

Studying Extensive Air Showers with **CONEX**

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CONEX is

- Easy to install
- Has just a few command line options
⇒ very straightforward to configure
- Very fast
- Produces small ROOT output files

Introduction

- CONEX is a small self-contained program with ROOT output. It can be obtained after request from `tanguy.pierog@kit.edu`
(version v2r2.3i is on the USB-sticks)
- The only required prerequisite is: `ROOT`
- Mix of fortran and C++

Installation

- Unpack: `tar xzvf conex2r2.3i.tgz`
- `cd conex2r2.3i`
- Compile: `make [opt]`

opt	Description
	QGSJet01, SIBYLL2.1
epos	QGSJet01, SIBYLL2.1, EPOS1.99
nexus	QGSJet01, SIBYLL2.1, EPOS1.99, QGSJetII.3, neXus3.97
qll	QGSJet01, SIBYLL2.1, EPOS1.99, QGSJetII.3
tutti	QGSJet01, SIBYLL2.1, EPOS1.99, QGSJetII.3, neXus3.97

Parameters and Options

- See README for all details
- **bin/conex2r -h**

```
conex2r -s [random seed] -S [autosave range] -a [alpha] -e [log10(emin/eV)] -E [log10(emax/eV)]  
-z [min zenith angle/deg] -Z [max zenith angle/deg] -i [azimuth angle/deg]  
-n [nShower] -p [0=gamma,100=p,5600=Fe,...] -x [prefix] -m [2=QGSJET, 5=Sibyll]  
-K [maxDetailLevel]
```

- The *.param files (e.g. conex_sibyll.param)

```
set xmaxp 2000 ! max slant depth  
set hacut1 1. ! cut for hadrons in GeV (not less than 1)  
set emcut1 0.001 ! cut for leptons in GeV (not less than 0.001)  
set hground 0. ! height of the observer in meter  
set altitude 0. ! height of the shower core (compared to hground) above obs. point  
set fehcut 0.05 ! relative threshold MC->CE for hadronic particle  
set feecut 0.005 ! relative threshold MC->CE for e/m particles  
set femcut 0.0005 ! relative threshold MC->CE for muons
```

CONEX searches for the param files at run-time at
\$CONEX_ROOT and \$PWD

Simulation Of Showers

- bin/conex2r -e 20 -E 20 -n 5 -p 100 -m 5
- ls -sh *.root

<2 min

85K conex_sibyll_179324753_100.root

- root conex_sibyll_179324753_100.root
- .ls

CONEX writes output files to \$ROOT_OUT or \$PWD

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CONEX writes output files to \$ROOT_OUT or \$PWD

You just simulated 5 air showers at 10^{20} eV in ~ 1 min.
File size: ~ 100 kBytes.

Description of Header

```
Seed1      : random seed1
Particle   : particle ID
Alpha       : spectral index
lgEmin     : log10 of the minimum energy in eV
lgEmax     : log10 of the maximum energy in eV
zMin       : minimum zenith angle in degree
zMax       : maximum zenith angle in degree
Version    : conex version
HEModel    : High Energy interaction model flag
              (1=NeXuS, 2= QGSJet, 4=Sibyll)
LEModel    : Low Energy (E < 80 GeV) model flag (only 3=Gheisha)
hadCut     : hadron and muon cutoff (minimum energy)
emCut      : e/m particles cutoff
              (minimum energy for electrons, positrons and gammas)
hadThr     : Emax(CE)/Emax(MC) for hadrons (threshold)
muThr     : Emax(CE)/Emax(MC) for muon (threshold)
emThr      : Emax(CE)/Emax(MC) for e/m particles (threshold)
SVNRevision : svn revision
OutputVersion : cxroot output version
```

Description of Shower

lgE : log10 of the primary energy in eV
zenith : zenith angle in degree
azimuth : azimuth angle in degree (0 = East)
Seed2 : random seed2 (number of random number generator calls below 1 billion)
Seed3 : random seed3 (number of billions of random number generator calls)
XfirstIn : inelasticity of first interaction ([0,1])
Xfirst : real first interaction point in slant depth (g/cm^2)
Hfirst : altitude of real first interaction point (m)
X0 : pseudo first interaction point for GH fit
Xmax : GH fit result for slant depth of the shower maximum (g/cm^2)
Nmax : Number of charged particles above cut-off at the shower maximum
p1 : first parameter for the polynomial function of the GH fit
p2 : second parameter for the polynomial function of the GH fit
p3 : third parameter for the polynomial function of the GH fit
chi2 : Chi squared / number of degree of freedom / sqrt (Nmax)
for the fit
 (small number not realistic because it's divided by sqrt (Nmax))
Xmx : Quadratic fit on real profile for slant depth of the shower maximum (g/cm^2)
Nmrx : Quadratic fit on real profile of the number of charged particles above cut-off at the shower maximum

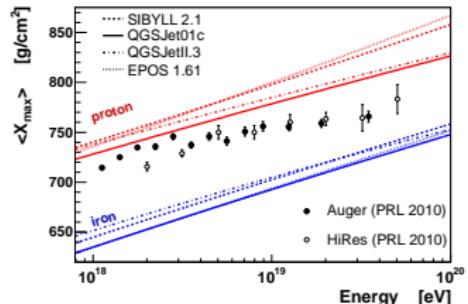
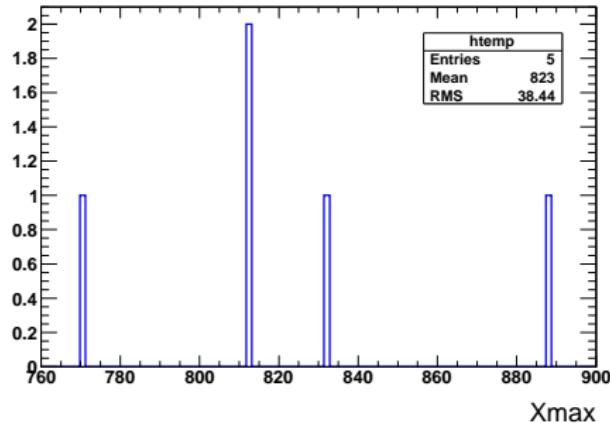
Description of Shower, Longitudinal Profiles

nX : number of points of X, N, H, D, dEdX, Mu, dMu,
Gamma, and Electrons array
X : slant depth array (g/cm²)
H : height array (m)
D : distance array (m)
N : array of number of charged particles above cut-off
crossing each X plane
dEdX : array of energy deposit (GeV/(g/cm²)) in the MIDDLE of
each X bin
Mu : array of number of muons above cut-off crossing each X plane
dMu : array of number of muons produce above cut-off in each bin
Electrons: array of number of e⁺ + e⁻ above cut-off crossing each X plane
Gamma : array of number of gammas above cut-off crossing each X plane
EGround : Energy of particles at maximum X (EGround[0]=e+gamma;
EGround[1]=hadrons; EGround[2]=muons)

Analysing the Data

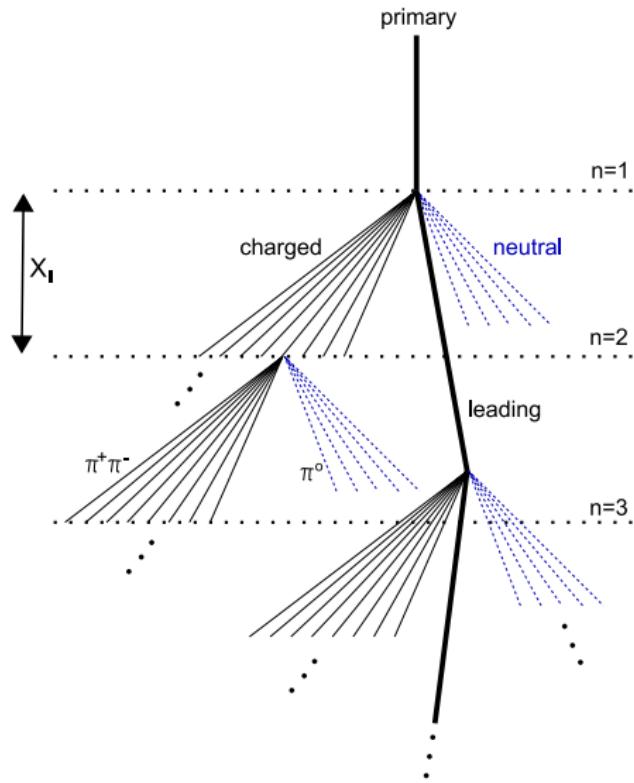
⇒ Depth of the Shower Maximum

Shower->Draw("Xmax")



Save Leading Interactions

- bin/conex2r -e 20 -E 20 -n 5 -p 100 -m 5 **-K 2**



Description of FirstInteraction Tree

nInt : Number of saved Interactions (size of array kinel, pId, pEnergy, mult, matg, depth) per shower

kinel : Inelasticity for each saved interactions

pId : Id of parent particle (CONEX Id)
(CORSIKA Id if LEADING_INTERACTIONS_CORSIKA is defined
in conexConfig.h)

pEnergy : Energy of parent particle for each saved interactions

mult : Multiplicity for each saved interactions

matg : Mass of target nucleus for each saved interactions

depth : Slant depth (g/cm²) of each saved interactions

height : altitude above see level (m) of each saved interactions

nPart : total number of recorded secondary particles per shower
(nPart=sum[nInt](mult))
(size of array Energy, px, py, pz, Type, and idInt)

idInt : Interaction number

Type : Id of secondary particles (CONEX Id)
(CORSIKA Id if -DLEADING_INTERACTIONS_CORSIKA is selected
in Makefile)

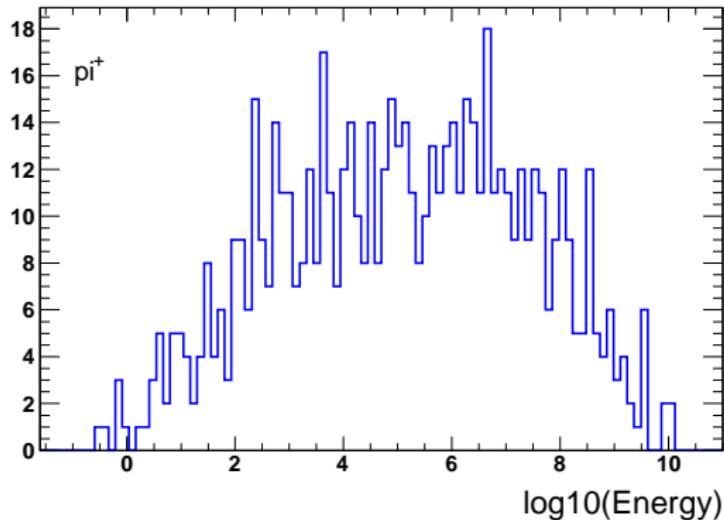
Energy : total energy (GeV) of secondary particles

px,py,pz : momentum (GeV/c) of secondary particles
(if -DLEADING_INTERACTIONS_TREE_EXT is used in Makefile)

Working with the FirstInteraction Data

⇒ Energy Distribution of π^+ in first interaction:

```
FirstInteraction->Draw("log10(Energy)", "Type==8&&idInt==0")
```



The End