

# **GEANT4 BEGINNERS COURSE**

Queen's University, Belfast (UK)

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## **Detector description: materials and geometry**

**Geant 4** tutorial course



# Part I: Materials

## Materials:

- The System of units & constants
- Definition of elements
- Materials and mixtures
- Some examples ...
- The NIST Data Base

# Units

- Geant4 has no default unit.
- To introduce the input data, unit **must** be “multiplied” to the number.
  - for example :  

```
G4double width = 12.5*m;
```

```
G4double density = 2.7*g/cm3;
```
  - Almost all commonly used units are available.
  - The user can define new units.
  - Refer to CLHEP: `SystemOfUnits.h`
- To output the data you can **divide** a variable by a unit you want to get.

```
G4cout << dE / MeV << “ (MeV)” << G4endl;
```

# System of Units

- System of units are defined in CLHEP, based on:
  - millimetre (**mm**), nanosecond (**ns**), Mega eV (**MeV**), positron charge (**eplus**) degree Kelvin (**kelvin**), the amount of substance (**mole**), luminous intensity (**candela**), radian (**radian**), steradian (**steradian**)
- All other units are computed from the basic ones
- Alternative way to output data: Geant4 can choose the most appropriate unit to use. Just specify the *category* for the data (**Length**, **Time**, **Energy**, etc...):

```
G4cout << G4BestUnit(StepSize, "Length") ;
```

**StepSize** will be printed in **km**, **m**, **mm** Or ... **fermi**, depending on its actual value

# Defining new Units

- New units can be defined directly as constants, or (suggested way) via `G4UnitDefinition`
  - `G4UnitDefinition` ( name, symbol, category, value )
- Example (speed):
  - `G4UnitDefinition` ("km/hour", "km/h",  
"Speed", km/3600\*s) ;
  - The new category "Speed" will be registered in the kernel in `G4UnitsTable`
- To print the list of units:
  - From the code  
`G4UnitDefinition::PrintUnitsTable()` ;
  - At run-time, as UI command:  
`Idle> /units/list`

# Definition of materials

- Different kinds of materials can be defined:
  - isotopes <> **G4Isotope**
  - elements <> **G4Element**
  - molecules <> **G4Material**
  - compounds and mixtures <> **G4Material**
- Attributes associated:
  - temperature, pressure, state, density
- **G4Isotope** and **G4Element** describe properties of the *atoms*:
  - Atomic number, number of nucleons, mass of a mole, shell energies, cross-sections per atoms, etc...
- **G4Material** describes the *macroscopic* properties of the matter:
  - temperature, pressure, state, density
  - Radiation length, absorption length, etc...
- **G4Material** is the only class used and visible to the toolkit: it is used by tracking, geometry and physics

# Elements and isotopes

- Isotopes can be assembled into elements

```
G4Isotope (const G4String& name,  
           G4int          z,      // atomic number  
           G4int          n,      // number of nucleons  
           G4double       a );   // mass of mole
```

- ... building elements as follows:

```
G4Element (const G4String& name,  
           const G4String& symbol, // element symbol  
           G4int          nIso );  // n. of isotopes  
G4Element::AddIsotope (G4Isotope* iso, // isotope  
                       G4double relAbund); // fraction of  
                                           // atoms  
                                           // per volume
```

# Materials of one element and molecules

- Single element material:

```
G4double z, a, density
density = 1.390*g/cm3;
a = 39.95*g/mole;
G4Material* lAr =
    new G4Material("liquidArgon", z=18, a, density);
```

- A molecule material (composition by number of atoms):

```
a = 1.01*g/mole;
G4Element* elH =
    new G4Element("Hydrogen", symbol="H", z=1., a);
a = 16.00*g/mole;
G4Element* elO =
    new G4Element("Oxygen", symbol="O", z=8., a);
density = 1.000*g/cm3;
G4Material* H2O =
    new G4Material("Water", density, ncomponents=2);
H2O->AddElement(elH, natoms=2);
H2O->AddElement(elO, natoms=1);
```



# Compound and mixture

- Mixture: composition by fraction of mass

```
a = 14.01*g/mole;  
G4Element* elN = new G4Element(name="Nitrogen",symbol="N",z= 7.,a) ;  
a = 16.00*g/mole;  
G4Element* elO = new G4Element(name="Oxygen",symbol="O",z= 8.,a) ;  
density = 1.290*mg/cm3;  
G4Material* Air = new G4Material(name="Air",density,ncomponents=2) ;  
Air->AddElement(elN, 70.0*perCent) ;  
Air->AddElement(elO, 30.0*perCent) ;
```

- Composition of compound materials

```
G4Element* elC = ...;    // define "carbon" element  
G4Material* SiO2 = ...;  // define "quartz" material  
G4Material* H2O = ...;   // define "water" material  
density = 0.200*g/cm3;  
G4Material* Aerog = new G4Material("Aerogel",density,ncomponents=3) ;  
Aerog->AddMaterial(SiO2,fractionmass=62.5*perCent) ;  
Aerog->AddMaterial(H2O ,fractionmass=37.4*perCent) ;  
Aerog->AddElement (elC ,fractionmass= 0.1*perCent) ;
```

# Example: gas

- It may be necessary to specify temperature and pressure
  - $(dE/dx)$  computation affected)

```
G4double density = 27.*mg/cm3;  
G4double temperature = 325.*kelvin;  
G4double pressure = 50.*atmosphere;  
G4Material* CO2 = new G4Material("CarbonicGas", density,  
    ncomponents=2, kStateGas, temperature, pressure);  
CO2->AddElement(C,natoms = 1);  
CO2->AddElement(O,natoms = 2);
```

- Absolute vacuum does not exist: gas at very low density !
  - Cannot define materials composed of multiple elements through **Z** or **A**, or with  $\rho=0$

```
G4double atomicNumber = 1.;  
G4double massOfMole = 1.008*g/mole;  
G4double density = 1.e-25*g/cm3;  
G4double temperature = 2.73*kelvin;  
G4double pressure = 3.e-18*pascal;  
G4Material* Vacuum = new G4Material("interGalactic",  
    atomicNumber,massOfMole, density, kStateGas,temperature, pressure);10
```

# NIST Material Data-Base in Geant4

- NIST database for materials is imported inside Geant4  
<http://physics.nist.gov/PhysRefData>
- Additional interfaces defined
- UI commands specific for handling materials
- The best accuracy for the most relevant parameters guaranteed:
  - Density
  - Mean excitation potential
  - Chemical bounds
  - Element composition
  - Isotope composition
  - Various corrections

Z	A	m	error	(%)	A <sub>eff</sub>
=====					
14	Si 22	22.03453	(22)		28.0855(3)
	23	23.02552	(21)		
	24	24.011546	(21)		
	25	25.004107	(11)		
	26	25.992330	(3)		
	27	26.98670476	(17)		
	28	27.9769265327	(20)	92.2297 (7)	
	29	28.97649472	(3)	4.6832 (5)	
	30	29.97377022	(5)	3.0872 (5)	
	31	30.97536327	(7)		
	32	31.9741481	(23)		
	33	32.978001	(17)		
	34	33.978576	(15)		
	35	34.984580	(40)		
	36	35.98669	(11)		
	37	36.99300	(13)		
	38	37.99598	(29)		
	39	39.00230	(43)		
	40	40.00580	(54)		
	41	41.01270	(64)		
	42	42.01610	(75)		

- *Natural isotope compositions*
- *More than 3000 isotope masses*

# NIST materials in Geant4

```
#####
### Elementary Materials from the NIST Data
Base
#####
```

Z	Name	ChFormula	density(g/cm^3)	I(eV)
1	G4_H	H_2	8.3748e-05	19.2
2	G4_He		0.000166322	41.8
3	G4_Li		0.534	40
4	G4_Be		1.848	63.7
5	G4_B		2.37	76
6	G4_C		2	81
7	G4_N	N_2	0.0011652	82
8	G4_O	O_2	0.00133151	95
9	G4_F		0.00158029	115
10	G4_Ne		0.000838505	137
11	G4_Na		0.971	149
12	G4_Mg		1.74	156
13	G4_Al		2.6989	166
14	G4_Si		2.33	173

```
#####
### Compound Materials from the NIST Data Base
#####
```

N	Name	ChFormula	density(g/cm^3)	I(eV)
13	G4_Adipose_Tissue		0.92	63.2
	1	0.119477		
	6	0.63724		
	7	0.00797		
	8	0.232333		
	11	0.0005		
	12	2e-05		
	15	0.00016		
	16	0.00073		
	17	0.00119		
	19	0.00032		
	20	2e-05		
	26	2e-05		
	30	2e-05		
4	G4_Air		0.00120479	85.7
	6	0.000124		
	7	0.755268		
	8	0.231781		
	18	0.012827		
2	G4_CsI		4.51	553.1
	53	0.47692		
	55	0.52308		

- NIST Elementary materials:
  - H -> Cf ( Z = 1 -> 98 )
- NIST compounds:
  - e.g. “**G4\_ADIPOSE\_TISSUE\_IRCP**”
- HEP and Nuclear materials:
  - e.g. **Liquid Ar, PbWO**
- It is possible to build mixtures of NIST and user-defined materials

# How to use the NIST DB

- No need to predefine elements and materials
- Retrieve materials from NIST manager:

```
G4NistManager* manager = G4NistManager::Instance();
```

```
G4Material* H2O = manager->FindOrBuildMaterial("G4_WATER");
```

- Some UI commands ...

```
/material/nist/printElement    ← print defined elements
```

```
/material/nist/listMaterials    ← print defined materials
```

# Part II: Geometry

## Geometry:

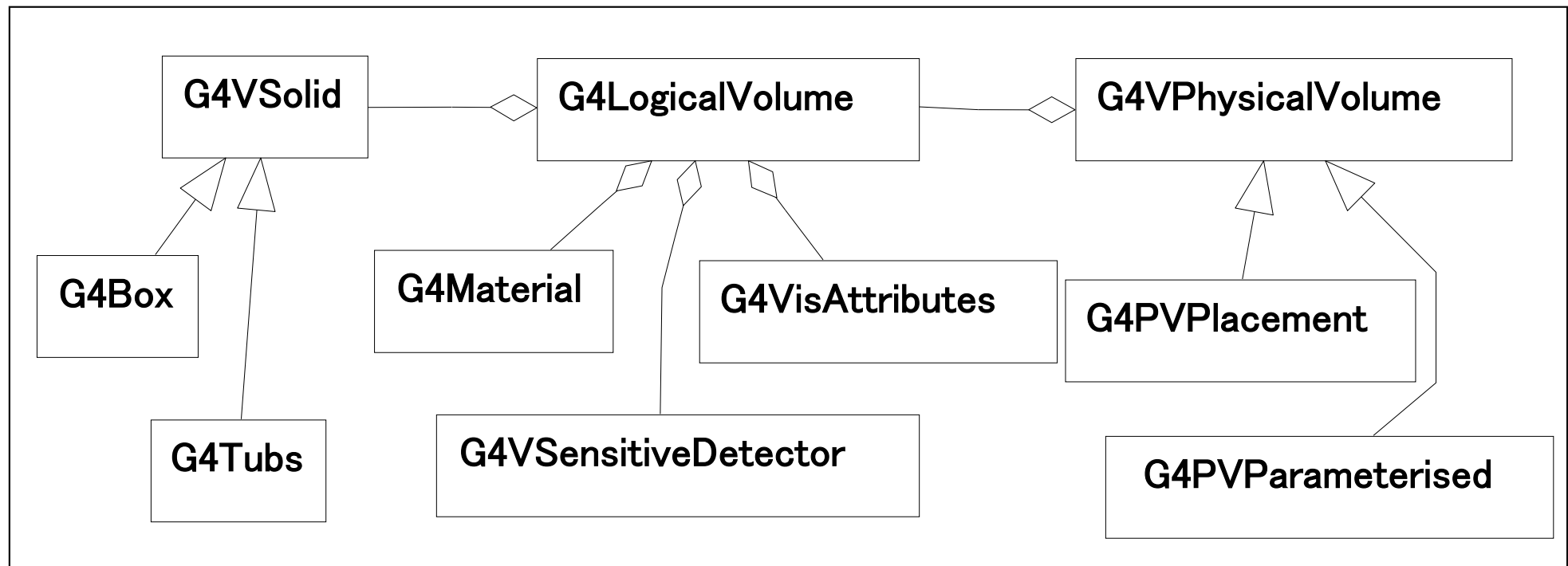
- Detector description: the basic
  - Detector geometry components
  - Define detector geometry
- Describing a detector
  - Solids
  - Logical volumes
  - Physical volumes
- Tools for geometry check

# Describe your detector

- A detector geometry is made of a number of volumes
- The largest volume is called **World** volume
  - It must contain all other volumes
- Derive your own concrete class from `G4VUserDetectorConstruction` abstract base class.
- Implementing the method `construct()`:
  - Modularize it according to each detector component:
    - Define shapes/solids required to describe the geometry
    - Construct all necessary materials
    - Construct and place volumes of your detector geometry
    - Define sensitive detectors and identify detector volumes which to associate them
    - Associate magnetic field to detector regions
    - Define visualization attributes for the detector elements

# Detector geometry components

- Three conceptual layers
  - **G4VSolid** -- *shape, size*
  - **G4LogicalVolume** -- *daughter physical volumes,*  
*material, sensitivity, magnetic field, etc.*
  - **G4VPhysicalVolume** -- *position, rotation*



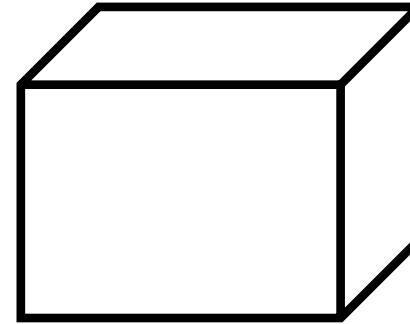


# Define detector geometry

- Basic strategy

Solid : shape and size

```
G4VSolid* pBoxSolid =  
    new G4Box("aBoxSolid",  
              1.*m, 2.*m, 3.*m);
```



## Step 1

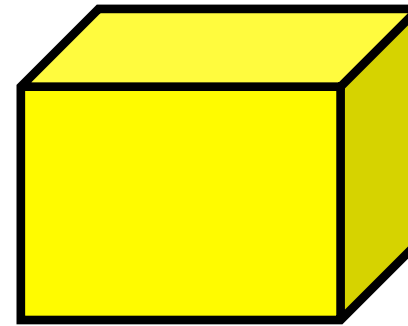
Create the  
geom. object :  
box

# Define detector geometry

- Basic strategy

```
G4VSolid* pBoxSolid =  
    new G4Box("aBoxSolid",  
              1.*m, 2.*m, 3.*m);  
  
G4LogicalVolume* pBoxLog =  
    new G4LogicalVolume( pBoxSolid,  
                          pBoxMaterial, "aBoxLog", 0, 0, 0);
```

Logical volume :  
+ material, sensitivity, etc.



## Step 1

Create the  
geom. object :  
box



## Step 2

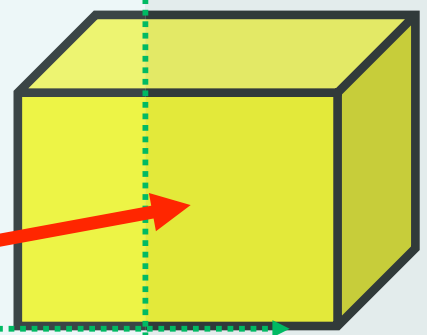
Assign properties  
to object :  
material

# Define detector geometry

- Basic strategy

Physical volume :  
+ rotation and position

```
G4VSolid* pBoxSolid =  
  
    new G4Box("aBoxSolid",  
              1.*m, 2.*m, 3.*m);  
  
G4LogicalVolume* pBoxLog =  
    new G4LogicalVolume( pBoxSolid,  
                          pBoxMaterial, "aBoxLog", 0, 0, 0);  
  
G4VPhysicalVolume* aBoxPhys =  
    new G4PVPlacement(pRotation,  
                      G4ThreeVector(posX, posY, posZ), pBoxLog,  
                      "aBoxPhys", pMotherLog, 0, copyNo);
```



## Step 1

Create the  
geom. object :  
box



## Step 2

Assign properties  
to object :  
material



## Step 3

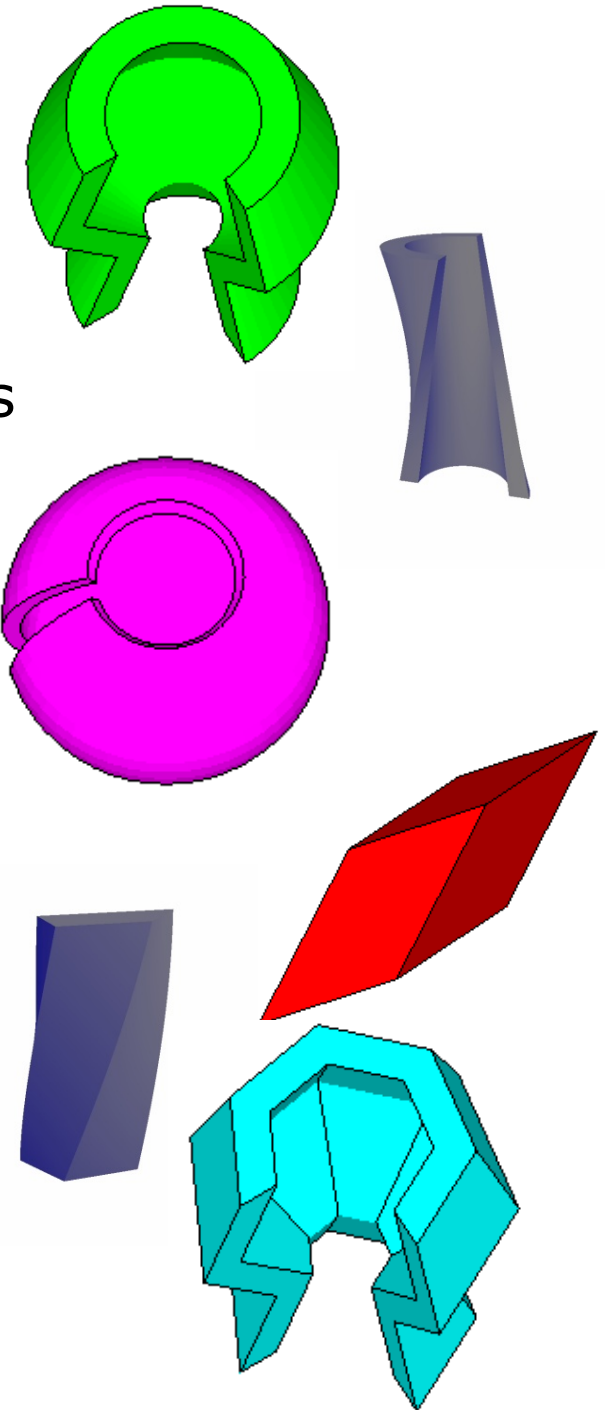
Place it in the  
coordinate  
system of  
mother volume

- A unique physical volume which represents the experimental area must exist and fully contains all other components

## ➤ The world volume

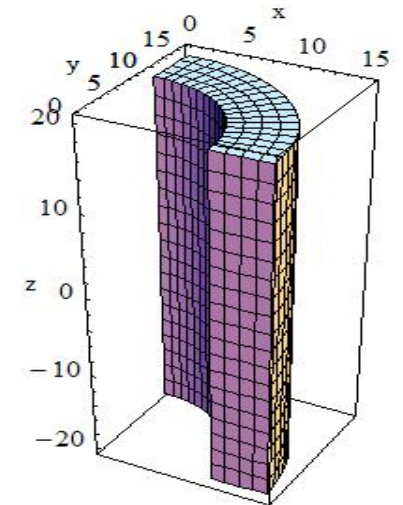
# Solids

- Solids defined in Geant4:
  - CSG (Constructed Solid Geometry) solids
    - G4Box, G4Tubs, G4Cons, G4Trd, ...
    - Analogous to simple GEANT3 CSG solids
  - Specific solids (CSG like)
    - G4Polycone, G4Polyhedra, G4Hype, ..
    - G4TwistedTubs, G4TwistedTrap, ...
  - BREP (Boundary REPresented) solids
    - G4BREPSolidPolycone, G4BSplineSurface, ...
    - Any order surface
  - Boolean solids
    - G4UnionSolid, G4SubtractionSolid, ...

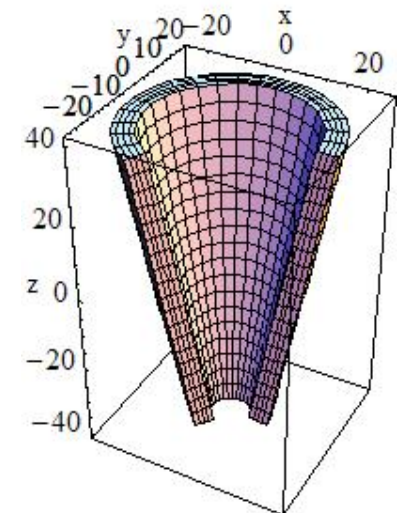


# CSG: G4Tubs, G4Cons

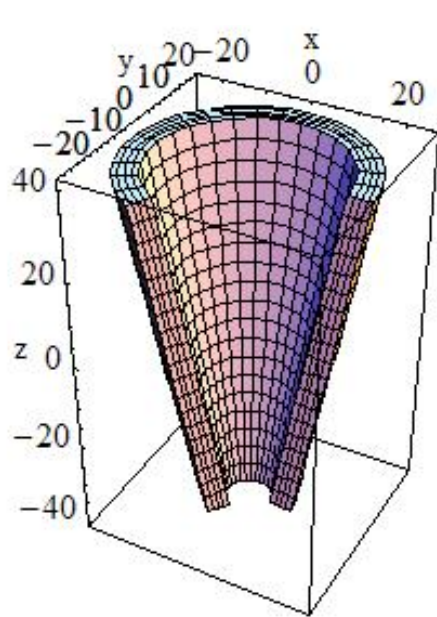
```
G4Tubs(const G4String& pname, // name
        G4double pRmin, // inner radius
        G4double pRmax, // outer radius
        G4double pDz, // Z half length
        G4double pSphi, // starting Phi
        G4double pDphi); // segment angle
```



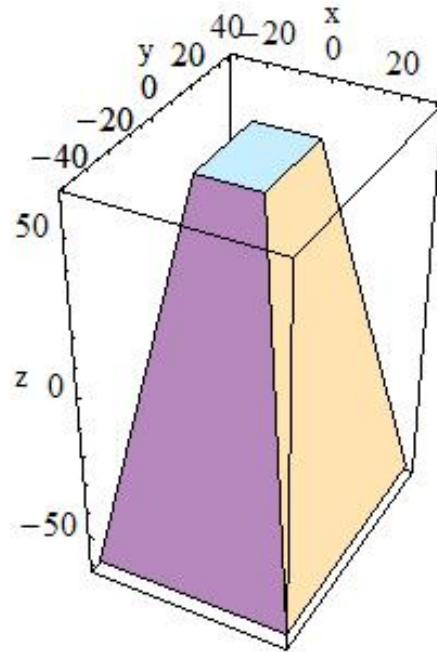
```
G4Cons(const G4String& pname, // name
        G4double pRmin1, // inner radius -pDz
        G4double pRmax1, // outer radius -pDz
        G4double pRmin2, // inner radius +pDz
        G4double pRmax2, // outer radius +pDz
        G4double pDz, // Z half length
        G4double pSphi, // starting Phi
        G4double pDphi); // segment angle
```



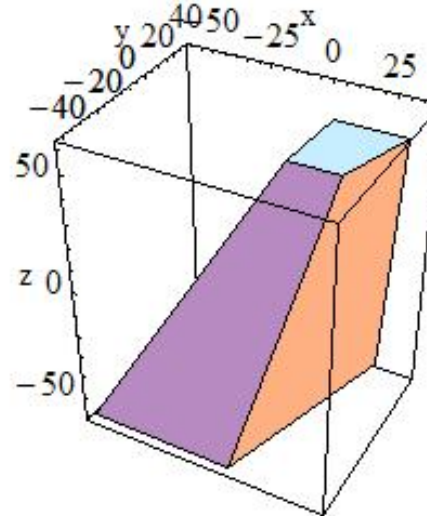
# Other CSG solids



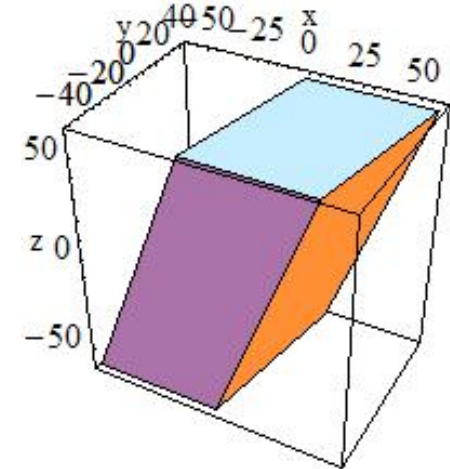
G4Cons



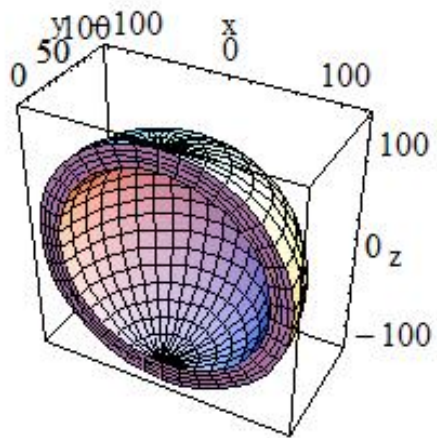
G4Trd



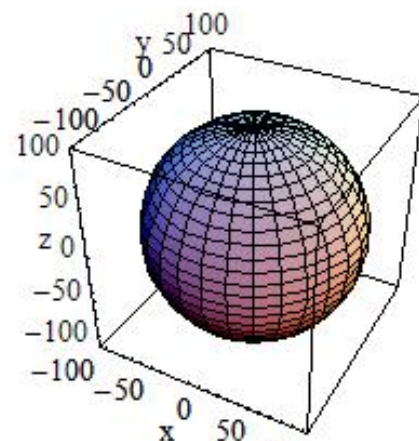
G4Trap



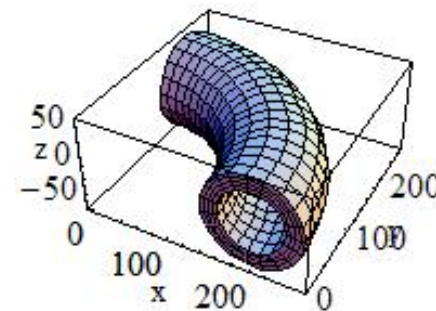
G4Para  
(parallelepiped)



G4Sphere



G4Orb  
(full solid sphere)

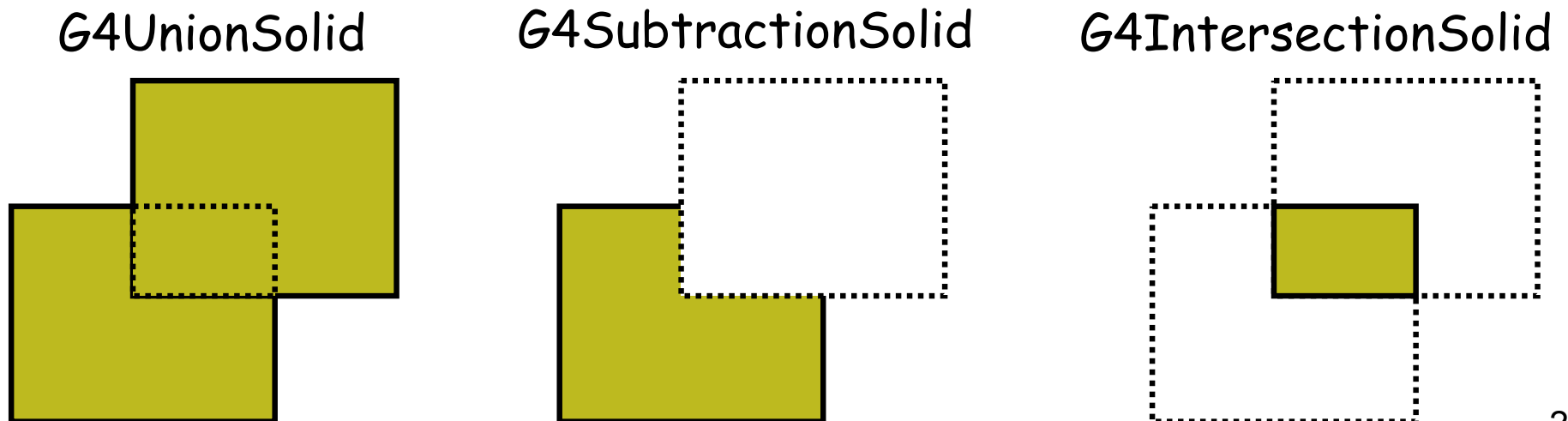


G4Torus

Consult to  
[Section 4.1.2 of Geant4 Application Developers Guide](#) for all available shapes.

# Boolean Solids

- ▶ Solids can be combined using boolean operations:
  - ▶ **G4UnionSolid, G4SubtractionSolid, G4IntersectionSolid**
  - ▶ Requires: 2 solids, 1 boolean operation, and an (optional) transformation for the 2<sup>nd</sup> solid
  - ▶ 2<sup>nd</sup> solid is positioned relative to the coordinate system of the 1<sup>st</sup> solid
  - ▶ Result of boolean operation becomes a solid. Thus the third solid can be combined to the resulting solid of first operation.
- ▶ Solids to be combined can be either CSG or other Boolean solids.
- ▶ **Note:** tracking cost for the navigation in a complex Boolean solid is proportional to the number of constituent CSG solids

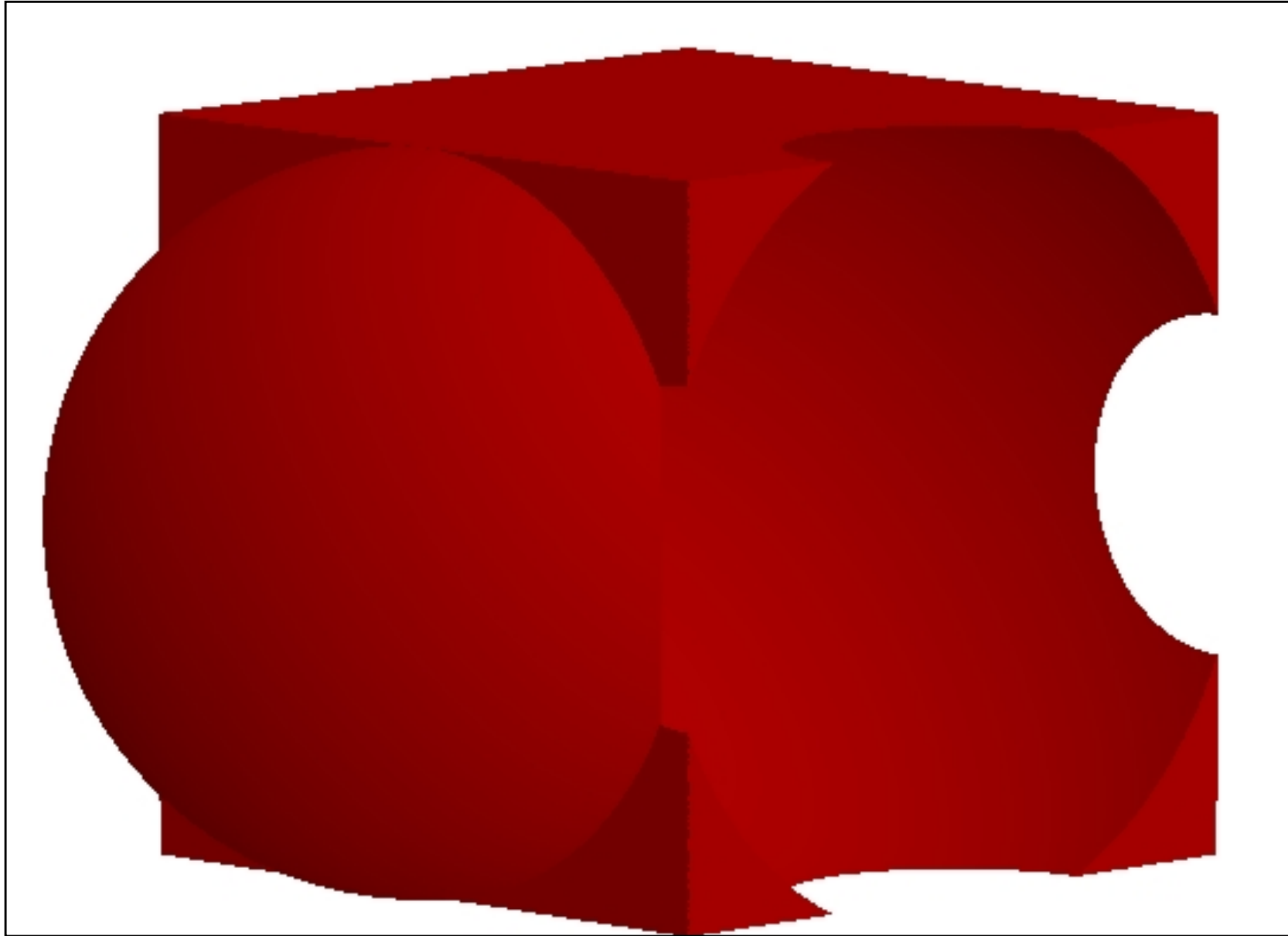


# Boolean Solids - example

```
G4VSolid* box = new G4Box("Box", 50*cm, 60*cm, 40*cm) ;  
G4VSolid* cylinder  
    = new G4Tubs("Cylinder", 0., 50.*cm, 50.*cm, 0., 2*M_PI*rad) ;  
  
G4VSolid* union  
    = new G4UnionSolid("Box+Cylinder", box, cylinder) ;  
  
G4VSolid* subtract  
    = new G4SubtractionSolid("Box-Cylinder", box, cylinder,  
        0, G4ThreeVector(30.*cm, 0., 0.)) ;  
  
G4RotationMatrix* rm = new G4RotationMatrix() ;  
rm->RotateX(30.*deg) ;  
G4VSolid* intersect  
    = new G4IntersectionSolid("Box&&Cylinder",  
        box, cylinder, rm, G4ThreeVector(0., 0., 0.)) ;
```



# Boolean Solids



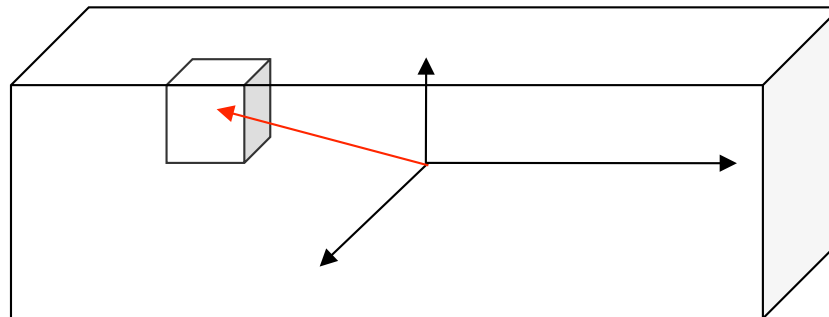
# G4LogicalVolume

```
G4LogicalVolume(G4VSolid* pSolid,  
                G4Material* pMaterial,  
                const G4String& name,  
                G4FieldManager* pFieldMgr=0,  
                G4VSensitiveDetector* pSDetector=0,  
                G4UserLimits* pULimits=0,  
                G4bool optimise=true);
```

- Contains all information of volume except position:
  - Shape and dimension (G4VSolid)
  - Material, sensitivity, visualization attributes
  - Position of daughter volumes
  - Magnetic field, User limits
- Physical volumes of same type can share a logical volume.
- The pointers to solid and material must be NOT null

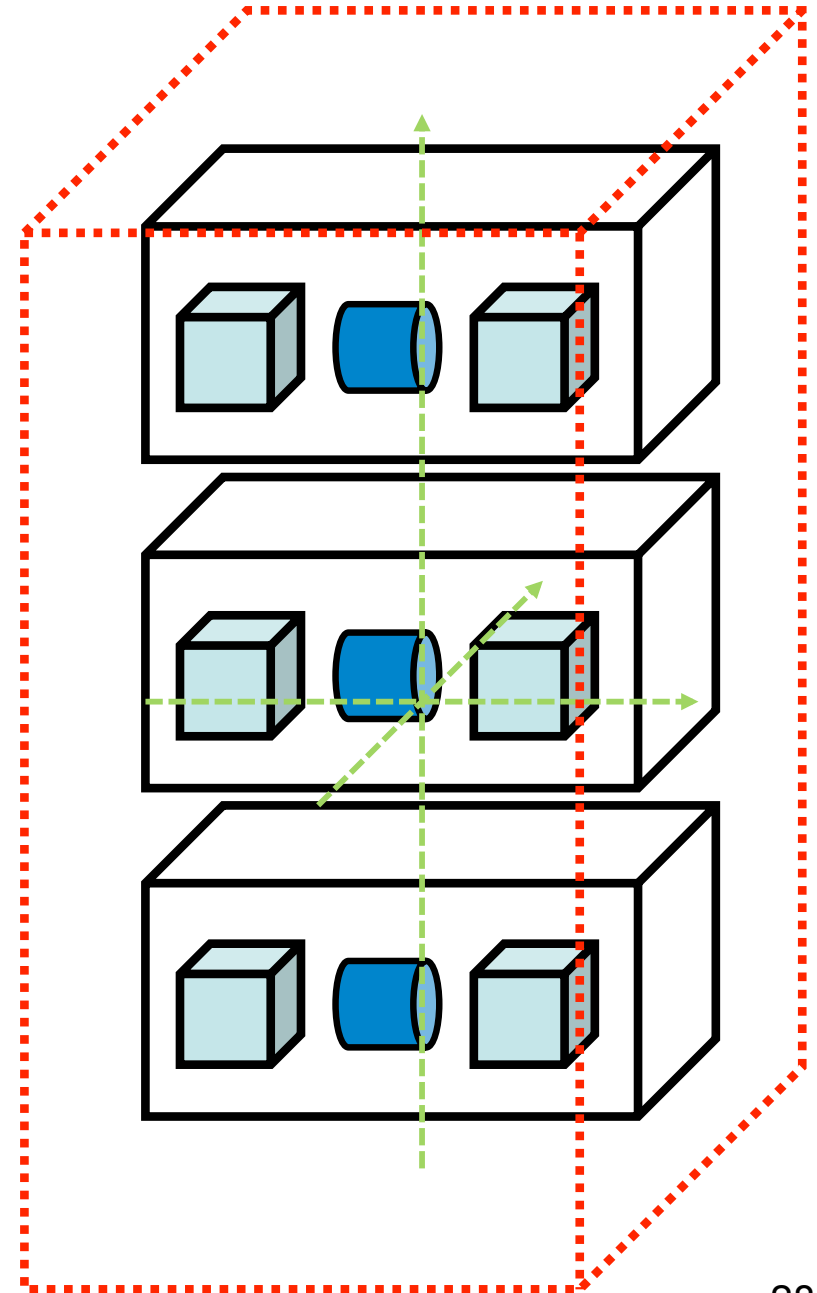
# Geometrical hierarchy

- Mother and daughter volumes
  - A volume is placed in its mother volume
    - Position and rotation of the daughter volume is described with respect to the local coordinate system of the mother volume
    - The origin of the mother's local coordinate system is at the center of the mother volume
    - Daughter volumes cannot protrude from the mother volume
    - Daughter volumes cannot overlap
  - The logical volume of mother knows the daughter volumes it contains
    - It is uniquely defined to be their mother volume



# Geometrical hierarchy

- One logical volume can be placed more than once. One or more volumes can be placed in a mother volume
- Note that the mother-daughter relationship is an information of **G4LogicalVolume**
  - If the mother volume is placed more than once, all daughters by definition appear in each placed physical volume
- The **world volume** must be a unique physical volume which fully contains with some margin all the other volumes (root volume of the hierarchy)
  - The world volume defines the **global coordinate system**. The origin of the global coordinate system is at the center of the world volume
  - Position of a track is given with respect to the global coordinate system

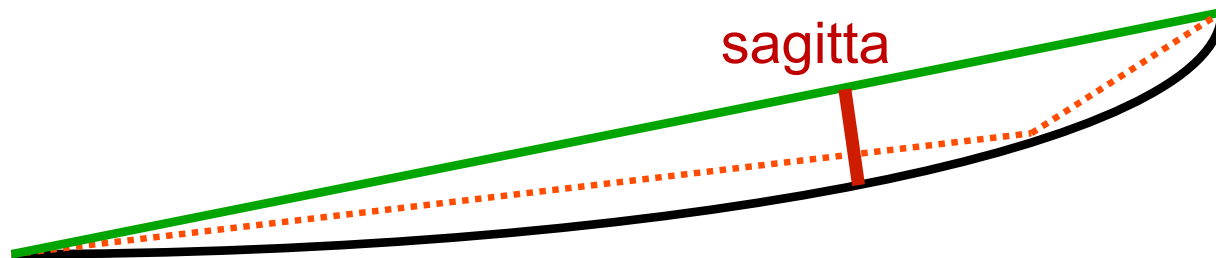


# Region

- A **region** may have its unique
  - Production thresholds (cuts)
    - If a region in the mass geometry does not have its own production thresholds, those of the default region are used (i.e., may not be those of the parent region).
  - User limits
    - Artificial limits affecting to the tracking, e.g. max step length, max number of steps, min kinetic energy left, etc.
    - You can set user limits directly to logical volume as well. If both logical volume and associated region have user limits, those of logical volume wins.
  - Fast simulation manager
  - Field manager
  - ...
- Please note :
  - World logical volume is recognized as **the default region**. User is **not** allowed to define a region to the world logical volume.

# Fields

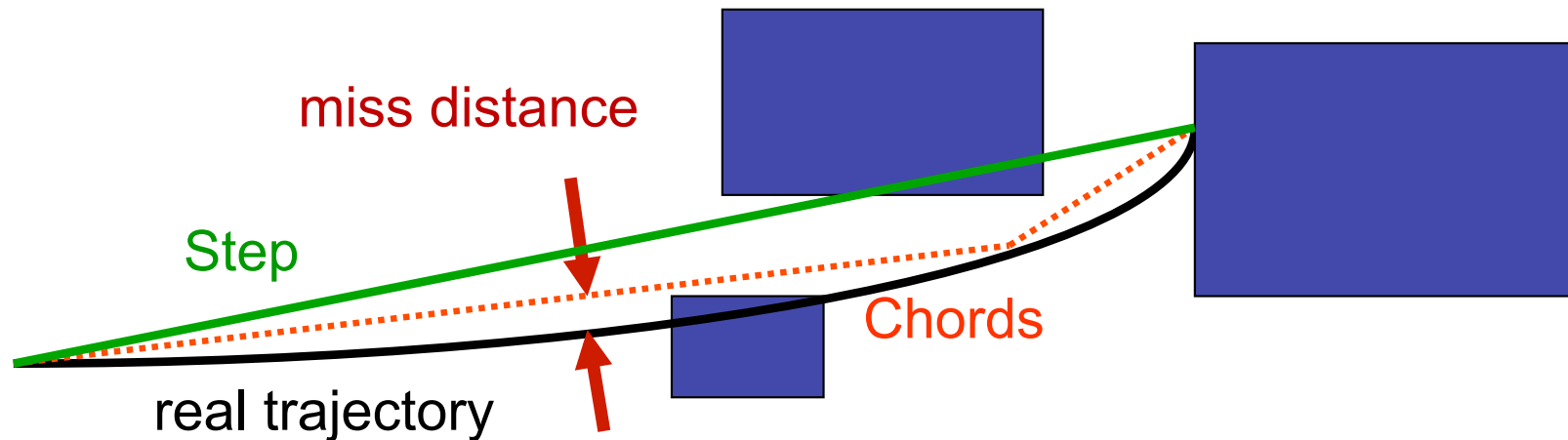
- In order to propagate a particle inside a field (e.g. magnetic, electric or both), we integrate the equation of motion of the particle in the field
- In general this is best done using a **Runge-Kutta** (RK) method for the integration of ordinary differential equations
  - Several RK methods are available
- In specific cases other solvers can also be used
- Once a method is chosen for calculating the track's motion in a field, Geant4 breaks up this curved path into linear chord segments



- The chord segments are determined so that they closely approximate the curved path; they're chosen so that their **sagitta** is small enough
  - The *sagitta* is the maximum distance between the curved path and the straight line

# Fields

- Chords are used to interrogate the Navigator
  - to see whether the track has crossed a volume boundary



- The accuracy of the volume intersection can be tuned
  - by setting a parameter called the “*miss distance*”:
    - Upper bound for the value of the sagitta
    - It provide a measure of the error resolution
    - Setting small *miss distance* may be highly CPU consuming
- One step can consist of more than one chord
  - In some cases, one step consists of several turns

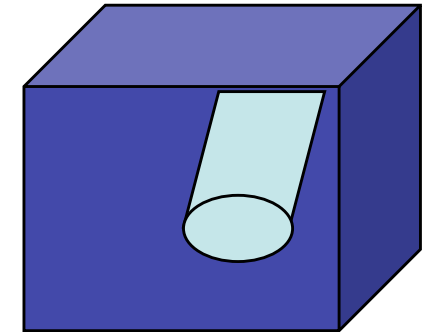
# Create a Magnetic Field

- How to create Uniform field ?
  - Use the constructor of `G4UniformMagField`
    - `G4ThreeVector fieldV ( 0.1 * Tesla, 1.0*Gauss, 0.0 );`
    - `G4MagneticField *myField = new G4UniformField( fieldV );`
- Non-uniform field
  - Concrete class derived from `G4MagneticField`
- A field is packaged together with properties and accuracy parameters into a **field manager**:
  - `G4FieldManager* localFieldMgr = new G4FieldManager(myField);`
- **One** field manager is associated with the 'world'
- Other volumes/regions in the geometry can override this
  - An alternative field manager can be associated with any logical volume:  
`G4FieldManager* localFieldMgr = new G4FieldManager(myField);`  
`logVolume->setFieldManager(localFieldMgr, true);`
  - The field is propagated to all the daughter volumes.

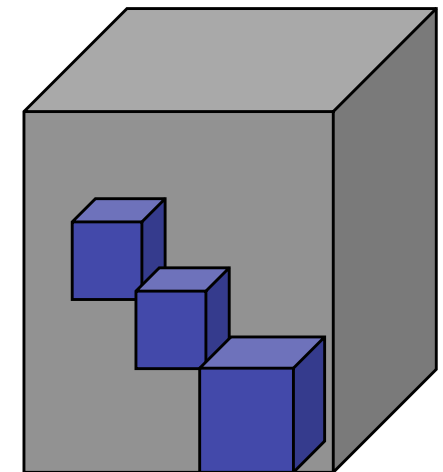


# Physical Volumes

- **Placement:** it is one positioned volume
  - can represent any number of volumes
  - reduces use of memory.
  - Replica
    - simple repetition, similar to G3 divisions
  - Parameterised
- **Repeated:** a volume placed many times



*placement*



*repeated*

# G4PVPlacement

```
G4PVPlacement(G4RotationMatrix* pRot,      // rotation of mother frame
               const G4ThreeVector& tlate, // position in rotated frame
               G4LogicalVolume* pCurrentLogical,
               const G4String& pName,
               G4LogicalVolume* pMotherLogical,
               G4bool pMany,                // not used. Set it to false...
               G4int pCopyNo,               // unique arbitrary index
               G4bool pSurfChk=false);      // optional overlap check
```

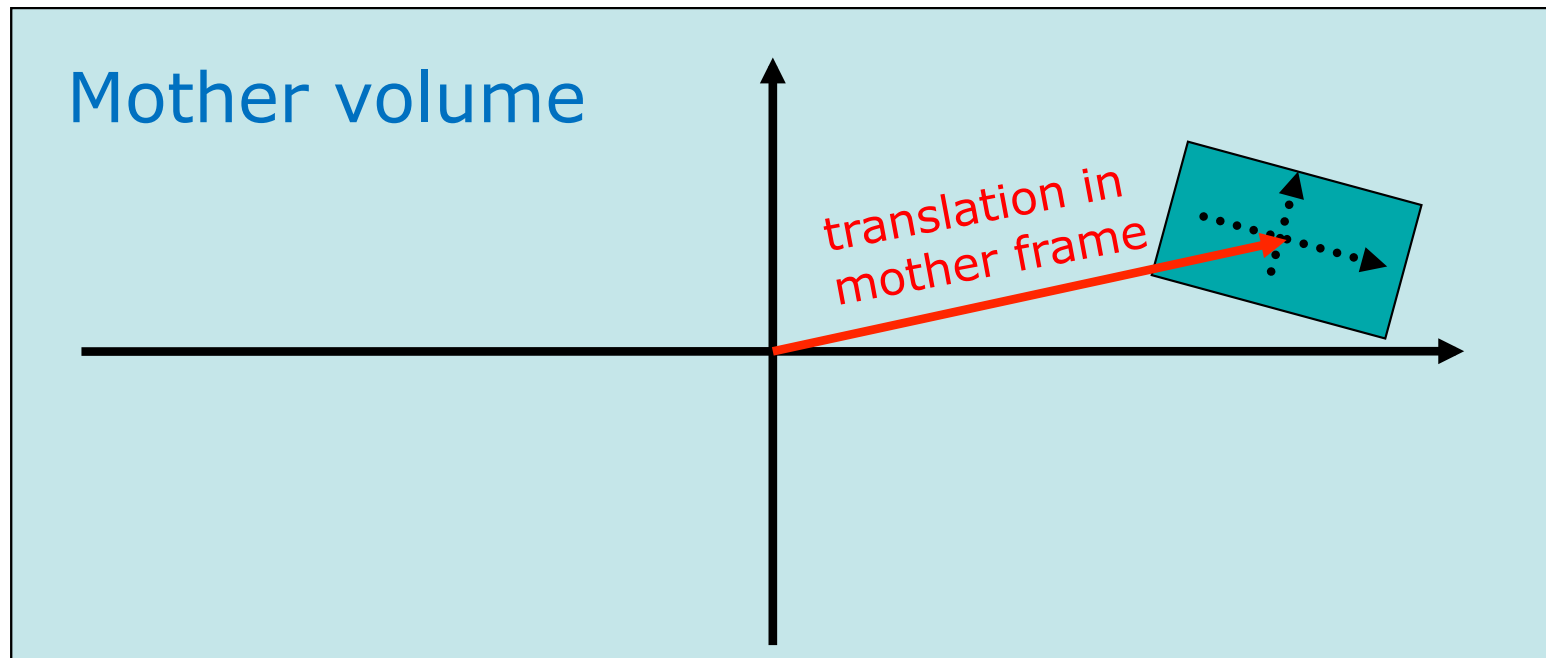
- Single volume positioned relatively to the mother volume
  - In a frame rotated and translated relative to the coordinate system of the mother volume
- Three additional constructors:
  - A simple variation: specifying the mother volume as a pointer to its physical volume instead of its logical volume.
  - Using `G4Transform3D` to represent the direct rotation and translation of the solid instead of the frame (*alternative constructor*)
  - The combination of the two variants above

# G4PVPlacement

## Rotation of mother frame ...

```
G4PVPlacement(G4RotationMatrix* pRot,          // rotation of mother frame
               const G4ThreeVector& tlate,    // position in mother frame
               G4LogicalVolume* pCurrentLogical,
               const G4String& pName,
               G4LogicalVolume* pMotherLogical,
               G4bool pMany,                  // not used. Set it to false..
               G4int pCopyNo,                 // unique arbitrary index
               G4bool pSurfChk=false );      // optional overlap check
```

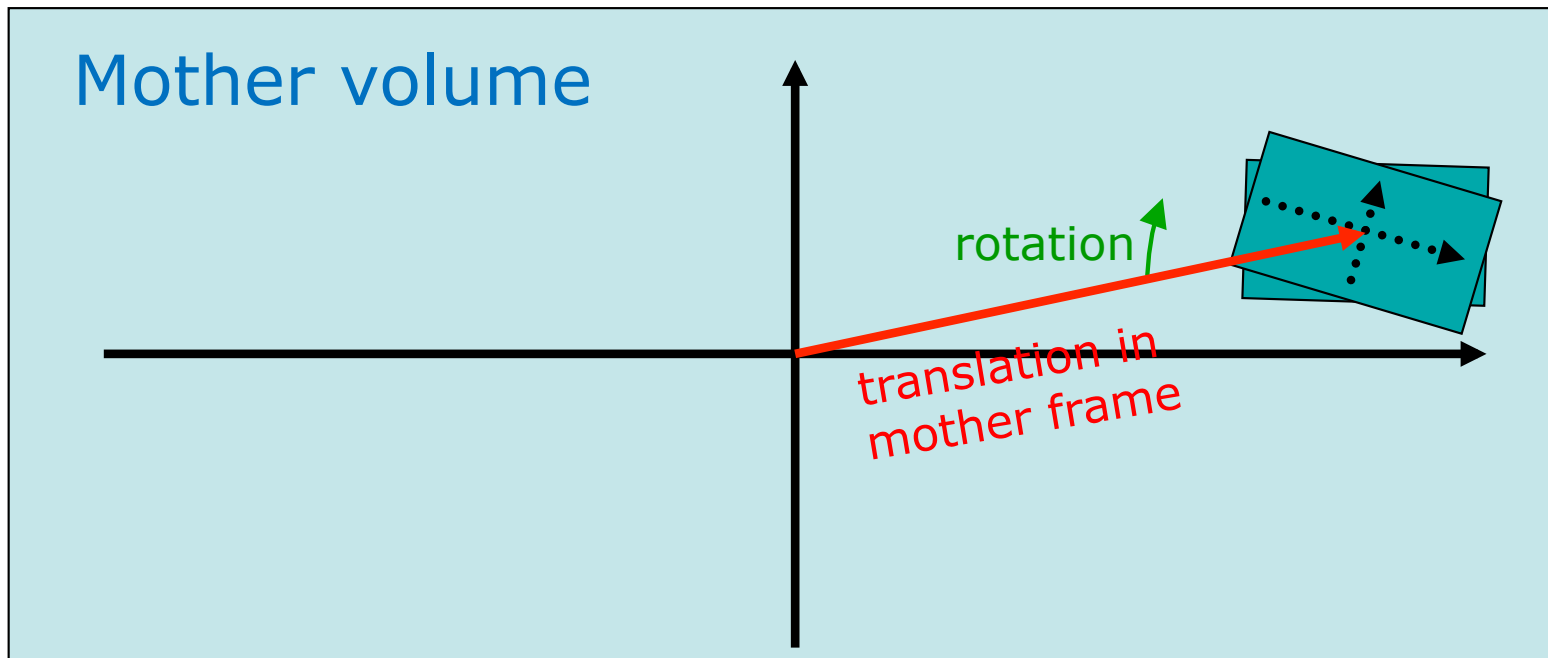
- Single volume positioned relatively to the mother volume



# G4PVPlacement

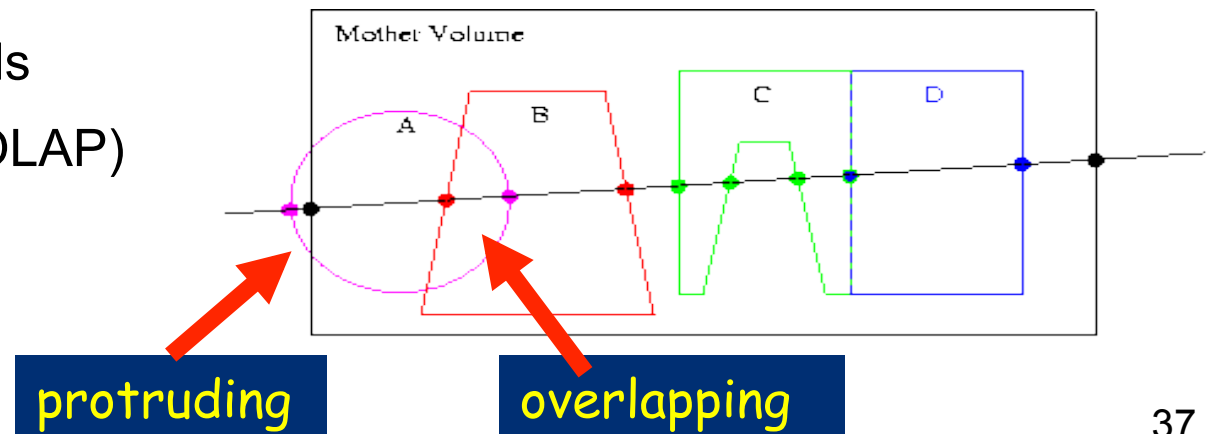
## Rotation in mother frame ...

```
G4PVPlacement( G4Transform3D( G4RotationMatrix &pRot,          // rotation of daughter frame
                             const G4ThreeVector &tlate), // position in mother frame
               G4LogicalVolume *pDaughterLogical,
               const G4String &pName,
               G4LogicalVolume *pMotherLogical,
               G4bool pMany,          // not used, set it to false...
               G4int pCopyNo,        // unique arbitrary integer
               G4bool pSurfChk=false ); // optional overlap check
```



# Tools for geometry check

- A **protruding** volume is a contained daughter volume which actually **protrudes** from its mother volume.
- When volumes in a common mother actually **intersect themselves** are defined as **overlapping**.
- Geant4 **does not allow** for malformed geometries, **neither protruding nor overlapping**.
  - The behavior of navigation is unpredictable for such cases.
- The problem of detecting overlaps between volumes is bounded by the complexity of the solid models description.
- Utilities are provided for detecting wrong positioning
  - Optional checks at construction
  - Kernel run-time commands
  - Graphical tools (DAVID, OLAP)



# Tools for geometry check

- Constructors of **G4PVPlacement** and **G4PVParameterised** have an optional argument “pSurfChk”.

```
G4PVPlacement(G4RotationMatrix* pRot,  
              const G4ThreeVector &tlate,  
              G4LogicalVolume *pDaughterLogical,  
              const G4String &pName,  
              G4LogicalVolume *pMotherLogical,  
              G4bool pMany, G4int pCopyNo,  
              G4bool pSurfChk=false);
```

If this flag is true, overlap check is done at the construction:

- some number of points are randomly sampled on the surface of creating volume.

*This check requires lots of CPU time, but it is worth to try at least once.*

- Built-in run-time commands to activate verification tests for the user geometry:
  - **geometry/test/run** or **geometry/test/grid\_test**  
*to start verification of geometry for overlapping regions based on a standard grid setup, limited to the first depth level*
  - **geometry/test/recursive\_test**  
*applies the grid test to all depth levels (may require lots of CPU time!)*
  - **geometry/test/line\_test**  
*to shoot a line along a specified direction and position*
  - ...

# Tools for geometry check

```
void MGManagerDetectorConstruction::CheckOverlaps()
{
    G4PhysicalVolumeStore* thePVStore =
    G4PhysicalVolumeStore::GetInstance();
    G4cout << thePVStore->size() << " physical volumes are
defined" << G4endl;
    G4bool overlapFlag = false;
    G4int res=1000;
    G4double tol=0.; //tolerance
    for (size_t i=0;i<thePVStore->size();i++)
    {
        overlapFlag = (*thePVStore)[i]->CheckOverlaps
(res,tol,fCheckOverlapsVerbosity) | overlapFlag;
    }
    if (overlapFlag)
        G4cout << "Check: there are overlapping volumes" <<
G4endl;
}
```

## Output example:

```
136 physical volumes are defined
Checking overlaps for volume BeamLineSupport ... OK!
Checking overlaps for volume BeamLineCover ... OK!
Checking overlaps for volume BeamLineCover2 ... OK!
Checking overlaps for volume VacuumZone ... OK!
Checking overlaps for volume FirstScatteringFoil ... OK!
...
...
...
----- WWWWW ----- G4Exception-START ----- WWWWW -
*** G4Exception : GeomVol1002 issued by :
    G4PVPlacement::CheckOverlaps()
    Overlap with volume already placed !
    Overlap is detected for volume BrassTube2
        with HoleNozzleSupport volume's
    local point (12.6381,12.8171,-25.1867), overlapping by
        at least: 3.5 mm
*** This is just a warning message. ***
----- WWWWW ----- G4Exception-END ----- WWWWW -
...
...
```

This method can be called at any point after run->Initialize();

**Thanks for your attention**