

Standard output, common errors and mistakes

Exploring the standard output | Handling errors | Common mistakes

Beginner online training, Fall 2020

Introduction

- We have already encountered most of the basic functions of FLUKA and introduced Flair as the main interface for setting up FLUKA simulations (building the geometry, defining materials and beams etc.)
- In the background, the code generates various output files that provide useful information

- The first objective of this lecture is to present the information that can be found in these files, focusing on the standard output file (.out) and, more briefly, on the .err and .log files
- In the second part, we shall present a non-exhaustive list of possible errors and, most importantly, how to use the available information in order to identify their cause



The standard output file

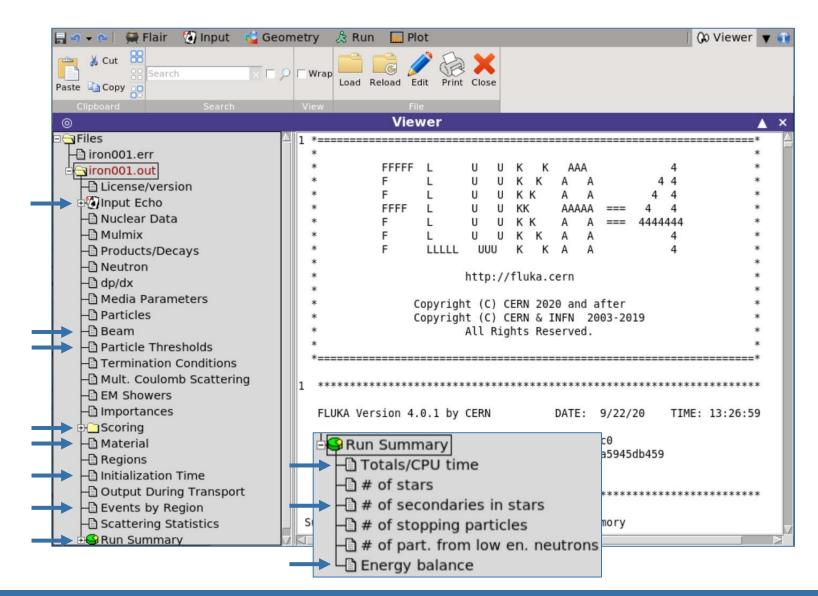
and the .err and .log files



- The standard output file (and other generated files) can be listed and viewed inside Flair via the Files menu under the Run tab
- It may be found in the working directory (i.e. the directory in which you launched the run) or in the temporary fluka_* subdirectory if the run is still ongoing or if it ended with errors
- Name: <inputname>###.out,
 e.g. myrun001.out
- 🚺 Input 🖨 Flair 🚴 Run Plot Ceometry 🔏 Cut 😽 a) →Ascii ... X Delete Filter Copy OD Runs Data Viewer Editor Paste Files **G** Refresh Run 0 Run Cycles Spawn File 001 <ex materials solution> raniron001 002 iron001.out iron 003 cheese iron001.log 004 iron001.err 005 iron001 fort.21 006 irco.out compile data input plot temporary
- One output file is produced per cycle (*001.out, *002.out etc.)
- Note: Not to be confused with the <inputname>.out file



- The .out file can be quite large
- The Flair viewer organises it into sections for easier navigation
- It contains a trove of useful information concerning the simulation
- Extremely useful for investigating strange results and errors
- Let's look at some of the contents!





Beam information

- Particle type, energy, source position and spatial profile
- (as long as you are not generating arbitrary distributions via source routine)

=== Output before the actual run - Beam properties ===

Fluka incident beam properties:

	Beam particle: PROTON Id: 1 (F Mass: 0.9383 (C		_	-	
	Average beam momentum :	Θ.	.644445 (GeV/c)	
	Average beam kinetic energy:	Θ.	.200000 (GeV)		
	Momentum deviation at FWHM (rectar	ngular):	0.0000000 (Ge	V/c)	
ſ	Beam hit position : 0.000000	0.000	90000 -5.	00000000	CM
	Beam direction cosines: 0.000000	0.000	00000 1.	000000000	
4	Beam spot FWHM X-width (Rectangula	ar): 0.0000	cm		
	Beam spot FWHM Y-width (Rectangula				
	Beam FWHM angular divergence (Rect	tangular):	0.0000	(mrad)	
	(Spatial distribution, polarization	on, and angular	direction and	distribution	
	are given in the beam frame of ref	ference)			
	Beam reference frame (world coordi	inates):			
	Beam X axis: 1.00000000	0.00000000	0.0000000		
	Beam Y axis: 0.00000000	1.00000000	0.0000000		
	Beam Z axis: 0.00000000	0.00000000	1.00000000		
	The nominal beam position belongs lattice cell: 0 ()	to region:	2 (VOID),	
	Possing constructions - Constructions - Constructions - Constructions				



Summary of transport thresholds

=== Particle transport thresholds:

100 keV (**PRECISION DEFAULTS**)

<u>Global cut-off kinetic energy for particle transport</u>: <u>1.000E-04 GeV</u> The cut-off kinetic energy is superseded by individual particle thresholds if set

Cut-off kinetic energy for 4-HELIUM transport: 1.000E-04 GeV Cut-off kinetic energy for 3-HELIUM transport: 1.000E-04 GeV Cut-off kinetic energy for TRITON transport: 1.000E-04 GeV Cut-off kinetic energy for DEUTERON transport: 1.000E-04 GeV Cut-off kinetic energy for PROTON transport: 1.000E-04 GeV Cut-off kinetic energy for APROTON transport: 1.000E-04 GeV Cut-off kinetic energy for ELECTRON transport defined in the Emfcut card Cut-off kinetic energy for POSITRON transport defined in the Emfcut card Cut-off kinetic energy for NEUTRIE transport: 0.000E+00 GeV Cut-off kinetic energy for ANEUTRIE transport: 0.000E+00 GeV Cut-off kinetic energy for PHOTON transport defined in the Emfcut card Cut-off kinetic energy for NEUTRON transport: 1.000E-14 GeV

Correspondence of regions	and EMF-FLUKA material num	bers
Region EMF	FLUKA	
1 BLKBODY	0 VACUUM	1 BLCKHOLE
2 VOID	0 VACUUM	2 VACUUM
3 TARGET	1 IRON 11	L IRON
Prompt:		
Ecut = 6.1100E-01 MeV,	Pcut = 3.3333E-02 MeV	
BIAS = F, Ray. = T,	S(q,Z) = T, $Pz(q,Z) = T$,	Acolin.= F
Delayed:		
Ecut = 6.1100E-01 MeV,	Pcut = 3.3333E-02 MeV	

In the absence of an **EMFCUT** card explicitly setting them, **electron/ positron** and γ transport thresholds are calculated internally



=== Material compositions: ===

Materials and material properties

- Includes FLUKA predefined materials and userdefined materials
- Basic material properties

 (In)elastic scattering length, radiation length for selected BEAM particle/energy

Material Number&Name	Atomic Number	Atomic Weight	Density	Inelastic Scattering Length for PROTON at Beam energy	Elastic Scattering Length for PROTON at Beam energy	Radiation Length
			g/cm**3	cm	cm	cm
1 BLCKHOL 2 VACUUM 3 HYDROGE	0.000	0.000 0.000 1.008	0.000 0.000 0.8370E-04	0.1000E+31 0.1000E+31 0.3641E+10	0.1000E+31 0.1000E+31 0.8848E+06	0.1000E+31 0.1000E+31 0.7532E+06
30 GRUYERE	3.294	5.975	0.9153	115.2	219.2	41.98
Materia	al Number	Atom content	Partial Densi	ties		
HYDROGE OXYGEN CARBON NITROGE CALCIUM PHOSPHO SODIUM POTASSI MAGNESI	8 6 2N 7 1 21 0 28 19 20	0.62121 0.17083 0.18469 0.19411E-01 0.15141E-02 0.11754E-02 0.95020E-03 0.11485E-03 0.99864E-04	0.95916E-01 0.41870 0.33981 0.41649E-01 0.92953E-02 0.55772E-02 0.33463E-02 0.68786E-03 0.37181E-03			
31 WATER	3.333	6.005	1.000	109.5	194.4	36.08
Materia	al Number	Atom content	Partial Densi	ties		
HYDROGE 0XYGEN	N 3 8	0.66667 0.33333	0.11190 0.88810			



Standard output, common errors and mistakes

Summary of requested scoring

Interpreted scoring cards

****** "usrbin" option: Cartesian binning n. 1 "Protons ", generalized particle n. 1 X coordinate: from -1.0000E+01 to 1.0000E+01 cm, 201 bins (9.9502E-02 cm wide) Y coordinate: from -1.0000E+01 to 1.0000E+01 cm, 201 bins (9.9502E-02 cm wide) Z coordinate: from -5.0000E+00 to 1.5000E+01 cm, 200 bins (1.0000E-01 cm wide) data will be printed on <u>unit -21</u> (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:

No user bdrx defined

******* "USRTRACK" option:

No user track-length estimator defined



Input echo

- Interpreted body and region definitions
- Very useful in case of complex inputs controlled by many different preprocessor directives, where the options may not be activated as intended due to a mistake
- The input echo will only contain the parts which were really active in the run

Initialisation time

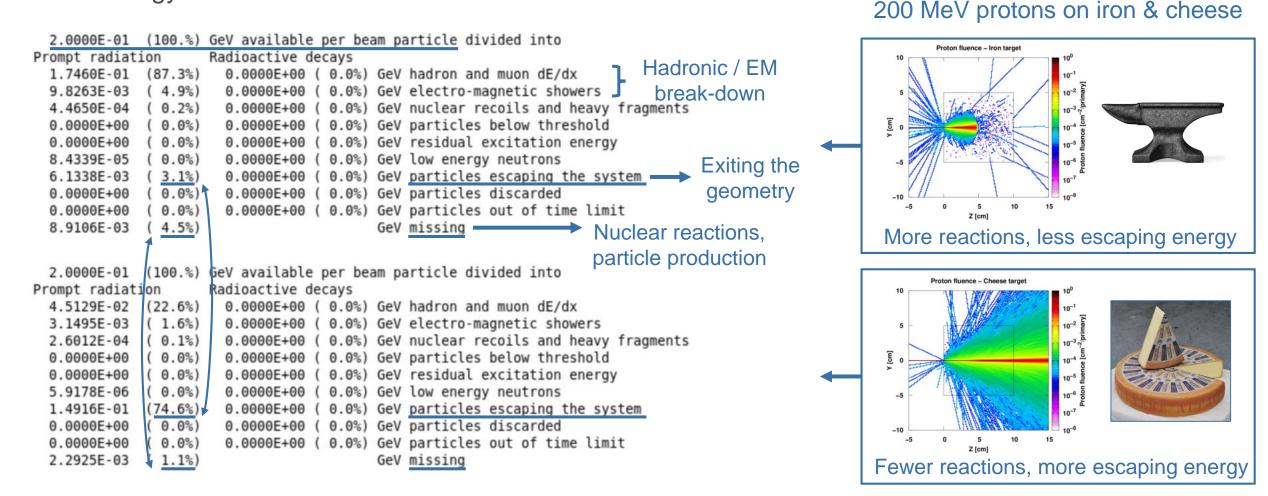
- Run initialisation happens at the beginning of each cycle
- It is generally quite fast (especially relative to the total duration of the run) but can reach up to several minutes for complex geometries, LATTICEs, scoring, preprocessor directives etc.

```
Total time used for initialization:0.539sBasic input templateTotal time used for initialization:82.1sComplex LHC simulation,<br/>34000 line input file
```



Run summary

Energy balance

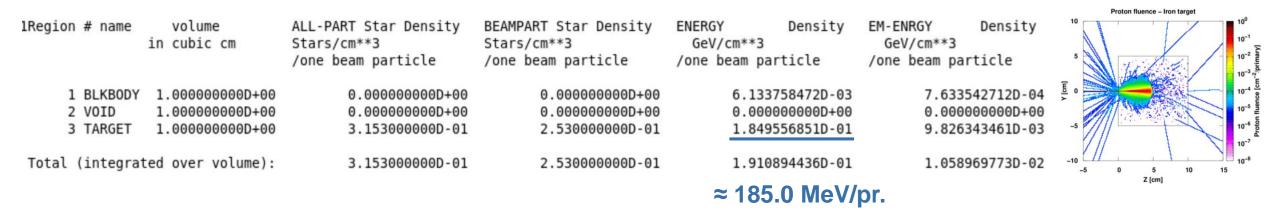


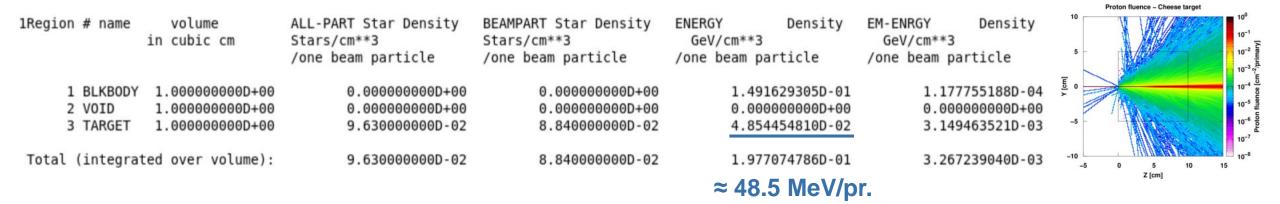


Recall the materials exercise:

Information per region

Energy deposition







Run summary

• Primary particles, number of interactions, CPU-time

	Total number of primaries run:10000 for a!!! Please remember that all results are normalizedThe main stack maximum occupancy was12 out	per unit weight !!!			
	Total <u>number of inelastic interactions</u> (stars): Total weight of the inelastic interactions (stars):	963 9.630000E+02			
	Total <u>number of elastic interactions</u> : Total weight of the elastic interactions: 6.490000E	649 +02			
	Total <u>number of low energy neutron interactions</u> : Total weight of the low energy neutron interactions:	1047 1.047000E+03			
From fractions of a millisecond to tens of minutes or more	Total CPU time used to follow all primary particles: Average CPU time used to follow a primary particle: Maximum CPU time used to follow a primary particle: Residual CPU time left:	6.594E-04 seconds of:			

of minutes or more

Standard output, common errors and mistakes

Run summary

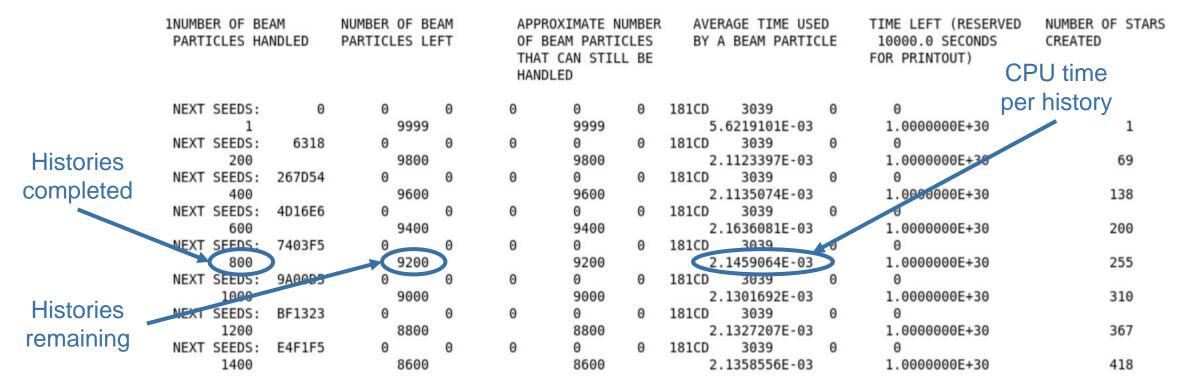
• Number and type of produced secondary particles

Number of secondaries Prompt radiation	generated in inelastic interactions per beam particle:
	0.0000E+00 (100.%)
	0.0000E+00 (0.0%) 4-HELIUM
	0.0000E+00 (0.0%) 3-HELIUM
and a second	0.0000E+00 (0.0%) TRITON
	0.0000E+00 (0.0%) DEUTERON
Eller and the construction of the second	0.0000E+00 (0.0%) HEAVYION
	0.0000E+00 (0.0%) OPTIPHOT
	0.0000E+00 (0.0%) RAY
5.1690E-01 (27.2%)	0.0000E+00 (0.0%) PROTON
0.0000E+00 (0.0%)	0.0000E+00 (0.0%) APROTON
0.0000E+00 (0.0%)	0.0000E+00 (0.0%) ELECTRON
0.0000E+00 (0.0%)	0.0000E+00 (0.0%) POSITRON
0.0000E+00 (0.0%)	0.0000E+00 (0.0%) NEUTRIE
0.0000E+00 (0.0%)	0.0000E+00 (0.0%) ANEUTRIE
7.0840E-01 (37.2%)	0.0000E+00 (0.0%) PHOTON
5.9500E-01 (31.3%)	0.0000E+00 (0.0%) NEUTRON



The .err file

- Note: The .err file is generated even for successful runs, its presence does not necessarily imply than any errors were encountered
- It contains information concerning the progress of the run
- Run-time error messages and warnings will also be written to this file





The .log file

 Critical errors, usually following a code abort will be written to the .log file, as in the following examples

Pro	gram received signal SIGFPE: Floating-point exception - erroneous arithmetic operation.
Bac	ktrace for this error:
#0	0x7f60e3aae3ff in ???
#1	0x7f60e3ab84fe in ???
#2	0x7f60e47b13a8 in read_real
	at /GCC/build/contrib/gcc-8.2.0/src/gcc/8.2.0/libgfortran/io/list_read.c:1872
#3	0x7f60e47b2fcd in list_formatted_read_scalar
	at /GCC/build/contrib/gcc-8.2.0/src/gcc/8.2.0/libgfortran/io/list_read.c:2180
#4	0x6b1859 in flkcgi_
	at comlat/flkcgi.f:556
#5	0x56d96e in geoinp_
	at geolat/geoinp.f:241
#6	0x40388e in flukam_
	at main/flukam.f:2442
#7	0x402100 in fluka
	at main/fluka.f:77
#8	0x402100 in main
	at /shared/src/usflmd.inc:15

STOP TOO MANY ERRORS IN GEOMETRY Note: The following floating-point exceptions are signalling: IEEE_UNDERFLOW_FLAG IEEE_DENORMAL STOP STOP: FLUKA ABORTED



User-defined messages

- Users may print information, warning or error messages to the .out and .err files from within their user routines, e.g.:
 - WRITE (LUNOUT, *) 'This is a message'
 - WRITE (LUNERR,*) 'This is a message'
- This can be useful when debugging user routines



Error handling and common mistakes

Progress

Status: Finished with ERRORS



Before running

- Have the latest FLUKA and Flair versions correctly installed, downloading the appropriate version for your system and following the instructions
- Try running a provided example to confirm that there are no installation errors
- Do not run inside the FLUKA directory!
- Avoid using file and directory names with spaces
- Flair will already detect and highlight many mistakes, such as nonsense inputs in certain fields, multiple **ASSIGNMAT** cards for a region etc.:

■ USRBIN Type: X-Y-Z ▼ Part: PROTON ▼	Xmin: -10.0 Ymin: 10.0 Zmin: -5.0	Unit: 21 BIN V Xmax: 10.0 Ymax: -10.0 Zmax: 15.0	Name: Protons NX: 201. NY: 201. NZ: 200.	>w> Warning:		card #19 IRON	region TARGET: TARGET
MATERIAL POTASSIU		#:	p: -0.862		ASSIGNMA	COPPER	TARGET
Z: 19	Am:	A:	dE/dx: 🔻				

• Build the geometry carefully, use the Geometry Editor to search for geometry errors



Click "Start"...

- Runs may fail during initialisation; this usually indicates the existence of a blatantly wrong setting in a card, some missing necessary setting, a mistyped expression or value, a missing file, #if directive without corresponding #endif, ...
- Crashes occurring once particle tracking has started are often attributable to geometry errors, numerical exceptions (often associated with user routines), ...



Crashes at initialisation

Low-energy neutron cross-sections not found

- As mentioned in the Materials lecture, low-energy neutron cross sections are associated to materials based on their names
- If a user-defined elemental material is named differently (1) than in the relevant library (2) (see section 10.4.1.2 of the FLUKA manual), then FLUKA will not be able to assign any neutron cross-sections



• This leads to a stop of the execution with the following message (found in the .out file)

**** Low energy neutron xsec not found for some media F



Crashes at initialisation

Missing random number file

- This error usually indicates that the previous cycle did not end successfully and therefore the random number file ran<inputname>### was not generated
- The relevant error messages are found in the .out, .err and .log files:

**** No Random file available !!!!!! **** Abort called from FLRM64 reason NO RANDOM FILE Run stopped! <inputname>###.out STOP NO RANDOM FILE

STOP NO RANDOM FILE STOP STOP: FLUKA ABORTED

<inputname>###.log

Abort called from FLRM64 reason NO RANDOM FILE Run stopped! <inputname>###.err STOP NO RANDOM FILE

• When a run fails, it is generally advisable to look at the output files of the first cycle, because all further cycles will fail with this error after a failed first cycle



Crashes at initialisation

Missing executable or external file

- If a custom executable or external auxiliary file (e.g. to be used for sampling by a source routine) is not found in the working directory, the code will stop
- An example of the error message in the <inputname>.out file:

```
*-* Running: errors
Dir: /errors
Cmd: /usr/bin/nohup /soft/fluka4-0.1/bin/rfluka -e /errors/myexe -M 5 errors
/usr/bin/nohup: ignoring input
```

FLUKA

Dir: /soft/fluka4-0.1 Data: /soft/fluka4-0.1/data Exec: /errors/myexe Input: /errors/errors.inp Error: /errors/myexe does not exist or it is not executable!

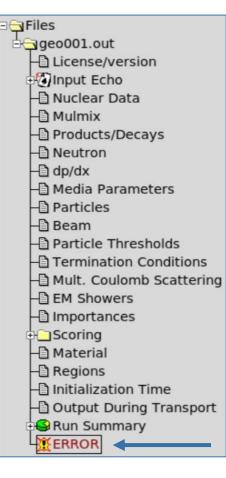
Input file formatting errors

- Syntax/alignment errors in manually edited input files
- Presence of hidden DOS characters in files edited in Windows (clean with dos2unix command)



Run-time errors

- Where to look?
 - The bottom of the .out and .err files, and the top of the .log file



When you open the standard output file in Flair, the navigation menu will point to the error message

- ...or, at least, to the bottom of the file; in case of an uncontrolled crash, the final information printed in the standard output file is less likely to be informative
- Instead, more meaningful information may be found in the .log and .err files



Run-time errors

Geometry and tracking accuracy errors

- A single geometry error or problem with tracking accuracy (often encountered when using LATTICEs) may not condemn a run, as the code attempts to recover the lost particle
- These attempts are recorded in the .err file ("geofar" errors)

NEXT SEEDS: F0A31F Θ 181CD 3039 Θ Θ Θ Θ 1100 1100 2.2694074E-04 1.000000E+30 23 3900 NEXT SEEDS: F25D77 Θ 181CD 3039 Θ Θ Θ 0 0 Geofar: Particle in region 3 (cell # 0) in position 9.208454725E-01 1.554251376E+00 1.303608751E+00 is now causing trouble, requesting a step of 5.206776825E+00 cm to direction 8.513230560E-02 3.069905258E-01 9.478973086E-01, error count: 0 R2: 1.806558531E+00 R3: 2.227790273E+00 cm X*U (2D): 5.555341454E-01 X*U (3D): 1.791221372E+00 cm X*UOLD(2D): 8.622110860E-01 X*UOLD(3D): 4.169256137E-01 cm 3, Irsav2: Kloop: 32194, Irsave: 3, error code: -33 Nfrom: 5000 old direction 9.398507183E-01 -2.088589279E-03 -3.415790758E-01, lagain, lstnew, lsense, lsnsct F F F T Particle index 7 total energy 4.547721732E-04 GeV Nsurf Θ Try again to establish the current region moving the particle of a 3.887637526E-08 long step We succeeded in saving the particle: current region is n. 3 (cell # 0)

• The run will stop if too many errors are encountered and/or the particle cannot be recovered

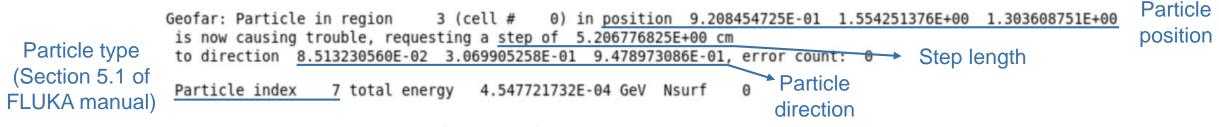
Abort called from FLKAG1 reason TOO MANY ERRORS IN GEOMETRY Run stopped! STOP TOO MANY ERRORS IN GEOMETRY



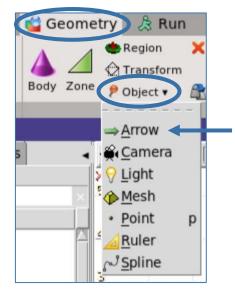
Run-time errors

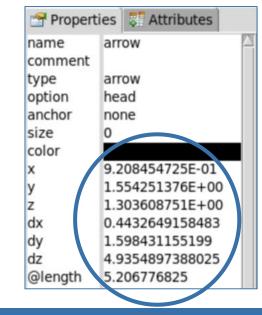
Geometry errors

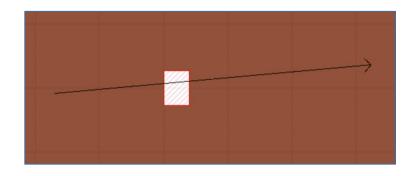
• These error messages offer valuable information that can help locate the problem



 In Flair, you can add a vector ("Arrow") setting its origin, direction and length according to the error message to help in visualising the geometry problem









Advanced debugging

 It can get worse: uncontrolled code aborts may lead to the creation of a core dump in the temporary FLUKA directory (fluka_*), with an error like the following to be found in the <inputname>.out file

- Such errors are usually caused by numerical issues, variable type discrepancies, attempts to access non-existent array elements etc., and are often associated with the use of user routines
- The core. * file contains information on the program state at the time of the crash and the relevant back-trace
- It can be accessed via the GNU debugger (gdb)
 - ...but this is a story for another course...



If the run ends without errors...

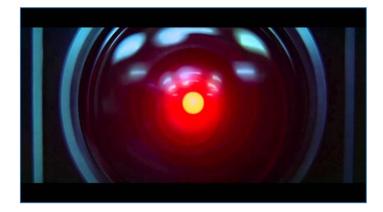
Progress Status: Finished OK

- Some mistakes will not lead to run-time crashes, but may have an impact on the results of the simulation
 - Overlapping geometry regions will not generate an error; during particle tracking the code will decide in which region the particle is and continue, potentially distorting the scoring of energy deposition, fluence etc.
 - Wrong beam energy assignment (e.g. MeV vs. GeV, energy vs. momentum, total HEAVYION energy in **BEAM** card instead of energy per nucleon, ...)
 - Wrong material densities
 - Wrong/multiple material assignments for a region
 - Inappropriate thresholds
 - A missing **RANDOMIZ** card will lead to zero statistical errors, as identical cycles are executed
- Such errors can be discovered upon inspection of the results which may completely deviate from expectations, display strange discontinuities etc.
- You should always critically assess your results based on your knowledge of the problem!



To conclude...

• Errors will always crop up, even for experienced users; do not despair!



"It can only be attributable to human error." -- HAL 9000

- Most of the time, the cause will be fairly mundane, and a hint on its origin will be found in the various error messages and output files
- If you are unable to discover the cause of the problem using the suggestions in this lecture, it's time to explore the FLUKA forum for similar issues and, finally, to post your question on the FLUKA forum
- Remember that meaningful debugging assistance generally requires that you at least provide your .flair file, and other relevant files, user routines etc.



