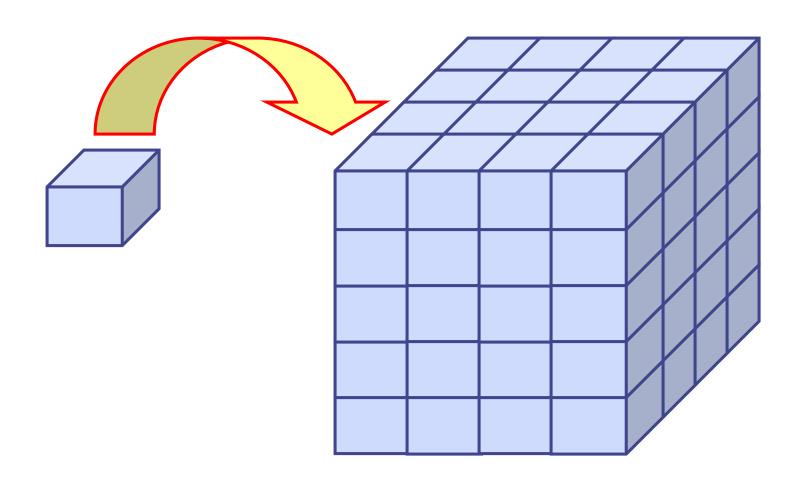


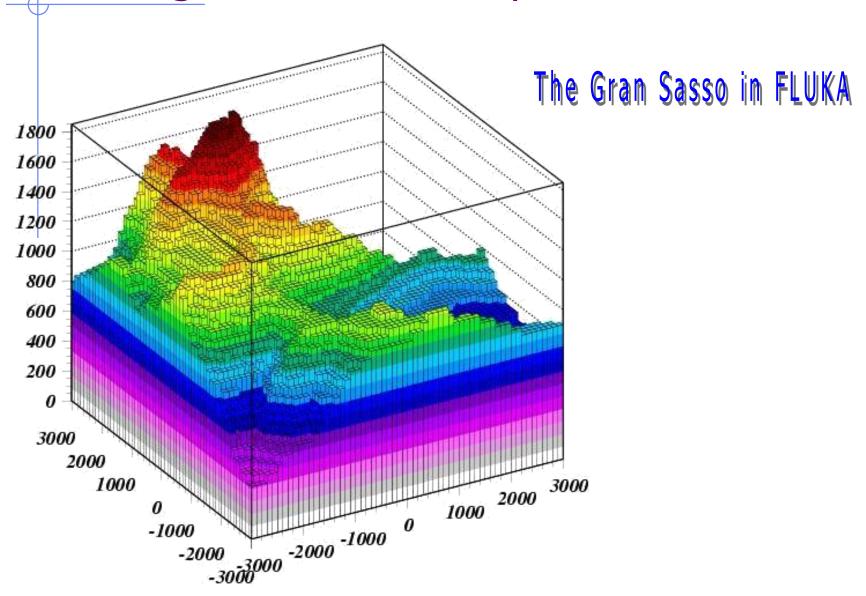
# **Voxels and Medical Applications**

FLUKA Beginners course

• It is possible to describe a geometry in terms of "voxels", i.e., tiny parallelepipeds (all of equal size) forming a 3-dimensional grid

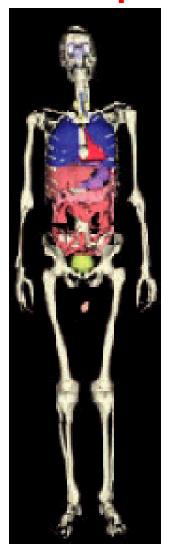


## Voxel geometries: examples



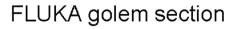
#### Voxel geometries: examples

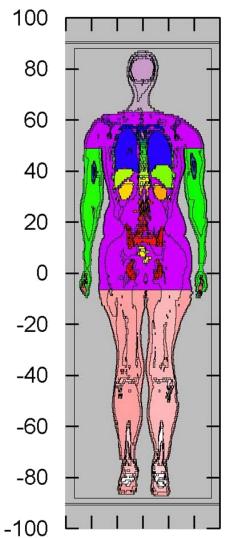
The anthropomorphic GOLEM phantom



Implementation in FLUKA (radioprotection applications)

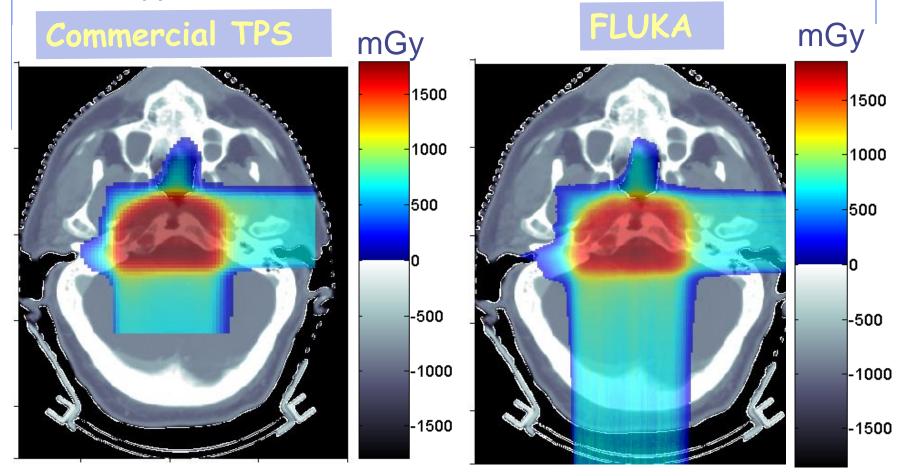
Petoussi-Henss et al, 2002





### Voxel geometries in medical applications

Voxel geometries are especially useful to import CT scan of a human body, e.g., for dosimetric calculations of the planned treatment in radiotherapy



Parodi et al., 2007

• The CT scan contains integer values "Hounsfield Unit" reflecting the X-ray attenuation coefficient  $\mu_x$ 

```
HU_x = 1000 \ (\mu_x - \mu_{H20}) \ / \ \mu_{H20} , typically -1000 \leq HU \leq 3500
```

- We will use loosely the word "organ" to indicate a group of voxels (or even more than one group) made of the same "tissue" material (same HU value or in a given HU interval)
- The code handles each organ as a CG region, possibly in addition to other conventional "non-voxel" regions defined by the user
- The voxel structure can be complemented by parts written in the standard Combinatorial geometry
- The code assumes that the voxel structure is contained in a parallelepiped. This RPP is automatically generated from the voxel information.

 To describe a voxel geometry, the user must convert his CT scan or equivalent data to a format understood by FLUKA. Starting from DICOM images, this is performed directly by FLAIR (see next slides).

#### This stage should :

- Assign an organ index to each voxel. In many practical cases, the user will have a continuum of CT values (HU), and may have to group these values in intervals
- Each organ is identified by a unique integer ≤32767. The organ numbering does not need to be contiguous i.e. "holes" in the numbering sequence are allowed.
- One of the organs must have number 0 and plays the role of the medium surrounding the voxels (usually vacuum or air).
- Assign to each NONZERO organ a voxel-region number. The voxel-region numbering has to be contiguous and starts from 1.

- The information is input to FLUKA through a special file \*vxl containing:
  - The number of voxels in each coordinate
  - The number of voxel-regions, and the maximum organ number
  - The voxel dimension in each coordinate
  - A list of the organ corresponding to each voxel
  - A list of the voxel-region number corresponding to each organ
  - (In the new release) definition of Regions of Interests (ROI)
    - A list of the ROIs for each voxel

## Input file

Prepare the usual FLUKA input file.

The geometry is written like a normal Combinatorial Geometry input, but in addition a VOXELS card must be inserted right after the GEOBEGIN card and before the Geometry title card

- WHAT(1), WHAT(2), WHAT(3) = x, y, z coordinates chosen as the origin of the "voxel volume", (i.e. of a region made of a single RPP body extending from WHAT(1) to WHAT(1) + NX\*DX, ...) which contains all the voxels
- WHAT(4) ROT-DEFI transformation applied to the whole voxel
- WHAT(5), WHAT(6): not used
- SDUM = name of the voxel file extension will be assumed to be .vxl)



x: -35.068359

y: -35.068359

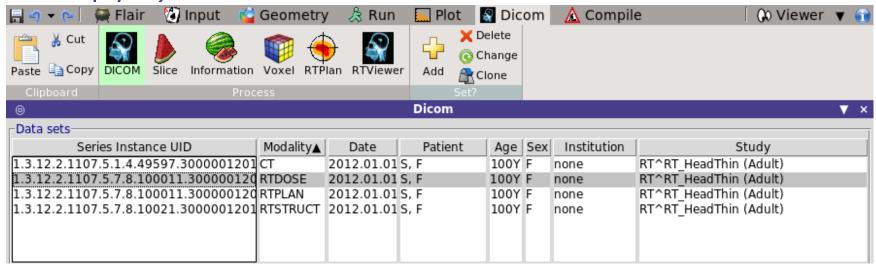
z: -88.6855

Trans: 🔻

Filename: VOXEL1 ▼

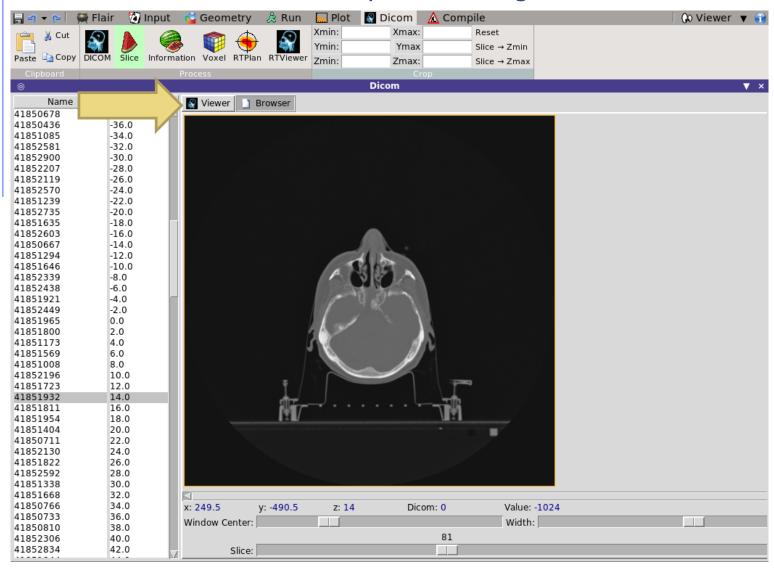
### Processing the **DICOM** files with FLAIR I

- DICOM = Digital Imaging and Comunications in Medicine is a medical standard for distributing any kind of medical image.
- FLAIR has a capability to process the DICOM files using the pydicom module and convert them to FLUKA VOXELS, USRBIN compatible files as well as providing input for Treatment Plan simulations
- First select the "Directory" where the DICOM data sets are located (if you have doubts press F1 and the flair manual will help you).

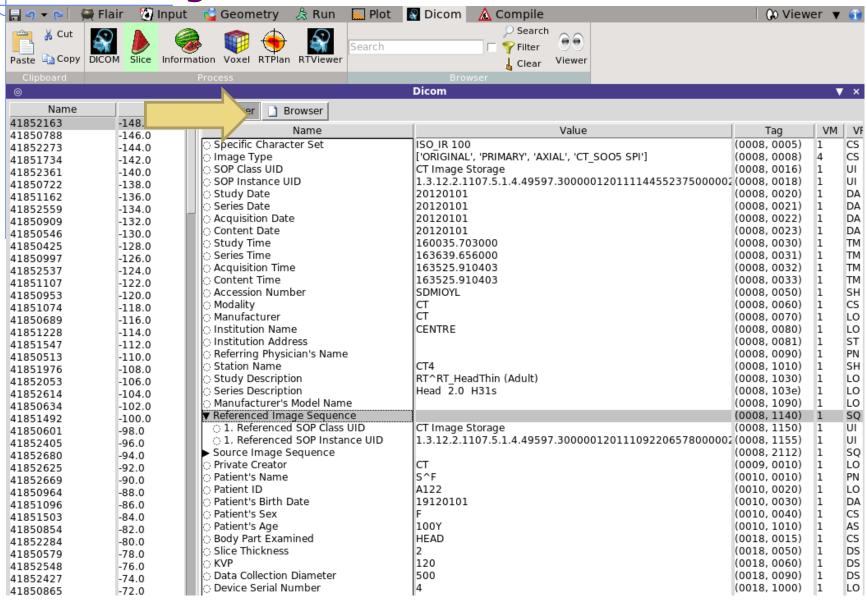


#### Processing the **DICOM** files with FLAIR II

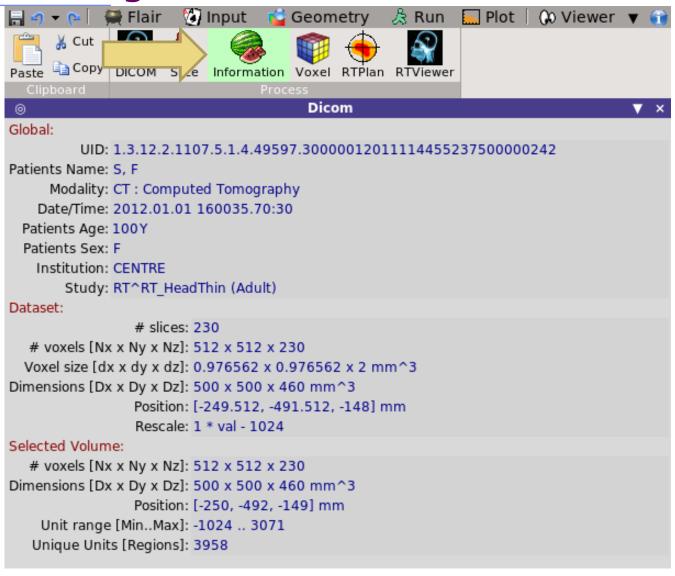
Select one "Data sets" and inspect the images.



### Processing the **DICOM** files with FLAIR III

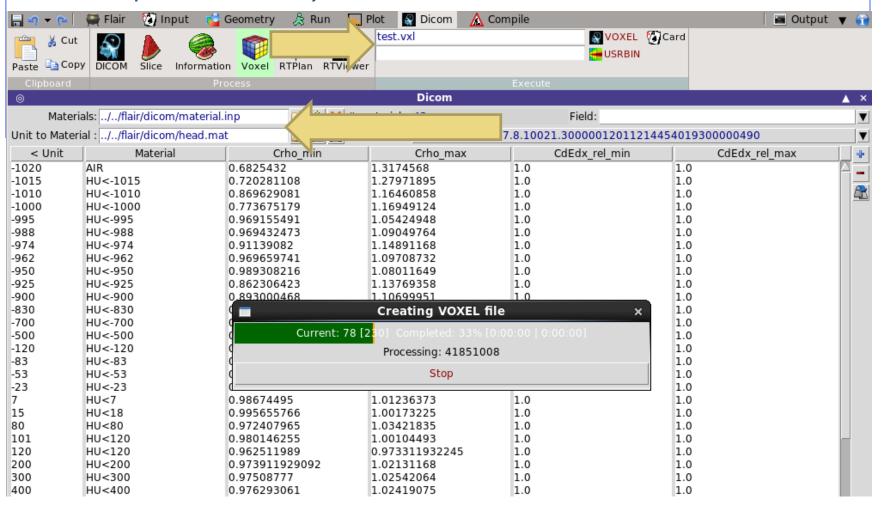


#### Processing the **DICOM** files with FLAIR IV



#### Processing the **DICOM** files with FLAIR V

The "Voxel" tab is used in order to convert the CT dataset to VOXELS. For the VOXEL geometry two additional files are needed (example: material.inp and head.mat).



#### Processing the **DICOM** files with FLAIR VI

- <= Unit: specify the upper limit of the range. Every entry will correspond to a range from the previous upper limit+1 until the current upper limit.
  </p>
- Material: select any of the predefined FLUKA materials defined previously.

Optionally you can specify correction factors for the density and dE/dx

- Crho\_min/Chro\_max: density correction factors to be applied on the lower/upper limit of the unit range (see next slides).
- CdEdx\_rel\_min/CdEdx\_rel\_max: relative correction factors on dE/dx for minimum/maximum unit in the range (see next slides).

≤ Unit	Material	Crho_min	Crho_max	CdEdx_rel_min	CdEdx_rel_max
-1020	AIR	0.6825432	1.3174568	1.0	1.0
-1015	HU<-1015	0.720281108	1.27971895	1.0	1.0
-1010	HU<-1010	0.869629081	1.16460858	1.0	1.0
-1000	HU<-1000	0.773675179	1.16949124	1.0	1.0
-995	HU<-995	0.969155491	1.05424948	1.0	1.0
-988	HU<-988	0.969432473	1.09049764	1.0	1.0
-974	HU<-974	0.91139082	1.14891168	1.0	1.0
-962	HU<-962	0.969659741	1.09708732	1.0	1.0
-950	HU<-950	0.989308216	1.08011649	1.0	1.0
-925	HU<-925	0.862306423	1.13769358	1.0	1.0
-900	HU<-900	0.893000468	1.10699951	1.0	1.0
-830	HU<-830	0.783902333	1.21609767	1.0	1.0
-700	HU<-700	0.75158871	1.24841129	1.0	1.0
-500	HU<-500	0.765689411	1.23431059	1.0	1.0
-120	HU<-120	0.734835247	1.26516475	1.0	1.0
-83	HU<-83	0.980501545	1.01835909	1.0	1.0
-53	HU<-53	0.98600717	1.01305997	1.0	1.0

#### Voxel Body

- The usual list of NB bodies, not including the RPP corresponding to the "voxel volume" (see VOXELS card above). This RPP will be generated and added automatically by the code as the (NB+1) th body, with one corner in the point indicated in the VOXELS card, and dimensions NX\*DX, NY\*DY and NZ\*DZ as read from the voxel file.
- The usual region list of NR regions, with the space occupied by body named VOXEL or numbered NB+1 (the "voxel volume") subtracted. In other words, the NR regions listed must cover the whole available space, excepted the space corresponding to the "voxel volume". This is easily obtained by subtracting body VOXEL or NB+1 in the relevant region definitions, even though this body is not explicitly input at the end of the body list.

VOXELS	x: -35.068359 Trans: ▼	y: -35.068359 Filename: VOXEL1 ▼	z: -88.6855
SPH BODY1	x: 0. R: 10000.	у: 0.	z: <b>0.</b>
SPH BODY2	x: 0. R: 1000.	у: 0.	z: <b>0.</b>
♦ END			
REGION REG1 expr: BODY1-BODY2		Neigh: 5	Volume:
REGION REG2 expr: BODY2-VOXEL		Neigh: 5	Volume:
♦ END			
TAT CECEND	_		

#### **Voxel Regions**

The code will automatically generate and add several regions:

NO additional regions, where NO = number of non-zero organs:

Name	Number	Description
VOXEL	NR+1	sort of a "cage" for all voxels. Nothing should ever be deposited in it. The user shall assign vacuum to it.
VOXEL001	NR+2	containing all voxels belonging to organ number 0. There must be at least 2 of such voxels, but in general they should be many more. Typical material assignment to this region is air
VOXEL002	NR+3	corresponding to organ 1
VOXEL003	NR+4	corresponding to organ 2
VOXEL###	NR+2+NO	corresponding to organ NO

#### Few remarks

- The assignment of materials is made directly by FLAIR. The user has to assign the materials to the regions defined by combinatorial geometry.
- The "head.mat" and "material.inp" files are examples, the user should update these files taking into account his calibration curves.

#### Practical issues for Medical Applications

#### General problems for MC calculations on CT scans

- How to assign realistic human tissue parameters (= materials) for MC Calculation ?
- How to find a good compromise between the number of different HU values (~ 3000-5000) and the materials to be considered in the MC?
  - (issues on memory and computation speed when attempting to treat each HU number as a different material !!!)
- How to preserve continuous, HU-dependent information when segmenting the HU numbers into intervals sharing the same "tissue" material?
  - (critical for ion range calculation in charged hadron therapy !!!)

### CT stoichiometric calibration (I)

CT segmentation into 27 materials of defined elemental composition (from analysis of 71 human CT scans)

Air, Lung,
Adipose tissue

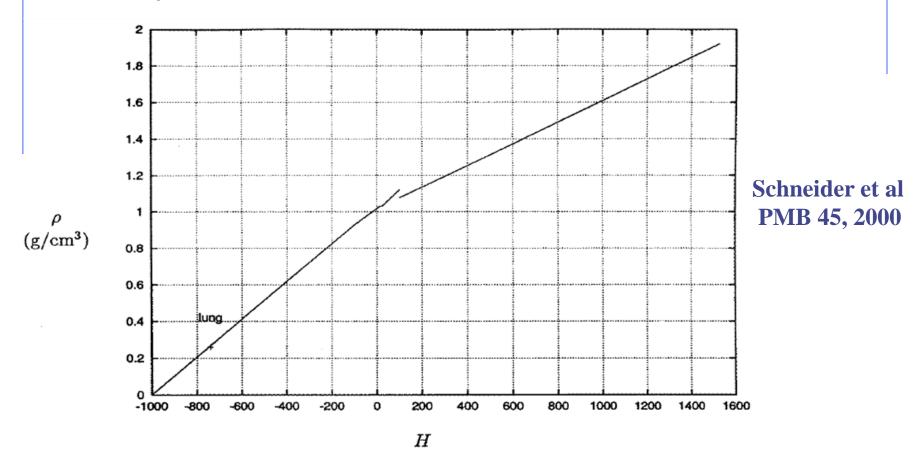
Soft tissue

Skeletal tissué

	$w_i(pp)$											
Н	Н	С	N	0	Na	Mg	P	S	C1	Ar	K	Ca
-1000950			75.5	23.2						1.3		
-950120	10.3	10.5	3.1	74.9	0.2		0.2	0.3	0.3		0.2	
-12083	11.6	68.1	0.2	19.8	0.1			0.1	0.1			
-8253	11.3	56.7	0.9	30.8	0.1			0.1	0.1			
-5223	11.0	45.8	1.5	41.1	0.1		0.1	0.2	0.2			
-22-7	10.8	35.6	2.2	50.9			0.1	0.2	0.2			
8-18	10.6	28.4	2.6	57.8			0.1	0.2	0.2		0.1	
19-80	10.3	13.4	3.0	72.3	0.2		0.2	0.2	0.2		0.2	
80-120	9.4	20.7	6.2	62.2	0.6			0.6	0.3			
120-200	9.5	45.5	2.5	35.5	0.1		2.1	0.1	0.1		0.1	4.5
200-300	8.9	42.3	2.7	36.3	0.1		3.0	0.1	0.1		0.1	6.4
300-400	8.2	39.1	2.9	37.2	0.1		3.9	0.1	0.1		0.1	8.3
400-500	7.6	36.1	3.0	38.0	0.1	0.1	4.7	0.2	0.1			10.1
500-600	7.1	33.5	3.2	38.7	0.1	0.1	5.4	0.2				11.7
600-700	6.6	31.0	3.3	39.4	0.1	0.1	6.1	0.2				13.2
700-800	6.1	28.7	3.5	40.0	0.1	0.1	6.7	0.2				14.6
800-900	5.6	26.5	3.6	40.5	0.1	0.2	7.3	0.3				15.9
900-1000	5.2	24.6	3.7	41.1	0.1	0.2	7.8	0.3				17.0
1000-1100	4.9	22.7	3.8	41.6	0.1	0.2	8.3	0.3				18.1
1100-1200	4.5	21.0	3.9	42.0	0.1	0.2	8.8	0.3				19.2
1200-1300	4.2	19.4	4.0	42.5	0.1	0.2	9.2	0.3				20.1
1300-1400	3.9	17.9	4.1	42.9	0.1	0.2	9.6	0.3				21.0
1400-1500	3.6	16.5	4.2	43.2	0.1	0.2	10.0	0.3				21.9
1500-1600	3.4	15.5	4.2	43.5	0.1	0.2	10.3	0.3				22.5

#### CT stoichiometric calibration (II)

Assign to each material a "nominal mean density", e.g. using the density at the center of each HU interval (Jiang et al, MP 2004)



But "real density" (and related physical quantities) varies continuously with HU value !!!

#### The region-dependent CORRFACT card

- "CORRFACT" card allows to alter material density for dE/dx and nuclear processes
- First two inputs specify a density scaling factor (restricted to the interval [2/3,3/2]) for charged particle ionization processes (WHAT(1)) and for all other processes (WHAT(2)) to the region(s) specified by the inputs WHAT(4-6) [cf. manual]
- This is especially important in ion beam therapy to force the MC to follow the same semi-empirical HU-range calibration curve as the Treatment Planning System (TPS) for dosimetric comparisons
- FLAIR automatically appends the CORRFACT cards calculated taking into account the calibration curves provided by the user at the end of the .vxl file.

#### How to account for HU-dependent dEdx

In the INPUT

 Let several regions share the same material composition and mean density according to CT segmentation (reduced number of materials to save memory / initialization time)

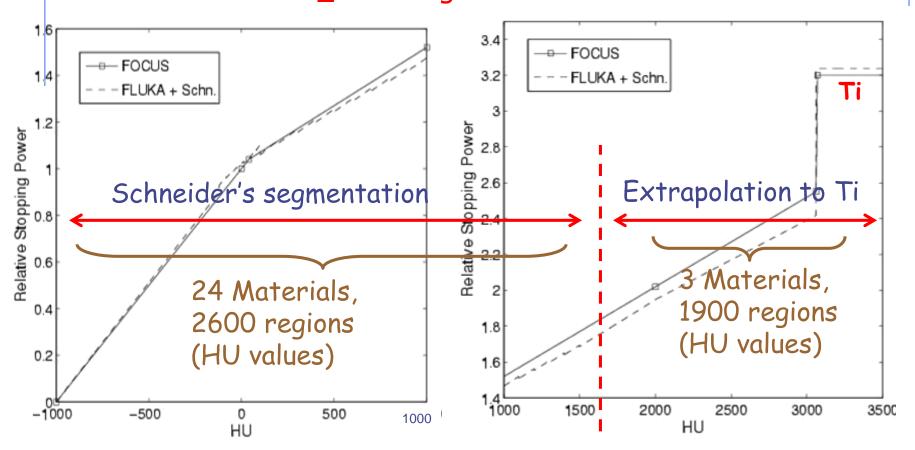
```
ASSIGNMA BONE VOXEL005 (region number 25)
ASSIGNMA BONE VOXEL016 (region number 31)
```

■ Use CORRFACT to impose the desired correction for stopping power (⇒ ion range!) in the regions KREG corresponding to different organs IO (i.e., different HU values) sharing the same MATERIAL assignment

CORRFACT	0.85	0.0	0.0	25	
CORRFACT	1.3	0.0	0.0	31	Region #25 corresponds to "softer" bone than #31

# Forcing FLUKA to follow the same range calibration cur as TPS for p @ MGH Boston

The CORRFACT ionization scaling factors were obtained from the dEdx ratio between TPS and FLUKA (+ Schneider "mass density") -> The user should update the "head.mat" file with his own calibration for CdEdX rel taking into account his TPS.



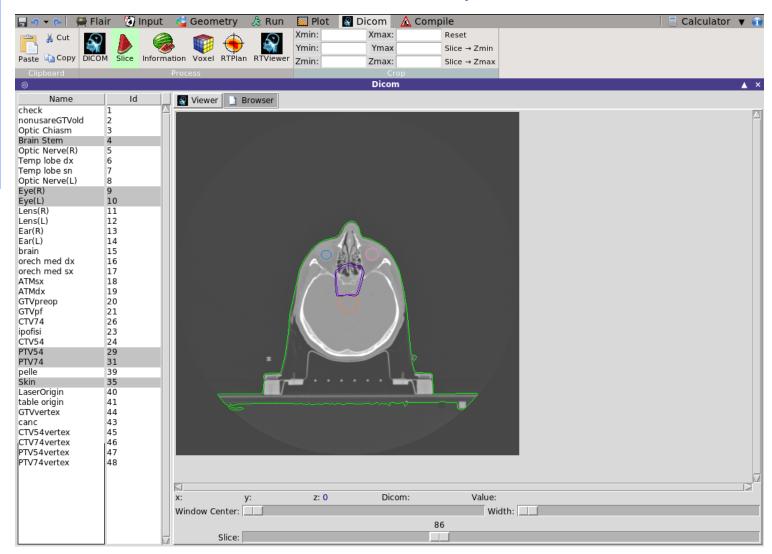
Parodi et al MP 34, 2007, Parodi et PMB 52, 2007

#### The FLUKA voxel with ROI (new release)

- The information for regions of interest (ROIs) and points of interests (eg. dose reference points) are included in RTSTRUCT file
- ROIs are represented as the points belonging to a closed polygon using 2D coordinates (not rounded to the pixel size of the corresponding CT image).
- User can embedded RTSTRUCT file into the VOXEL
  - For each voxel Flair identifies list of ROIs and creates ROI to voxel correspondence matrix
  - Voxel with RT STRUCT information can be currently used as an input for DVH calculations, however simulations will be available with the next FLUKA release
  - Flair provides some checks on the structures, like calculating volumes using the true polygonal information or the discretiza- tion to voxels.
     Differences up to few percent can be visible due to quantification process

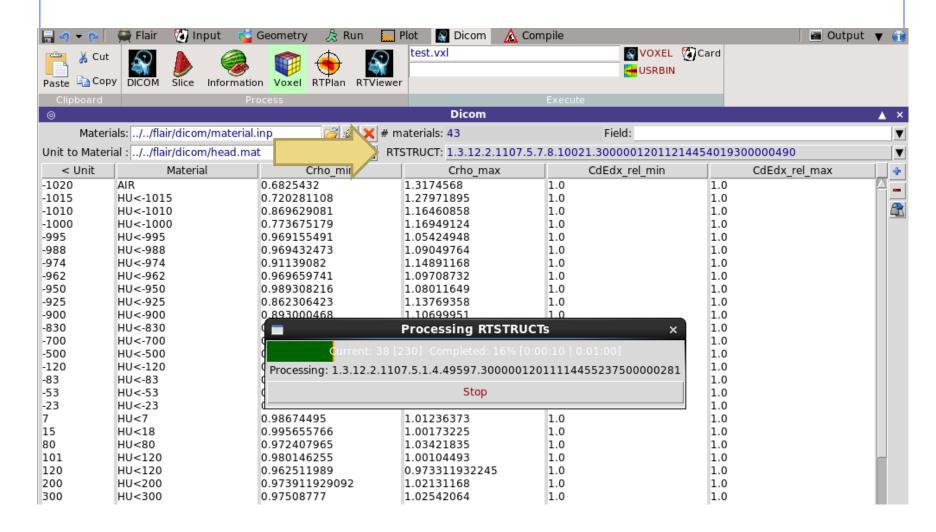
#### Processing the **DICOM** files with FLAIR VII

Select RTSTRUCT "Data sets" and inspect the ROI's



#### Processing the **DICOM** files with FLAIR VIII

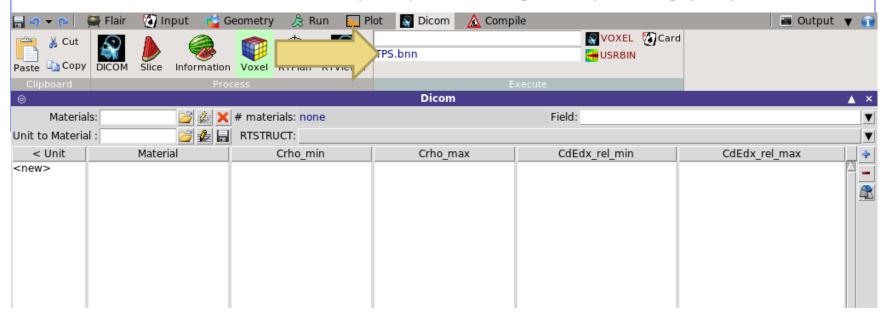
User can include ROIs imported from RTSTRUCT. Selection of corresponding file is done in VoxelTab.



#### Processing the **DICOM** files with FLAIR IX

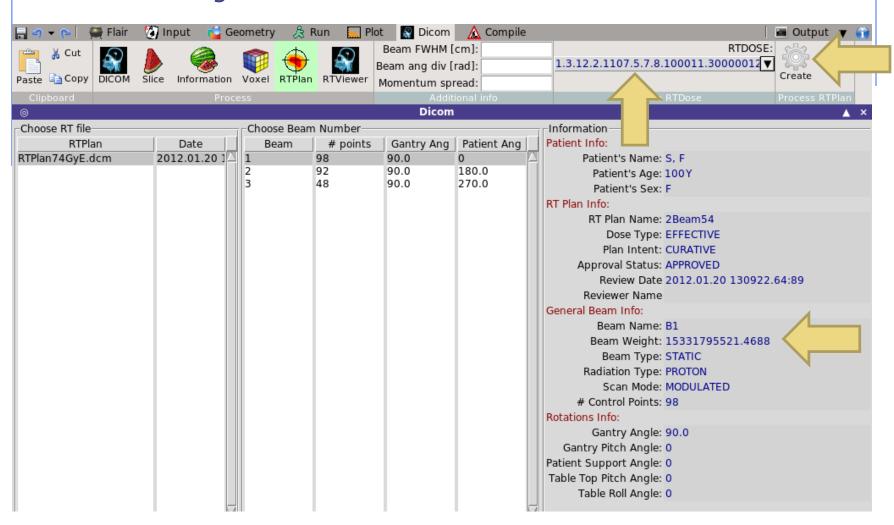
The "VoxelTab" is also for converting the RTDOSE dataset to FLUKA understandable format – USRBIN.

This can be further used for postprocessing and plotting purposes.



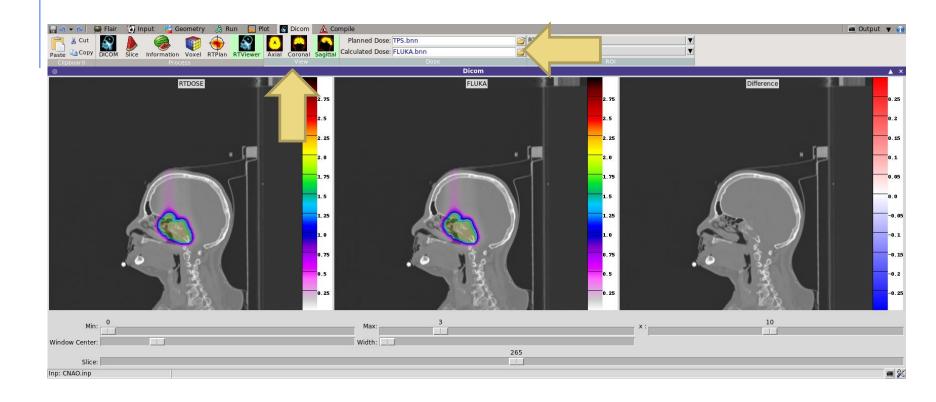
#### Processing the **DICOM** files with FLAIR X

The "RTPlan" tab is used to modify input file for Treatment Plan Simulations using data from RTPLAN and RTDOSE



#### Processing the **DICOM** files with FLAIR XI

- The "RTViewer" tab provides graphical comparison of TPS and FLUKA calculations using USRBIN files
- Available views for Axial, Coronal and Sagittal planes



### Processing the **DICOM** files with FLAIR XII

FLAIR
allows to
Dose-Volume
plots

postproce of the creation of t

1

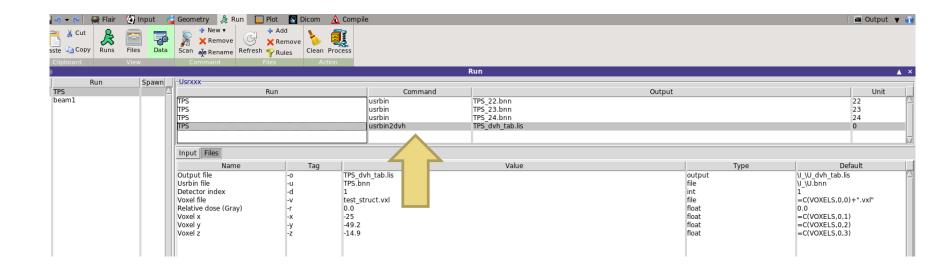
1.5

2

2.5

0.5

TPS dvh plot



10

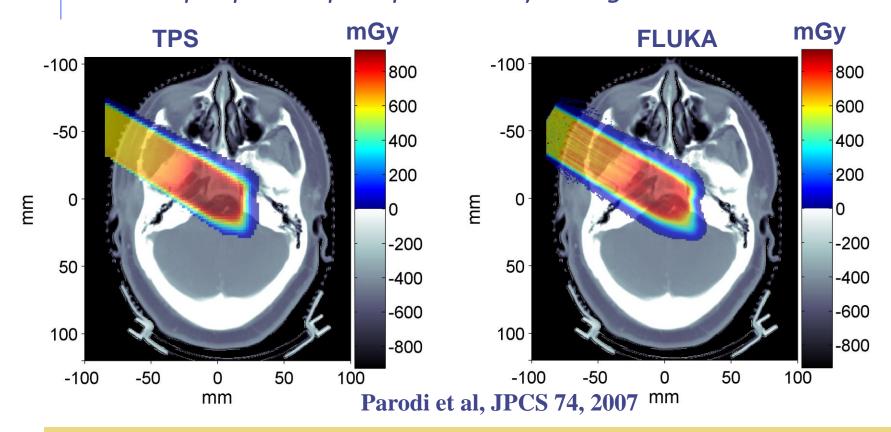
0.526245, 108.930

#### Few remarks

- Updated input file from RTPlan Tab uses #include cards in order to import pencil beam parameters.
   Due to typical vast number of described beams user needs to activate/deactivate these cards .
- Simulated and processed FLUKA files have to be weighted according to the information from RTPlan usrwei routine available
- DVH plots requires \*.vxl file with RTSTRUCT information.
- In RTViewer graphical comparison is available only for the \*.bnn files with the same binnings and positions.
   Editing input file by RTPlan Tab allows to provide correct parameters – user chooses corresponding RTDOSE file.

### Applications of FLUKA to p therapy @ MGH

Input phase-space provided by H. Paganetti, MGH Boston

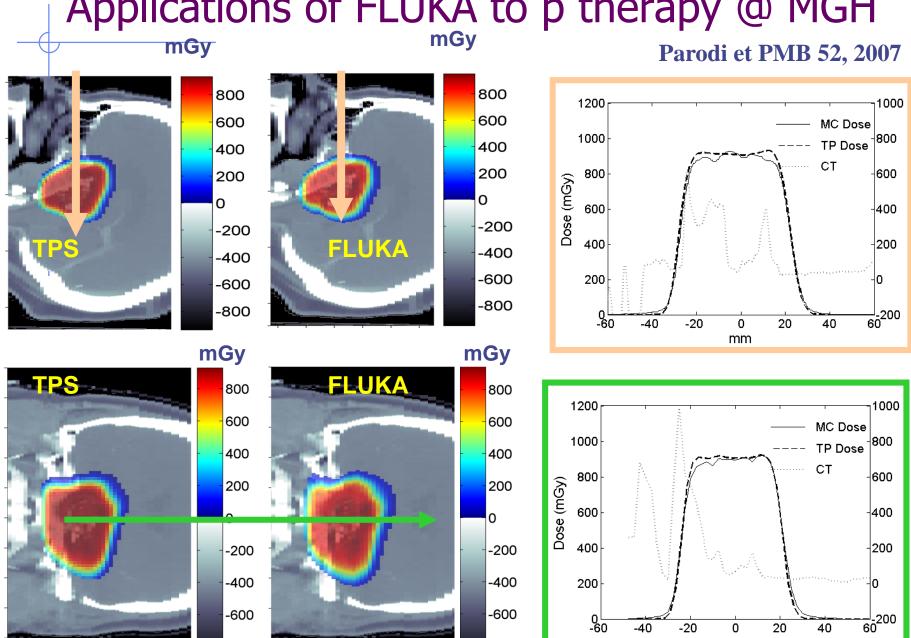


Prescribed dose: 1 GyE

 $MC: \sim 5.5 \cdot 10^6$  protons in 10 independent runs

(11h each on Linux Cluster mostly using 2.2GHz Athlon processors)

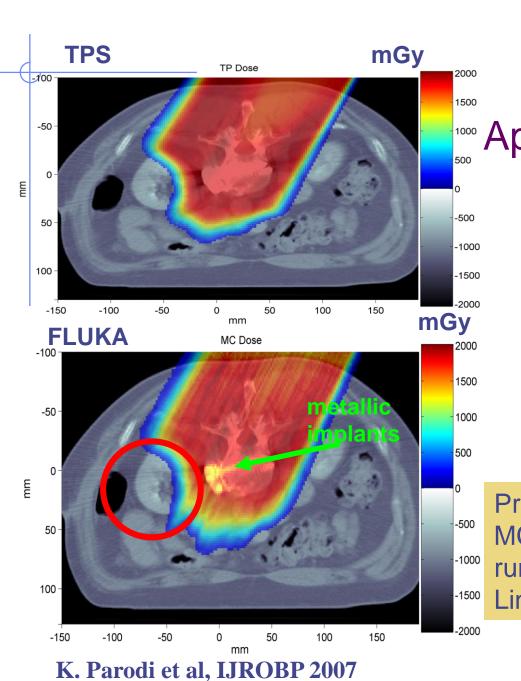
# Applications of FLUKA to p therapy @ MGH



-800

-800

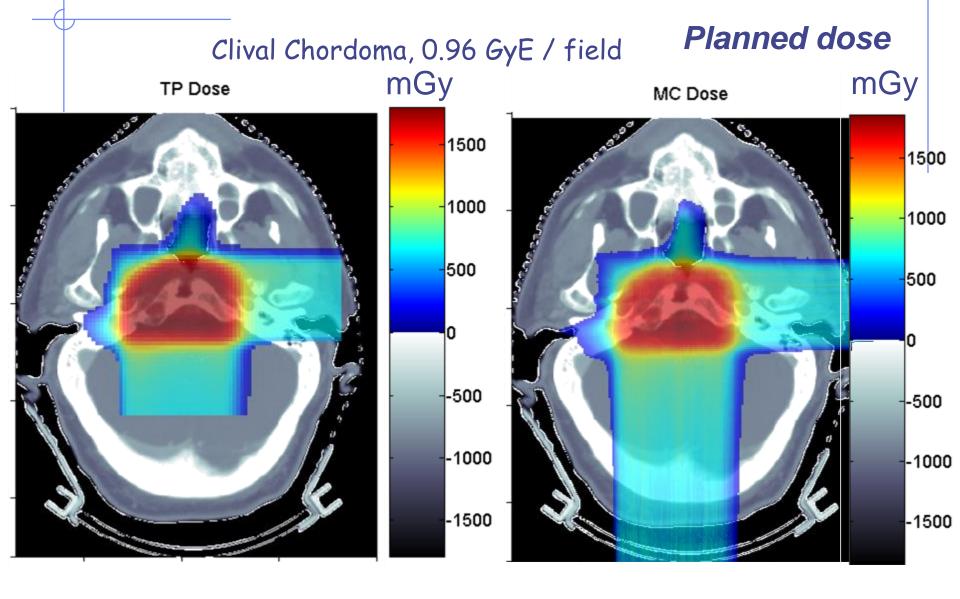
mm



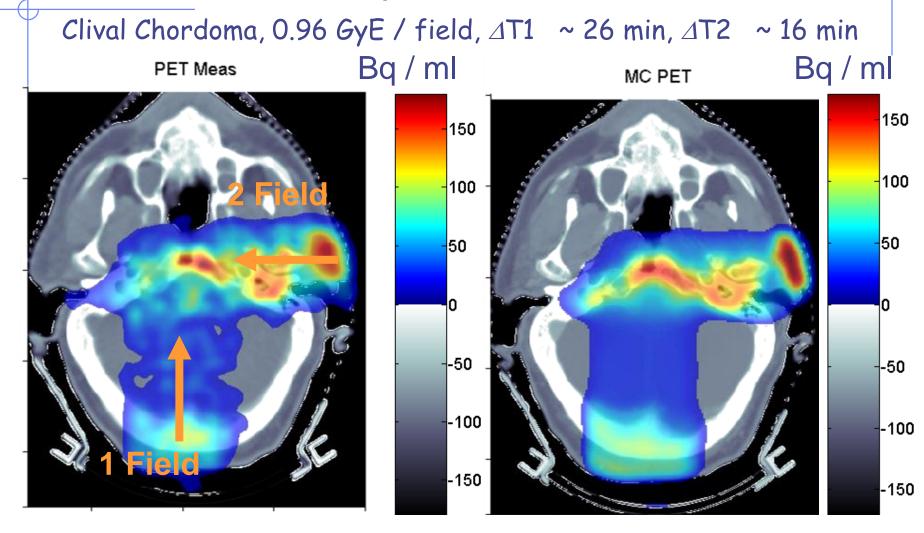
## Applications of FLUKA to p therapy @ MGH

Prescribed dose: 2 GyE MC:  $\sim 7.4 \ 10^7 p$  in 12 independent runs (~ 130h each on 2.2 GHz Linux cluster)

## Applications of FLUKA to p therapy @ MGH



# Post-radiation PET/CT @ MGH Average Activity



K. Parodi et al, IJROBP 2007

... and FLUKA-voxel functionalities being also used at HIT and CNAO ...

#### Additional material – user routine

- In current version user needs to compile special routine for defining pencil beam source parameters.
- In the next FLUKA release new cards will allow user to simulate RTPLAN without external routines.

