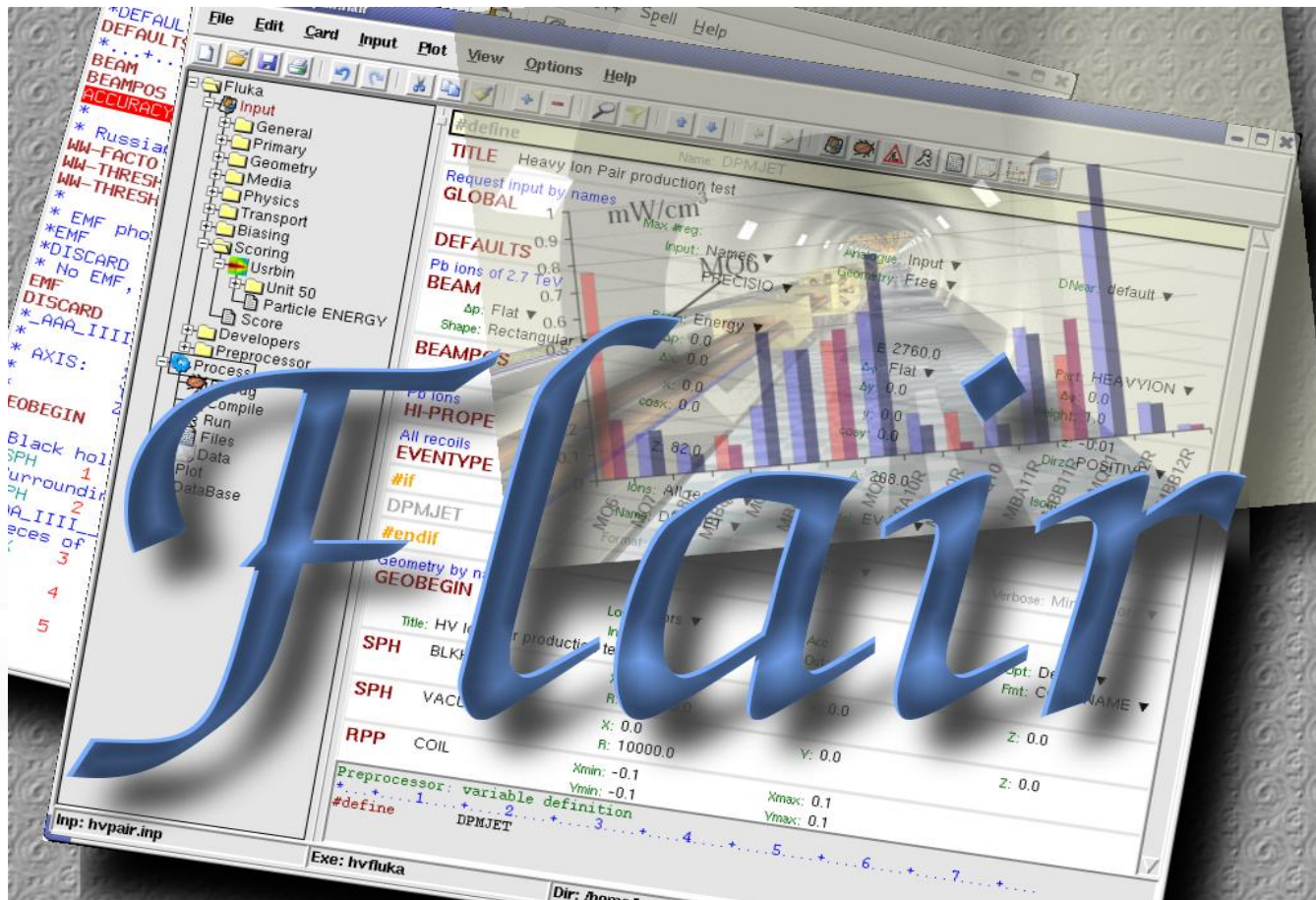




# Flair Advanced Features

Advanced FLUKA Course

# About



**fleɪ(r)** n [U,C] natural or instinctive ability (to do something well, to select or recognize what is best, more useful, etc.  
[Oxford Advanced Dictionary of Current English]

# What is flair [1/2]

## **FLUKA Advanced Interface** [<http://www.fluka.org/flair>]

- **All-in-one** User friendly graphical Interface;
- Minimum requirements on additional software;
- Working in an intermediate level

**Not hiding the inner functionality of FLUKA**

### **Front-End interface:**

- Fully featured **Input file Editor**
  - Mini-dialogs for each card, allows easy and almost error free editing
  - Uniform treatment of all FLUKA cards
  - Card grouping in categories and card filtering
  - Error checking and validation of the input file during editing
- **Geometry:** interactive visualization editing, transformation, optimizations and debugging (tomorrows talk);
- **Compilation** of the FLUKA Executable;
- **Running** and **monitoring** of the status of a/many run(s)

# What is flair [2/2]

## Back-End interface:

- Inspection of the output files (core dumps and directories)
- Output file(s) viewer dividing into sections
- Post processing (merging) the output data files
- Plot generation through an interface with **gnuplot**;

## Other Goodies:

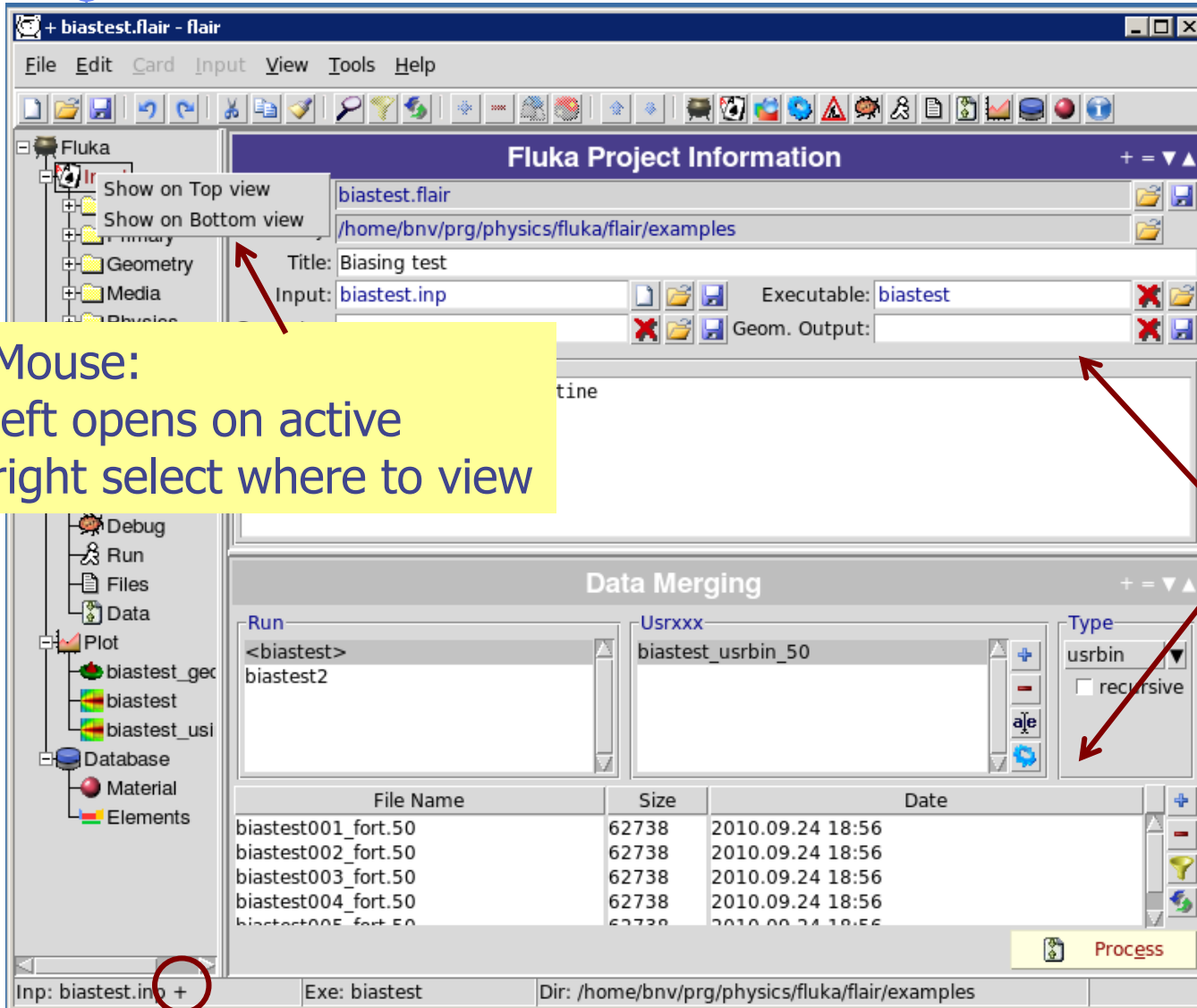
- Access to FLUKA manual as hyper text
- Checking for release updates of FLUKA and flair
- Nuclear wallet cards
- Library of materials
- Database of geometrical objects (Not yet completed)
- Programming python **API**
- Everything is accessible with keyboard shortcuts

# Concepts: Flair Project

- Store in a **single file** all relevant information:
  - Project notes
  - Links to needed files: **input file**, **source routines**, **output files** ...
  - **Multiple runs** from the same input file, as well running status
  - Procedures on how to **run the code**
  - **Rules** on how to perform **data merging**
  - Information on how to post process and **create plots** of the results
- You can consider Flair as an **editor** for the project files.
- Can handle any FLUKA input format (reading & writing), but internally it works using the **names format** for the input, **free with names** for the geometry (Recommended way of working)
- The format is plain ASCII file with extension: **.flair**

**Note:** If you want to copy a project you need to copy also all linked files especially the input and source routines!

# Interface



active

- + vertical/horizontal
- = equalize
- ▼ minimize
- ▲ maximize

Mouse:  
left opens on active  
right select where to view

2 working frames

inactive  
click to activate

input modified and not saved

# Command line options

Usage: flair [options] <filename | filename.flair | filename.inp>

Options:

- -d/D Activate/Deactivate the beta-development features
- -e exe Use exe as fluka executable
- **-g** Open geometry editor window
- -i inputfile Fluka input file (w/o the .inp extension)
- **-r** Load most recent project
- -R # Load recent project (number 1..10 or filename)
- -l List recent projects
- -x Run through an xterm (default)
- **-X** Run without an xterm (**useful in case of start failure**)
- -1 Load the first flair file in the folder

# Interface

## Keyboard:

Almost everything is possible with the keyboard see manual for shortcuts

**Ctrl-Enter:** Execute most important action

**Ins/Del:** Add or Delete

## Mouse:

**right-click** anywhere to get a popup menu

## Listboxes:

all listboxes are searchable. Typing only the characters (A-Z) and numbers (0-9) all other are ignored

## LabelFrames:

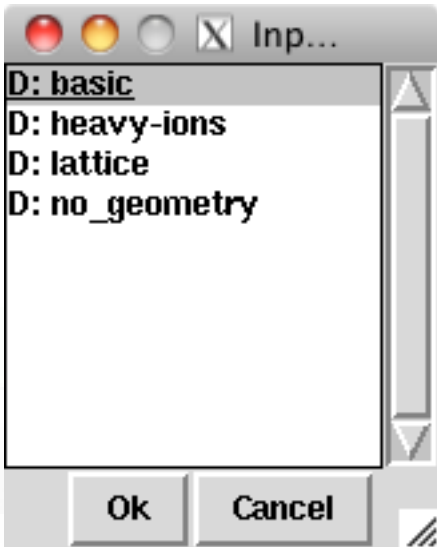
can be collapsed/expanded by clicking on the label



# Input Templates

Default template: **basic.inp**

- When requesting a new input or a new project flair will prompt to select an input template:



```

TITLE
GLOBAL                                1.0      1.0
DEFAULTS
BEAM
BEAMPOS
GEOBEGIN                                COMBNAME
      0      0
* Black body
SPH blkbody      0.0 0.0 0.0 10000000.0
* Void sphere
SPH void         0.0 0.0 0.0 1000000.0
* Cylindrical target
RCC target      0.0 0.0 0.0 0.0 0.0 10.0 5.0
END
* Black hole
BLKBODY      5 +blkbody -void
* Void around
VOID         5 +void -target
* Target
TARGET      5 +target
END
GEOEND
* .+.+.1+.+.2+.+.3+.+.4+.+.5+.+.6+.+.7..
ASSIGNMA      BLCKHOLE      BLKBODY
ASSIGNMA      VACUUM        VOID
ASSIGNMA      COPPER        TARGET
RANDOMIZ      1.0
START
STOP
    
```

- Flair default templates are prefixed with "D:"
- User templates will be prefixed with "U:"

**The user can create his own set of input templates. They are normal FLUKA input files and they have to be placed in the directory `~/flair/templates` (create the directory if not existing)**

# Card Categories

For easier access, cards are groups in the following categories:

- **General** General purpose (TITLE, DEFAULTS, GLOBAL...);
- **Primary** Definition of the primary starting particles;
- **Geometry** Cards related to the definition of the geometry bodies/regions/lattices plotting and rotations/translations;
  - **Bodies** Subcategory containing only the bodies definition;
  - **Transformations** Subcategory containing only the geometrical directives;
- **Media** Definition and assignment of materials;
- **Physics** Setting physics properties of the simulation;
- **Transport** Modify the way particles are transported in FLUKA;
- **Biasing** Cards for importance biasing definition;
- **Scoring** Cards related to scoring;
- **Flair** flair special cards;
- **Preprocessor** Definitions for creating conditional input files.

# Concepts: Extended Cards [1/2]

- Flair is treating the input file as a **list of extended cards**;
- Each extended card contains:
  - **Comment**: All commented lines preceding the card(s) as well the inline comments;
  - **Tag**: The 8 character word identifying the card. All tags not recognized by flair will be converted to **#error**;
  - **WHATs**: Multiple number of **WHATs** (0=sdum, 1-6 first line, 7-12 continuation line...)
  - **Extra**: multi line string of extra information for special cards like **REGION, TITLE, PLOTGEOM** etc.
  - **State** (Enable/Disable);
- Flair recognize automatically (and separates them from the comments) all the disabled valid FLUKA cards;

# Concepts: Extended Cards [2/2]

- The region definition in the in geometry is emphasized by the presence of a card named "REGION";
- All the COMPOUND cards related to one material are joined in one card;
- Cards are edited with the flair editor through the use of the mini-dialogs, forcing the user to enter the correct information (default);
- The user can nevertheless gain full control of the card using the Edit dialog (*Ctrl-E*);
- Flair will try to find the best floating point representation of each number, to ensure the maximum accuracy; number of digits that fits in the specific width (10 for the fixed format, 22 for the free format).
- Function evaluation: a field value starting with = will force flair to evaluate its content as a function e.g.

BEAMPOS    x: =2\*10+length

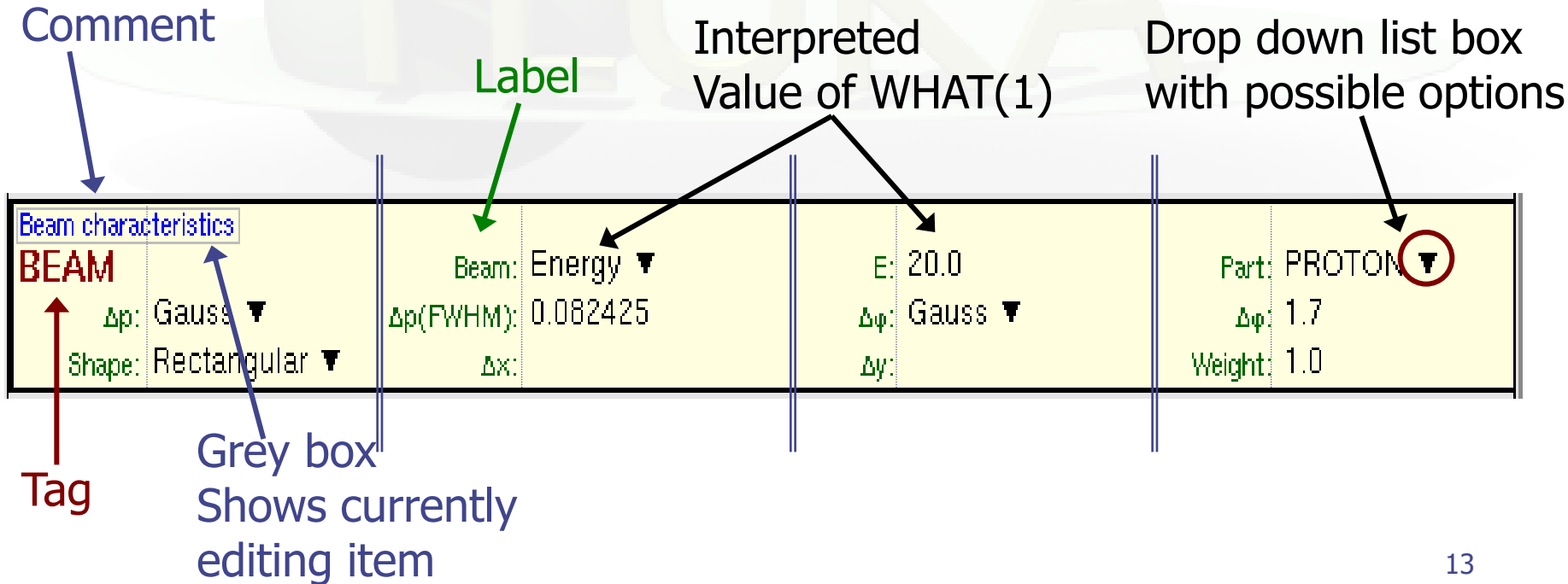
Flair will create a valid fluka input containing the evaluation of the formula and keep the formula inside the comments as

\*@what.1 =2\*10+length

# Anatomy of a card mini-dialog [1/2]

- For each extended card flair has a mini dialog (currently in 4 columns), interpreting all information stored in the card

```
* Beam characteristics
BEAM          -20.0 -0.082425      -1.7          1.0PROTON
```



# Anatomy of a card mini-dialog [2/2]

\* Energy deposition in 3D binning

```

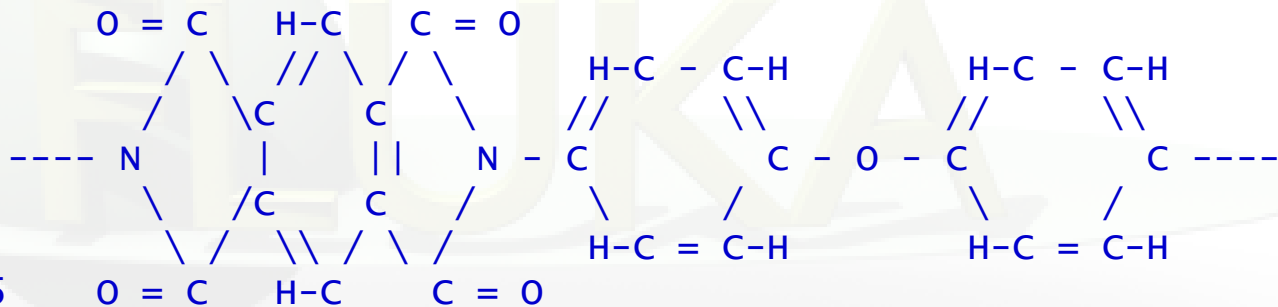
USRBIN          10.0      ENERGY      -50.0        45.0        54.0        36.0EneDep
USRBIN          -45.0     -54.0     -33.0        100.0       100.0       100.0&
    
```

<b>USRBIN</b>	Unit: 50 BIN ▼	Name: EneDep
Type: X-Y-Z ▼	Xmin: -45.0	NX: 100.0
Part: ENERGY ▼	Ymin: -54.0	NY: 100.0
	Zmin: -33.0	NZ: 100.0
	Xmax: 45.0	
	Ymax: 54.0	
	Zmax: 36.0	

\* Polypyromellitimide Polyimide, Kapton

\* Chemical

\* Formula



```

MATERIAL          1.43
COMPOUND          10.0  HYDROGEN          22.0  CARBON          2.0  NITROGEN
COMPOUND          5.0  OXYGEN
    
```

<b>MATERIAL</b>	Name: Polyimid	#	p: 1.43
Z:	Am:	A:	dE/dx: ▼

<b>COMPOUND</b>	Name: Polyimid ▼	Mix: Atom ▼	Elements: 6 ▼
f1: 10.0	M1: HYDROGEN ▼	f2: 22.0	M2: CARBON ▼
f3: 2.0	M3: NITROGEN ▼	f4: 5.0	M4: OXYGEN ▼
f5:	M5: ▼	f6:	M6: ▼

# Input Editor - 1

```

#define BIAS
TITLE Biasing test
GLOBAL Max #reg: Analogue: DNear:
           Input: Names Geometry: Free
DEFAULTS NEW-DEFA
BEAM Beam: Energy E: 0.005 Part: NEUTRON
        Δp: Flat Δp: Δφ: Isotropic
        Shape: Rectangular Δx: Δy: Weight:
BEAMPOS x: y: z:
           cosx: cosy: Type: POSITIVE
GEOBEGIN Log: Acc: Opt:
            Inp: Out: Fmt: COMBNAME
            Title:
Black body
SPH blkbody x: 0.0 y: 0.0 z: 10
           R: 10000000.0
Void sphere
SPH void x: 0.0 y: 0.0 z: 10
           R: 1000000.0
Cylindrical target
RPP target Xmin: -100. Xmax: 100.
           Ymin: -100. Ymax: 100.
           Zmin: -100. Zmax: 100.
Black hole
REGION BLKBODY Neigh: 5 Volume:
        Expr: +blkbody -void
Void around
REGION VOID Neigh: 5 Volume:
        Expr: +void -target
*.....1.....2.....3.....4.....5.....6.....7.....
SPH blkbody 0.0 0.0 10. 10 00000.0
    
```

highlight differences during editing

# Input Editor - 2

- Drag'n'drop from the TAG of the cards
- Double click on card TAG to select all similar cards
- Editing multiple cards: select cards and modifying the value in one card will propagate the change to all similar selected cards
- Ctrl-Double-Click Show/Hide selected cards
- #if..#endif, \$transform, \$translat or \$expand flair will enclose the selected cards with the #if #endif, or \$start\_xxx, \$end\_xxx transformation cards
- Popup Balloon tooltip displays short help:
  - for every option on every card
  - body description in the REGION expression
- Right-click: shows popup-menu
  - Quick filtering by REGION, MATERIAL, scoring etc...
- Easter Eggs: AWARI by Double-Right-Click on dialog showing the card representation as text at the bottom of the screen



# Input Editor - 3

- Automatic indentation of nested `#if..#endif` and `$start..$end` directives.
- To refresh the display press **Ctrl-R**
- Each **REGION** can be split into many cards if needed to be used with preprocessor commands.
- Use as a name **"&"**

Void around

**REGION VOID**

Neigh: 5

Volume:

Expr: +void -target

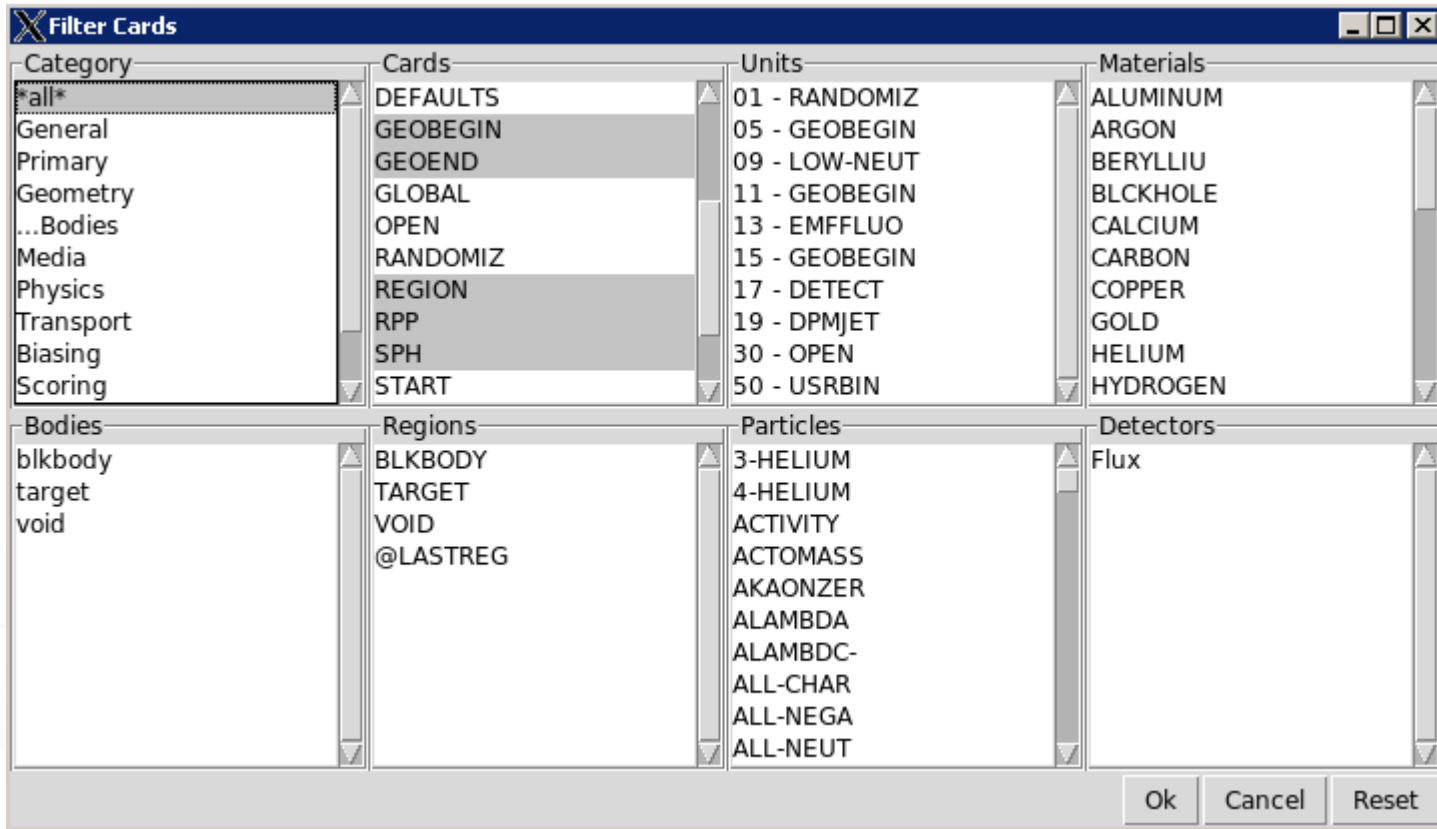
**#if** BIAS ▼

**REGION &**

cont: -bias

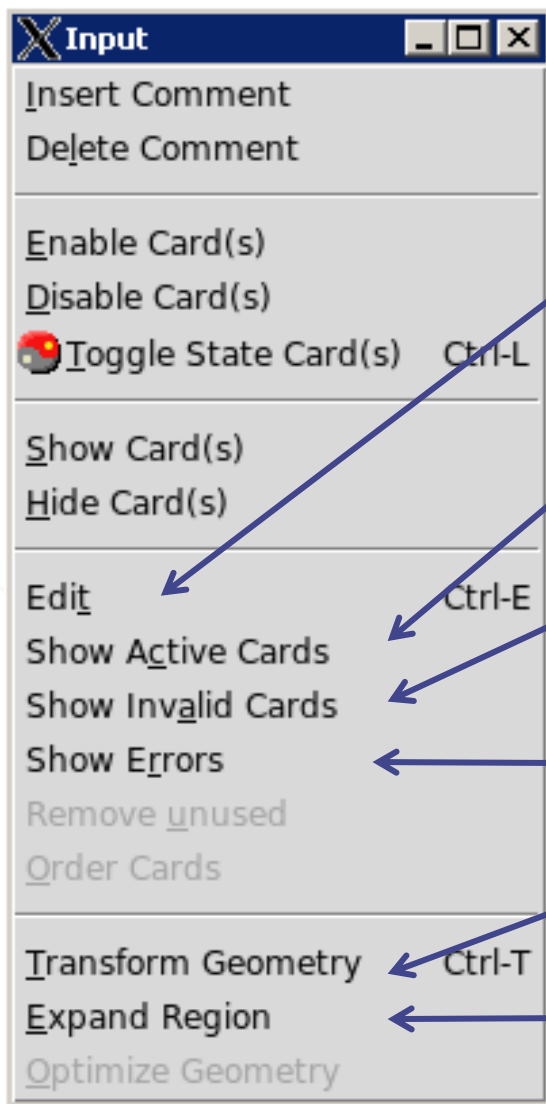
**#endif**

# Input Card Filtering



- Filter Cards dialog allows a more advanced selection of cards to be displaced, by showing only the cards that match the selected options

# Input Menu



Manual editing of the card

Scan input and display only active cards (not excluded by the preprocessor)

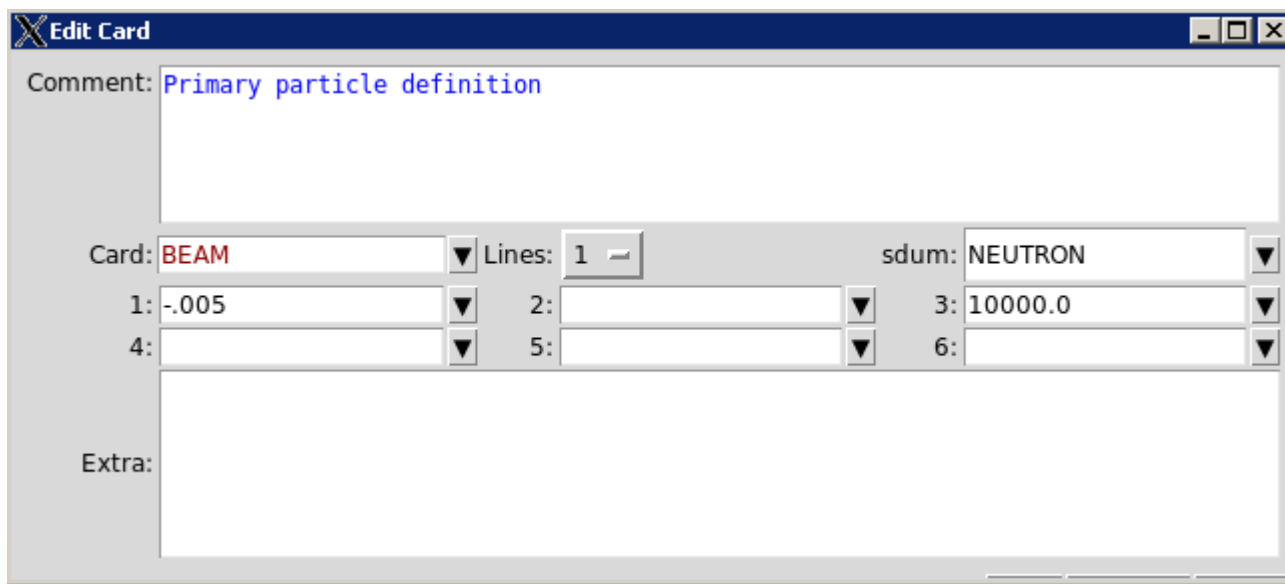
Show cards containing problems/errors

Display a message with the errors identified in the cards

Open the geometry transformation dialog

Expand parenthesis in the region (only logical optimization will performed)

# Manual Card Editing



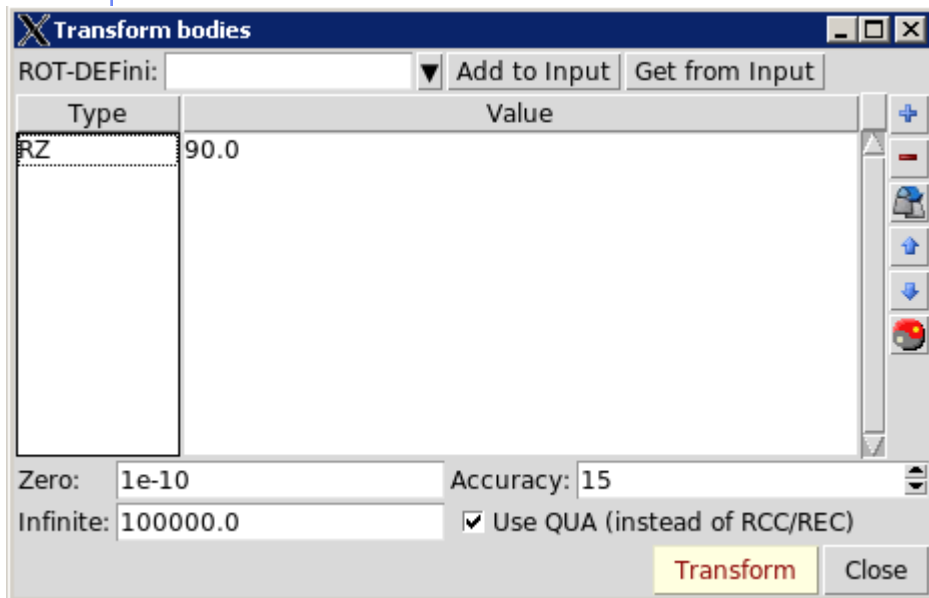
Accessible: **Ctrl-E** or **right-click** → **Edit, Menu** → **Input** → **Edit**

Lines: Number of lines the card extends

Extra: additional information for a card like title string for TITLE, or region expression for REGION

Dropdown box: shows with categories all items defined in the input (bodies, regions, materials, particles...)

# Bodies Transformation



## Transformation Types:

- T translate along a vector
- TX TY TZ translate along axis
- RX RY RZ **axis** rotation (degrees)
- S scaling

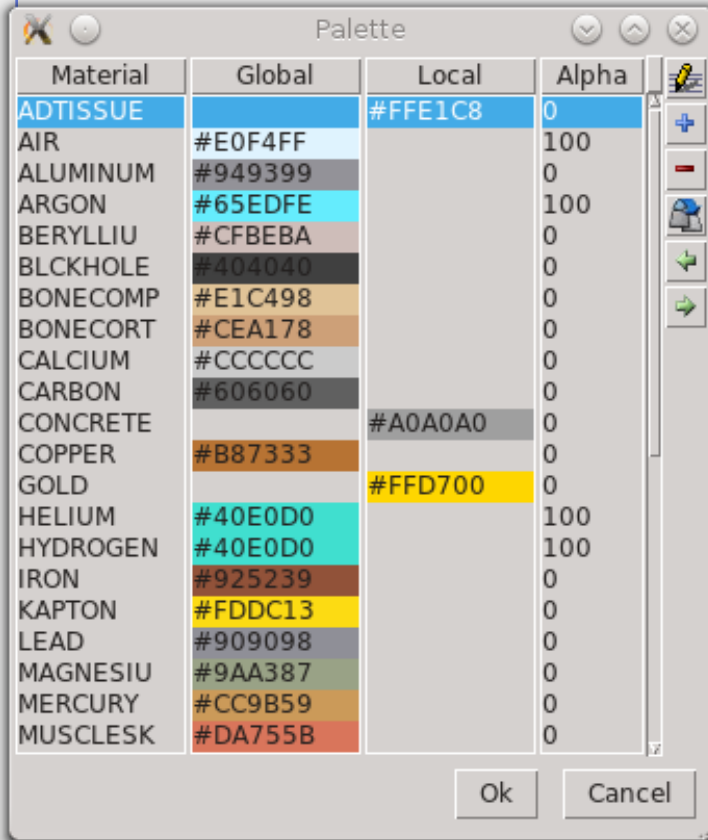
- Applies a user transformation to the selected bodies on the input editor.
- Convert transformations to/from **ROT-DEFini** cards
- **Zero**: limit below which to be considered as zero
- **Accuracy**: Numeric digits
- **Infinite**: infinite bodies when converted to which size to use
- Use **QUA**: convert infinite cylinders to infinite QUAdrics

## Remember:



When transforming bodies for use with **LATTICE** card, use the maximum precision

# Color Palette



Accessible: Menu → View → Palette

- Edit colors used for material display in **Geometry plots** and **GeometryEditor**
- **Global** colors are saved inside flair.ini and are shared between all projects
- **Local** colors are initially randomly assigned and saved inside the project file
- **Alpha** channel for setting transparency on materials. Used on the 3D raytracing plotting.  
0 = opaque  
100 = transparent

# Compiling

Compile Executable

File	Size	Date
usimbs.f	4856	2010.09.24 18:57
opncrd.f	5751	2010.09.24 18:40
oauxfi.f	6059	2010.09.24 18:49

automatic selecting needed routines from usermvax/

FLUKA User routines

File	Size	Date	Desc
pshckp.f	1274	2005.06.02 13:16	
queffc.f	1605	2005.03.24 10:40	quantum efficiency (for optical photons)
rflctv.f	1469	2005.03.24 10:40	reflectivity (for optical photons)
rfrndx.f	1469	2005.03.24 10:40	refraction index (for optical photons)
soevsv.f	2507	2005.06.17 16:13	saving source events
source.f	7327	2009.09.09 16:08	to generate any distribution for source particles
stupre.f	4223	2005.03.24 10:40	set user variables (electrons and photons)
stuprf.f	1981	2005.07.25 13:43	set user variables (hadrons, muons and neutrinos)
ubsset.f	5585	2005.03.24 10:40	to override input biasing parameters
udcdrl.f	2425	2005.03.24 10:40	decay direction biasing
usimbs.f	3262	2008.10.30 11:56	user-defined importance biasing
usrein.f	1553	2005.03.24 10:40	event initialisation
usreou.f	1480	2005.03.24 10:40	post-event output

Link: lfluka Exe: biastest  Default main:  D Line  Bound Check

Options:

**Build** Compile Clean

## Filetypes accepted:

- Fortran: .f, .F, .for, .FOR
- C/C++: .c, cpp, .cxx, .cc
- Libraries: .a, .so

Automatic scanning of necessary user routines and copying them to project folder.

**Build:** behaves like a "makefile" compiles based on files timestamp when are newer

**Compile:** Forces compile of the selected files

**Clean:** cleanup of all produced files

When you are unsure, click on "Clean" before "Build"

# Running

Run / Input

- <rdsourc>
- rdsourc\_first
- rdsourc\_second
- rdsourcPrecision\_first
- rdsourcPrecision\_secon

Override Options

Title

Primaries 0 Rnd 0

Time 0 Exe rdsourc

Defines

Sel	Name	Value
[ ]	FIRST	
[X]	SECOND	
[X]	ANALOGUE	
[X]	ENERGY	=10*MeV

Custom preprocessor #defines even with value/function

Cycles: Continue Previous 0 No. Cycles 5 Last 5

Run Stop Cycle Stop Run Kill Attach Refresh Queue \*Default

Progress

Status: Finished OK Input: rdsourcPrecision\_second Dir:

Started: ETA: Time/prim:

Elapsed: Cycle: Run:

Cycles:

Primaries:

<inputname> refers to the input file AS IT IS in the input editor.

Create additional runs based on the same input file by overriding:

- Title
- Preprocessor definitions
- Random number seed
- Starting particles
- Execution timeout
- Executable

- Monitors the status of the run by inspecting the FLUKA output files. If **timeout** occurs try to re-**Attach** to the running process.
- The timeout is user-definable in the **Preferences** dialog



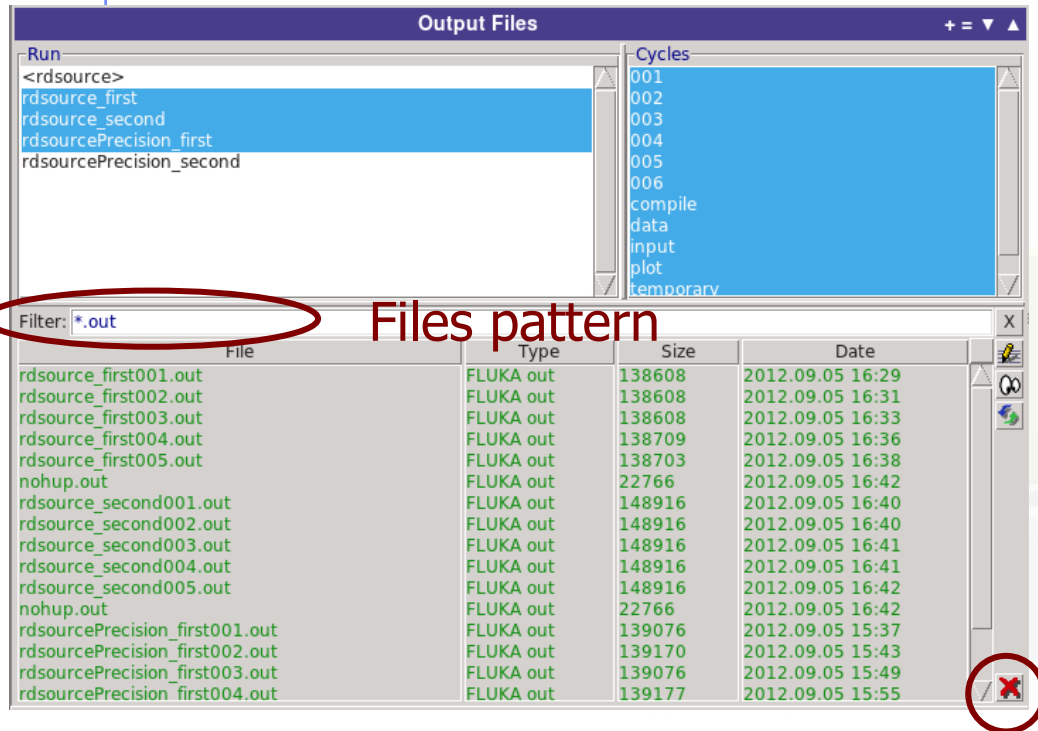
# Running: How to use multicore CPU's

- Create clones of the current input e.g. **test.inp** named: test**1**.inp, test**2**.inp, test**3**.inp ...
- Assign a **different random number seed** on each run (Rnd entry)
- Select all in the listbox and click Run

## Multiple Selection:

- To modify **many runs** at the same time, select them in the listbox
- The options will be "*disabled*"
- **Right-click** on the options you want to **enable** and modify them
- Modify the filters in Data processing for summing up all cycles from all runs (see later)

# Output Files



Files pattern

Delete selected files

Inspect Output files generated by FLUKA classified per:

Run/Cycle

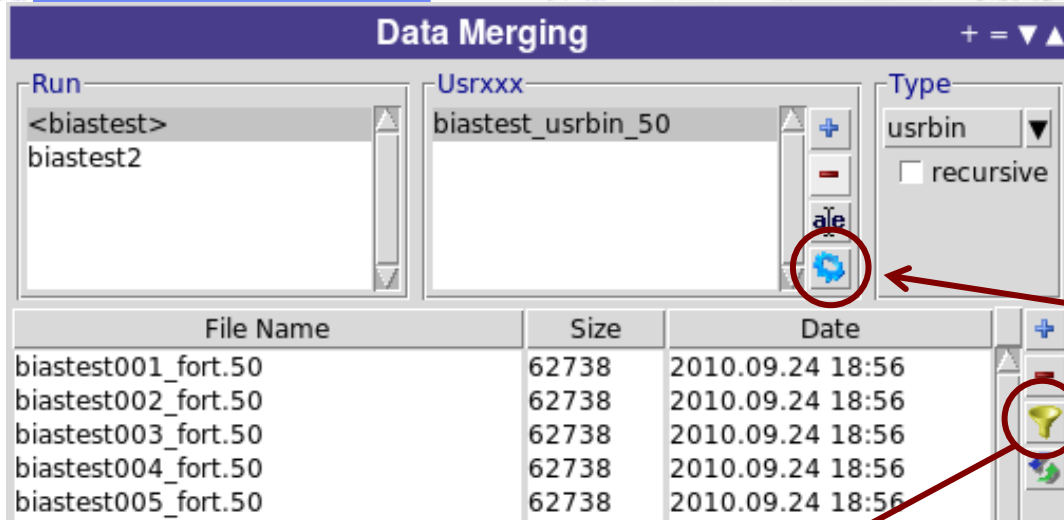
As well special output files from compilation data processing plotting and temporary

Double clicking opens:

- Files in the file Viewer
- coredumps in debugger

Right click can convert USBIN's from formatted to unformatted

# Data Processing



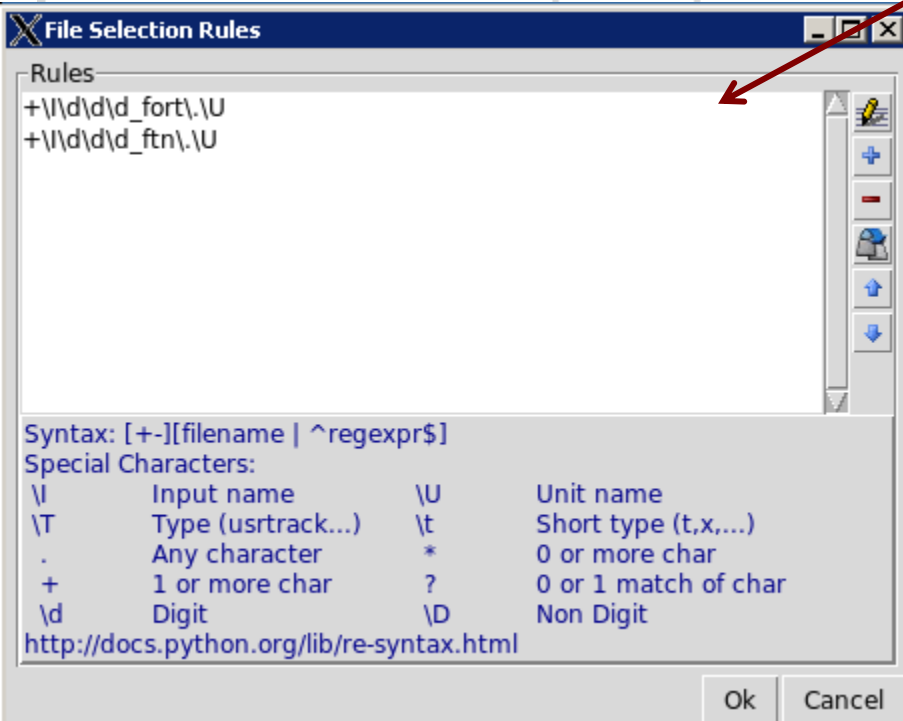
Process all scoring BINARY output files for each Run.

Name rules are defined in Preferences

Automatically scan input for scoring cards

+/- Modify file list by adding / removing items

Dialog for editing scanning rules for files.



Use the rules to merge from multiple runs. e.g. add a `\d` in the target like `+\\I\\d\\d\\d\\d_fort\\.\\U`

To modify the rules for multiple scoring cards, select all Usrxxx before

The default rules can be modified in the Preferences Dialog

# Plot List

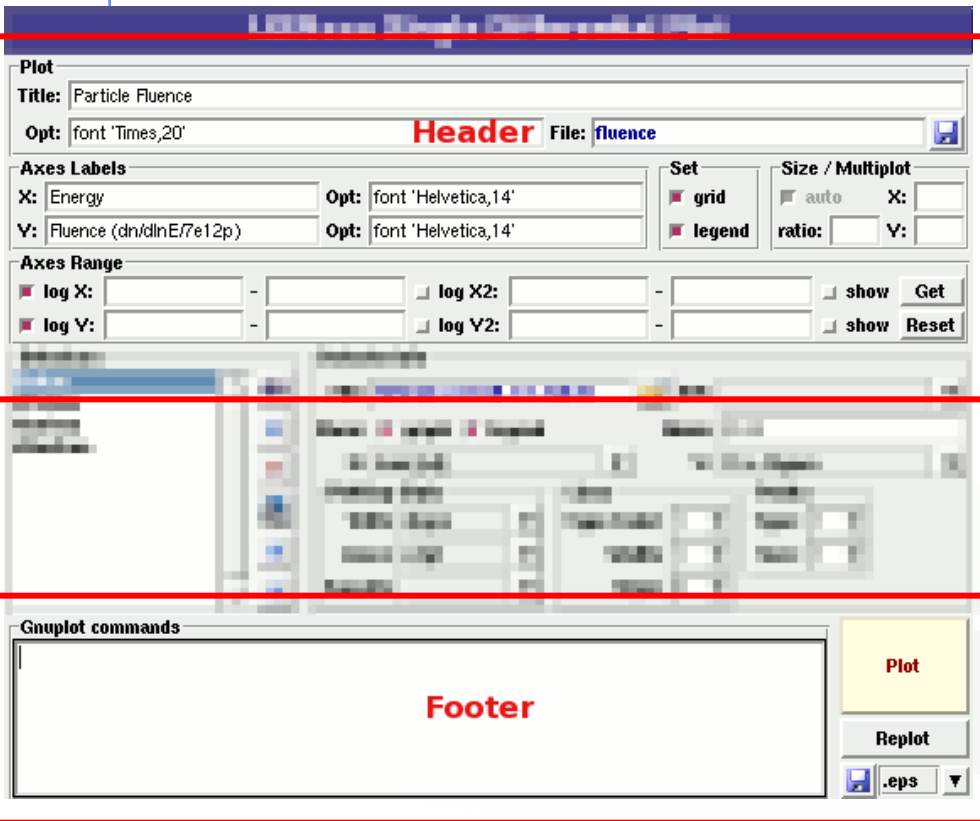
File	Title	Type
geometry	nTOF Target Geometry	Geometry
enedep	Deposited Energy	USRBIN
Fluence	Particle Fluence	USR-1D
resnuc	Residual Nuclei	RESNUCLE

- 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 (the extension will change)
- The Filter button creates automatically one plot for each processed unit (From the default input file)
- Hit Enter or click the Edit icon to display the plotting dialog
- Fast Double click on item to open the corresponding dialog
- Slow Double click to modify the value

## 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 (ToDo)

# Plotting Frames



- 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.
- **Re-Plot** will force the creation of all files regardless their state
- 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'



Look in the flair manual for a short reference of gnuplot commands

# General Tips

- In the Configuration Dialog you can set global commands to execute before or after any plot
- The **output window** displays all the commands that are sent to gnuplot. As well as the errors. In case of problem always consult the output window!
- In the **Gnuplot commands** you can fully customize the plot by adding manually gnuplot commands:
- Special commands:
  - **plot, splot** with no options, defines the order where flair should insert the plot or splot command.
  - **replot <plot-cmd>** append extra plots to the one generated by flair

# USRBIN Plots - 1

**Binning Detector**

File:  Title:

Cycles:  Primaries:  Weight:  Time:

**Binning Info**

Det:  X:  Min:

Type:  Y:  Max:

Score:  Z:  Int:

**Projection & Limits**

Type:

<input type="radio"/> X: <input type="text" value="-36"/>	<input type="text" value="1"/>	<input type="text" value="25.6"/>	<input type="text" value="Get"/>
<input type="radio"/> Y:	<input type="text" value="1"/>		<input type="checkbox"/> swap
<input checked="" type="radio"/> Z:	<input type="text" value="1"/>		<input checked="" type="checkbox"/> errors
Norm:			<input checked="" type="checkbox"/> log

Color Band: Min:  Max:

CPD:  Colors:

Palette:   Round

Geometry: Use:  Pos:  Axes:

**Gnuplot commands**

Rebinning

Swap axes

Get limits from gnuplot  
using right-mouse

Draw errors. (combined with log)  
Correct only if one slice is used

# USRBIN Plots - 2

**Binning Detector**

File:  Title:

Cycles:  Primaries:  Weight:  Time: \*\*\*\*\* Sum file \*\*\*\*\*

**Binning Info**

Det:  X:  Min:

Type:  Y:  Max:

Score:  Z:  Int:

**Projection & Limits**

<input type="radio"/> X: <input type="text" value="-36"/>	<input type="text" value="1"/>	<input type="text" value="25.6"/>	<input type="button" value="Get"/>
<input type="radio"/> Y: <input type="text" value=""/>	<input type="text" value="1"/>	<input type="text" value=""/>	<input type="checkbox"/> swap
<input checked="" type="radio"/> Z: <input type="text" value=""/>	<input type="text" value="1"/>	<input type="text" value=""/>	<input checked="" type="checkbox"/> errors

Norm:

log

**Type:**

**Color Band**

Min:  Max:

CPD:  Colors:

Palette:   Round

**Geometry**

Use:

Pos:

Axes:

**Gnuplot commands**

Normalization could be used as:

- number or expression evaluating in a number  $65e-3/2.7$
- function with x as variable. e.g  $E2T(x*65e-3/2.7)-293$   
with the function defined in the Gnuplot commands  
 $E2T(x) = ((3.00629e-08*x-0.000108436)*x+1.01097)*x+311.839$



# USRBIN Plots - 3

**Binning Detector**

File:  Title:

Cycles:  Primaries:  Weight:  Time: \*\*\*\*\* Sum file \*\*\*\*\*

**Binning Info**

Det:	<input type="text" value="1 EneDep"/>	X: <input type="text" value="[-40 .. 40] x 100 (0.8)"/>	Min: <input type="text" value="1.95034673E-07"/>
Type:	<input type="text" value="10: X-Y-Z"/>	Y: <input type="text" value="[-40 .. 40] x 100 (0.8)"/>	Max: <input type="text" value="0.0254351143"/>
Score:	<input type="text" value="ENERGY"/>	Z: <input type="text" value="[-30 .. 35] x 100 (0.65)"/>	Int: <input type="text" value="11.0419018"/>

**Projection & Limits**

<input type="radio"/> X:	<input type="text" value="-36"/>	<input type="text" value="1"/>	<input type="text" value="25.6"/>	<input type="button" value="Get"/>
<input type="radio"/> Y:	<input type="text" value=""/>	<input type="text" value="1"/>	<input type="text" value=""/>	<input type="checkbox"/> swap
<input checked="" type="radio"/> Z:	<input type="text" value=""/>	<input type="text" value="1"/>	<input type="text" value=""/>	<input checked="" type="checkbox"/> errors

Norm:  log

Type:

**Color Band**

Min:  Max:

CPD:  Colors:

Palette:   Round

**Geometry**

Use:

Pos:

Axes:

**Gnuplot commands**

Normalization could be plotted:

- 2D projection, 1D projection
- Trace of the maximum
- Full width at half maximum

# USRBIN Plots - 4

**Binning Detector**

File:  Title:

Cycles:  Primaries:  Weight:  Time: \*\*\*\*\* Sum file \*\*\*\*\*

**Binning Info**

Det: <input type="text" value="1 EneDep"/>	X: <input type="text" value="[-40 .. 40] x 100 (0.8)"/>	Min: <input type="text" value="1.95034673E-07"/>
Type: <input type="text" value="10: X-Y-Z"/>	Y: <input type="text" value="[-40 .. 40] x 100 (0.8)"/>	Max: <input type="text" value="0.0254351143"/>
Score: <input type="text" value="ENERGY"/>	Z: <input type="text" value="[-30 .. 35] x 100 (0.65)"/>	Int: <input type="text" value="11.0419018"/>

**Projection & Limits**

<input type="radio"/> X: <input type="text" value="-36"/>	<input type="text" value="1"/>	<input type="text" value="25.6"/>	<input type="button" value="Get"/>
<input type="radio"/> Y: <input type="text" value=""/>	<input type="text" value="1"/>	<input type="text" value=""/>	<input type="checkbox"/> swap
<input checked="" type="radio"/> Z: <input type="text" value=""/>	<input type="text" value="1"/>	<input type="text" value=""/>	<input checked="" type="checkbox"/> errors

Norm:   log

Type:

**Color Band**

Min:  Max:

CPD:  Colors:

Palette:   Round

**Geometry**

Use:

Pos:

Axes:

**Gnuplot commands**

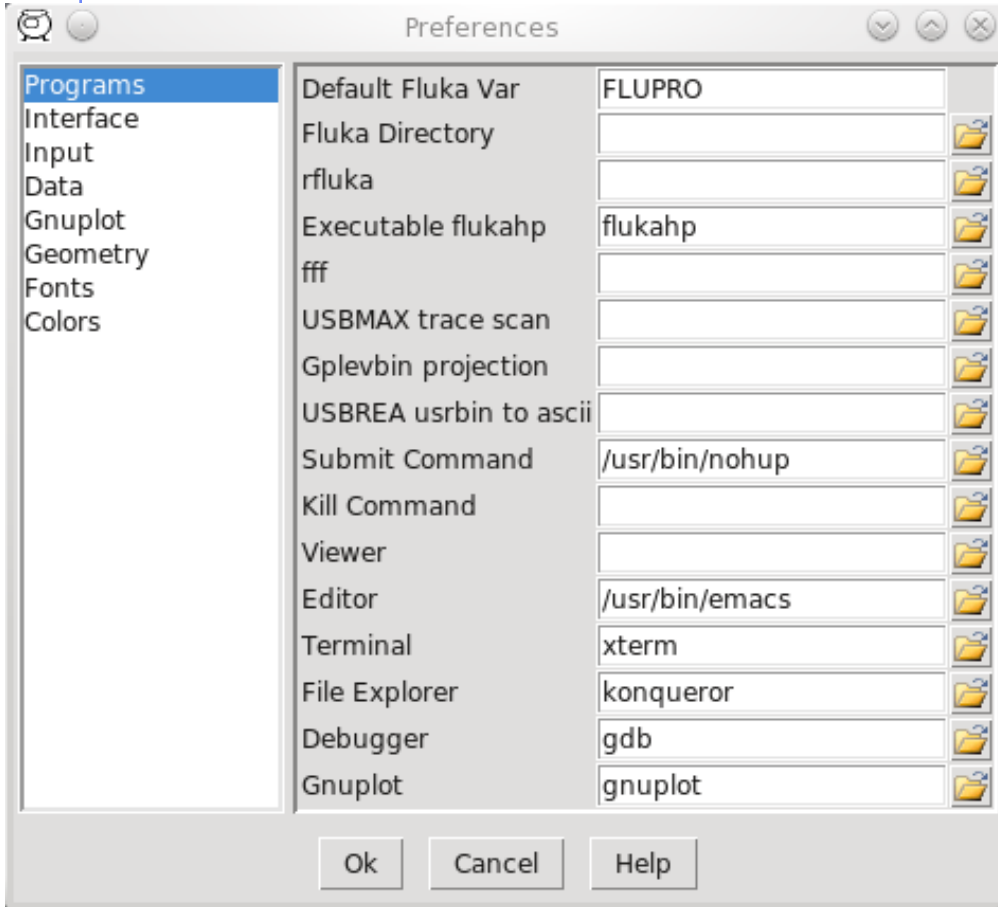
Geometry plot overlay (useful for LATTICE's):

**-Auto-** generates automatically from FLUKA a geometry at the middle position of the projection

otherwise you can use **any existing geometry plot** from the drop down list.

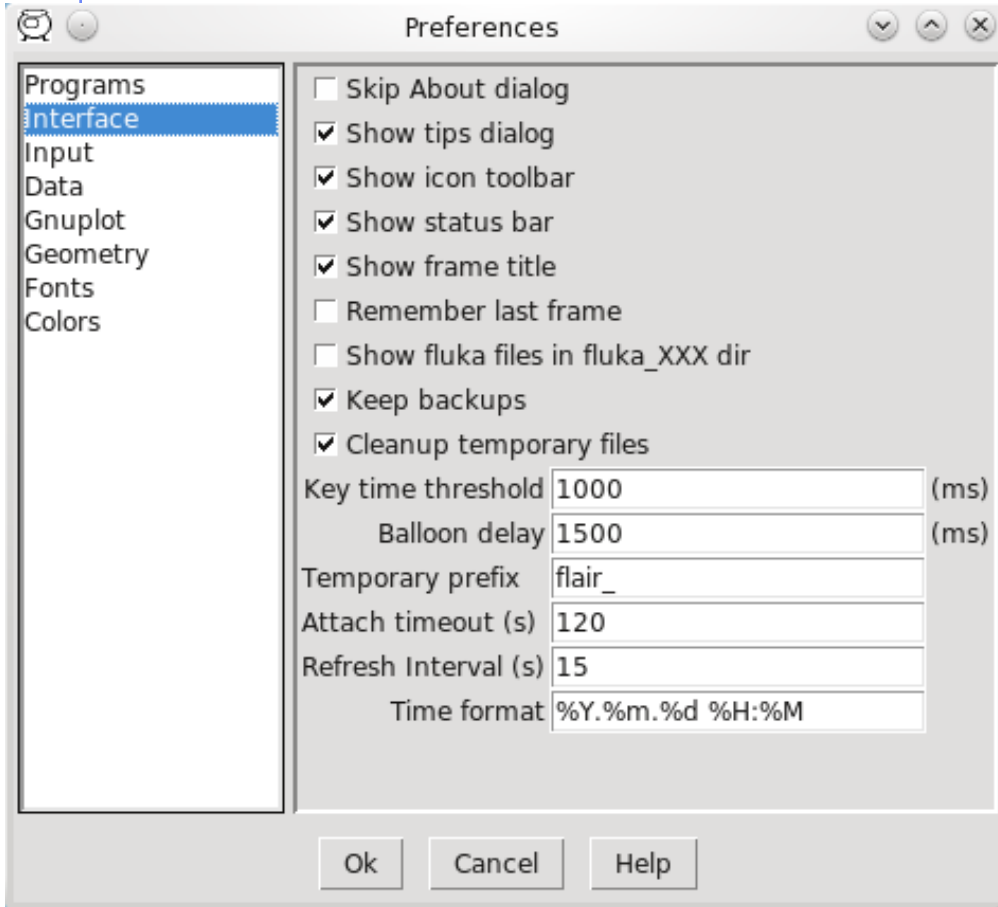
Be careful to properly match the axes that you are using

# Configuration Dialog: Programs



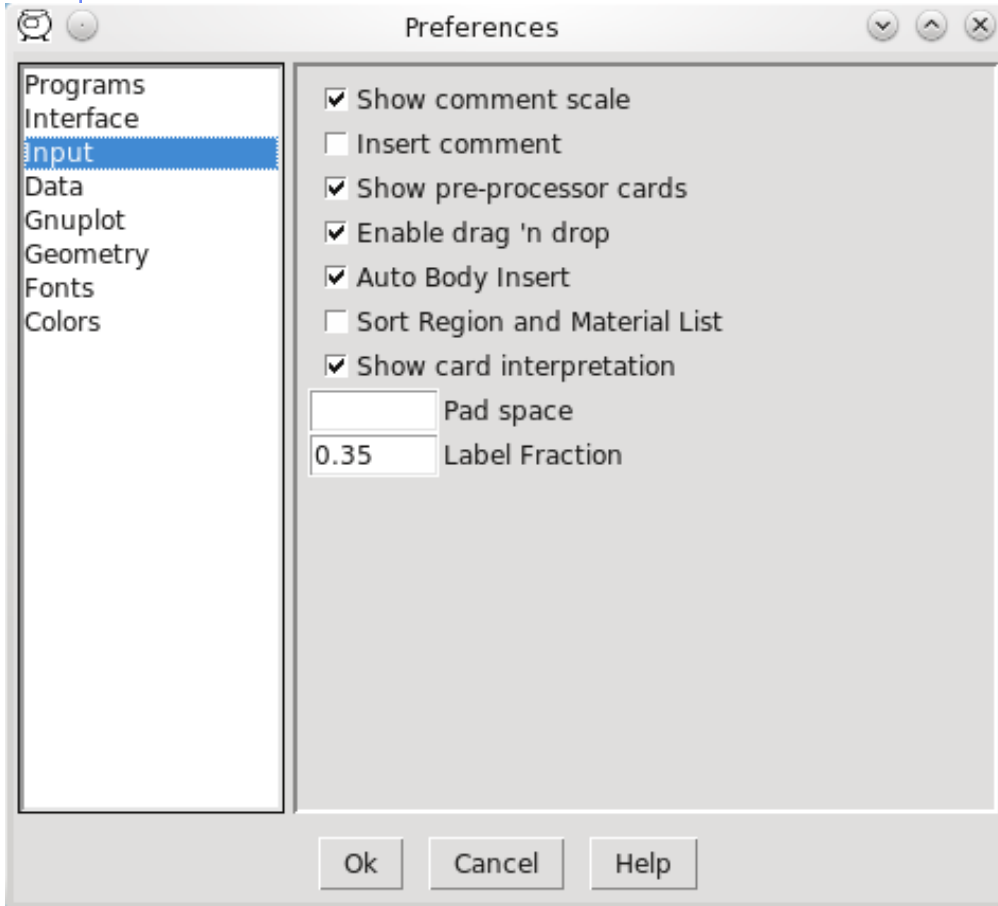
- Set FLUKA directory
- Override default programs to use
- Processing programs are in the "Data" section

# Configuration Dialog: Interface



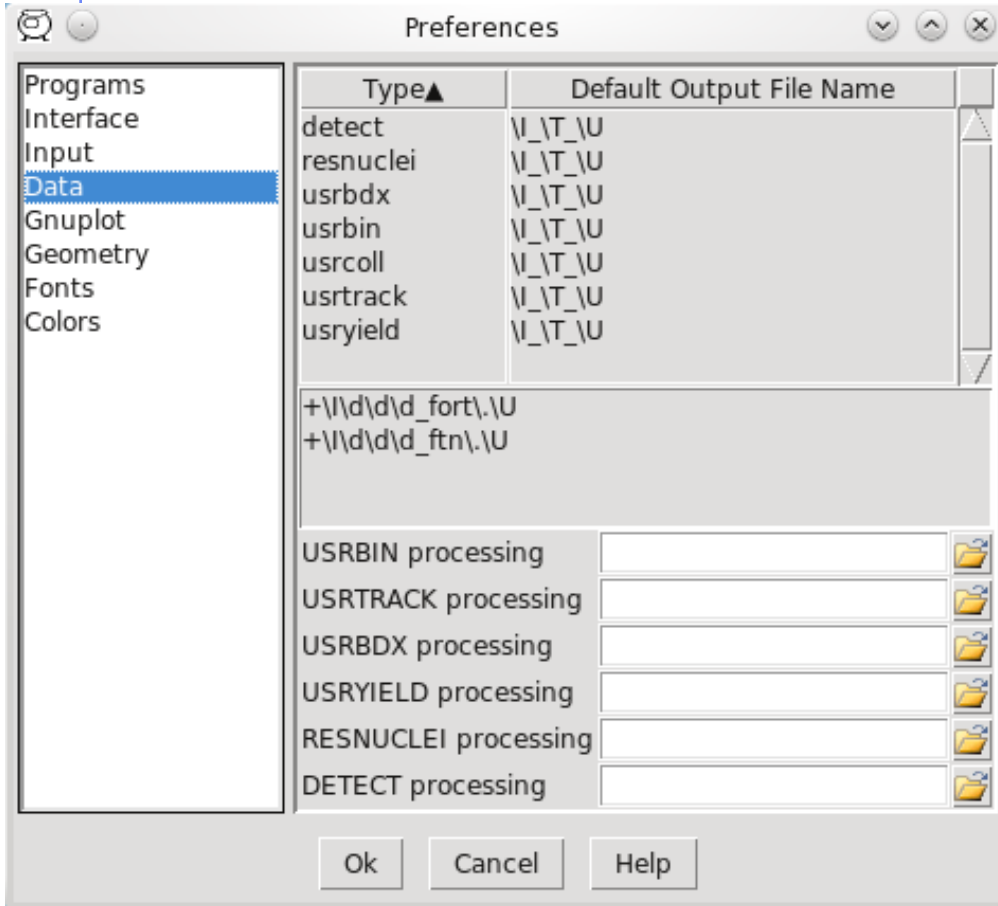
- General interface settings
- Keep backups when files are saved as (file~)
- Automatically Cleanup temporary files. Disable only if you want to inspect files after Debug or Plot when an error occurs
- Key time to reset the type-in search in listboxes
- Balloon delay time
- Time format for files (follows python&C syntax)
- Time out to attach to a running simulation
- Automatic refresh interval of information

# Configuration Dialog: Input Editor



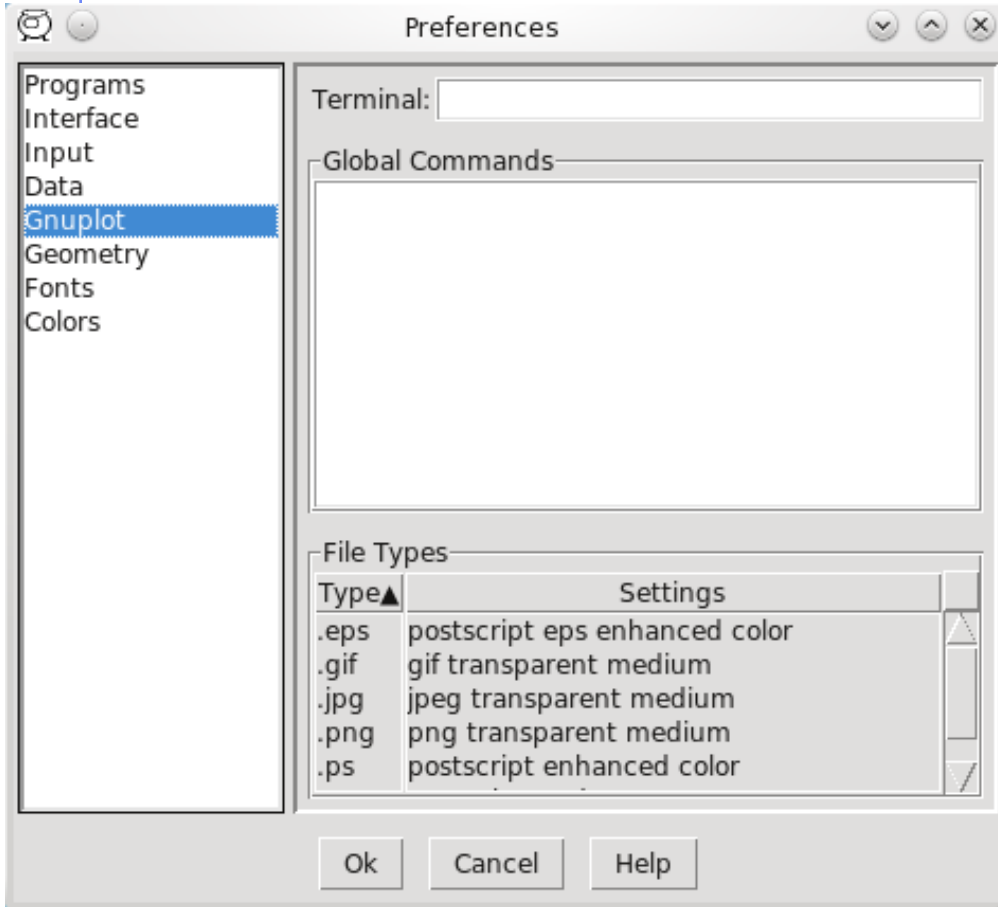
- Show alignment scale
- Automatically insert comment
- Always display preprocessor cards
- Enable drag'n'drop
- Automatic body insertion while editing the region expression
- Sort the region and material list
- Display card interpretation at the bottom of the screen

# Configuration Dialog: Data



- Define how to generate the automatic filenames
- \I will be replaced by input  
\T by card name  
\t by card character
- |           |   |
|-----------|---|
| usrbdx    | x |
| usrbin    | b |
| usrcoll   | c |
| usrtrack  | t |
| usryield  | y |
| resnuclei | r |
- \U the abs(unit-number)

# Configuration Dialog: Gnuplot



## Terminal:

additional options to supply to default terminal

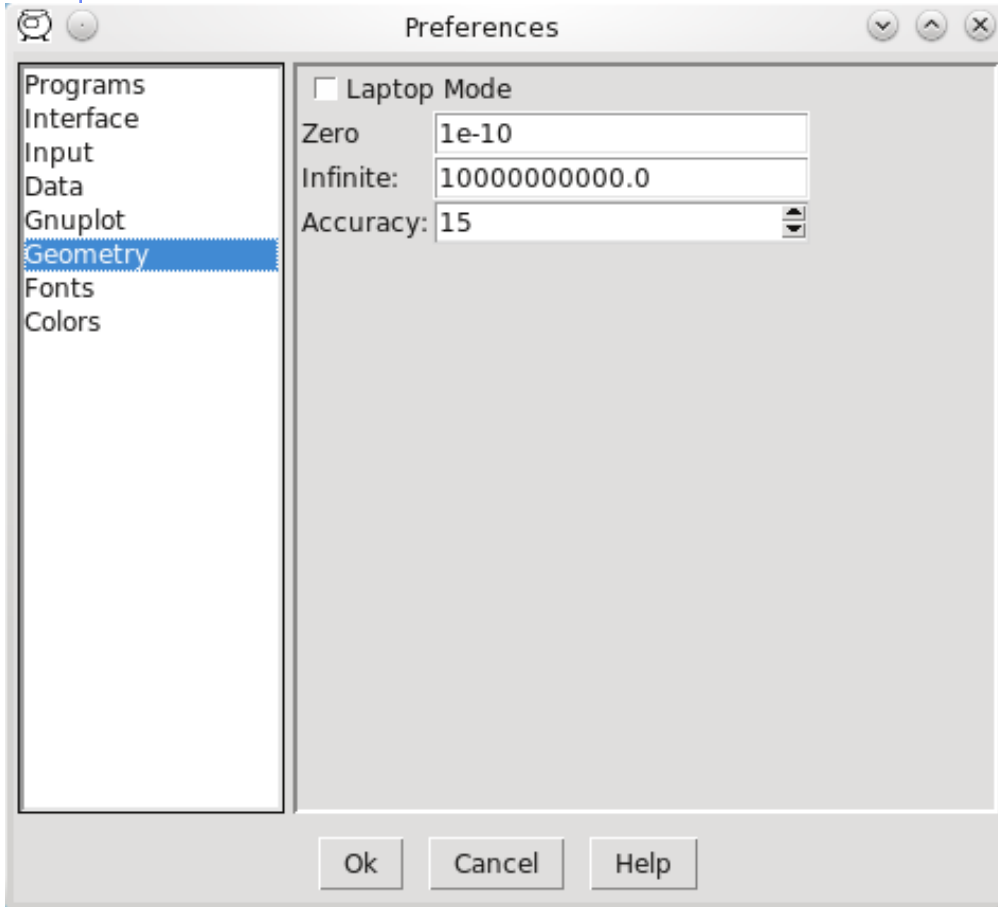
## Global Commands:

gnuplot commands to be executed before any plot

## File Types:

Right-click: to Add/Delete/Modify file types.

# Configuration Dialog: Geometry



## Laptop Mode:

check to swap middle with right mouse buttons. Middle button is used in GeometryEditor for panning, zooming, rotating etc...

## Zero:

## Infinite:

## Accuracy:

same as in the Bodies Transformation dialog



# Materials Database

Material Database

Search:

Group

- Biological
- Elements
- General
- ICRU
- Implantation
- Liquids / Gases
- Metal Alloys
- Plastics / Polymers
- Targets

Material List

Material	Density	Stoichiometry
Mercury	13.546	Hg
728 Cyclohexanone	0.9478	H-10, C-6, O-1
Skeletal Muscle (W&W type 1)	1.05	H-10.1, C-17.1,
Lead	11.35	Pb
Thallium	11.72	Tl
Cyclobutane	0.00125	H-8, C-4
1-Chlorobutane	0.8862	H-9, C-4, Cl-1
Sodium nitrate Na_N_O3	2.261	N-16.5, O-56.5,

Material Properties

Title: Mercury

Notes:

Names: MERCURY

Stoichiometry Properties

Composition: mass liquid 13.546 Group: Elements

Z	A	El	Name	Frac
80		Hg	Mercury	1.0

search database

insert material to input  
add/del material  
edit material

add names to be used  
by FLUKA

Modify Stoichiometry  
and properties of material

**WARNING:** When modifying the database a local copy will be created in `~/flair` folder!!!

# Periodic Table

**Table of Elements**

Table	List																		
Group	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	
Period																			
1	1 H																	2 He	
2	3 Li	4 Be											5 B	6 C	7 N	8 O	9 F	10 Ne	
3	11 Na	12 Mg											13 Al	14 Si	15 P	16 S	17 Cl	18 Ar	
4	19 K	20 Ca	21 Sc	22 Ti	23 V	24 Cr	25 Mn	26 Fe	27 Co	28 Ni	29 Cu	30 Zn	31 Ga	32 Ge	33 As	34 Se	35 Br	36 Kr	
5	37 Rb	38 Sr	39 Y	40 Zr	41 Nb	42 Mo	43 Tc	44 Ru	45 Rh	46 Pd	47 Ag	48 Cd	49 In	50 Sn	51 Sb	52 Te	53 I	54 Xe	
6	55 Cs	56 Ba	* 71 Lu	72 Hf	73 Ta	74 W	75 Re	76 Os	77 Ir	78 Pt	79 Au	80 Hg	81 Tl	82 Pb	83 Bi	84 Po	85 At	86 Rn	
7	87 Fr	88 Ra	** 103 Lr	104 Rf	105 Db	106 Sg	107 Ns	108 Hs	109 Mt	110 Ds	111 Rg	112 Cn	113 Nh	114 Fl	115 Mc	116 Lv	117 Ts	118 Og	
* Lanthanides			* 57 La	58 Ce	59 Pr	60 Nd	61 Pm	62 Sm	63 Eu	64 Gd	65 Tm	66 Yb	67 Lu						
** Actinides			** 89 Ac	90 Th	91 Pa	92 U	93 Np	94 Pu	95 Am	96 Cm	97 Bk	98 Cf	99 Es	100 Fm	101 Md	102 No	103 Lr		

<span style="color: blue;">H</span> - gas	Li - solid	Br - liquid
<span style="background-color: #90EE90;"> </span> Non-Metals	<span style="background-color: #ADD8E6;"> </span> Transition Metals	<span style="background-color: #ADD8E6;"> </span> Rare Earth
<span style="background-color: #FFD700;"> </span> Alkali Metals	<span style="background-color: #ADD8E6;"> </span> Alkali Earth Metals	<span style="background-color: #FFB6C1;"> </span> Other Me

**Element: 80-Hg Mercury**

Hg

Z: 80

Atomic Weight: 200.59 (2)

Density: 13.546 c

Melting: -38.83

Boiling: 356.73

Oxidation: +1,+2

A	Jπ	Δ (MeV)	T1/2, Γ, Abundance	Decay Mode
175		-8.2s	20 ms (+40-13)	A
176	0+	-11.80	34 ms (+18-9)	A ~100%
177		-12.7	0.130 s (5)	A 85%, EC 15
178	0+	-16.32	0.254 s (19)	A ~ 70%, EC
179		-17.0s	1.09 s (4)	A ~ 53%, EC
180	0+	-20.2s	3.0 s (2)	EC 51%, A 49
181	1/2(-)	-20.7s	3.6 s (3)	EC 64%, A 36
182	0+	-23.5s	10.83 s (6)	EC 84.8%, A 1
183	1/2-	-23.9s	9.4 s (7)	EC 74.5%, A 2
184	0+	-26.2s	30.6 s (3)	EC 98.89%, A

c) At 20 C.

# Import / Export

## Importing

- **Input:** merge parts or entire input file with the current
- **Mcnp:** import mcnp geometry into FLUKA. (experimental)

## Exporting

- **Gnuplot:** save active plot to a gnuplot script
- **Makefile:** create a makefile for compiling the executable
- **Mcnp:** save input in MCNP format: Geometry, Materials, Importances
- **Povray:** save geometry into povray 3D format

# Programming Interface: API

There is work presently going on to decouple the functionality from the interface, some of the basic classes can be used to input processing

file: **Input.py** - to manipulate input files

```
import Input
```

```
Input.init([database])
```

 to initialize the database of cards

Most commonly used classes:

Card containing the description of each card

Input manipulating the FLUKA input file

file: **Project.py** - to manipulate project files

# API: class Card

Constructor: `Input.Card(tag, what [,comment [,extra]])`  
what is a list starting with `what[0]=sdum`

## Important Methods:

<code>setWhat(n, value)</code>	set value to <code>what#n</code>
<code>nwhats()</code>	return number of whats
<code>what(n)</code>	return value of <code>what#n</code>
<code>numWhat(n)</code>	return numeric value of <code>what#n</code>
<code>intWhat(n)</code>	return integer value of <code>what#n</code>
<code>clone()</code>	return a copy of the card
<code>setEnabled(e)</code>	enable/disable card

# API: class Input

Constructor: `Input.Input()`

initialize the structure to hold an input file

## Important Variables:

cardlist

a list with pointers to cards

cards

a dictionary with pointers to cards grouped per tag

## Important Methods:

`read(filename)`

read input from file

`write(filename)`

write input to filename

`addCard(card,pos)`

add card to position pos (or end of file)

`delCard(pos)`

delete card from position pos

`preprocess()`

preprocess input to check for active cards

`setEnabled(e)`

enable/disable card

# API: class Project

Constructor: `Project.Project()`

initialize the structure to hold a project file

## Important Methods:

`clear()` to re-initialize project

`load(filename)` load project from file filename

`save([filename])` save project to filename

`runCmd(run)` create run command

# API: example

Read an input file and modify the random number seed

```
import Input
Input.init()
input = Input.Input()
input.read("test.inp")
try:
    rndcard = self.cards["RANDOMIZ"][0]
    rndcard.setWhat(2,5723)
except:
    print "No RANDOMIZE card found"
    sys.exit(0)
input.write("test2.inp")
```



# API: .flair file structure

# comments

Variable: Value

Notes:

multi-line values are terminated with  
Ctrl-L

Run: name

...

Block of Run related information

Data:

...

... Including Data processing information

EndData

EndRun

Plot: name

...

Plot related informations

EndPlot