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# Highlight on Advanced Topics

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FLUKA Beginner's Course

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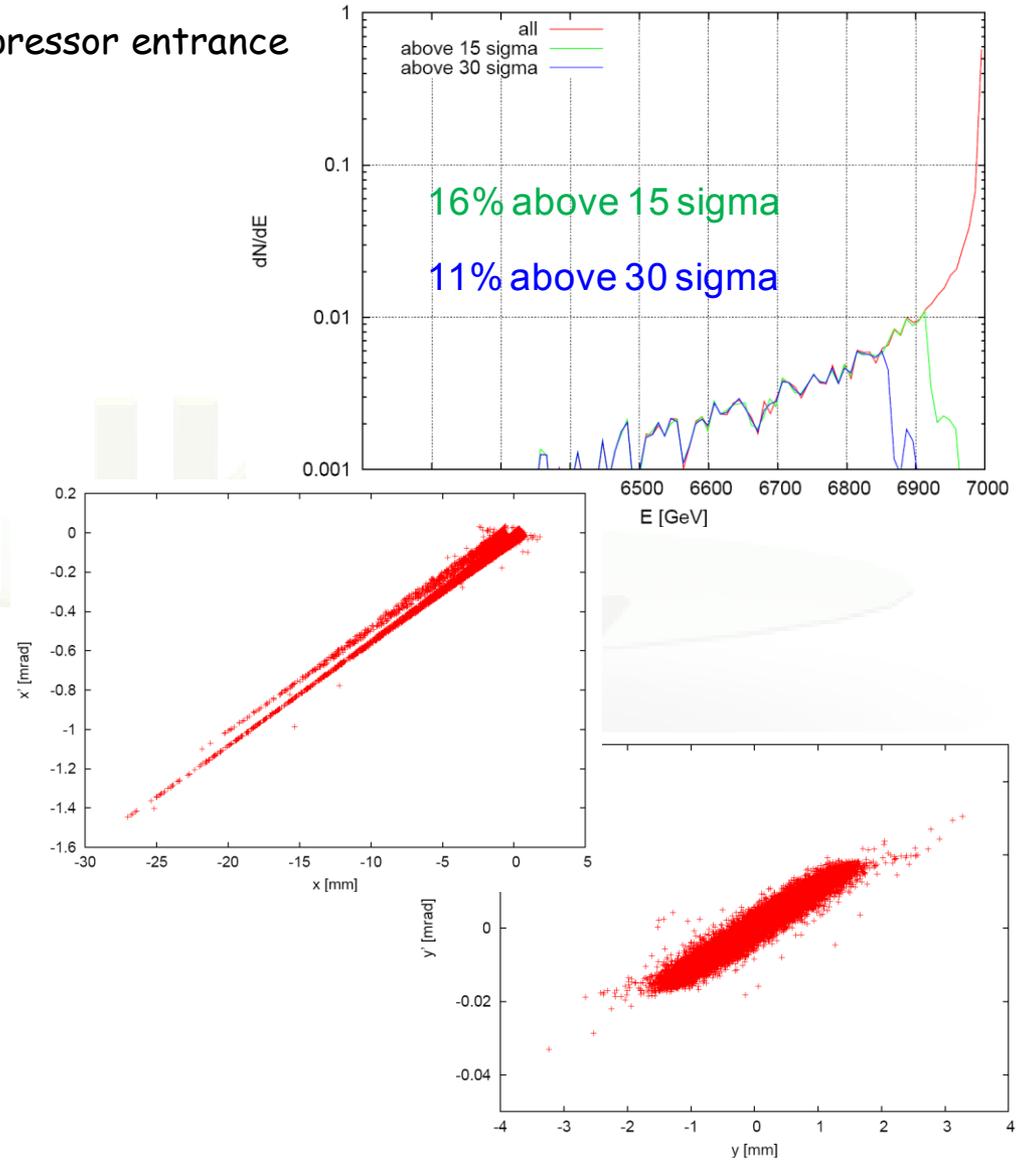
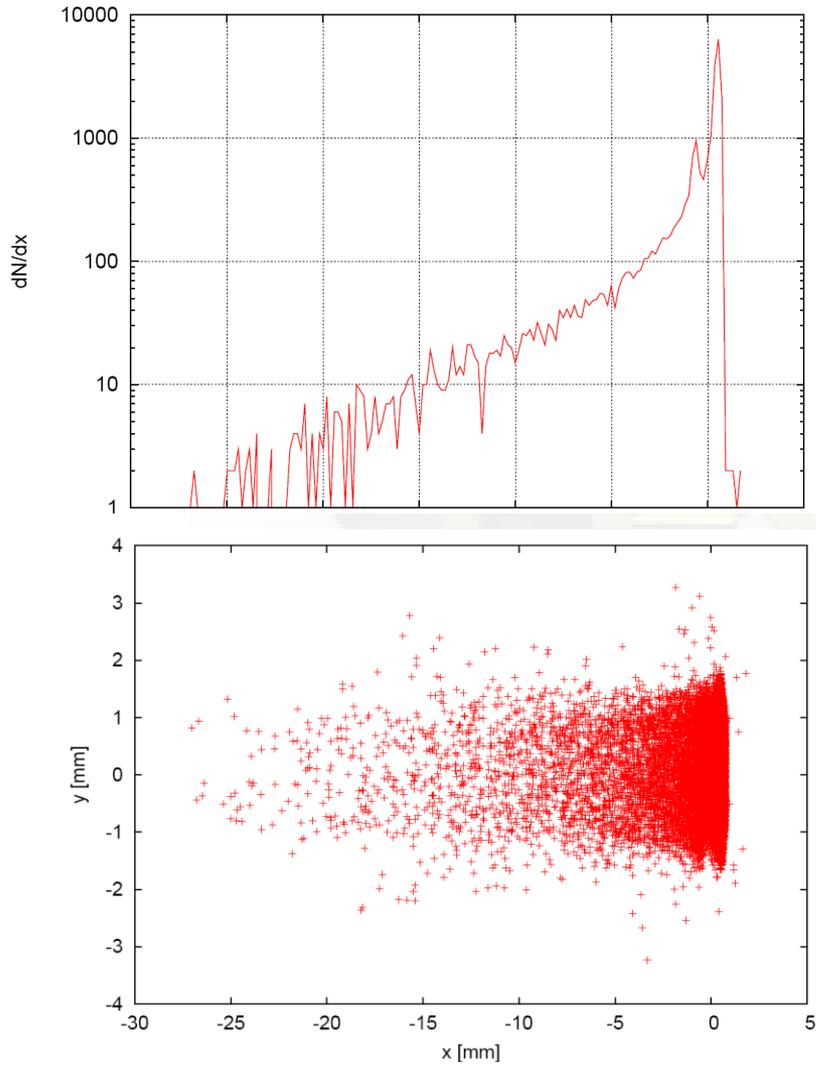
## This is not the end ...

- As said in the first day, most applications can be run through data-cards only, exploiting the FLUKA built-in capacities
- Sometimes, more is needed...
- A set of templates User Routines is provided in the `$FLUPRO/usermvax` directory
- All common and default parameters are located in `$FLUPRO/flukapro`
- These can be modified by the user to fit his input/output needs
- We'll give here some hints of what can be done, more can be found in the manual, in the fluka-discuss archive, or

at the second FLUKA advanced course and workshop  
to be held in Vancouver, Canada, on Sep 15-20 2012

# Implementing customized beam distributions - 1

LHC proton beam at the Dispersion Suppressor entrance



# Implementing customized beam distributions - 2

Input card: **SOURCE**

User routine: [\\$FLUPRO/usermvax/source.f](#)  
to be linked in one's own executable !

In source.f

you can sample beam particle position, direction, and energy  
from an external file or any (analytic or numerical) distribution

you can assign different weights to primary particles

you can load reaction products in the same primary history

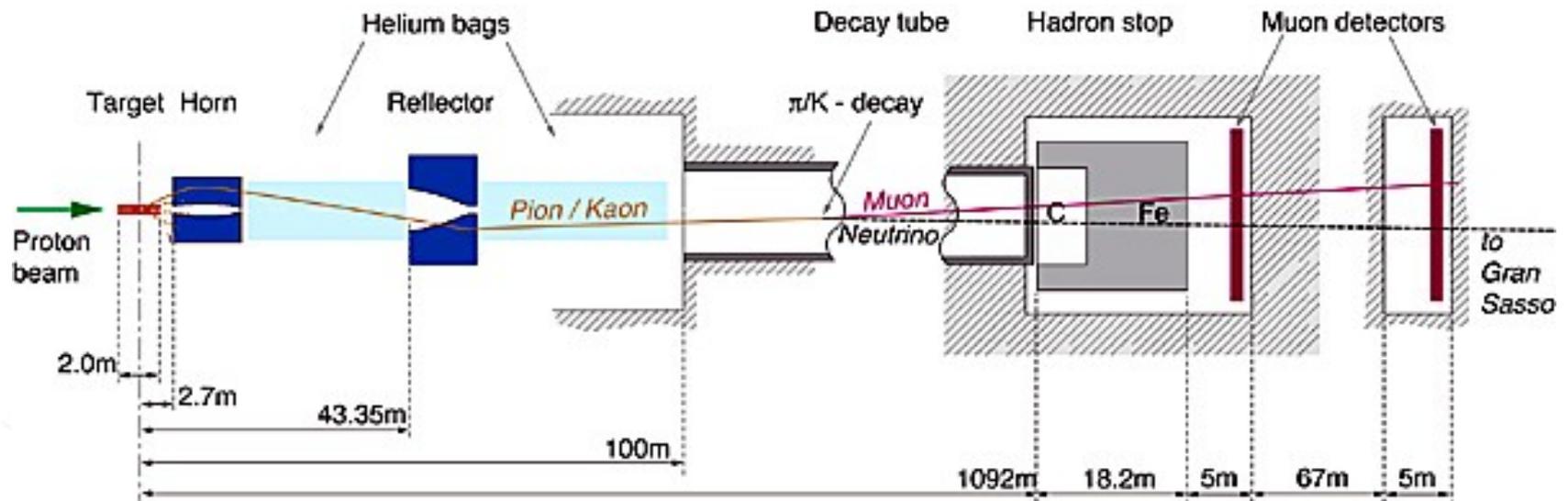
several sampling routines already exist in the FLUKA code

parameters input in the **SOURCE** card are available

a **BEAM** card with a momentum/energy higher than the maximum one is still needed  
for initialization purposes (to define the tabulation limit)

# Implementing magnetic field - 1

CERN Neutrino to Gran Sasso



The two magnetic lenses (blue in the sketch) align positive mesons towards the Decay tunnel, so that neutrinos from the decay are directed to GranSasso, 730~km away  
Negative mesons are deflected away  
The lenses have a finite energy/angle acceptance

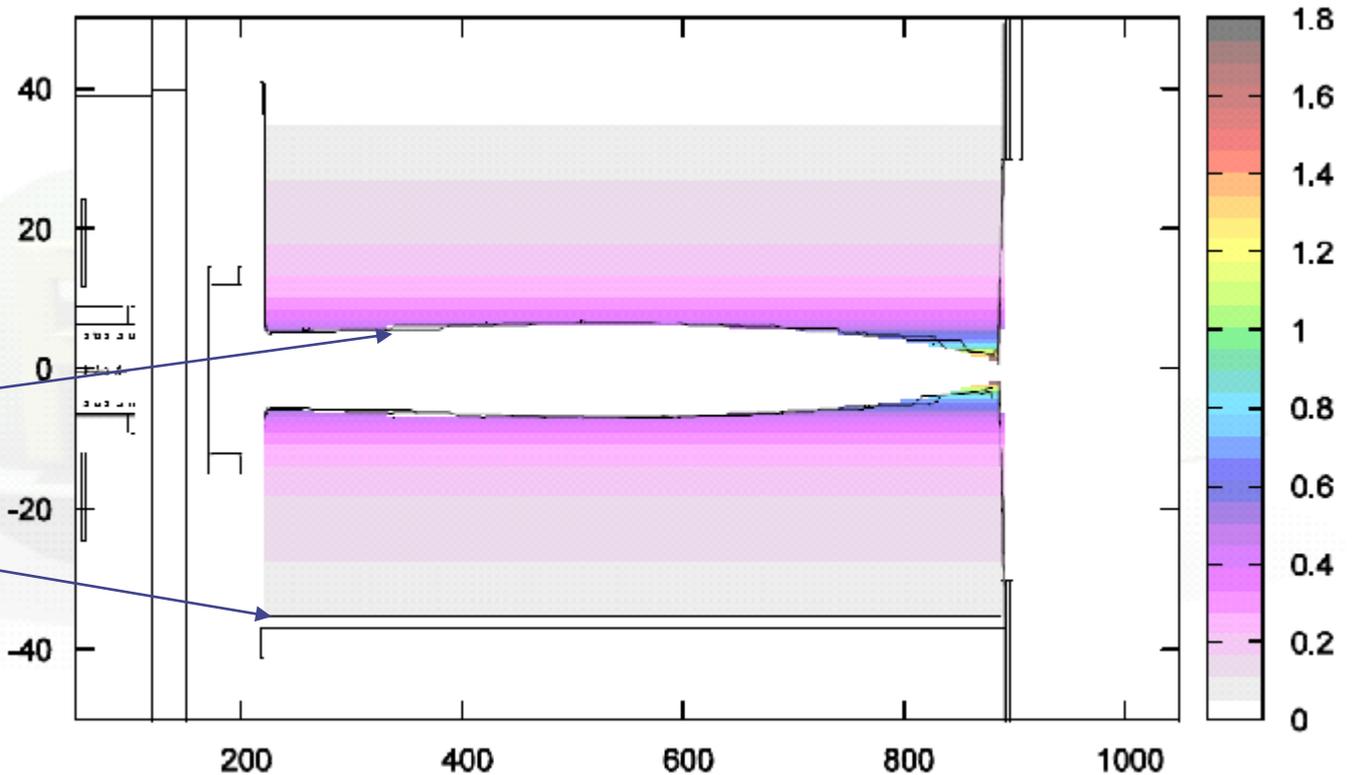
# Implementing magnetic field - 2

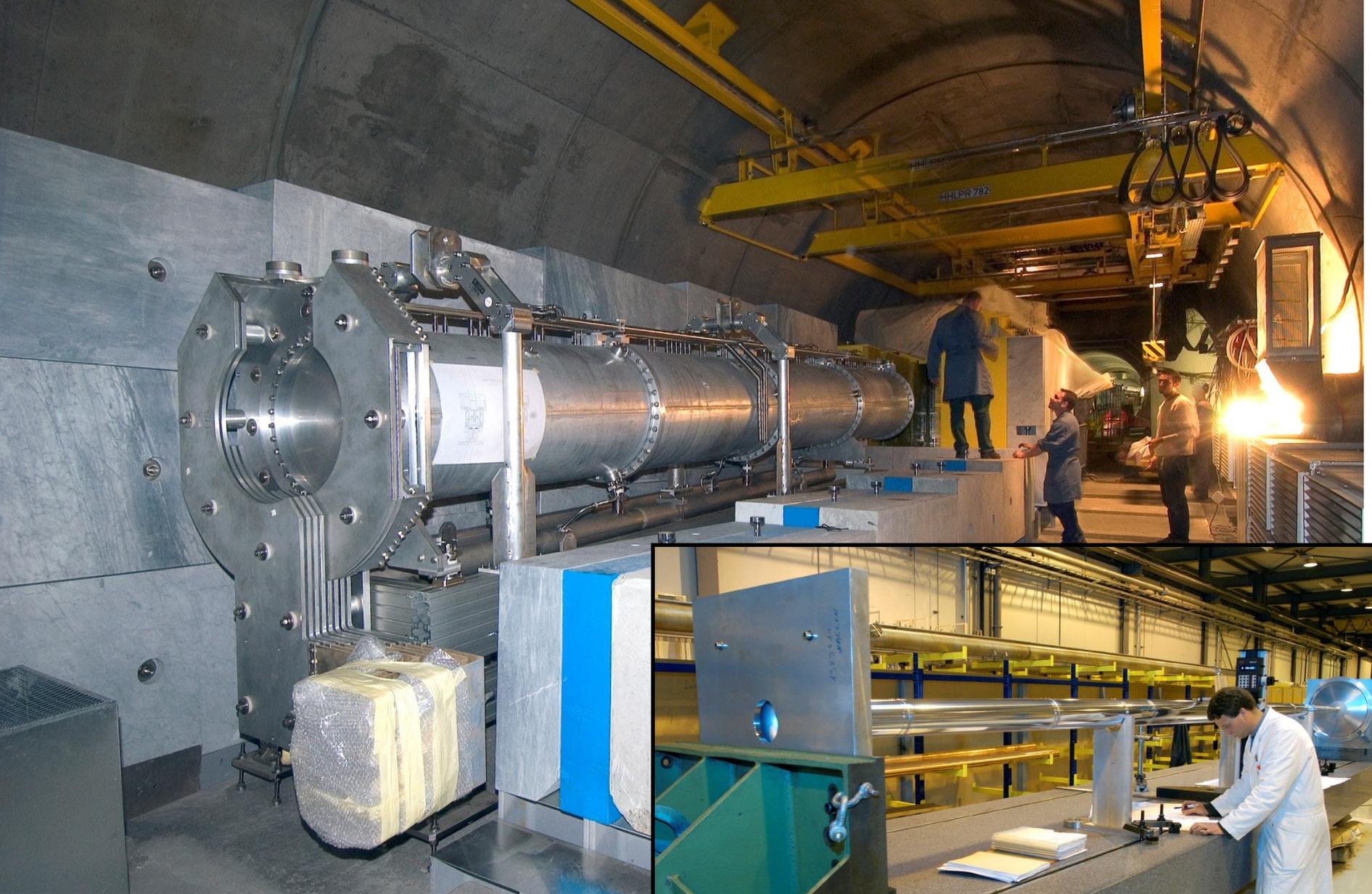
Input card: **ASSIGNMA**

User routine: **`$FLUPRO/usermvax/magfld.f`**  
to be linked in one's own executable !

Magnetic field intensity in the CNGS horn

A current  $\approx 150\text{kA}$ ,  
pulsed, circulates through the  
Inner  
and  
Outer  
conductors  
The field is toroidal,  
 $B \propto 1/R$





# Implementing magnetic field - 4

```
SUBROUTINE MAGFLD ( X, Y, Z, BTX, BTY, BTZ, B, NREG, IDISC )
```

```
IF ( NREG .EQ. NRHORN ) THEN
```

```
  RRR = SQRT ( X**2 + Y**2 )
```

```
  BTX = -Y / RRR
```

```
  BTY = X / RRR
```

```
  BTZ = ZERZER
```

```
  B = 2.D-07 * CURHOR / 1.D-02 / RRR
```

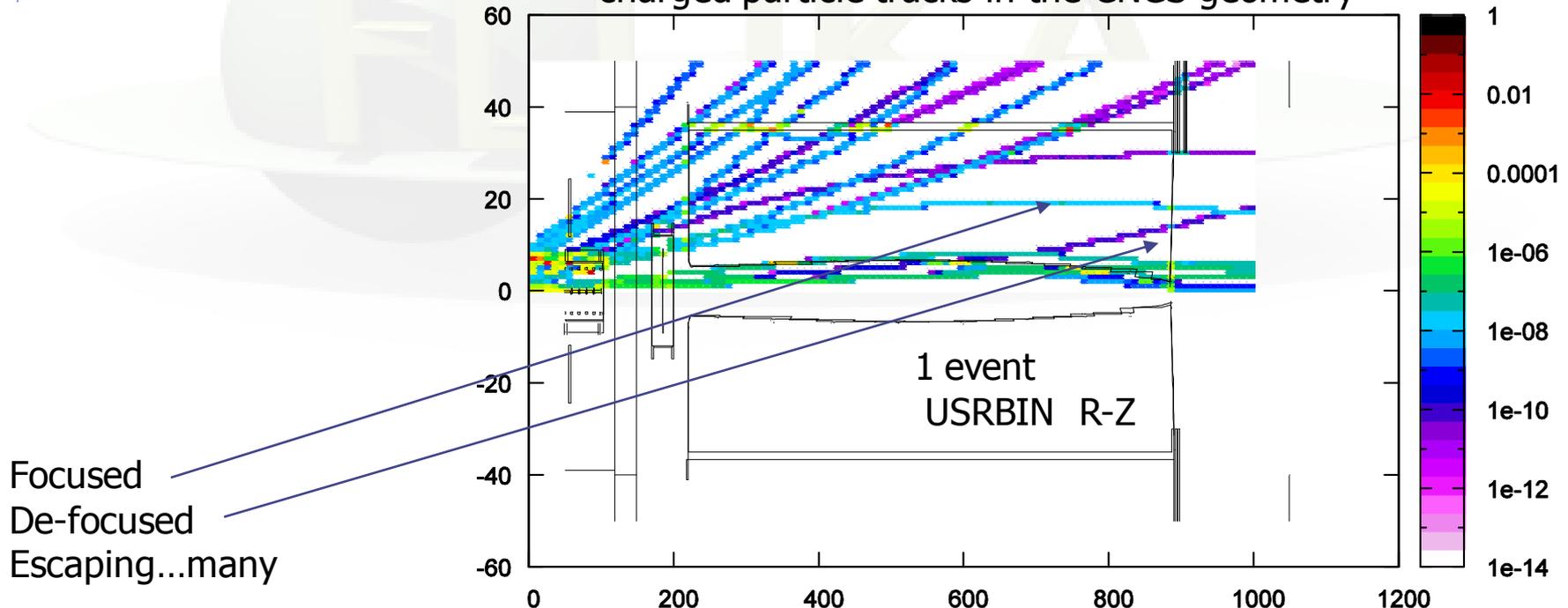
```
END IF
```

In magfld.f

you can define the magnetic field  
analytically or  
by interpolating an external map

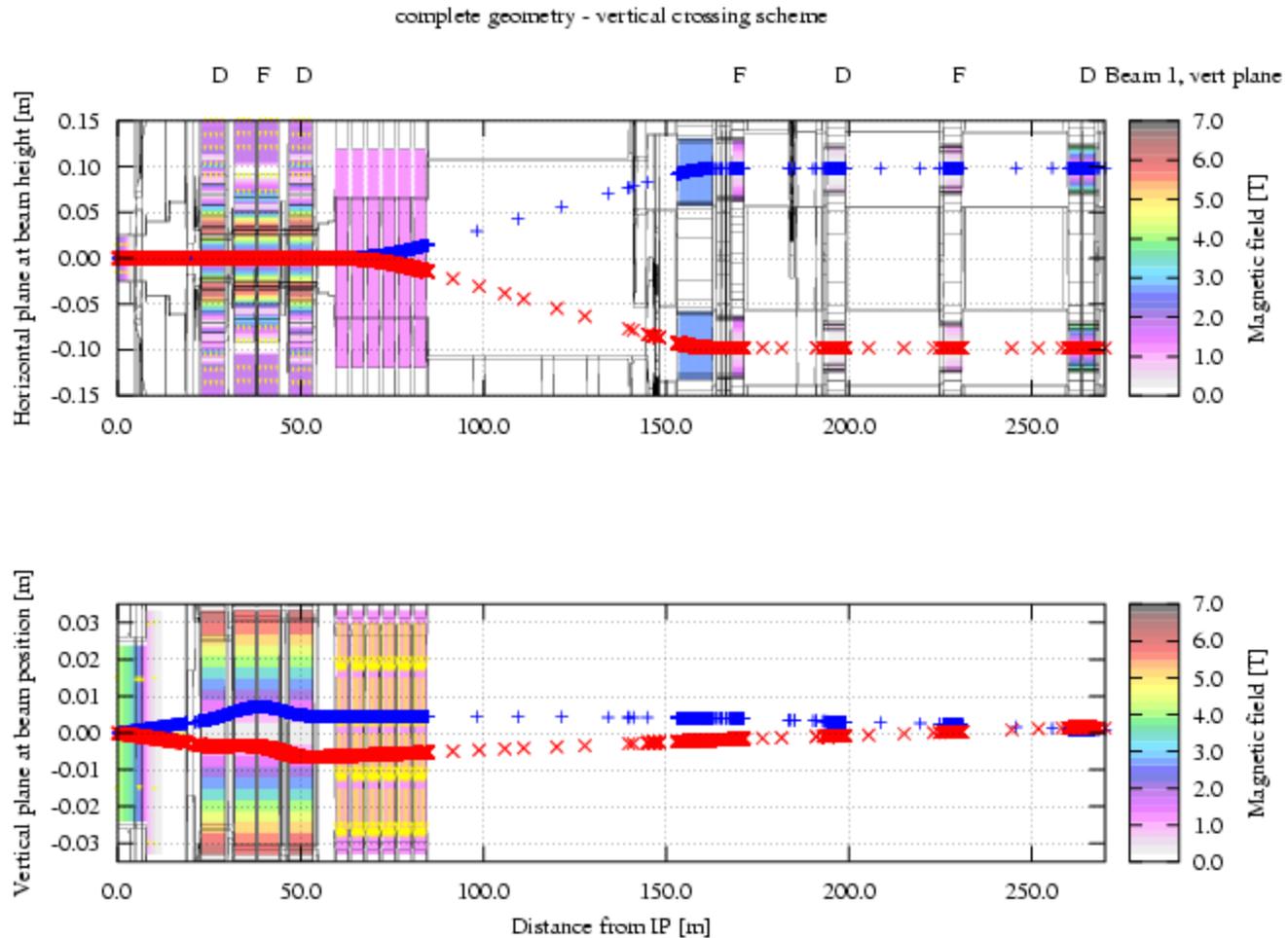


charged particle tracks in the CNGS geometry



# Implementing customized scoring - 1

beam crossing in ATLAS



## Implementing customized scoring - 2

Input card: **USERDUMP**

User routine: **[\\$FLUPRO/usermvax/mgdraw.f](#)**  
to be linked in one's own executable !

In mgdraw.f

you can get particle trajectories and (continuous and local) energy losses

*you can go over reaction products*

you can access information at each boundary crossing, particle step,  
energy deposition event, *interaction*

In mdstck.f (for which no activating card is needed) as well,  
with the additional – dangerous – possibility of influencing  
the subsequent transport of secondaries

# Implementing customized scoring-3

Input card: **USERWEIG**

User routines: [\\$FLUPRO/usermvax/comscw.f](#)  
[\\$FLUPRO/usermvax/fluscw.f](#)

to be linked in one's own executable !

In comscw.f

you can apply a user defined weight (even discard) on deposited energy, stars or residual nuclei

you can extract information (and dump it on a file) about the involved particles

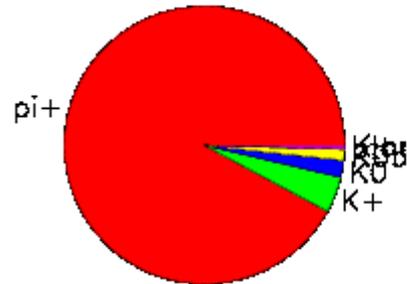
In fluscw.f

same as comscw.f for fluence scoring

# Implementing customized flags

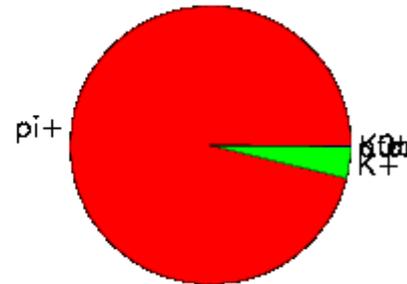
identification of ancestors of muon-neutrinos reaching Gran Sasso

out of target



ancestor as exiting the target

decay parent



parent decaying in  $\nu_\mu$

No input card

User routines: [\\$FLUPRO/usermvax/stuprf.f](#)  
[\\$FLUPRO/usermvax/stupre.f](#)

to be linked in one's own executable !

In `stuprf.f`

you can keep track of the particle origin

`stupre.f` applies in case of electrons, photons and positrons

# Implementing region independent importance biasing

Input card: **BIASING**

User routines: **\$FLUPRO/usermvax/usimbs.f**

to be linked in one's own executable !

usimbs.f

is called for every particle step in user selected regions in order to return the change in the importance from the beginning to the end of the step, in a region-independent way!

It doesn't produce as accurate results as the manual region importance biasing, but it is a great time saver as segmenting a complex geometry can be a cumbersome task