



# Estimators and Scorings

21<sup>st</sup> FLUKA Beginner's Course  
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# FLUKA Scoring & Results - Estimators

- It is often said that Monte Carlo (MC) is a “mathematical experiment”  
The MC equivalent of the result of a real experiment (*i.e.*, of a measurement) is called an estimator.
- Just as a real measurement, an estimator is obtained by sampling from a statistical distribution and has a statistical error (and in general also a systematic one).
- There are often several different techniques to measure the same physical quantity: in the same way the same quantity can be calculated using different kinds of estimators.
- FLUKA offers numerous different estimators, *i.e.*, directly from the input file the users can request scoring the respective quantities they are interested in.
- As the latter is implemented in a very complete way, users are strongly encouraged to preferably use the built-in estimators with respect to user-defined scoring
- For additional requirements FLUKA user routines are provided

# Built-In and User Scoring

- Several **pre-defined estimators** can be activated in FLUKA.
- One usually refers to these estimators as "**scoring**" capabilities
- Users have also the possibility to build their own scoring through **user routines**, HOWEVER:
  - **Built-in scoring** covers most of the **common needs**
  - **Built-in scoring** has been **extensively tested**
  - **Built-in scoring** takes BIASING **weights automatically into account**
  - **Built-in scoring** has **refined algorithms** for track subdivision
  - **Built-in scoring** comes with **utility programs** that allow to evaluate statistical errors
- Standard scoring can be weighted by means of **simple user routines** (fluscw, comscw)

# FLUKA Scoring & Results

**What?** Energy deposition, and derivatives (dose), fluence or current versus energy/ angle/ other kinematical variables, time, DPA.....

**Where?** In regions, on boundaries, on region-independent grids

**When?**

At the end of each cycle

At each event

**Output?** Files `in $nnn$ _fort.##` where ## is logical unit number chosen by user

**Results?**

Postprocessing utilities (in `$FLUPRO/flutil`) merge cycles, calculate average and rms, provide data files for plotting

User code needed

Results normalized  
**per primary**

Data merging and plotting available in **FLAIR**

# The FLUKA Output Files

The FLUKA output consists of:

- A **main (standard) output**, written on logical output unit **LUNOUT** (predefined as 11 by default) [**\*.out**]
- A file with the last random number seeds, unit **LUNRAN** (2 by default) [**ran\***]
- A file of error messages, unit **LUNERR** (15 by default) [**.err**]
- Any number (including zero) of **estimator output files**. Their logical unit number is defined by the user [**\*\_fort.xx**]
- The available range of logical output numbers is: 21-99
- Generally, the user can choose between **formatted (ASCII) and unformatted (binary)** scoring (negative or positive sign in the logical unit number). Unformatted scoring is mandatory for the use of provided post-processing utilities.
- Several estimators can be output on the same file (same logical unit) **provided they are of the same type**
- Possible **additional output generated by the user** in any user routine

# Before giving details

- A couple of slides on important definitions / quantities:
- Reaction rate and cross section
- Fluence and the difference wrt current
- Fluence as the correct quantity to evaluate the potential effect of a particle field on a “detector”
- Different methods to evaluate fluence

# Definitions

- $N$ : number of identical particles
- $N_0$ : number of atoms per unit volume
- $\lambda[cm]$ : mean free path, i.e. average distance travelled by a particle in a material before an interaction. It depends on the material, particle type and energy.
- $l$ : total distance travelled
- $v$ : average particle velocity

# Cross Section

- $\Sigma[cm^{-1}] = 1/\lambda[cm]$  : **macroscopic cross section**, i.e. probability of interaction per unit distance. It depends on the material, particle type and energy.
- $\sigma = \frac{\Sigma}{N_0} = \text{atom effective area}$  , [barn =  $10^{-24}cm^2$ ] : **microscopic cross section**, i.e.
  - i) the **area of an atom weighted with the probability of interaction** (hence the name “cross section”);
  - ii) or the **probability of interaction per unit length**, with the length measured in **atoms/cm<sup>2</sup>** (the number of atoms contained in a cylinder with a 1 cm<sup>2</sup> base).

Microscopic and macroscopic cross sections have a similar physical meaning of “probability of interaction per unit length”, with length measured in different units. Thus, the number of interaction can be obtained by multiplying both by the corresponding particle track-length.

# Reaction Rate and Fluence

- $R = \Sigma\Phi V$  : number of reactions in a given time interval inside the volume  $V$  (where  $\Phi$  is the fluence and the product  $\Sigma\Phi$  is integrated over energy or velocity)
- $\dot{R} = N \frac{dl}{dt} \Sigma = Nv\Sigma$  : **reaction rate**
- $\frac{d\dot{R}}{dV} = \frac{dN}{dV} v \Sigma = n(\mathbf{r}, v)v\Sigma$  : reaction rate inside the volume element  $dV$
- $\Phi(\mathbf{r}, v) = n(\mathbf{r}, v)dl$  ,  $[cm^{-2}]$  : **fluence**, i.e. time integral of the flux density
- $\dot{\Phi}(\mathbf{r}, v) = n(\mathbf{r}, v)v$  ,  $[cm^{-3}cms^{-1}] = [cm^{-2}s^{-1}]$  : **fluence rate** or **flux density**.
- Fluence is measured in **particles per  $cm^2$**  but in reality it describes the **density of particle tracks**

# Fluence vs Current (1/2)

## Surface crossing estimation

- Imagine a surface having an infinitesimal thickness  $dt$ .

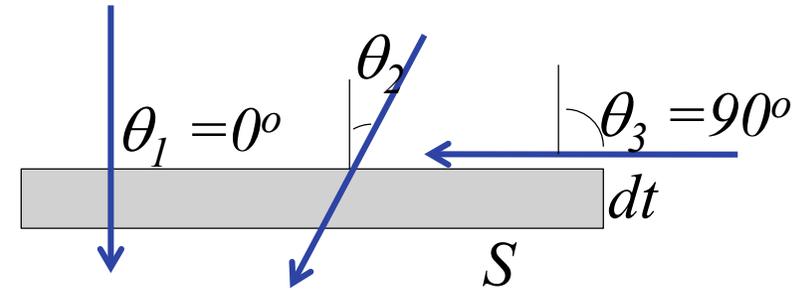
A particle incident with an angle  $\theta$  with respect to the normal of the surface  $S$  will travel a segment  $dt/\cos\theta$ .

- Therefore, we can calculate an **average surface fluence** by adding  $dt/\cos\theta$  for each particle crossing the surface, and dividing by the volume  $S dt$ :

$$\Phi = \lim_{dt \rightarrow 0} \frac{\sum_i \frac{dt}{\cos \theta_i}}{S dt}$$

- While the **current**  $J$  counts the number of particles crossing the surface divided by the surface:

$$J = dN/dS$$



# Fluence vs Current (2/2)

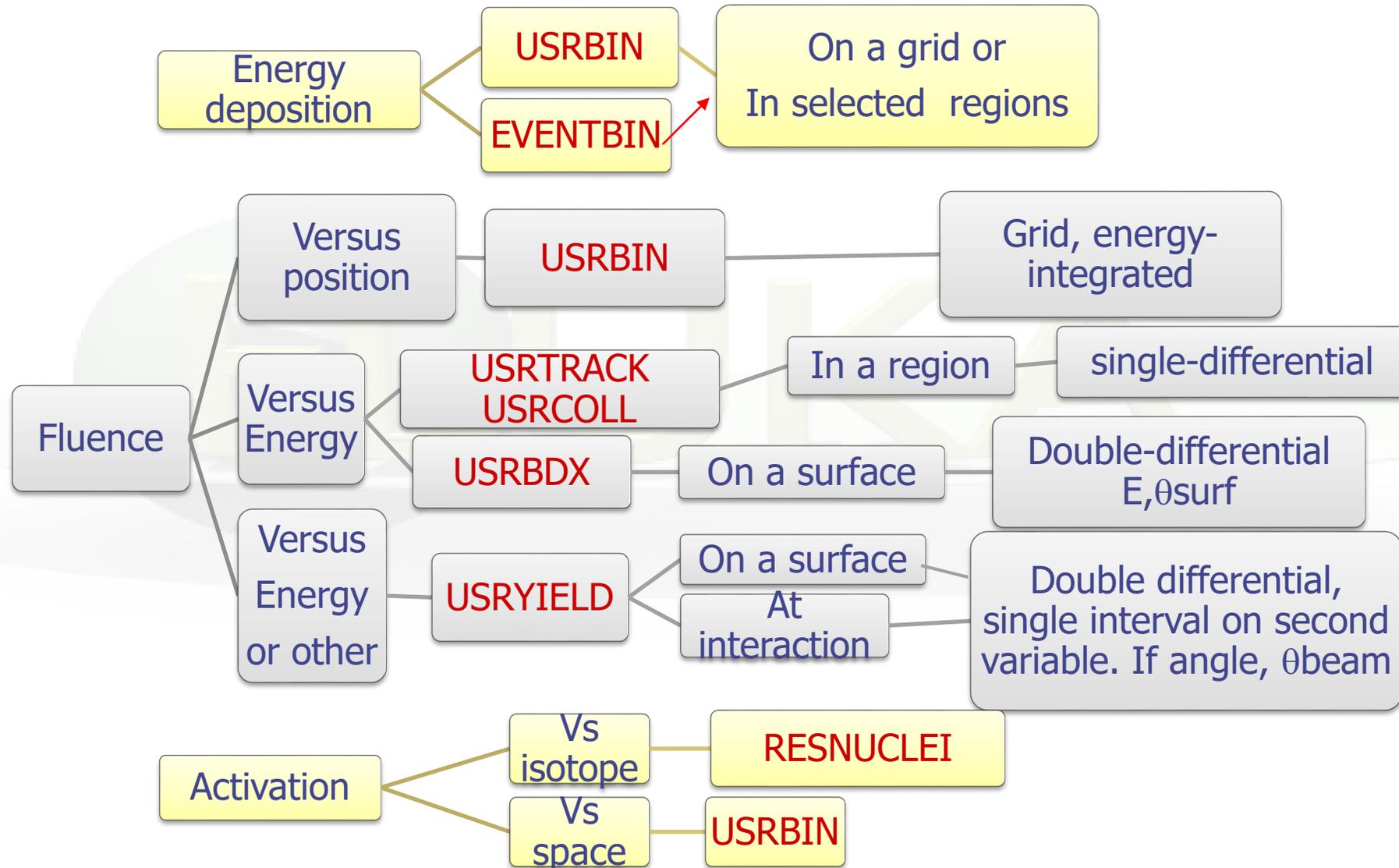
The **fluence** is independent from the orientation of **surface**  $S$ ,  
while the **current** is **NOT!**

In an isotropic field can be easily seen that on a flat surface  $J = \Phi/2$

Current is *meaningful* if particles are only counted (e.g. signal trigger).

But if someone estimates **dose, activation, radiation damage, instrument response** the quantity to be used is **fluence**.

# FLUKA estimators (1/2)



# FLUKA estimators (2/2)

- **USRTRACK**, **USRCOLL** score average  $d\Phi/dE$  (differential fluence) of a given type or family of particles in a given region (see next slide);
- **USRBDX** scores average  $d^2\Phi/dEd\Omega$  (double-differential fluence or current) of a given type or family of particles on a given surface;
- **USRBIN** scores the spatial distribution of energy deposited, or total fluence (or star density, or momentum transfer) in a regular mesh (cylindrical, Cartesian or per region) described by the user;
- **USRYIELD** scores a double differential yield of particles escaping from a surface. The distribution can be with respect to energy and angle, but also other more “exotic” quantities;
- **SCORE** scores energy deposited (or star density) in all regions;
- The output of SCORE will be printed in the main (standard) output, written on logical output unit **LUNOUT** (pre-defined as **11** by default)
- All other detectors write their results into logical output units assigned by the user (the unit numbers must be **>20**)

# Fluence estimation

- Track length estimation:

USRTRACK

$$\dot{\Phi}(v) dt = n(v) v dt = \frac{dN(v)}{dV} \frac{dl(v)}{dt} dt = \lim_{\Delta V \rightarrow 0} \frac{\sum_i l_i(v)}{\Delta V}$$

- Collision density estimation:

USRCOLL

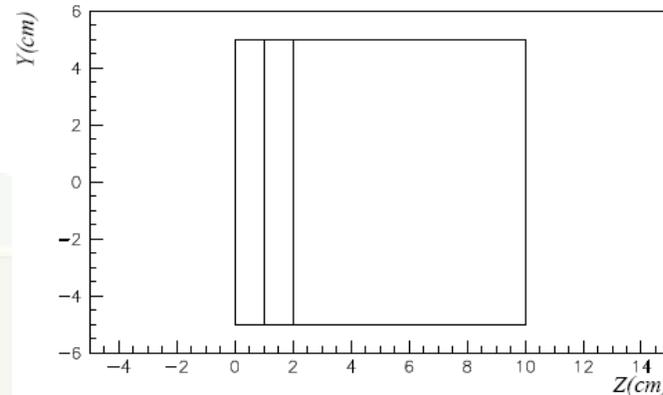
$$\dot{\Phi}(v) = \frac{\dot{R}(v)}{\sigma(v) N_o \Delta V} = \frac{\dot{R}(v)}{\Sigma(v) \Delta V} = \frac{\dot{R}(v) \lambda(v)}{\Delta V}$$

# More “Special” Scoring

- **RESNUCLEi** scores residual nuclei in a given region (more details are given in the respective lecture on activation);
- **DETECT** scores energy deposition in coincidence or anti-coincidence with a trigger, separately for each “event” (primary history);
- **EVENTBIN** is like **USRBIN**, but prints the binning output after each event instead of an average over histories;
- **ROTPRIBIN** sets the storage precision (single or double) and assigns rotations/translations for a given user-defined binning (**USRBIN** or **EVENTBIN**). Useful in case of LATTICES;
- **TCQUENCH** sets scoring time cut-offs and/or Birks quenching parameters for binnings (**USRBIN** or **EVENTBIN**) indicated by the user;
- **USERDUMP** defines the events to be written onto a “collision tape” file;
- **AUXSCORE** defines filters and conversion coefficients.
- **DCYSCORE** assigns cooling times (see lecture on activation)

# Extending the example with Scoring

- Cylinder along Z, filled by water-aluminum-lead and surrounded by Air



- The **USRBIN** command allows to define a **3-D grid**, either cartesian, cylindrical (R-Z- $\Phi$ ) or by region.
- On this grid, one can score energy deposition, particle fluence (total or by particle type), as well as the density of interactions;
- An equivalent **EVENTBIN** command outputs the same quantities event-by-event;
- Using **USERWEIG** the results can be weighted by the **comscw.f** or **fluscw.f** external routines (Advanced).

# Example: USRBIN

- Start from the solution of ex\_Geometry1 ( Either copy your .inp and .flair files and rename them to example\_score, or download from the site & rename):

```
mkdir example_score ;  
cp ex_Geometry1/ex_Geometry1_final.inp example_score/example_score.inp ;  
cd example_score
```

If you had some problem or you could not complete the ex\_Geometry1, copy the solution ex\_Geometry1\_final.inp

Check that there are **no spurious STOP** in the input file (inserted for the debugging test during the exercise Geometry1).

# Example: USRBIN

- Start from the solution of ex\_Geometry1 ( Either copy your .inp and .flair files and rename them to example\_score, or download from the site & rename):

```
mkdir example_score ;  
cp ex_Geometry1/ex_Geometry1_final.inp example_score/example_score.inp ;  
cd example_score
```

- Open in flair or with your preferred editor
- Add USRBIN to:
  - 1) score **ENERGY** on a **CYLINDRICAL GRID** ( $r, z, \phi$ ) covering the target and surroundings:  $0 < r < 10$  cm,  $-5$  cm  $< z < 15$  cm, with cells having  $\Delta r = \Delta z = 1$  mm,  $\Delta \phi = 2\pi$ . Output unit = 40 BIN
  - 2) score **Neutron Fluence** on the same grid as above. Output unit = 41 BIN
  - 3) score **Charged Hadron Fluence** on the same grid. Output unit = 42 BIN
- Run 5 cycles, 1000 primaries each

# USRBIN - WHATs

**WHAT(1)** : code indicating the type of binning selected.

Example : Usrbin WHAT(1) = 11.0 → Mesh: R-Z or R-Phi-Z, no symmetry

**WHAT(2)** : particle (or particle family) type to be scored

**WHAT(3)** = logical output unit

The first three fields have similar meanings for all estimators.

The other WHATs will contain the limits or regions or boundaries, and the number of intervals for cells/histograms

**SDUM** (optional) : name of the estimator

## **For Ursbin, r-phi-z:**

**WHAT(4)**= For R-Z and R-Phi-Z binning: Rmax

**WHAT(5)**= For Cartesian binning: Ymax For R-Z and R-Phi-Z binning: Y coordinate of the binning axis

..... And so on

# USRBIN (1)

1) Add USRBIN to score **ENERGY** on a **CYLINDRICAL GRID** ( $r, z, \phi$ ) covering the target and surroundings:  $0 < r < 10$  cm,  $-5$  cm  $< z < 15$  cm, with cells having  $\Delta r = \Delta z = 1$  mm,  $\Delta \phi = 2\pi$

\* Energy deposition [GeV/cm<sup>3</sup>]

*	+	1	+	2	+	3	+	4	+	5	+	6	+	7	+
<b>USRBIN</b>				11.0	<b>ENERGY</b>			-40.0		10.0				15.0	<b>TargEne</b>
<b>USRBIN</b>				0.0				-5.0		100.0				200.0	&

 <b>USRBIN</b>	Unit: 40 BIN ▼	Name: TargEne
Type: R-Φ-Z ▼	Rmin: 0	NR: 100
Part: ENERGY ▼	X:	NΦ:
	Zmin: -5.0	NZ: 200
	Y:	
	Zmax: 15.0	

- This is an R-Z-Φ binning (what(1)=11), scoring energy density (generalized particle ENERGY, or 208), writing the unformatted output on unit 40, spanning  $0 < R < 10$  cm in 100 bins,  $0 < \Phi < 2\pi$  in 1 bin (default),  $-5$  cm  $< z < 15$  cm in 200 bins.

# USRBIN (2-3)

2) Add USRBIN to score **Neutron Fluence** on the same grid as before

\* Neutron fluence [ $\text{cm}^{-2}$ ]

*	+	1	+	2	+	3	+	4	+	5	+	6	+	7	+
USRBIN				11.0	NEUTRON			-41.0		10.0				15.0	TargNeu
USRBIN				0.0				-5.0		100.0				200.0	&

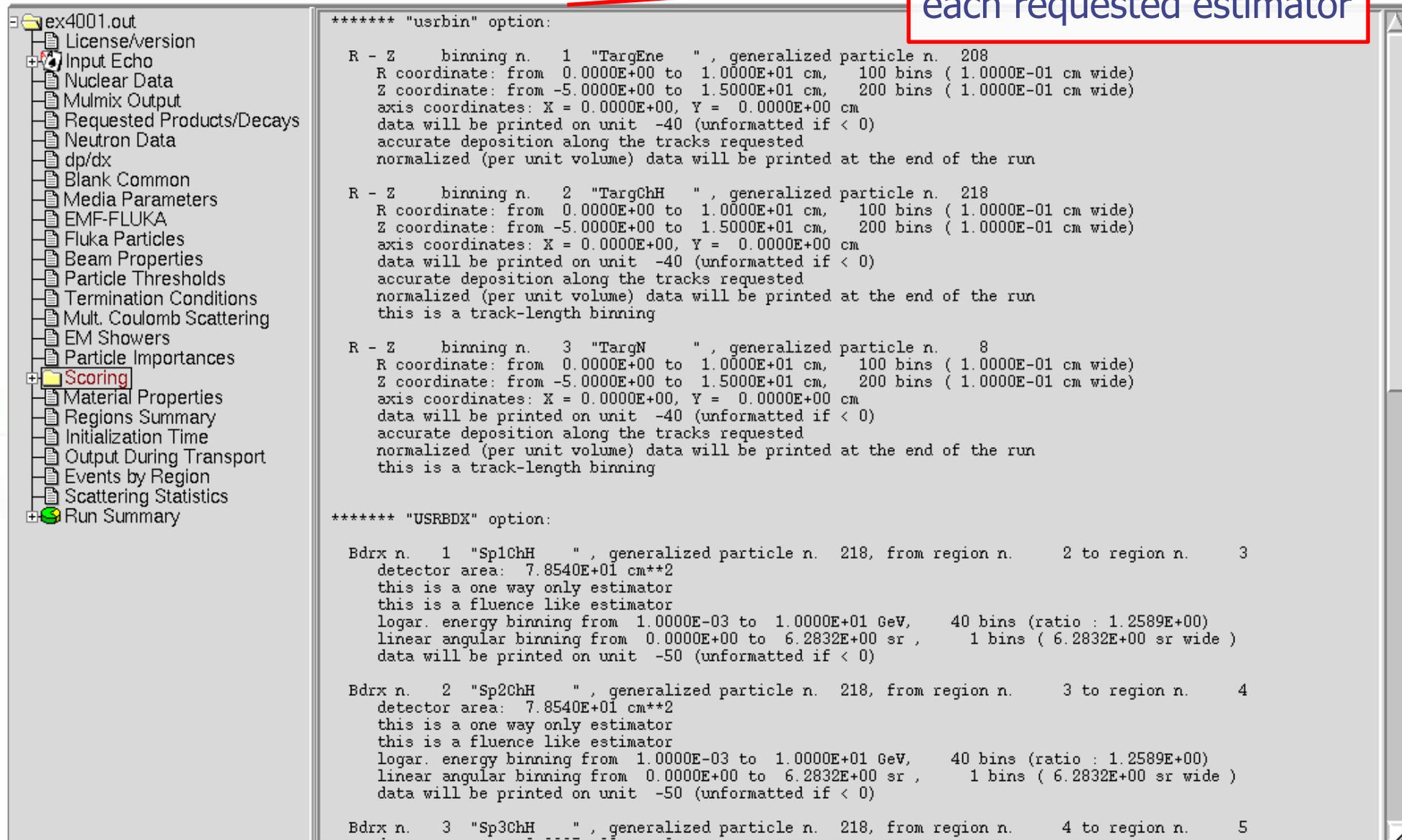
 <b>USRBIN</b>	Unit: 41 BIN ▼	Name: TargNeu
Type: R- $\Phi$ -Z ▼	Rmin: 0	NR: 100
Part: NEUTRON ▼	X:	N $\Phi$ :
	Zmin: -5.0	NZ: 200
	Rmax: 10	
	Y:	
	Zmax: 15.0	

- This is a R-Z- $\Phi$  binning (what(1)=11), scoring neutron fluence, writing the unformatted output on unit 40, spanning  $0 < R < 10$  cm in 100 bins,  $0 < \Phi < 2\pi$  in 1 bin (default),  $5 \text{ cm} < z < 15 \text{ cm}$  in 200 bins.

3) Add USRBIN to score **Charged Hadron Fluence** (HAD-CHAR) on the same grid as before.  
Output unit = 42 BIN

# Check in the .out file

Complete description of each requested estimator



The image shows a file explorer on the left and a terminal window on the right. The file explorer displays a directory tree for 'ex4001.out', with 'Scoring' highlighted. The terminal window shows the output of the simulation, including options for 'usrbin' and 'USRBDX', and detailed descriptions of three different estimators (Bdrx n. 1, 2, and 3).

```
***** "usrbin" option:  
  
R - Z      binning n. 1 "TargEne  ", generalized particle n. 208  
R coordinate: from 0.0000E+00 to 1.0000E+01 cm, 100 bins ( 1.0000E-01 cm wide)  
Z coordinate: from -5.0000E+00 to 1.5000E+01 cm, 200 bins ( 1.0000E-01 cm wide)  
axis coordinates: X = 0.0000E+00, Y = 0.0000E+00 cm  
data will be printed on unit -40 (unformatted if < 0)  
accurate deposition along the tracks requested  
normalized (per unit volume) data will be printed at the end of the run  
  
R - Z      binning n. 2 "TargChH  ", generalized particle n. 218  
R coordinate: from 0.0000E+00 to 1.0000E+01 cm, 100 bins ( 1.0000E-01 cm wide)  
Z coordinate: from -5.0000E+00 to 1.5000E+01 cm, 200 bins ( 1.0000E-01 cm wide)  
axis coordinates: X = 0.0000E+00, Y = 0.0000E+00 cm  
data will be printed on unit -40 (unformatted if < 0)  
accurate deposition along the tracks requested  
normalized (per unit volume) data will be printed at the end of the run  
this is a track-length binning  
  
R - Z      binning n. 3 "TargN     ", generalized particle n. 8  
R coordinate: from 0.0000E+00 to 1.0000E+01 cm, 100 bins ( 1.0000E-01 cm wide)  
Z coordinate: from -5.0000E+00 to 1.5000E+01 cm, 200 bins ( 1.0000E-01 cm wide)  
axis coordinates: X = 0.0000E+00, Y = 0.0000E+00 cm  
data will be printed on unit -40 (unformatted if < 0)  
accurate deposition along the tracks requested  
normalized (per unit volume) data will be printed at the end of the run  
this is a track-length binning  
  
***** "USRBDX" option:  
  
Bdrx n. 1  "Sp1ChH  ", generalized particle n. 218, from region n. 2 to region n. 3  
detector area: 7.8540E+01 cm**2  
this is a one way only estimator  
this is a fluence like estimator  
logar. energy binning from 1.0000E-03 to 1.0000E+01 GeV, 40 bins (ratio : 1.2589E+00)  
linear angular binning from 0.0000E+00 to 6.2832E+00 sr , 1 bins ( 6.2832E+00 sr wide )  
data will be printed on unit -50 (unformatted if < 0)  
  
Bdrx n. 2  "Sp2ChH  ", generalized particle n. 218, from region n. 3 to region n. 4  
detector area: 7.8540E+01 cm**2  
this is a one way only estimator  
this is a fluence like estimator  
logar. energy binning from 1.0000E-03 to 1.0000E+01 GeV, 40 bins (ratio : 1.2589E+00)  
linear angular binning from 0.0000E+00 to 6.2832E+00 sr , 1 bins ( 6.2832E+00 sr wide )  
data will be printed on unit -50 (unformatted if < 0)  
  
Bdrx n. 3  "Sp3ChH  ", generalized particle n. 218, from region n. 4 to region n. 5
```

# Standard Postprocessing Programs (1/2)

- To analyse the results of the different scoring options, several programs are made available
- The most natural ones are kept in `$FLUPRO/flutil`.
- They assume that the **estimator files are unformatted**, and can calculate standard deviations and average values over many cycles:
  - `ustsuw.f` to analyze **USRTRACK** and **USRCOLL** outputs
  - `usxsuw.f` to analyze **USRBDX** outputs
  - `usysuw.f` to analyze **USRYIELD** outputs
  - `usbsuw.f` to analyze **USRBIN** outputs
  - `ursuw.f` to analyze **RESNUCLEi** outputs

Example in the terminal:

```
$FLUPRO/flutil/name.f
```

```
Type the input file: inputname###_fort.nn
```

```
Type the input file: inputname###_fort.nn
```

```
...
```

```
Type the input file:
```

```
Type the output file: outputname
```

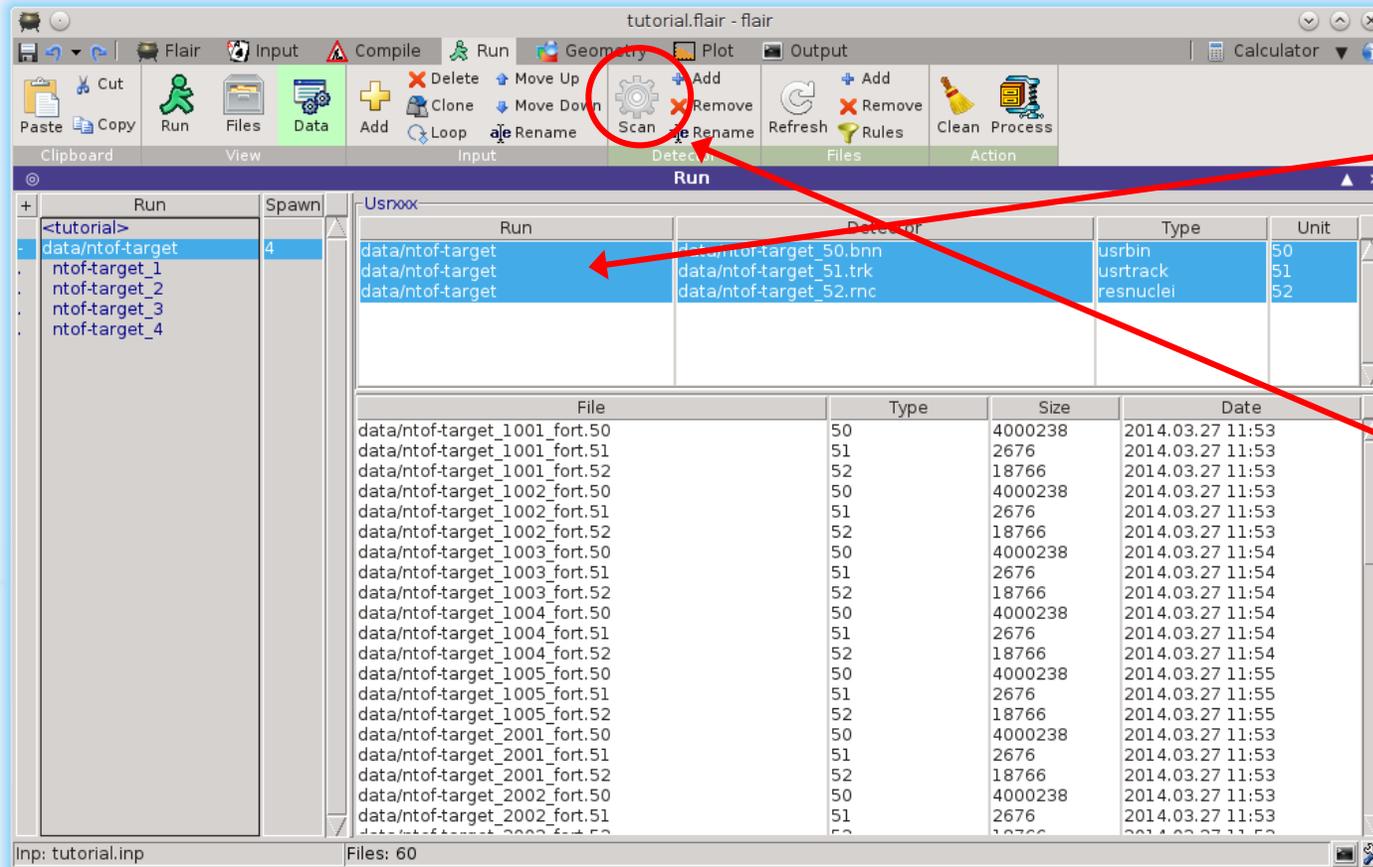
# Standard Postprocessing Programs (2/2)

- `usbsuw` produces a `_.bnn` file (binary), which can then be read by Flair for plotting
- All other programs (except `usbsuw`) produces three files:
  - a text file with extension `_sum.lis` which contains averaged distributions, standard deviations, **cumulative (integral)** quantities
  - an unformatted file which can replace the  $N$  unformatted estimator files in further postprocessing
  - a text file with extension `_tab.lis` to be easily readout by graphics codes

[Simpler programs are also provided in the manual, as guide for users who would like to write their own analysis program].

# Flair: Data Processing

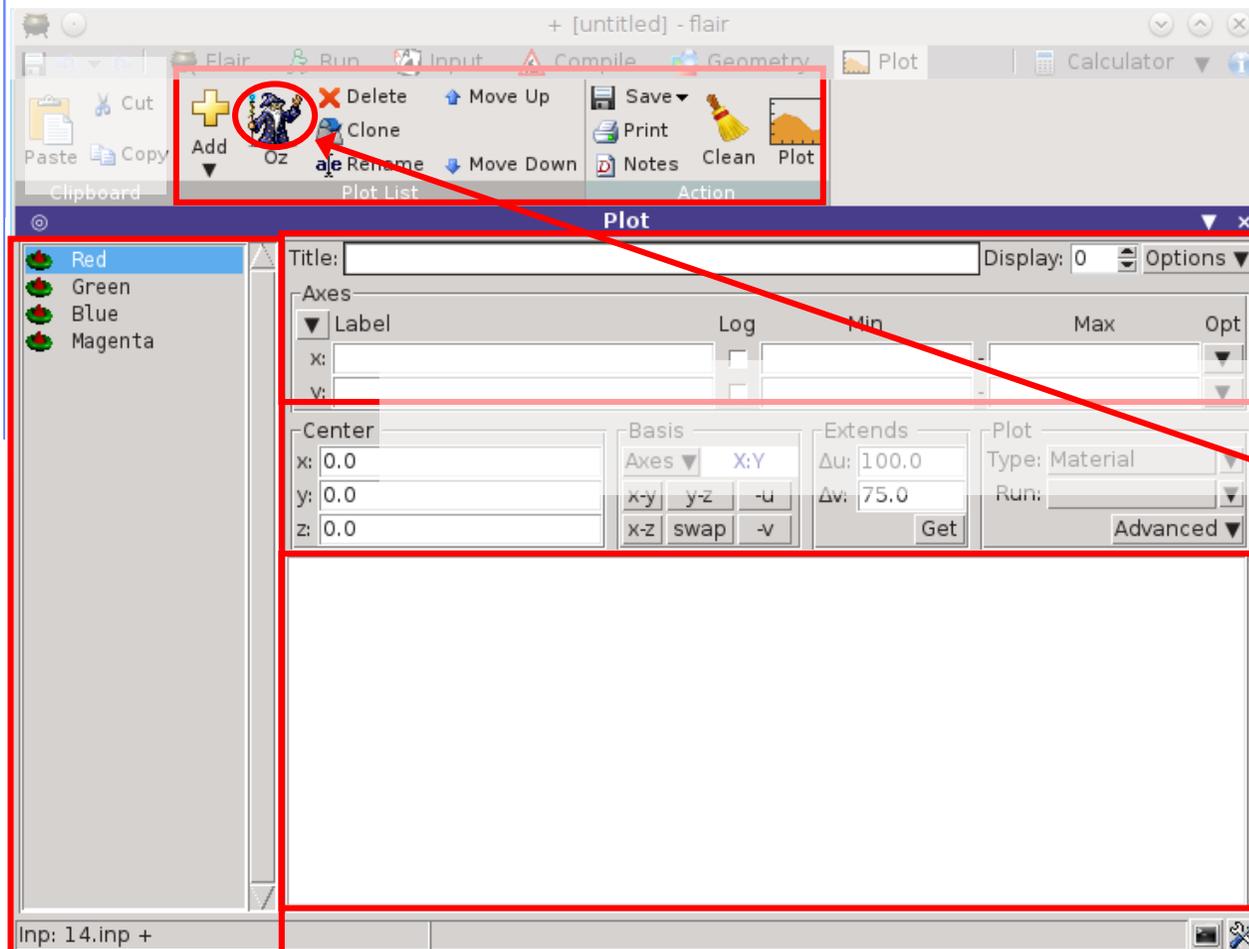
In the run tab



- Flair **initially** scans the input for possible unformatted output data for each scoring card. It **creates automatic rules for processing** (merging).
- If in the mean time you have modified the input click the **"automatic" scan**
- The default names are generated by the rules specified in the preference dialog, can be changed if desired (advanced)

FLAIR uses the auxiliary programs available in \$FLUPRO/flutil to merge output data. For USRBIN: the **usbsuw** code

# Plot List

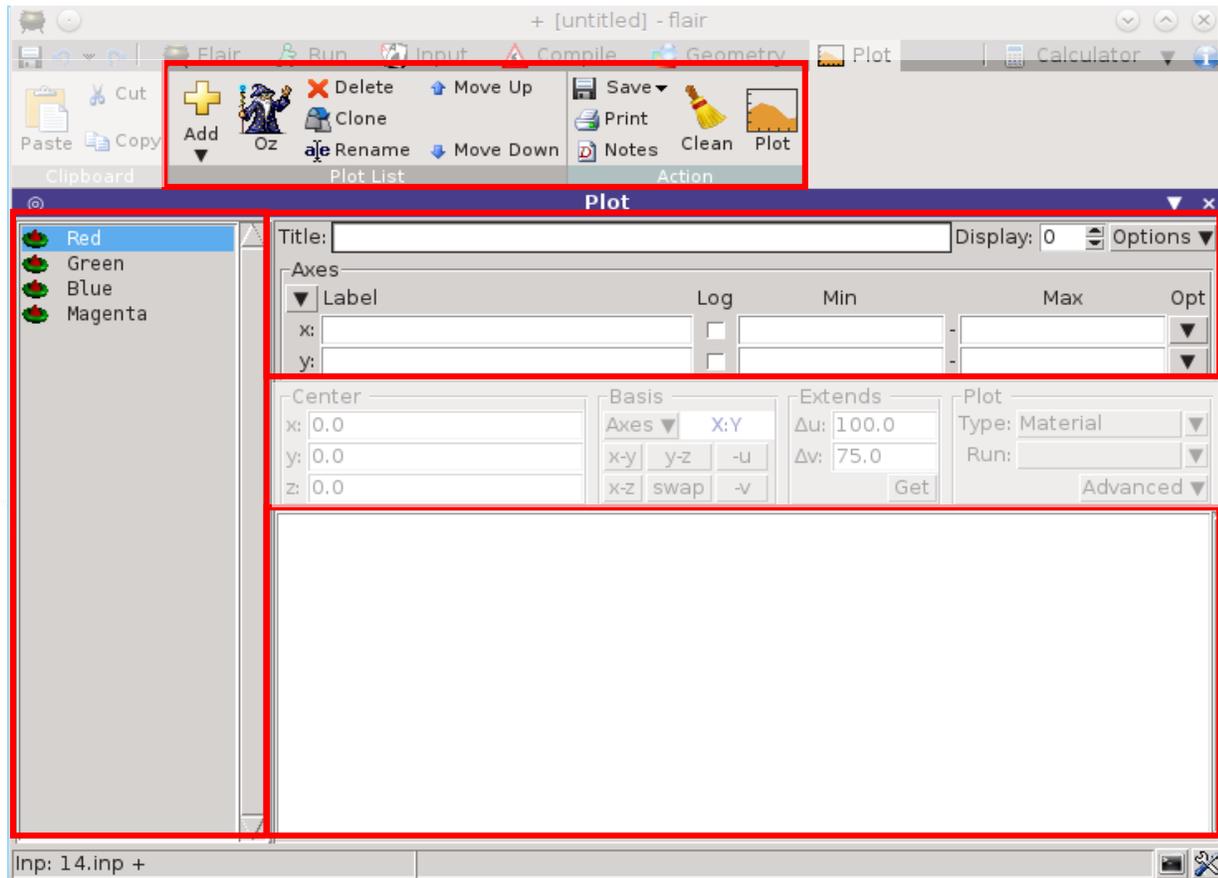


- Plots can be created in the "Plot" list frame. Either Add new plots or Clone from existing ones.
- It is important to set a unique filename for each plot.  
This filename will be used for every auxiliary file that the plot needs (with different extensions)
- The Wizard button creates automatically one plot for each processed unit
- Double click on a plot, or hit **Enter** or click the **Edit icon** to display the plotting dialog

## Plot Types

- Geometry For geometry plots
- USRBIN For plotting the output of USRBIN
- USR-1D To plot single differential quantities from cards  
**USRBDX, USRTRACK, USRCOLL, USRYIELD**
- USR-2D To plot double differential from **USRBDX**
- RESNUCLE To plot 1d or 2d distributions of **RESNUCLEi**
- USERDUMP To plot the output of USERDUMP, Useful for visualizing the source distribution.

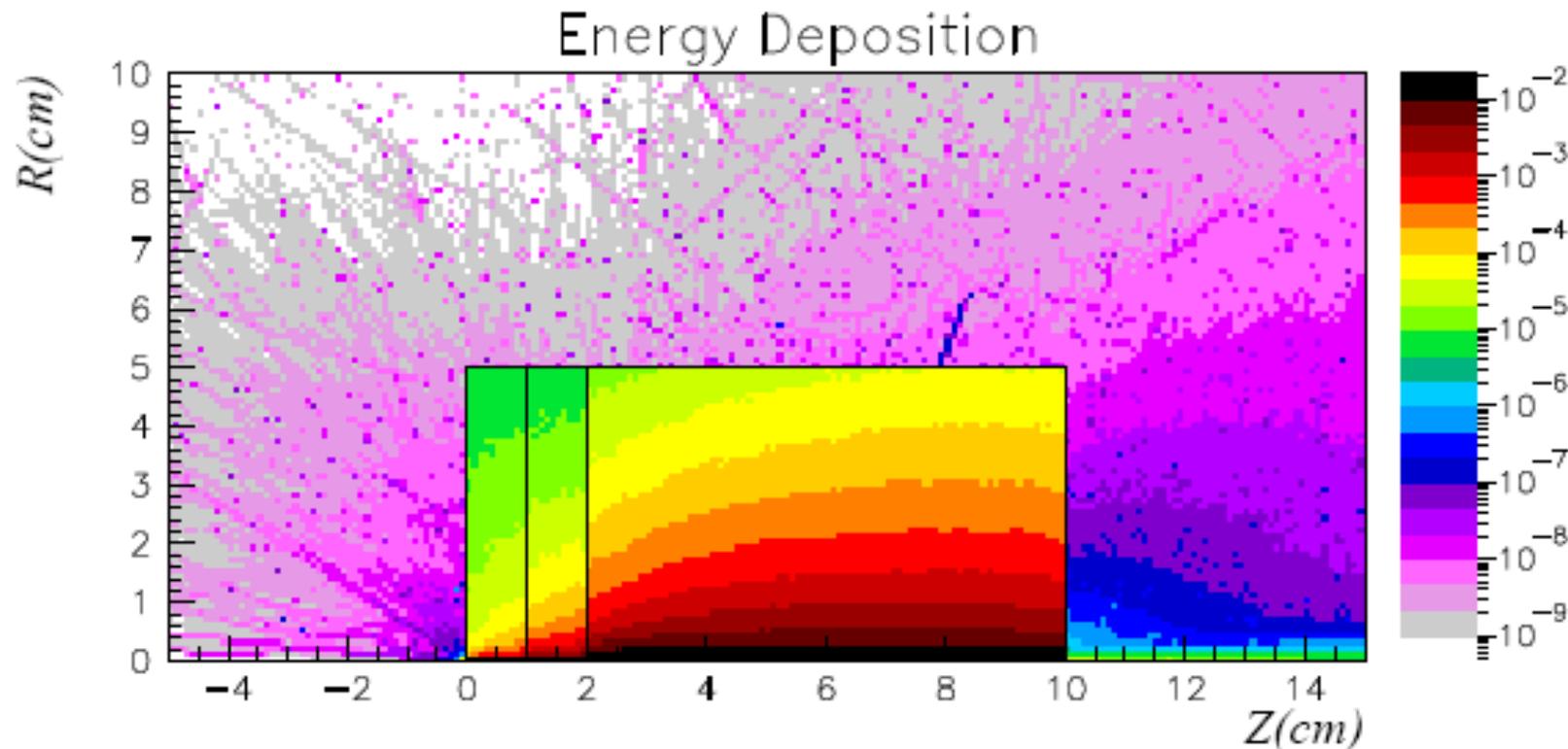
# Plotting Frames



- **FLAIR uses gnuplot to produce plots**
- All plot types share some common fields: Title + options, Filename, Axis Labels, Legends (Keys) and Gnuplot Commands.
- **Plot** button (Ctrl-Enter) will generate all the necessary files to display the plot, **ONLY** if they do not exist.
- All plots are listed in the **Plot List**
- **Clean** button will remove all files generated by Flair during plotting process.
- Check the Gnuplot manual to provide additional customization commands: e.g. To change the title font to Times size=20, add in the Opt: field the command: font 'Times,20'

# USRBIN → The Result

**WHAT(2) = ENERGY** : Energy deposition from a 3.5 GeV proton beam hitting at [0.,0.,0.] directed along z. Results are normalized to **GeV/cm<sup>3</sup> per primary**

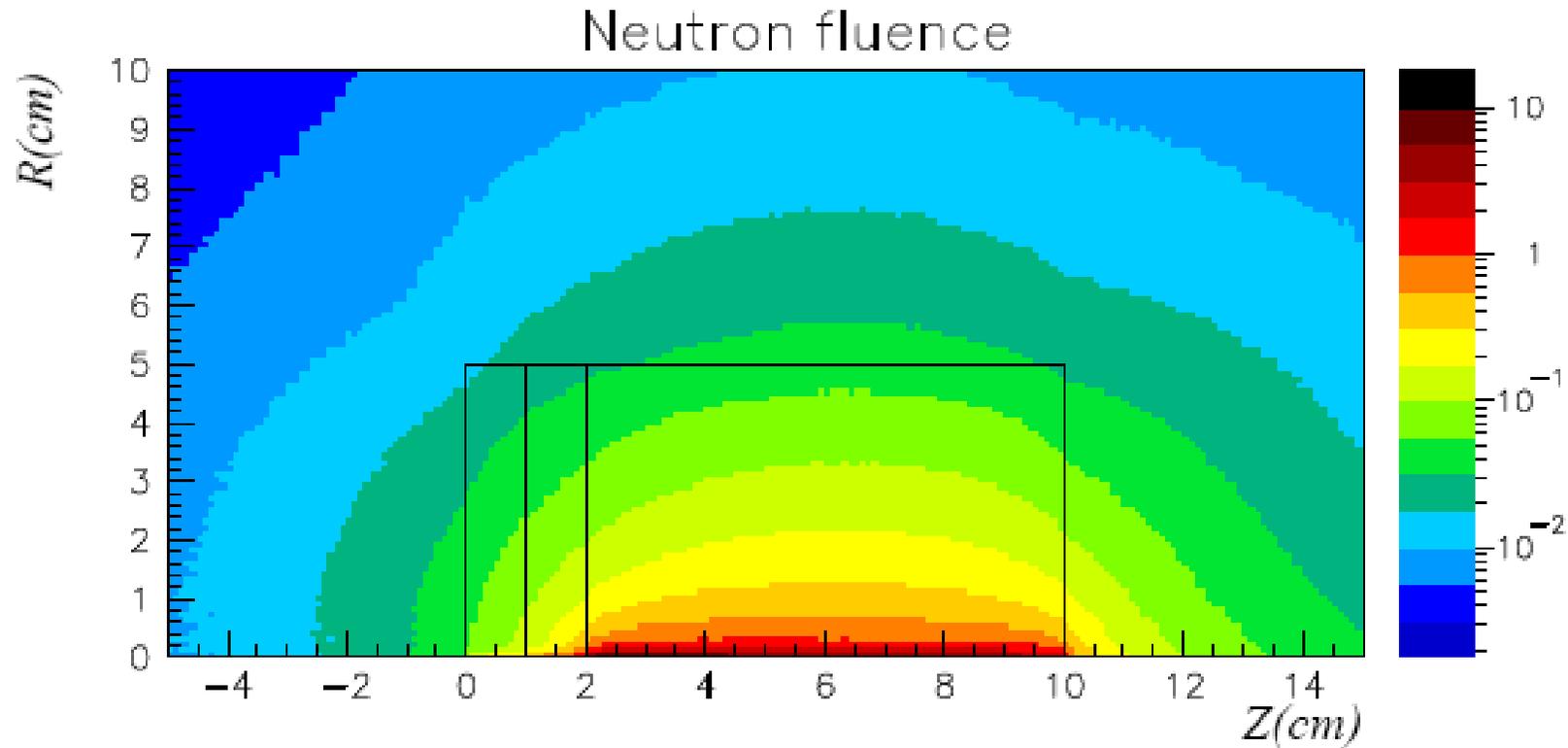


This plot is a 2D projection of a 3D structure → the result is the **AVERAGE** over the 3<sup>rd</sup> coordinate. Projection limits can be set in FLAIR

# USRBIN → The Result

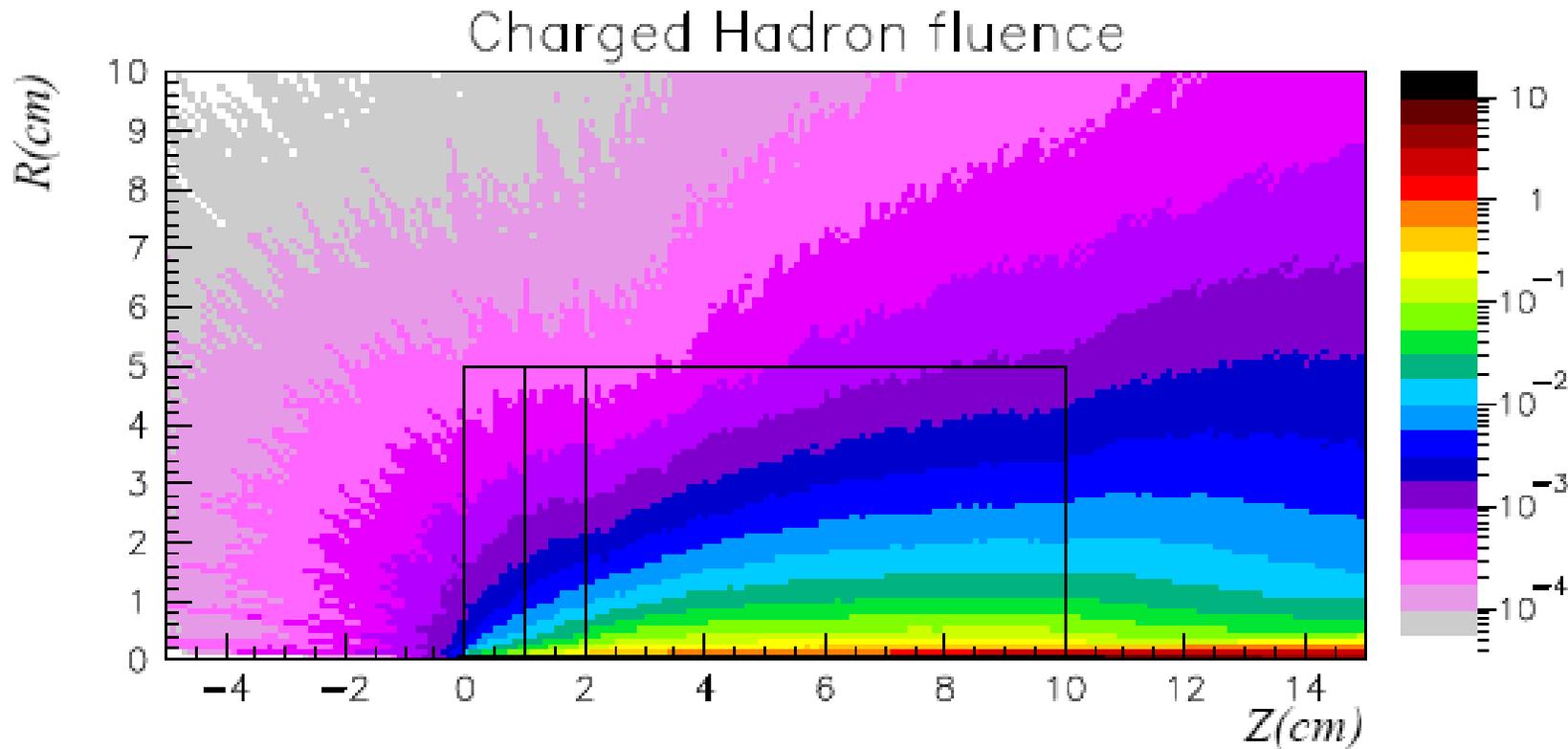
Same, **WHAT(2)**= NEUTRON to get neutron fluence.

Results are normalized to **particles/cm<sup>2</sup> per primary**



# USRBIN → The Result

Same, **WHAT(2)**= HAD-CHAR to get charged hadron fluence.  
Results are normalized to **particles/cm<sup>2</sup> per primary**



# USRBIN – more quantities

**USRBIN** can score Particle fluence as well as “Generalized particles”, either fluence-like or energy-like, for instance with **what(2)** =

- **DOSE:** Energy/unit mass (GeV/g)
- **DPA-SCO:** Displacements per atom ( see the lecture on Ionization and transport)
- **X-MOMENT:** x-component of momentum transfer (GeV/c)
- **ACTIVITY:** activity per unit volume (Bq/cm<sup>3</sup>) (see lecture on radioactivity)
  
- ... and more (see in the manual)

# USRBDX

**USRBDX** scores double differential (energy and angle) particle distributions across a boundary surface. The **angle** is with respect to the normal of the surface.

**WHAT(1)** : the distribution can be fluence or current, one-way or two-ways, according to the value of **WHAT(1)** (see the manual)

**WHAT(2)** = particle type to be scored

**WHAT(3)** = logical output unit

**WHAT(4)** = first region defining the boundary

**WHAT(5)** = second region defining the boundary

**WHAT(6)** = area of the detector in  $\text{cm}^2$

**SDUM** = detector identifier

**Continuation card:**

**WHAT(1)** = maximum kinetic energy for scoring (GeV)

**WHAT(2)** = minimum kinetic energy for scoring (GeV)

**WHAT(3)** = number of energy intervals

**WHAT(4)** = maximum solid angle in sr

**WHAT(5)** = minimum solid angle (if linear angular binning) or solid angle (if logarithmic angular binning)

**WHAT(6)** = number of angular bins

# Example : USRBDX (1)

\* in this case post-processed results are single differential (already integrated over the solid angle)

1) Score **CHARGED HADRONS** at the outer surface of the lead segment (from TARGS3 to INAIR). **WHAT(1)=99** means: fluence, one-way only, log. intervals in energy. From 1 MeV to 10 GeV in 40 intervals, and *one angular interval (default)\**. **WHAT(6)** is a normalization factor: setting it equal to the surface area provides results normalized to  $\text{cm}^{-2}$  (fluence unit)  $\text{GeV}^{-1} \text{sr}^{-1}$ . Output to unformatted unit 50

```

*   +   1   +   2   +   3   +   4   +   5   +   6   +   7   +
* out from lead
USRBDX          99.0  HAD-CHAR          -50.  TARGS3      INAIR      329.87Sp3ChH
USRBDX          10.0    0.001          40.                                &
  
```

<b>▲ USRBDX</b>	Unit: 50 BIN ▼	Name: Sp3ChH
Type: Φ1,LogE,LinΩ ▼	Reg: TARGS3 ▼	Area: 329.87
Part: HAD-CHAR ▼	Emin: 0.001	Ebins: 40
	Ωmin:	Ωbins:
	Emax: 10.	
	Ωmax:	
	to Reg: INAIR ▼	

Repeat the same between **TARGS1** and **TARGS2**, and between **TARGS2** and **TARGS3** (use the right normalization factor!).

# Example : USRBDX (2)

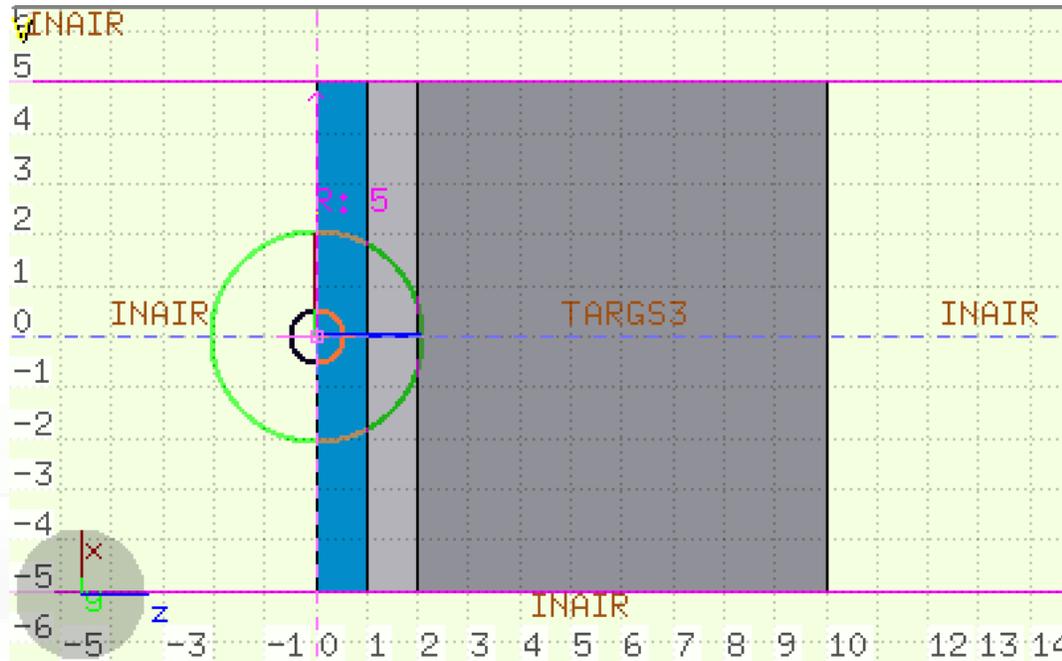
\* in this case post-processed results are single differential (already integrated over the solid angle)

2) Score at the surface between 2<sup>nd</sup> and 3<sup>rd</sup> target section, same as before but in 3 angular bins.

*	+	1	+	2	+	3	+	4	+	5	+	6	+	7	+
USRBDX				99.0	HAD-CHAR			-54.	TARGS2		TARGS3		78.5398	Sp2ChHA	
USRBDX				10.0		0.001		40.						3.0	&

<b>USRBDX</b>	Unit: 54 BIN ▼	Name: Sp2ChHA
Type: $\Phi_1, \text{LogE}, \text{Lin}\Omega$ ▼	to Reg: TARGS3 ▼	Area: 78.5398
Part: HAD-CHAR ▼	Emin: 10.	Ebins: 40
Reg: TARGS2 ▼	$\Omega$ min:	$\Omega$ bins: 3

# Example : Normalization



$R_{\text{TARG}} = 5 \text{ cm}$   
 $Z_{\text{TARGS1}} = 1 \text{ cm}$   
 $Z_{\text{TARGS2}} = 1 \text{ cm}$   
 $Z_{\text{TARGS3}} = 8 \text{ cm}$

Area between TARGS2 and TARGS3 =  $\pi R_{\text{TARG}}^2 = 78.5398$

Area between TARGS3 and INAIR =  $2\pi R_{\text{TARG}} Z_{\text{TARGS3}} + \pi R_{\text{TARG}}^2 = 329.87$

# Standard Postprocessing Programs

The FLUKA postprocessing program for USRBDX, `usxsuw.f`, generates files with the extension **`_sum.lis`** and **`_tab.lis`**. Example of `sum.lis` from USRBDX:

```
Total primaries run:          3000
Total weight of the primaries run: 3000.00000
Detector n:          1 (      1 ) Sp2ChHA
(Area:              78.5398026 cmq,
distr. scored:      218 , Charged Hadrons
from reg.           3 to      4 ,
one way scoring,
fluence scoring scoring)

Tot. resp. (Part/cmq/pr) 1.4218389E-02 +/- 1.486696 %
( --> (Part/pr)      1.116709 +/- 1.486696 % )
```

User provided the area in what(6), thus normalization is /cm<sup>2</sup>

\*\*\*\* Different. Fluxes as a function of energy \*\*\*\* (integrated over solid angle) \*\*\*\*

.....

\*\*\*\* Cumulative Fluxes as a function of energy \*\*\*\* (integrated over solid angle)

.....

\*\*\*\* Double diff. Fluxes as a function of energy \*\*\*\*

# Standard Postprocessing Programs

- Example of **tab.lis** for USRBDX

# Detector n: 1 Sp2ChHA (integrated over solid angle)

# N. of energy intervals 40

<b>Emin</b>	<b>Emax</b>	<b>Result</b>	<b>Error (%)</b>
1.000E-02	1.259E-02	5.049E-03	6.420E+01
1.259E-02	1.585E-02	1.115E-03	6.934E+01
1.585E-02	1.995E-02	2.826E-03	4.675E+01
1.995E-02	2.512E-02	2.356E-03	3.866E+01
2.512E-02	3.162E-02	6.437E-03	2.857E+01

...

# double differential distributions

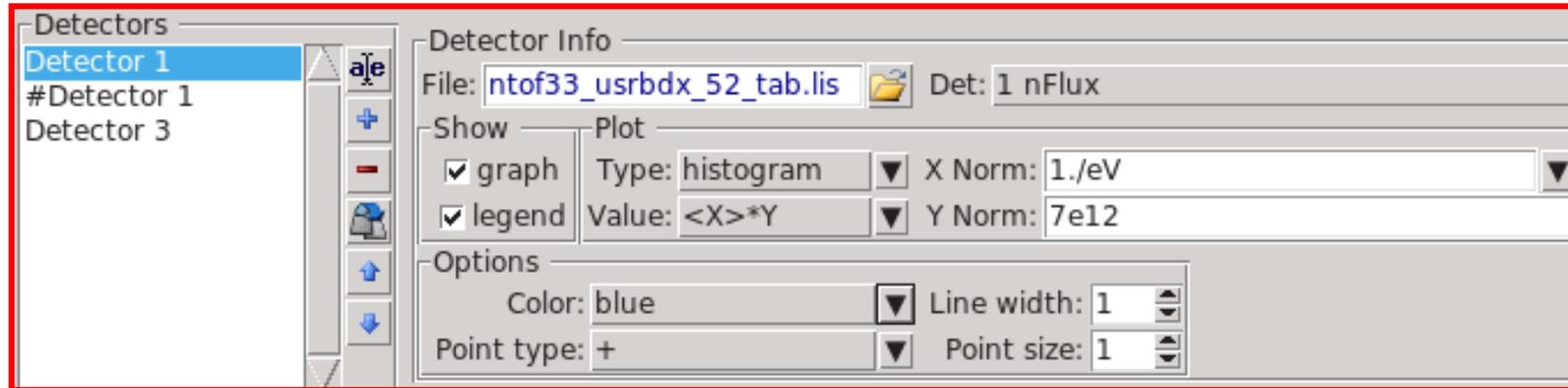
# number of solid angle intervals 3

# Block n: 1 0.00000000 : 2.09439516

1.000E-02	1.259E-02	2.980E-04	9.900E+01
1.259E-02	1.585E-02	1.981E-04	9.900E+01
1.585E-02	1.995E-02	3.866E-04	6.700E+01
1.995E-02	2.512E-02	7.171E-04	3.402E+01
2.512E-02	3.162E-02	5.544E-04	4.550E+01

# FLAIR USR-1D Single Differential Plot

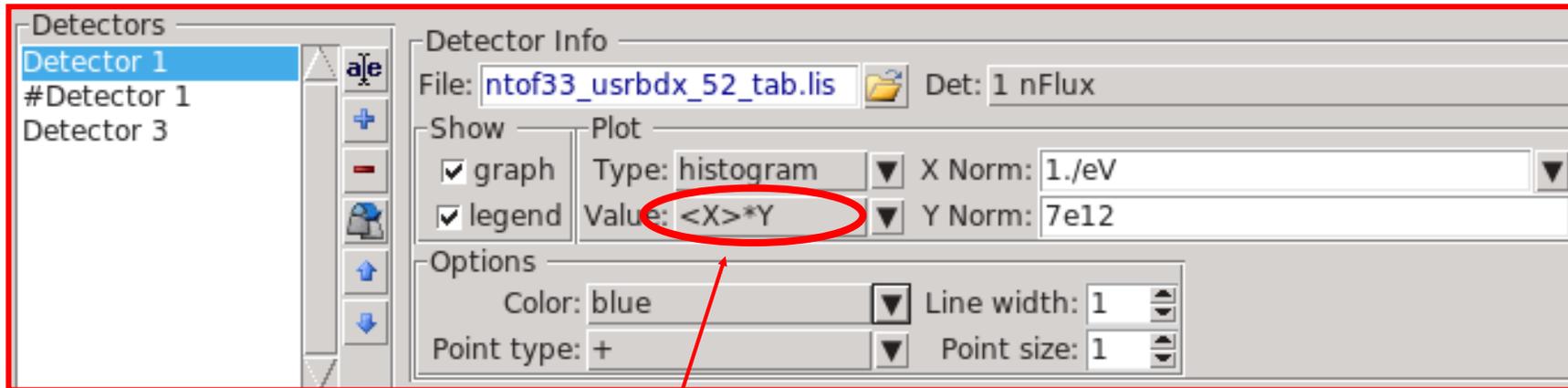
Plot → Add → USR-1D (or using the Wizard button)



- USR-1D is able to plot the 1D single differential information from the USRBDX, USRCOLL, USRTRACK and USRYIELD cards (The 2D information is not handled).
- The file type in use should have the extension **\_tab.lis** (generated by the FLUKA data merging tools)
- You can superimpose many scoring output in a single plot.
- Error bars can be plotted (for instance using histerr instead of histogram)

# FLAIR USR-1D Single Differential Plot

Plot → Add → USR-1D (or using the Wizard button)



**WARNING:  
use Lethargy plot!!!**

# Lethargy plots

- What is Lethargy? Why to use it?

Case 1: Linear plot  $F(x)$  vs  $x$ .

The integral between  $x_1$  and  $x_2$  is

$$I = \int_{x_1}^{x_2} F(x) dx$$

Comparing the  $F(x)$  value in each bin one has a direct, immediate feeling which bins contribute more to the integral.

# Lethargy plots

- What is Lethargy? Why to use it?

Case 2: x-logarithmic plot  $F(x)$  vs  $\log(x)$

$$z = \log(x) \quad , \quad dz = \frac{dx}{x} \quad \Rightarrow \quad dx = x dz$$

Making a change of variable in  $I = \int_{x_1}^{x_2} F(x) dx$  one gets:

$$I = \int_{\log(x_1)}^{\log(x_2)} x F(z) dz = \int_{\log(x_1)}^{\log(x_2)} x F(\log(x)) d(\log(x))$$

$$\neq \int_{\log(x_1)}^{\log(x_2)} F(\log(x)) d(\log(x))$$

Comparing the  $x * F(\log(x))$  value in each bin one has a direct, immediate feeling which bins contribute more to the integral.

# Lethargy plots

- What is Lethargy? Why to use it?
  - If  $E$ =particle energy and  $N$ =number of particles, a plot in  $E \cdot dN$  vs  $\log(E)$  (in FLAIR :  $y^* <x>$  plot option) is called **Lethargy plot**. It allows giving the correct importance to different energy ranges in the plot:
    - i.e. the area of each bin is proportional to the corresponding integral flux, and comparing the bins you have immediately the feeling in which energy bin you have more/less particles.
  - Lethargy plots must be used in general for any particle type with an energy spectrum extended over many orders of magnitude.
  - It is the standard representation for Low Energy Neutrons spectra, also because of the  $1/E$  dependence of moderated n spectra (see lecture on Low-Energy Neutrons).

# USRBDX -> The Result

This is true only if the surface area is explicitly given

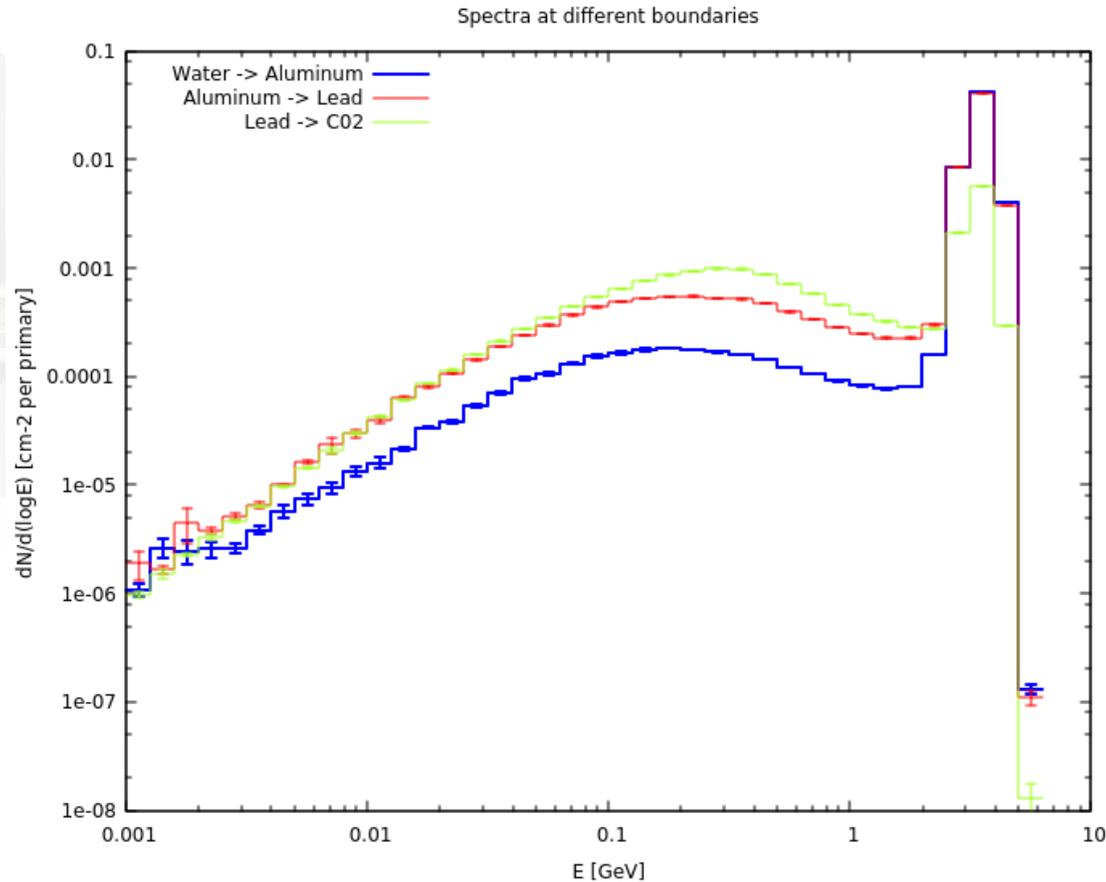
Evolution of charged hadron spectra at the various surfaces  
post-processed results are normalized to  $\text{GeV}^{-1} \text{cm}^{-2}$  per primary

From post-processing we get **single** differential spectra since we asked for one angular bin only

$$y = \frac{dN}{d(\log E)} = E \frac{dN}{dE}$$

Value:

Lethargy plot



# USRBDX ->The Result

This is true only if the surface area is explicitly given

Evolution of charged hadron spectra at the various surfaces  
post-processed results are normalized to  $\text{GeV}^{-1} \text{cm}^{-2}$  per primary

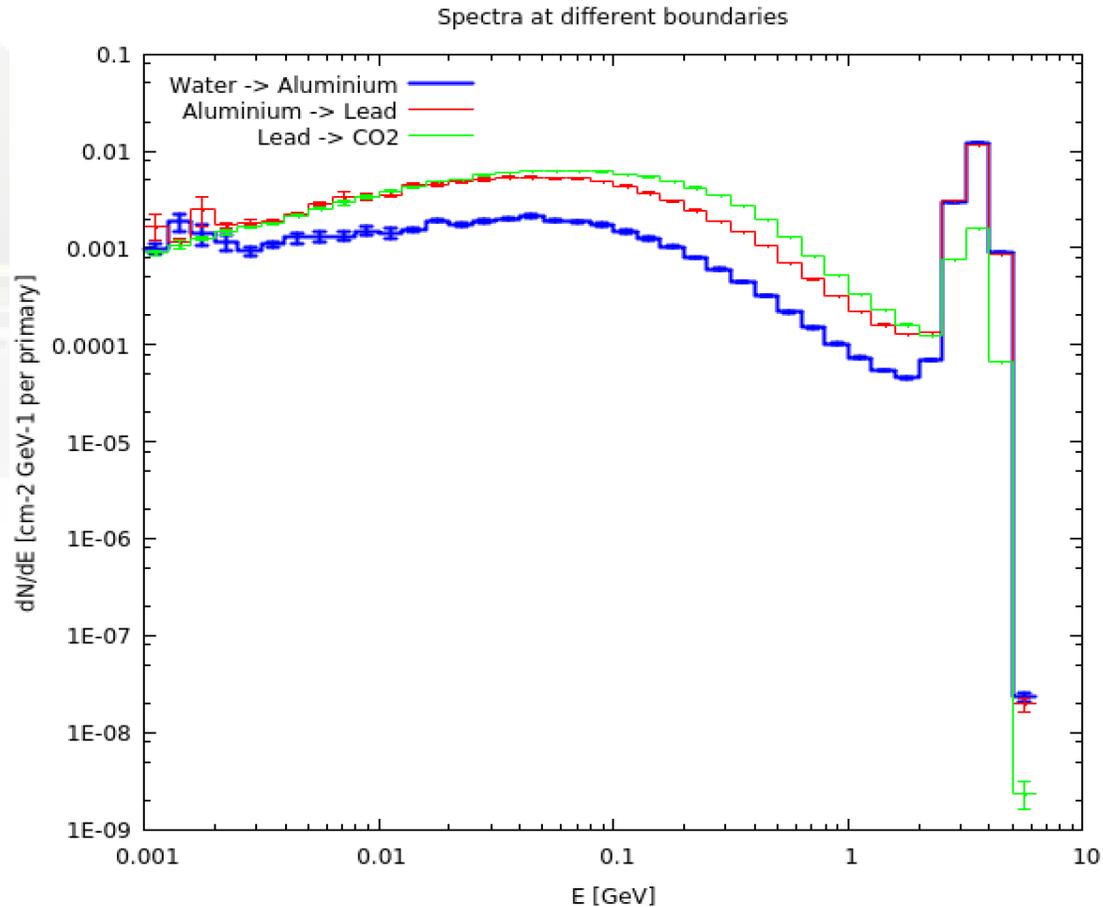
$$y = \frac{dN}{dE}$$

Value: Y ▼

Not Lethargy plot.

The low energy area  
"seems" to  
contribute more.  
This may be  
misleading...  
Be careful!

From post-processing we get **single** differential spectra  
since we asked for one angular bin only

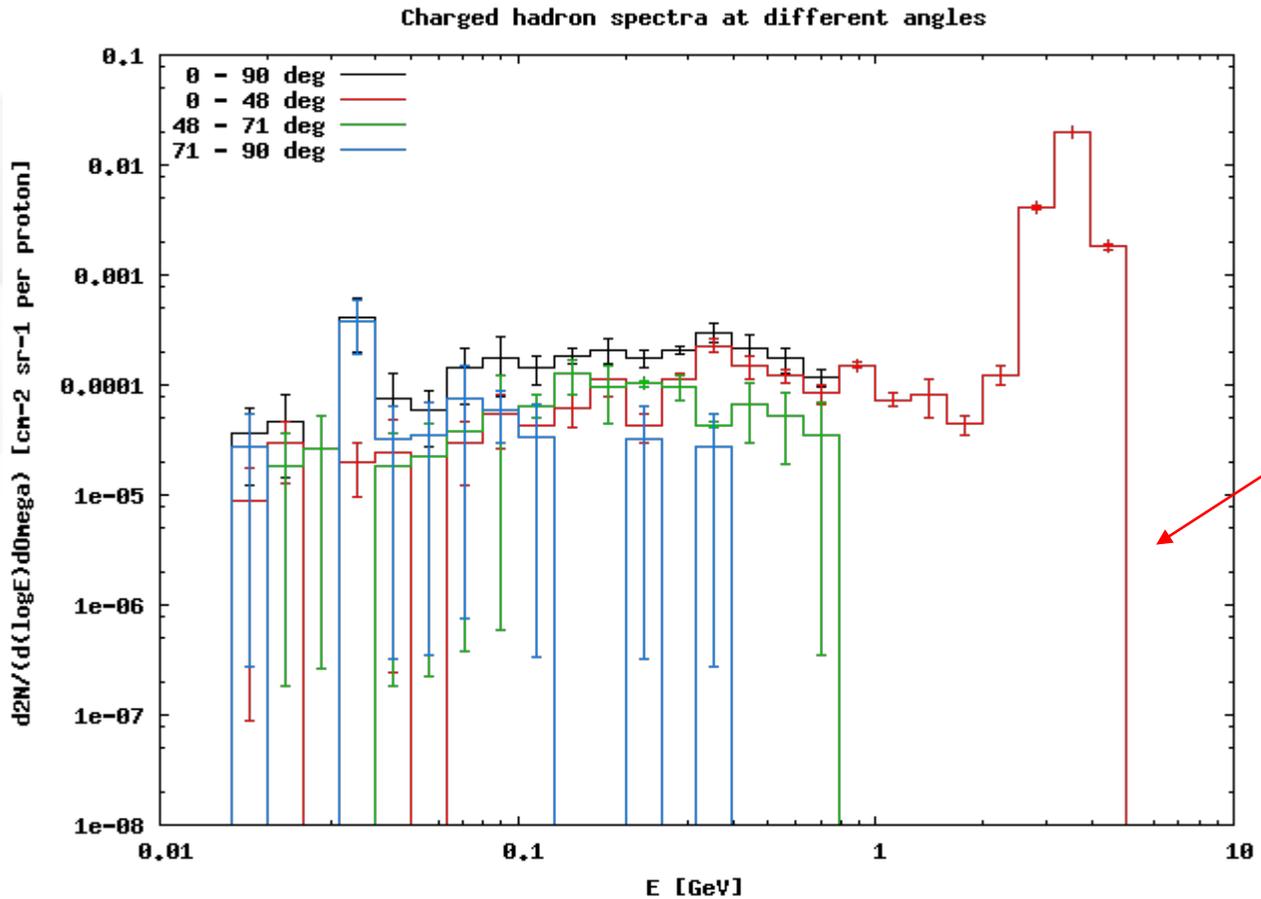


# USRBDX -> The Result

Double differential charged hadron spectra for consecutive solid angle portions results are normalized to  $\text{GeV}^{-1} \text{sr}^{-1} \text{cm}^{-2}$  per primary

From post-processing we get **double** differential spectra since we asked for more than one angular bin, but the angle-integrated spectrum is provided as well on top

This is true only if the surface area is explicitly given



Forward is high energy

# USRTRACK

- Calculates **differential fluence** as a function of **energy** by scoring track-length in a given **region**. Results are normalized to  $GeV^{-1} cm^{-2}$  per primary if the region volume is provided (otherwise should be intended as  $GeV^{-1} cm$  per primary, i.e. differential track-length)

For all SDUMs except EN-NUCL and ENERGY:

WHAT(1) : linear (=1) or logarithmic (= -1) binning energy

WHAT(2) = particle type to be scored

WHAT(3) = logical output unit

WHAT(4) = region defining the detector (if >0) or all regions (if = -1)

WHAT(5) = volume of the detector in  $cm^3$

WHAT(6) = number of energy bins

SDUM = track-length detector identifier

Continuation card:

WHAT(1) = maximum kinetic energy for scoring (GeV)

WHAT(2) = minimum kinetic energy for scoring (GeV)

For SDUM= EN-NUCL: The energy scale will be changed from energy to energy per nucleon. For SDUM= ENERGY: the energy scale will be changed to the default, that is total kinetic energy.

# USRTRACK

- Score track-length of charged hadrons in TARGS3, in a logarithmic binning energy (WHAT(1)=-1) using 40 energy bins, between 1 MeV and 10 GeV. Normalize to the region volume in order to have the results in  $[GeV^{-1} cm^{-2} per primary]$

```

*   +   1   +   2   +   3   +   4   +   5   +   6   +   7   +
*           log   partype  out.unit  region  volume  #bins
*           Emax   Emin
USRTRACK   -1.0  HAD-CHAR  -55.    TARGS3  628.31  40.TrChH
USRTRACK   10.0   0.001
  
```

---

```

USRTRACK                               Unit: 55 BIN ▼           Name: TrChH
Type: Log ▼                               Reg: TARGS3 ▼           Vol: 628.31
Part: HAD-CHAR ▼                           Emin: 0.001           Emax: 10.           Bins: 40.
  
```

**Remember:** USRTRACK scores differential fluence in a **region**, USRBDX scores fluence or current on a **surface**, and USRBIN scores e.g. fluence in **volumes** and gives no differential information.

# USRYIELD

- Scores a **double-differential particle yield** across a boundary surface or at interaction points. Angles wrt **beam direction**. Only 1 interval in the “second” quantity
- “Energy-like” quantities

Kinetic energy , total momentum , total energy , longitudinal momentum in the lab frame ,  
longitudinal momentum in the c.m.s. frame **LET**

- “Angle-like” quantities

Rapidity in the lab frame , rapidity in the c.m.s. frame , pseudorapidity in the lab frame ,  
pseudorapidity in the c.m.s. frame , Feynman-x in the lab frame ,  
Feynman-x in the c.m.s. frame , transverse momentum , transverse mass ,  
polar angle (\*) in the lab frame , polar angle (\*) in the c.m.s. frame ,  
square transverse momentum , **charge** , weighted angle in the lab frame ,  
weighted transverse momentum

# USRYIELD

**WARNING!! calculating a cross section has no meaning in case of a thick target.**

- While option USRBDX calculates angular distributions **WITH RESPECT TO THE NORMAL** to the boundary at the point of crossing, **USRYIELD's** distributions are calculated **WITH RESPECT TO THE BEAM DIRECTION**, as defined by BEAMPOS (or a different direction specified with **SDUM=BEAMDEF**).

**WHAT(1)** : the two physical quantities with respect to which the double-differential yield is calculated.

**WHAT(2)** = particle type to be scored

**WHAT(3)** = logical output unit

**WHAT(4)**= first region defining the boundary

**WHAT(5)**= second region defining the boundary

**WHAT(6)** = normalization factor

**SDUM** = detector name

Continuation card:

**WHAT(1)**= upper limit for the first quantity

**WHAT(2)**= lower limit for the first quantity

**WHAT(3)** = number of scoring intervals for the first quantity

**WHAT(4)**= upper limit for the second quantity

**WHAT(5)**= lower limit for the second quantity

... (see the manual)

# USRYIELD

- Score plain double differential yield (continuation card WHAT(6)=3) of pions, with the first quantity being the polar angle (degree) and the second quantity being the kinetic energy (WHAT(1)=124), between TARGS3 and INAIR, between 0 and 180 deg in 18 bins and between 0 and 10 GeV.

```

*      124 = 24 + 1 * 100 => polar angle (in degrees) and kinetic energy
*
*      out.unit      Reg1      Reg2      Norm
*      Amax      Amin      #Abins      Emax      Emin      dbl.differential
*      +      1      +      2      +      3      +      4      +      5      +      6      +      7      +
USRYIELD      124.0      PIONS+-      -57.      TARGS3      INAIR      1.0YieAng
USRYIELD      180.0      0.0      18.      10.0      0.0      3.0&
  
```

<b>⚠ USRYIELD</b>	Type: Yield ▼	Unit: 57 BIN ▼	Name: YieAng
ie: Polar θ lab deg ▼	ie: Ekin GeV ▼	Log: Linear ▼	Norm: 1.0
Part: PIONS+- ▼	Yield: ▼	Reg: TARGS3 ▼	to Reg: INAIR ▼
Min1: 0.0	Max1: 180.0	Nbins1: 18.0	
Min2: 0.0	Max2: 10.0	Kind: d2N/dx1dx2 ▼	Mat: ▼

- Only one interval is possible for the second variable, BUT results are normalized as Double Differential (in this case, charged pions GeV<sup>-1</sup> sr<sup>-1</sup> per primary)

**WARNING!!**  
 Use WHAT(6) = 3 for plain double differential yield, the DEFAULT is plain double-differential cross section !! 50

# Standard Postprocessing Programs

The FLUKA postprocessing program for USRYIELD, `usysuw.f`, generates files with the extension **`_sum.lis`** and **`_tab.lis`**. Example of `sum.lis` from USRYIELD

Detector n: 1 YieAng

(User norm: 1.

sigma: 1. mb

distr. scored: 209 , **PIONS+-**

from reg. 4 to reg. 5 )

linear 1st variable (x1) binning from 0.0000E+00 to 3.1416E+00 18 bins  
( 1.7453E-01 wide)

2nd variable (x2) ranges from 0.0000E+00 to 1.0000E+01

1st variable (x1) is: **Laboratory Angle (radians)**

2nd variable (x2) is: **Laboratory Kinetic Energy**

The scored double differential yield is (normalized per primary particle):

**plain d2 N / dx1 dx2** where x1, x2 are the first and second variables

Tot. response (integrated over x1) 2.6339998E-02 +/- 3.883959 %

**WARNING!! The Tot. response is NOT integrate over the second quantity !!**

in this case it turns out to be *particles/GeV per primary*  
(to be multiplied by the energy width interval of 10GeV)

# Standard Postprocessing Programs

- Example of `tab.lis` for USRYIELD

# Detector n: 1 YieAng

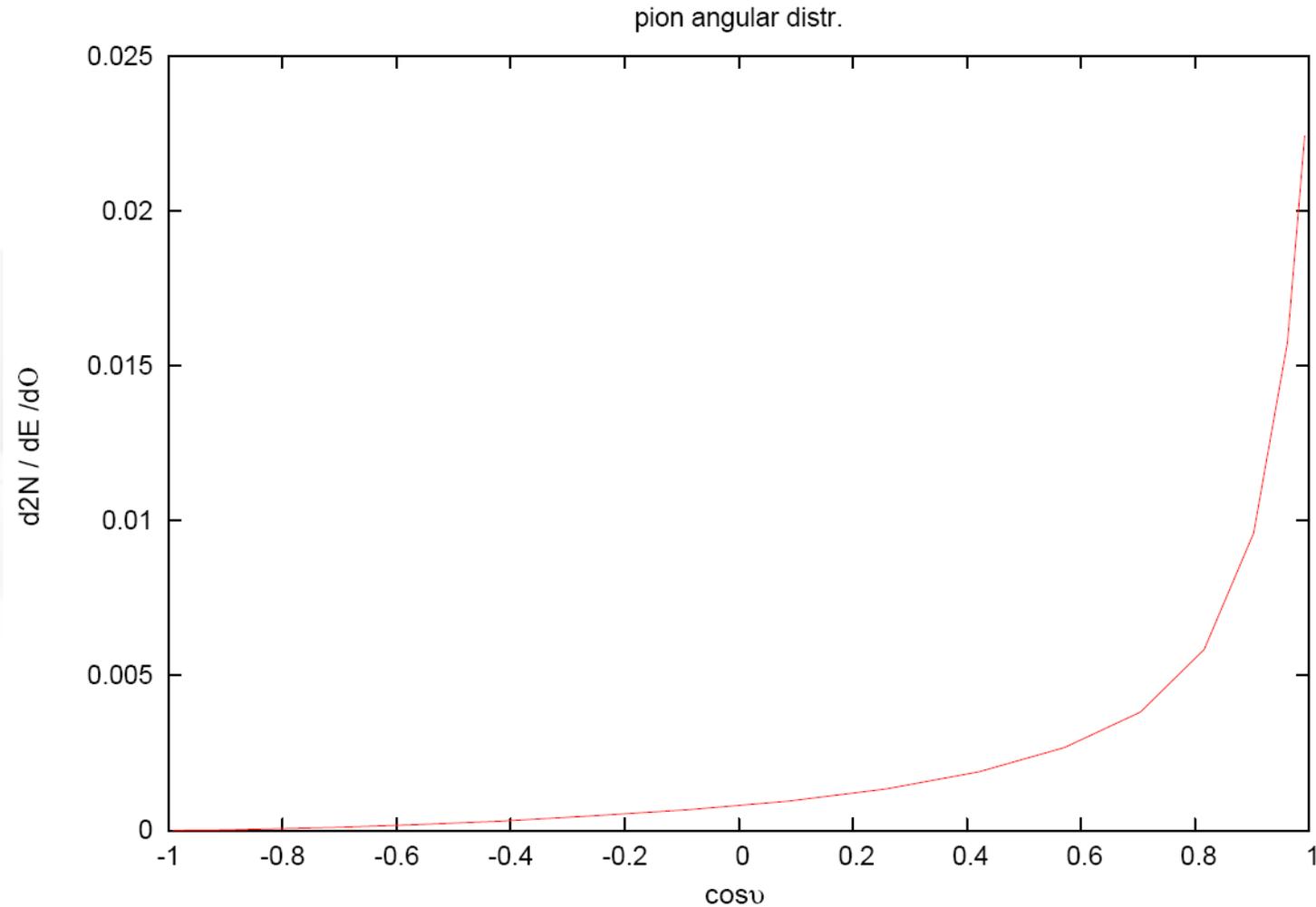
# N. of x1 intervals 18

Thetamin	Thetamax	Result	Error (%)
0.000000	0.1745329	2.0742605E-02	10.87912
0.1745329	0.3490658	1.4463779E-02	10.65940
0.3490658	0.5235988	9.8084798E-03	7.649231
0.5235988	0.6981317	5.8580171E-03	4.966214
0.6981317	0.8726646	3.8220894E-03	10.60832
0.8726646	1.047198	2.6973977E-03	5.450788

...

# USRYIELD -> The Result

- pion angular distribution



# Bins, units etc

- Results from USRBDX, USRYIELD, USRTRACK, USRCOLL are given as **DIFFERENTIAL** distributions of fluence in energy, in units of (cm<sup>-2</sup>) **GeV<sup>-1</sup>** per primary. Thus,
- A) results are independent on the chosen binning
- B) to obtain INTEGRAL results (fluence in cm<sup>-2</sup> PER ENERGY INTERVAL per primary) one must multiply the value of each energy bin by the width of the bin (even for logarithmic binning):

$$N = \int \frac{dN}{dE} dE$$

- When scoring **Neutrons** (see lecture), the E bin limits below 20 MeV are automatically set==transport groups limits
- The normalization is per cm<sup>-2</sup> ONLY if the user provides the surface area or region volume. FLUKA is not able to calculate areas/volumes
- The same if USRBIN is used by region
- Instead, the normalization for USRBIN on grids is automatically in cm<sup>-2</sup> ( or cm<sup>-3</sup> for deposited energy etc)

# FILTERS : AUXSCORE

There is the possibility to **filter** the estimators, restricting the scoring to a selected subset of particles.

For instance: : score the yield of 56-Iron ions (very useful: there is no separate name for each ion specie, except light ones. HEAVYION score all isotopes heavier than alpha's together!)

*	+	1	+	2	+	3	+	4	+	5	+	6	+	7	+
USRYIELD				124.0		ALL-PART		-87.		TARGS3		INAIR		1.0	Fe56
USRYIELD				180.0		0.0		18.		10.0		0.0		3.0	&
AUXSCORE				USRYIELD		-5602600.				Fe56		Fe56			

<b>USRYIELD</b> ie: Polar θ lab deg ▼ Part: ALL-PART ▼ Mint: 0.0 Minz: 0.0	Type: Yield ▼ ia: Ekin GeV ▼ Yield: ▼ Max1: 180.0 Max2: 10.0	Unit: 87 BIN ▼ Log: Linear ▼ Reg: TARGS3 ▼ Nbins1: 18.0 Kind: d2N/dx1dx2 ▼	Name: Fe56 Norm: 1.0 to Reg: INAIR ▼ Mat: ▼
<b>AUXSCORE</b> Delta: ▼	Type: USRYIELD ▼ Z: 26 Det: Fe56 ▼	Part: ▼ A: 56 to Det: Fe56 ▼	Set: ▼ Isomer: 0 Step:

The requested ion is coded in what(2) according to its **A**, **Z** and (optionally) isomeric state **m**:

$$\text{what}(2) = - (100 * \mathbf{Z} + 100000 * \mathbf{A} + \mathbf{m} * 100000000)$$

Z,A,m=0 means all , e.g. -2600 == all Iron isotopes

# Warning on AUXSCORE

- To be used with care, or **NOT** used at all for energy deposition scoring.
- Why? Because **In real world, energy is eventually deposited mostly by electrons only.**
- In Monte Carlo, part of the energy is deposited “by other particles” as continuous energy deposition or point energy deposition, depending on the delta-ray threshold, production threshold, transport threshold (see lecture on transport)
- **Any filtering done with AUXSCORE on energy deposition will depend on the adopted thresholds and settings. (see lecture on transport)**