



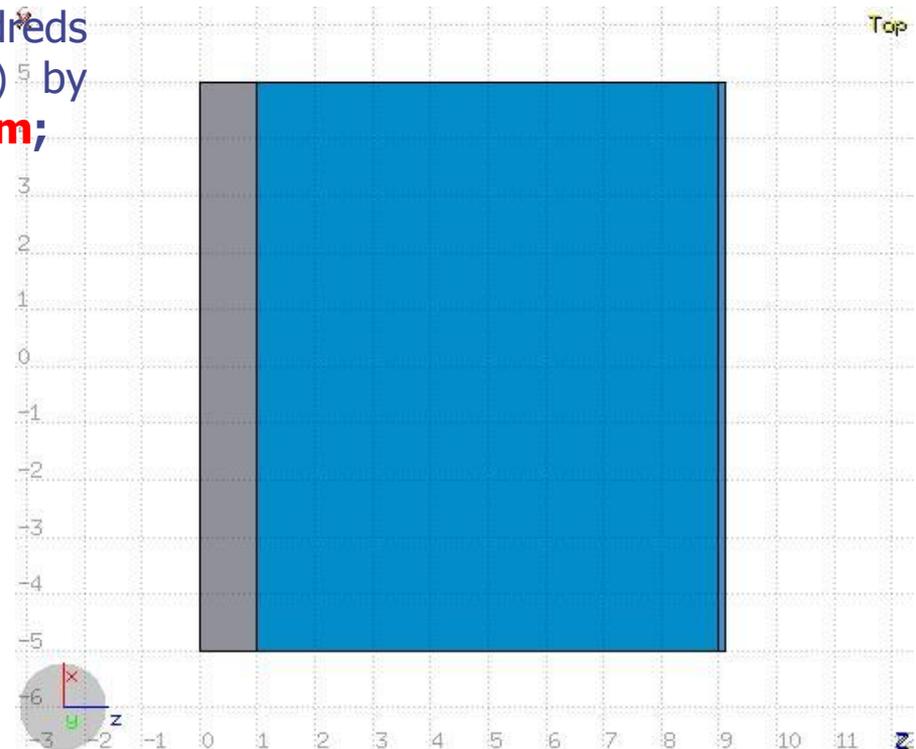
Exercise 7: Low Energy Neutrons

Beginners' FLUKA Course

Exercise: Low Energy Neutrons

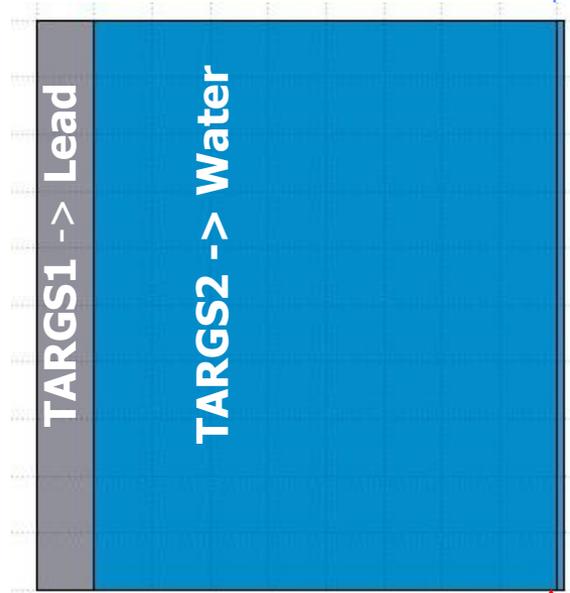
GOAL: Simulate the effect of the thermal neutron cutoff of a thin Cd foil.

- Create a folder called **ex7** and start there a new flair project based on the **course** template (as before)
- Save the input as **ex7.inp**
- Modify the geometry:
 - Increase the **TARGS2** dimensions by moving the **T2seg** plane to **z=9 cm**
 - Thin out the **TARGS3** to few hundreds of microns (start with 100 um) by moving plane **ZThigh** to **z=9.01 cm**;



Exercise: Low Energy Neutrons

- Change the materials as follows:
 - **TARGS1** -> **Lead**
 - **TARGS2** -> **Water**
 - **TARGS3** -> **Iron/Cadmium**
(between **#if ... #else ... #endif** statement)
- **Note: Cadmium** is not a **FLUKA** default element so you must define the material card first (you can try to use flair to add it)



```

END
GEOEND
Materials definition
-----

define Cadmium
MATERIAL                                Name: CADMIUM                #                p: 8.65
    Z: 48.                               Am:                       A:                dE/dx:
define a flag
#define flagIRON                          :
if "flagIRON" flag is set assign "iron" to the TARGS3 region
if not assign Cadmium instead
#if flagIRON
ASSIGNMA                                Mat: IRON                    Reg: TARGS3        to Reg:
    Mat(Decay):                          Step:                  Field:
#else
ASSIGNMA                                Mat: CADMIUM                 Reg: TARGS3        to Reg:
    Mat(Decay):                          Step:                  Field:
#endif
ASSIGNMA                                Mat: BLKHOLE                 Reg: BLKHOLE       to Reg:
    
```

TARGS2 -> Iron/Cadmium

Exercise: Low Energy Neutrons

- Add a boundary crossing estimator to score the neutron *fluence* from the thin foil region **TARGS3** to the region **INAIR**:
 - Use logarithmic energy binning down to the group of lowest energy
 - Write the output unformatted to unit 56
- Run for Iron and Cadmium 5 cycles of 20000 primaries and plot the results as a lethargy spectrum (x-axis: GeoMean, y-axis: $Y^* < X_{geo} >$, both axis logarithmically)
- For the **Iron** case: Identify the peak in thermal part of spectrum (note the automatic matching of neutron group structure)
- Compare the results with the **Cadmium** case

Note: not to overwrite the results when running the second time you can create two runs in flair and run them independently

