**User’s guide for MINOS simulations**

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**The main file**

MINOS\_simulation: MINOS\_sim.cc and executable MINOS\_sim

The main files allow constructing the different classes and the visualization window.

**The different folders**

The folder src/ contains the source files of the classes.

The folder include/ contains the header files of the classes.

The folder result/ contains the result files of the simulation.

The folder Inputs/ contains the files used to define some characteristics of the setup and read by the libraries.

The folder lib/ contains some libraries used by the simulation.

**The Inputs files**

ConfigFileBeam.txt: definition of the characteristics of the beam which is sent in the setup in terms of energy, momentum, position, charge and mass number.

ConfigFileSetup.txt: definition of the dimensions of the elements constituting the setup

ConfigFileRings.txt: definition of all the parameters used in the second step of the simulations for the drift of ionization electrons

**The libraries**

For each library one has the header file .hh, the source file .cxx and the GNUmakefile.

The library used to define the beam

libExN03BeamIn.so: this library allows to determine the characteristics of the beam by reading the ConfigFileBeam.txt file.

ExN03BeamIn.cxx: source file defining the function ReadConfigFile() which allows to read ConfiFileBeam.txt

ExN03BeamIn.hh: header file defining the beam parameters

|  |  |
| --- | --- |
| MeanX | central position of the beam in X (mm) |
| SigmaX | standard deviation of the position in X (mm) |
| MeanY | central position of the beam in Y (mm) |
| SigmaY | standard deviation of the position in Y (mm) |
| Z0 | Z position of emission of the beam (mm) |
| MeanMomentumX | mean momentum along X |
| SigmaMomentumX | standard deviation in momentum along X |
| MeanMomentumY | mean momentum along Y |
| SigmaMomentumY | standard deviation in momentum along Y |
| MomentumZ0 | momentum along Z |
| MeanEnergy | mean total kinetic energy (MeV) |
| SigmaEnergy | standard deviation in energy (MeV) |
| Z | charge |
| A | mass |

The library used to define the setup

libExN03Setup.so: this library allows to determine the characteristics of the setup by reading the ConfigFileSetup.txt file.

ExN03Setup.cxx: source file defining the function ReadConfigFile() which allows to read ConfiFileSetup.txt

ExN03Setup.hh: header file defining the setup parameters

|  |  |
| --- | --- |
| TargetRadius | target radius in mm |
| TargetLength | target length in mm |
| ChamberInnerRadius | chamber inner radius in mm |
| ChamberThickness | chamber thickness in mm |
| ChamberLength | chamber length in mm |
| InnerRohacellThickness | inner Rohacell thickness in mm |
| KaptonThickness | Kapton thickness in mm |
| OuterRohacellThickness | outer Rohacell radius in mm |
| TPCRadiusExt | TPC outer radius in mm |
| WindowThickness | Mylar window thickness in mm |

The output library

libExN03datai.so: this library allows to define observables which will be saved by the simulation; for example energy loss as a function of the position in the different materials, total energy loss in the different materials, vertex information in terms of energy, momentum, position, charge and mass of particles created, detection in the TPC.

ExN03Datai.cxx: source file defining the function ClearEvent() which allows to clear all the observables at the beginning of a new event simulated

ExN03DataoRings.hh: header file defining the observables

|  |  |
| --- | --- |
| Z, A | charge and mass of particles |
| energy0 | kinetic energy of each particle at vertex (MeV) |
| theta0, phi0 | θ and φ angle of each particle at vertex (deg) |
| x0, y0, z0 | vertex coordinates (mm) of each particle |
| detection | Boolean of energy deposited by each particle in the TPC |
| Et\_tar | energy in eV deposited by each particle in the target |
| Et\_win | energy in eV deposited by each particle in the Mylar window |
| Et\_ch | energy in eV deposited by each particle in the chamber |
| Et\_InnerRohacell | energy in eV deposited by each particle in the inner Rohacell |
| Et\_Kapton | energy in eV deposited by each particle in the Kapton |
| Et\_OuterRohacell | energy in eV deposited by each particle in the outer Rohacell |
| Et\_tpc | energy in eV deposited by each particle in the TPC |
| Et\_tpc\_tot | total energy in eV deposited by all particles in the TPC |
| x\_tar, y\_tar, z\_tar, e\_tar | position (mm) and energy (eV) deposited step by step in the target |
| x\_win, y\_win, z\_win, e\_win | position (mm) and energy (eV) deposited step by step in the Mylar window |
| x\_ch, y\_ch, z\_ch, e\_ch | position (mm) and energy (eV) deposited step by step in the chamber |
| x\_InRoh, y\_InRoh, z\_InRoh, e\_InRoh | position (mm) and energy (eV) deposited step by step in the inner Rohacell |
| x\_Kap, y\_Kap, z\_Kap, e\_Kap | position (mm) and energy (eV) deposited step by step in the Kapton |
| x\_OutRoh, y\_OutRoh, z\_OutRoh, e\_OutRoh | position (mm) and energy (eV) deposited step by step in the outer Rohacell |
| x\_tpc, y\_tpc, z\_tpc, e\_tpc | position (mm) and energy (eV) deposited step by step in the TPC |
| MomentumXAfterTPC | momentum of particle in the X direction |
| MomentumYAfterTPC | momentum of particle in the Y direction |
| MomentumZAfterTPC | momentum of particle in the Z direction |

The drift library

libExN03DataoRings.so: this library is used for the second step of the simulations to achieve the drift of ionization electrons. It defines a concentric matrix of several rings divided in several pads. Three main aspects:

* Allows to read the output file of the simulation
* Allows to define parameters used for the drift of electrons by reading the ConfigFileRings.txt file
* Allows to define observables saved by the program of drift, especially number of electrons deposited on the different pads of the Micromegas as a function of the position of creation

ExN03DataoRings.cxx: source file defining the two following functions:

* The function ClearEvent() allows to clear all the observables at the beginning of a new event treated
* The function ReadConfigurationFile() allows to read the parameters given in ConfigFileRings.txt

ExN03DataoRings.hh: header file defining the observables/parameters/functions

|  |  |
| --- | --- |
| datai | class ExN03Datai |
| x\_pad | X coordinate of the pads (mm) |
| y\_pad | Y coordinate of the pads (mm) |
| t\_pad | time coordinate of the pads (ns) |
| q\_pad | charge deposited on the pads (number of electrons) |
| dt | difference between time fitted and real time of pads |
| Et\_pad | total charge deposited on the detection matrix |
| nb\_pads | number of pads triggered (number of electrons) |
| NRings | number of rings of the detection matrix |
| NSegments | number of pads per ring |
| sigL | longitudinal diffusion in the gas ( mm/√(cm) ) |
| sigT | transverse diffusion in the gas ( mm/√(cm) ) |
| driftV | drift speed of electrons in the gas (mm/ns) |
| Ionis | mean energy of creation of an electron-ion pair (eV) |
| TimeBinSize | sampling time of the electronics (ns) |
| ShapingTime | shaping time of the electronics (ns) |
| Threshold | detection threshold of the electrons |
| Gain | mean gain of the Micromegas in the Polya distribution |
| Theta | θ parameter in the Polya distribution |
| NoiseRMS | σrms of the noise (number of electrons) |
| FirstRing | distance from the center of the first ring |
| LastRing | distance from the center of the last ring |

**Description of the classes:**

The class ExN03DetectorConstruction

This class allows to define the experimental setup, more precisely by the following way.

The constructor allows to define:

* the dimensions of the materials by calling the class ExN03Setup
* several materials by the function DefineMaterials()
* the material constituting each element of the setup by the functions SetXMaterial(“Y”) where X will be filled by the material Y

The function Construct() allows to build the setup by first calling the ComputeMinosParameters() function defined in the corresponding header file, only used to define the size of the “world volume” from the size of the TPC. Take care of giving dimensions of the world big enough to contain all the elements of the detector. Elements defined from inner to outer are:

* the target (LH2)
* the Mylar window (entrance window, exit window and tube)
* the chamber (Inox)
* a first Rohacell, Kapton and a second Rohacell for the field cage
* the TPC (“mix” gas for a T2K mix)

Moreover, in the function Construct(), we define the region TPCLog used by ExN03PhysicsList and we define the magnetic field too.

The class ExN03PhysicsList

This class allows to define the standard electromagnetic physics list. In the constructor we define the default cut value for tracking. The function SetCuts() allows to assign this cut value for electrons, positrons and gammas and a minimum energy value for the creation of particles. Moreover, a more accurate value for tracking can be added for the region TPCLog corresponding to the TPC.

The same parameters can be defined and modified in the QGSP\_INCL\_ABLA physics list in the file /source/physics\_lists/lists/include/QGSP\_INCL\_ABLA.cc

The class ExN03EventAction

This class defines some functions at the beginning (BeginOfEventAction which calls ExN03ROOTuple::RecordBeginOfEvent) and at the end (EndOfEventAction which calls ExN03ROOTuple::RecordEndOfEvent) of each event generated.

The class ExN03RunAction

This class defines functions called at the beginning (BeginOfRunAction which calls ExN03ROOTuple::RecordBeginOfRun) and the end (EndOfRunAction which calls ExN03ROOTuple::RecordEndOfRun) of the simulation.

The class ExN03TrackingAction

This class defines functions called at the beginning (PreUserTrackingAction which calls ExN03ROOTuple::RecordBeginOfTrack) and the end (PostUserTrackingAction which calls ExN03ROOTuple::RecordEndOfTrack) of tracking a particle.

The class ExN03SteppingAction

This class defines the function UserSteppingAction which allows to call the functions ExN03ROOTuple::RecordStepDExxx at each step of the tracking of particles where xxx is the material encountered.

The class ExN03ROOTuple

This class defines user’s functions called by the previous classes. The main functions are as follows:

* RecordBeginOfRun: used to define the output file and the output tree containing the observables defined in libExN03Datai.so
* RecordEndOfRun: used to write tree and close output file
* RecordBeginOfEvent: clears the observables at each beginning of event
* RecordEndOfEvent: used to fill the output tree at the end of each event
* RecordBeginOfTrack: used to initialize some observables calculated for each tracked particle like the energies loss in materials, the detection in the TPC.
* RecordEndOfTrack: used to save the characteristics of tracked particles (initial conditions of emission, charge, mass, energy loss in materials, detection in the TPC)
* RecordStepDExxx: used to save at each step of tracking the energy loss and the position of particles. For some materials, only the total energy loss in materials is incremented.

The class ExN03PraimaryGeneratiorAction

* The function SetDefaultPrimaryParticle defines the default parameters of the particle which is to be sent in the setup in terms of position, momentum, charge, mass, energy. If an input file is needed, it is opened here.
* The function GeneratePrimaries generates randomly the particle(s) sent in the setup.

The class MyMagneticField

This class allows by the constructor to define the magnetic field.

**To run a simulation**

1. Change the paths inside the codes. Compile the libraries in ./lib using the corresponding GNUmakefileXXX

*cd MINOS\_simulation/lib; make –f GNUmakefileXXX*

Add MINOS\_simulation/lib to the LD\_LIBRARY\_PATH

1. Change the paths inside the codes. Generate the Makefile for the main program using the CMakeList.txt file. Simulation is tested with Geant4.10.00.

*cd MINOS\_simulation; cmake .*

1. Run the simulation; root output file is written in ./results

The simulations need the file vis.mac to define the visualization window (can be modified to visualize particles; see inside the file) or run\_nothing.mac if only root file is needed as output.

**Drift of ionization electrons**

The second step of the simulations needed to achieve the drift if electrons and obtain the signal on pads is situated in the folder Drift/. The source file is DriftRings.cxx compiled with GNUMakefileRings to generate the executable DriftRings.

This program simulates a concentric matrix of rings divided in several segments. The characteristics of the gas, the amplification stage, the AGET electronics and the noise are taken into account. The program uses the library libExN03DataoRings.so. Parameters for the drift simulation are defined in ConfigFileRings.txt. Results of drift are in the folder Drift/result/.

More explanations are described in the source file itself of the folder MINOS\_beam.

**Calculation of the resolution**

The program used to calculate the resolution is in Analysis/. Resolution is calculated along the beam axis. The histogram of the difference between reconstructed vertex and real vertex is in the folder result/ and can be fitted by a Gaussian or Lorentz function. Tracks of protons are fitted with parametric lines and with a helicoid (preliminary, without energy loss taken into account along the track) if a magnetic field is applied. Note that this simulation uses the library which defines the beam, so the beam used for the Geant4 simulation has to be considered.

The source file is tpcFit.cxx, compiled with GNUMakefileResol, which generates the executable tpcFit.

More explanations can be found in the source files of MINOS\_beam and MINOS\_Bfield.

A better algorithm based on Hought transorm was developed. This is in Analysis\_Clem/.

To run it:

1. convert file from Drift to format suitable for Hough transform analysis

.L convtree.C+; convtree(“inputfile”);

1. compile and run minos\_simu\_Newclusterfind.cpp
2. plotvertexHoughNew.C