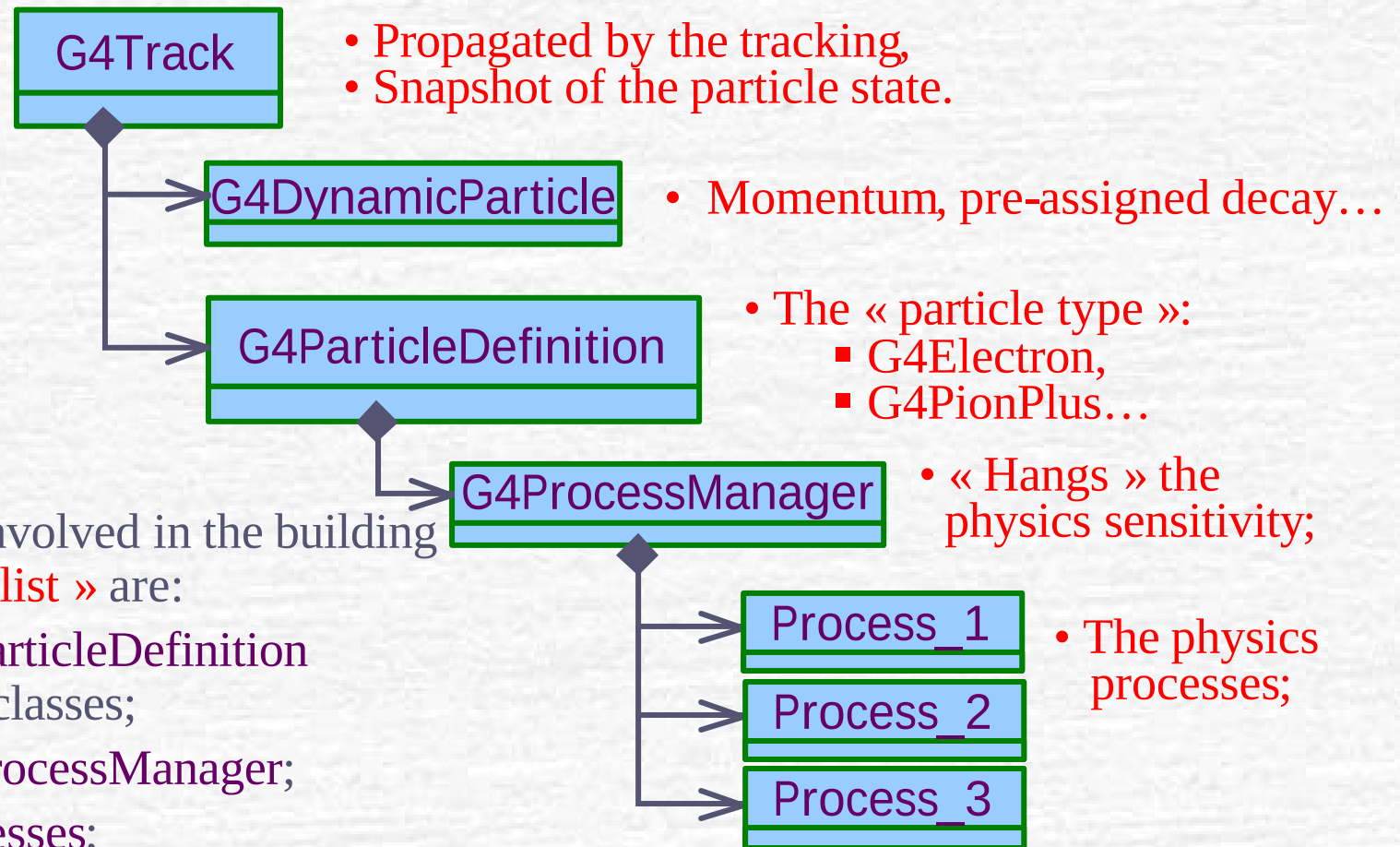
G4DynamicParticle

- ☛ G4DynamicParticle describes the purely dynamic part (ie no position, nor geometrical information...) of the particle state:
 - Momentum, energy, polarization;
 - It hangs a G4ParticleDefinition pointer;
 - Retains eventual pre-assigned decay informations:
 - decay products;
 - lifetime;
- ☛ Class defined in `source/particles/management`;

G4Track

- ✦ G4Track defines the class of objects **propagated by the GEANT4 tracking**;
 - Class defined in **source/track**;
- ✦ The G4Track represents a **« snapshot »** of the particle state;
- ✦ A G4Track object aggregates:
 - A G4ParticleDefinition;
 - A G4ParticleDynamics;
 - **Geometrical informations**:
 - Position, current volume ...
 - Track ID, parent ID;
 - process which created this G4Track;
 - weight, used for event biasing technic;
 - ...
- ✦ A G4Track is tracked from its birth until:
 - **It is killed**:
 - By an interaction
 - Or because it comes to rest, and is stable;
 - Or, by a user's action (under his responsibility !).
 - **Or, it exits the world volume**;
- ✦ Class users need to be familiar with !

Summary view of « What is tracked in GEANT4 »



- The classes involved in the building the « physics list » are:

- The G4ParticleDefinition concrete classes;
- The G4ProcessManager;
- The processes;



II. The process interface;

Speak about:

G4VProcess;
The Stepping;



G4VProcess (1)

- 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:

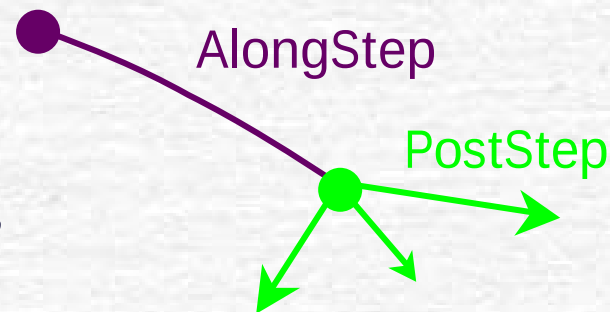
- Decay, e^+ annihilation ...

- **AlongStep** actions:

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

- **PostStep** actions:

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



G4VProcess (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 size***:

- 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.

G4VProcess (3)

The « action » methods are thus:

- `AtRestGetPhysicalInteractionLength()`, `AtRestDoIt()`;
- `AlongStepGetPhysicalInteractionLength()`, `AlongStepDoIt()`;
- `PostStepGetPhysicalInteractionLength()`, `PostStepDoIt()`;

A set of processes implementing given combinations of actions exists:

- `G4VDiscreteProcess`: only PostStep actions;
- `G4VContinuousDiscreteProcess`: AlongStep + PostStep actions;
- ...

G4VProcess also defines the method:

- `G4bool IsApplicable(const G4ParticleDefinition &);`
- which returns « true » if the process is applicable to the given particle type;

G4VProcess & G4ProcessManager

- In practice the G4ProcessManager retains **three vectors of actions**:
 - One for the **AtRest** methods of the particle;
 - One for the **AlongStep** ones;
 - And one for the **PostStep** actions.
- These are those vectors the user sets up in the « physics list » and which are used by the tracking.

How the Stepping handles processes

- ☛ The stepping treats processes **generically**:
 - The stepping does not know what processes it is handling;
- ☛ The stepping imposes on the processes to
 - Cooperate in their AlongStep actions;
 - Compete for PostStep and AtRest actions;
- ☛ Processes can optionally emit also a «signal» to require particular treatment:
 - **notForced**: «standard» case;
 - **forced**: PostStepDoIt action is applied anyway;
 - **conditionallyForced**: PostStepDoIt applied if AlongStep has limited the step;
 - More on this will be said in the session «Adding a new process»;

Stepping Invocation Sequence of Processes for a particle travelling

1. At the beginning of the step, determine the step length:
 - Consider all processes attached to the current `G4Track`;
 - Define the step length as the smallest of the lengths among:
 - All `AlongStepGetPhysicalInteractionLength()`
 - All `PostStepGetPhysicalInteractionLength()`
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`.
3. Apply `PostStepDoIt()` action(s) « sequentially », as long as the particle is alive:
 - Apply `PostStepDoIt()` of process which proposed the smallest step length;
 - apply « forced » and « conditionally forced » actions.

Stepping Invokation Sequence of Processes for a Particle at Rest

1. If the particle is at rest, *is stable and can't annihilate*, it is killed by the tracking:
 - To be more accurate: if a particle **at rest** has no « **AtRest** » actions defined, it is killed.
2. Otherwise determine the lifetime:
 - Take the smallest time among:
 - All **AtRestGetPhysicalInteractionLength()**
 - Called «physical interaction length» but returns a time!
3. Apply the **AtRestDoIt()** action of the process which returned the smallest time.

Processes ordering

☛ The Ordering of processes matters!

☛ Ordering of following processes is **critical**:

- Assuming n processes, the ordering of the **AlongGetPhysicalInteractionLength** of the last processes should be:

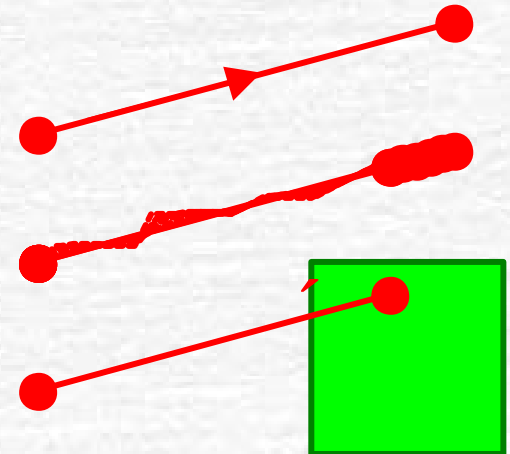
[$n-2$] ...

[$n-1$] multiple scattering

[n] transportation

☛ Why ?

- Processes return a « true path length »;
- The **multiple scattering** « virtually folds up » this true path length into a **shorter** « geometrical » path length;
- Based on this new length, the **transportation** can geometrically limit the step.



☛ Ordering of most other process does not matter.



III. The production cuts;

Speak about:

Why production cuts are needed;

The cuts scheme in GEANT4



The cuts in GEANT4

- ✦ In GEANT4 there is **no tracking cut**:
 - Particles are tracked down to a **zero range**/kinetic energy;
- ✦ Only **production cuts** exist;
 - ie cuts allowing a particle to be born or not;
- ✦ Why are production cuts needed ?
- ✦ Some electromagnetic processes involve infrared **divergences**:
 - This leads to an **infinity[huge number]** of smaller and smaller energy photons[electrons] (like in bremsstrahlung, δ -ray productions);
 - Production cuts limit this production to particles above the threshold;
 - The remaining, divergent part is treated as a « **net** » **continuous effect** (ie « AlongStep » action);
- ✦ For other processes, production cuts can be an « option » to speed-up the simulation.

Range versus Energy production cuts

- The production of a secondary particle is relevant if it can be « visible » in the detector:
 - I.e. produce a signal -say an energy deposition- visible compared to the signal of the primary alone;
- Range cut allows to easily define such visibility:
 - « I want to produce particles able to travel at least 1 mm; »
 - Criteria which can be applied uniformly accross the detector;
- A cut of the same energy would lead to very different ranges:
 - For the same particle type, depending on the material;
 - For the same material, depending on particle type;
- Range cut has been adopted by GEANT4;
- Actual input to cross-section is the energy threshold, but the conversion range-energy is done automatically in GEANT4;

«Violations» of the production threshold

- ✎ In some (many) cases, particles are produced, **even if they are below the production threshold**;
- ✎ This is intended to let the processes doing the « best » they can;
- ✎ This happens typically for:
 - Decays;
 - Positrons production:
 - In order to simulate the subsequent γ from the annihilation;
 - Hadronic processes:
 - Since no infrared divergences affect the cross-sections;
- ✎ Note: these are not « hard-coded » exceptions, but is a sophisticated, **generic**, mechanism of the tracking;

How GEANT4 produces the production cuts

- ☛ The user specifies a **range cut** for each particle 'type'
 - In the « **physics list** »;
- ☛ This range cut is converted into energy cuts:
 - Each particle -**G4ParticleWithCut**- converts the range cut into an energy cut, for each material;
- ☛ **Physics processes** can then compute the cross-section based on those energy cuts;
- ☛ Done at initialization time;

Relaxing the 'unique' range-cut

- ☛ Today each particle type has a **unique** range-cut
 - ◆ One for electrons, ◆ one for protons,
 - ◆ one for γ , ◆ one for other particles
- ☛ This ensures **consistency**
 - ◆ but is not optimal for applications in which the accuracy of energy deposition varies greatly between regions of the setup/detector.
- ☛ We are developing functionality to relax the restriction of a single cut
 - ◆ Allowing each range cut to be set for a **region**
 - ◆ In addition to the 'global'/default range-cut



IV. Building the « physics list »;

Speak about:

G4VUserPhysicsList;
Concrete physics lists;



G4VUserPhysicsList

- ☛ It is one of the « mandatory user classes »;
 - Defined in **source/run**
- ☛ Defines the **three pure virtual methods**:
 - ConstructParticles();
 - ConstructProcesses();
 - SetCuts();
- ☛ Need to **inherit** from G4VUserPhysicsList to implement a physics list;
- ☛ (Note that a G4UserPhysicsListMessenger allows to control interactively the physics list.)

ConstructParticles() (1)

- To get particle G4XXX, you need to invoke the static method XXXDefinition() in your ConstructParticles() method:

```
void MyPhysicsList::ConstructParticles()
{
    G4XXX::XXXDefinition();
}
```

- For example, to have electrons, positrons and gammas only:

```
void MyPhysicsList::ConstructParticles()
{
    G4Electron::ElectronDefinition();
    G4Positron::PositronDefinition();
    G4Gamma::GammaDefinition();
}
```

ConstructParticles() (2)

☞ Alternatively, some **helper classes** are provided:

- G4BosonConstructor, G4LeptonConstructor
- G4MesonConstructor, G4BaryonConstructor
- G4IonConstructor, G4ShortlivedConstructor

☞ You can use as:

```
G4BaryonConstructor baryonConstructor;  
baryonConstructor.ConstructParticle();
```

☞ Those helper classes are defined in **source/particles/**

- bosons, leptons
- hadrons/mesons, hadrons/barions
- hadrons/ions, shortlived

ConstructProcesses() the hard way

- The class heavily used there is the **G4ProcessManager**:
 - Defined in **source/processes/management**
 - It is used to attach processes to particles;
 - And set their ordering;
- Several ways to « add » a process:
 - **AddProcess**
 - **AddRestProcess**, **AddDiscreteProcess**, **AddContinuousProcess**
- And to order AtRest/AlongStep/PostStep actions of processes:
 - **SetProcessOrdering**
 - **SetProcessOrderingToFirst**, **SetProcessOrderingToLast**
 - (This is the ordering for the **DoIt()** methods, the **GetPhysical-InteractionLength()** ones have the **reverse** order.)
- Please review those **G4ProcessManager** methods !
- Show now various examples.

Example 1: G4VUserPhysicsList::AddTransportation()

☞ Helper method to add the transportation process:

void G4VUserPhysicsList::AddTransportation()

```
{
    G4Transportation* theTransportationProcess= new G4Transportation();
    // loop over all particles in G4ParticleTable
    theParticleIterator->reset();
    while( (*theParticleIterator)() ){
        G4ParticleDefinition* particle = theParticleIterator->value();
        G4ProcessManager* pmanager = particle->GetProcessManager();
        if (!particle->IsShortLived()) {
            // Add transportation process for all particles other than "shortlived"
            if ( pmanager == 0 ) {
                // Error !! no process manager
                G4Exception("G4VUserPhysicsList::AddTransportation : no process manager!");
            } else {
                // add transportation with ordering = ( -1, "first", "first" )
                pmanager->AddProcess(theTransportationProcess);
                pmanager->SetProcessOrderingToFirst(theTransportationProcess, idxAlongStep);
                pmanager->SetProcessOrderingToFirst(theTransportationProcess, idxPostStep);
            }
        } else {
            // shortlived particle case
        }
    }
}
```

Example 2: EM processes for gamma

- Simple example of « **discrete** » processes: ie only PostStep actions;
 - Show usage of helper function **AddDiscreteProcess**;
 - pmanager** is the **G4ProcessManager** of the gamma;
 - Assume the transportation has been set by **AddTransportation**;
- Code sample:
// Construct processes for gamma:
pmanager->AddDiscreteProcess(new G4GammaConversion());
pmanager->AddDiscreteProcess(new G4ComptonScattering());
pmanager->AddDiscreteProcess(new G4PhotoElectricEffect());
- Simple case, where we don't have to deal with processes ordering (except for the transportation which has been set to « first » elsewhere).
- A more complicated case now...

Example 3: EM processes for positrons

// Construct processes for positron

```
G4VProcess* theeplusMultipleScattering = new G4MultipleScattering();
```

```
G4VProcess* theeplusIonisation = new G4eIonisation();
```

```
G4VProcess* theeplusBremsstrahlung = new G4eBremsstrahlung();
```

```
G4VProcess* theeplusAnnihilation = new G4eplusAnnihilation();
```

// add processes

```
pmanager->AddProcess(theeplusMultipleScattering);
```

```
pmanager->AddProcess(theeplusIonisation);
```

```
pmanager->AddProcess(theeplusBremsstrahlung);
```

```
pmanager->AddProcess(theeplusAnnihilation);
```

// set ordering for AtRestDoIt

```
pmanager->SetProcessOrderingToFirst(theeplusAnnihilation, idxAtRest);
```

// set ordering for AlongStepDoIt

```
pmanager->SetProcessOrdering(theeplusMultipleScattering, idxAlongStep, 1);
```

```
pmanager->SetProcessOrdering(theeplusIonisation, idxAlongStep, 2);
```

// set ordering for PostStepDoIt

```
pmanager->SetProcessOrdering(theeplusMultipleScattering, idxPostStep, 1);
```

```
pmanager->SetProcessOrdering(theeplusIonisation, idxPostStep, 2);
```

```
pmanager->SetProcessOrdering(theeplusBremsstrahlung, idxPostStep, 3);
```

```
pmanager->SetProcessOrdering(theeplusAnnihilation, idxPostStep, 4);
```

An alternative way to implement particles and processes

- ☞ It exists the `G4VModularPhysicsList` class:
 - Defined in `source/run`;
 - Which **inherits** from `G4VUserPhysicsList`;
- ☞ Which makes use of a set of `G4VPhysicsConstructor`:
 - Defined in `source/run`;
- ☞ `G4VPhysicsConstructor` defines the pure virtual methods:
 - `ConstructParticle()`;
 - `ConstructProcess()`;
 - It is a kind of « sub-physics list », each of those implementing - say- the EM physics only, the hadronics physics only, etc...
- ☞ Allows to avoid all the physics definition in a single class;
- ☞ Please see `example/novice/N04`

For the best and newest way to create the UserPhysicsList

- ✓ See the presentation later today on Hadronics Physics Lists,
- ✓ It builds on the Modular Physics lists
 - ‘Builders’ modularise further the creation of Physics Lists
 - Accumulating physics models and processes for a particular use case.

SetCuts() (1)

- ✎ This pure virtual method is used to define the cut range;
- ✎ I will here talk only about the **recommended way** of setting cuts:
 - I.e: same cut range for all particles;
 - Setting particle dependent cuts is possible but might be reserved to **advanced** (perverted ? ;->) users.
- ✎ The **G4VUserPhysicsList** base class has the protected member:
 - protected:**
 - G4double** defaultCutValue;
 - Which is set to **1.0*mm** in the constructor;
- ✎ You can eventually change this value in an implementation of **SetCuts()**;
- ✎ The helper **G4VUserPhysicsList::SetCutsWithDefault** method implements the machinery to set the cuts using this **defaultCutValue** value;

SetCuts() (2)

- A typical implementation of `SetCuts()` is then simply:

```
void MyPhysicsList::SetCuts()
{
    defaultCutValue = 1.0*mm;
    SetCutsWithDefault();
}
```



V. User-defined limits;

Speak about:

G4UserLimit;

G4UserSpecialCuts process;



G4UserLimit

- ☛ This class allows to define the following limits in a given **G4LogicalVolume**:
 - **Maximum step size**;
 - Maximum track length;
 - Maximum track time;
 - Minimum kinetic energy;
 - Minimum range;
 - Class defined in **source/global/management**
- ☛ The user can inherit from **G4UserLimit**, or can instantiate the default implementation;
- ☛ The object has then to be set to the **G4LogicalVolume**;

G4UserSpecialCuts

- How to activate G4UserLimit ?
- The maximum step size is **automatically taken into account by the stepping**;
 - This is only the case for this G4UserLimit's attribute;
- For the other limits, the G4UserSpecialCuts process (discrete process) can be set in the physics list;
 - Defined in **source/processes/transportation**
- Or, a simple implementation of discrete process can be done to deal with only the cuts the user is interested in;

Conclusion

- ✦ Creating the « physics list »
 - ✦ exposes, deliberately, the **user** to the **choice** of **physics** (particles + processes) **relevant** to his/her application;
 - ✦ is a critical & complex task
- ✦ To assist users
 - ✦ The **examples** have been used as a starting point;
 - ✦ Then a **modular** structure was created;
 - ✦ Now a better, more structured, **builder** solution has been created; see later talk (J.P. Wellisch).
- ✦ Hypernews, email etc..., remain important to exchange experiences and expertise !