

A³MCNP Computer Exercises

by

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Using A³MCNP

Here, we will utilize the simplified VENUS problem to demonstrate the A³MCNP code.

- Copy the necessary input files

```
cp /usr/local/workshop/a3/cpa3mcpn .  
./cpa3mcpn  
cd a3
```

- Discussion on the new input information for A³MCNP

```
vi venus.a31  
vi venus.a32  
vi venus.a33  
vi sailor96_zaid.in (given as part of the code)
```

COMPUTER EXERCISE #1

1. Run A³MCNP

```
runa3_1 venus sailor96  
runa3_2 venus  
runa3_3 venus
```

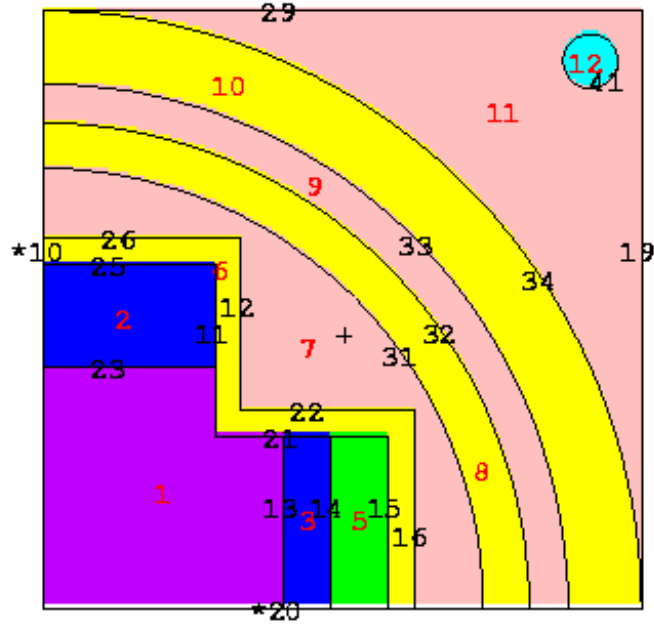
2. Run MCNP4C2

```
run3_0 venus
```

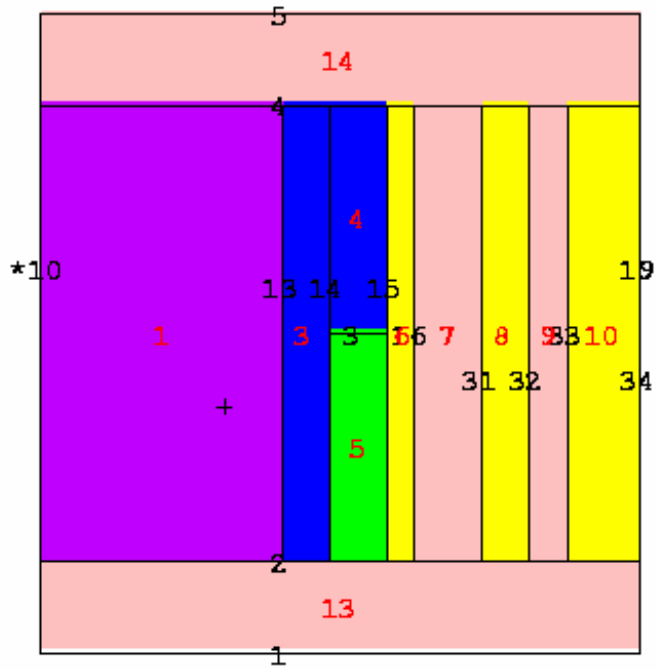
3. Compare the results

```
vi venus_3o (A3MCNP output)  
vi venus_0o (MCNP4C2 output)
```

SIMPLIFIED VENUS-3 MODEL



15



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file: venus

message: DATAPATH=/usr/mcnpxs

TITLE: Simplified VENUS-3 Model

```
c
c      10th International Training Course/Workshop on
c      Methodologies for Particle Transport Simulation...
c
c CELL CARDS
c core region
1  1 7.16E-02 (2 -4 10 -13 20 -21)
                : (2 -4 10 -11 21 -23)          $ fuel1
2  2 7.16E-02 (2 -4 10 -11 23 -25)          $ fuel2a
3  2 7.16E-02 (2 -4 13 -14 20 -21)          $ fuel2b
4  2 7.16E-02 (3 -4 14 -15 20 -21)          $ fuel2c (above PLSA)
5  3 9.34E-02 (2 -3 14 -15 20 -21)          $ PLSA
6  4 -7.90    (2 -4 10 -12 25 -26)          $ baffle1
                : (2 -4 11 -12 21 -25)          $ baffle2
                : (2 -4 12 -15 21 -22)          $ baffle3
                : (2 -4 15 -16 20 -22)          $ baffle4
7  5 -1.00    (2 -4 10 -12 26 -31)          $ water (baffle-barrel)
                : (2 -4      12 22 -31)
                : (2 -4 16 -31 20 -22)
8  4 -7.90    2 -4 10 20 31 -32            $ barrel
9  5 -1.00    2 -4 10 20 32 -33            $ water (barrel-vessel)
10 4 -7.90    2 -4 10 20 33 -34            $ vessel
11 5 -1.00    2 -4 10 20 34 -19 -29 #12    $ outside vessel
12 6 -2.70    2 -4 -41                    $ dosimeter
13 5 -1.00    1 -2 10 -19 20 -29          $ bottom reflector
14 5 -1.00    4 -5 10 -19 20 -29          $ top reflector
15 0          (-1:5:-10:19:-20:29)        $ outside
```

```
c
c SURFACE CARDS
c
c NOTE: * indicates reflective boundaries
```

```
c top and bottom model boundaries
1 pz 95.00      $ bottom of model
2 pz 105.00     $ fuel bottom
3 pz 130.00     $ fuel midplane
4 pz 155.00     $ fuel top
5 pz 165.00     $ top of model
```

```
c symmetry boundaries/outer boundaries
*10 px 0.0
19 px 65.573
*20 py 0.0
29 py 65.573
```

```
c planes (normal to x-axis)
11 px 18.90     $ inner square/baffle in
12 px 21.758    $ inner square/baffle out
13 px 26.46     $ fuel1/fuel2 boundary
14 px 31.50     $ fuel/PLSA boundary
15 px 37.80     $ core/baffle boundary
16 px 40.658    $ baffle outer boundary
```

```
c planes (normal to y-axis)
21 py 18.90     $ inner square/baffle in
22 py 21.758    $ inner square/baffle out
23 py 26.46     $ fuel1/fuel2 boundary
25 py 37.80     $ core/baffle boundary
26 py 40.658    $ baffle outer boundary
```

```
c outside of core
31 cz 48.283    $ barrel IR
32 cz 53.273    $ barrel OR
33 cz 57.545    $ vessel IR
```

```

34  cz      65.573      $ vessel OR
c
c  cavity dosimeter at 45 deg
41  c/z    60.0 60.0 3.00 $ (c/z - cylinder parallel to z-axis)
c  segmenting planes
42  pz 115
43  pz 125
44  pz 135
45  pz 145.

c
c  DATA CARDS
mode n          $ neutron only
imp:n 1 13r 0   $ set all cell importances to unity
c
c  TALLY CARDS
c  default energy bins (SAILOR/BUGLE group structure) for tallying
e0  1.003E+00  1.353E+00  1.653E+00  1.920E+00  2.231E+00  2.346E+00
    2.365E+00  2.466E+00  2.725E+00  3.012E+00  3.679E+00  4.966E+00
    6.065E+00  7.408E+00  8.607E+00  1.000E+01  1.221E+01  1.419E+01
    1.733E+01
c
f14:n (1 2 3 4)
fc14  volume flux in the core
fm14  1.271E+13 $ Source normalization factor
c
f24:n 10
fc24  volume flux in vessel
sd24  38814.0 $ volume of pv (1/4)*(Ro^2 - Ri^2)*pi*h
fm24  1.271E+13 $ Source normalization factor
c
f32:n 33
fc32  surface flux, vessel inner surface
sd32  4519.57 $ inner area
fm32  1.271E+13 $ Source normalization factor
c
f42:n 34
fc42  surface flux, vessel outer surface
sd42  5150.09 $ outer area
fm42  1.271E+13 $ Source normalization factor
c
f104:n 12
fc104 volume flux in cavity dosimeter
fq104 e s
fm104 1.271E+13 $ Source normalization factor
c
f114:n 12
fc114 Cu(n,a) reaction rate (BUGLE-96) in cavity dosimeter
fm114 1.271E+13 $ Source normalization factor
c  Note: response cross sections for the Cu(n,a) reaction are from
c  the BUGLE-96 multigroup library (Ref: RSICC DLC-185)
c  energy multiplier card
em114 0.0000E-00  1.6179E-08  2.4778E-07  6.7643E-07  8.1765E-07  9.4358E-07
      2.9883E-06  8.2245E-06  4.8528E-05  5.9165E-04  3.3924E-03  1.0490E-02
      1.8464E-02  2.7224E-02  3.7670E-02  4.3444E-02  3.2358E-02
e114  1.653E+00  1.920E+00  2.231E+00  2.346E+00  2.365E+00  2.466E+00
      2.725E+00  3.012E+00  3.679E+00  4.966E+00  6.065E+00  7.408E+00
      8.607E+00  1.000E+01  1.221E+01  1.419E+01  1.733E+01
c
f124:n 12
fc124 Cu(n,a) reaction rate (CE ENDF/B-VI) in cavity dosimeter
c  Note: response cross sections for the Cu(n,a) reaction are from
c  the MCNP continuous energy (CE) ENDF/B-VI library (Ref: RSICC DLC-189)
c  energy multiplier card
fm124 1.271E+13 $ Source normalization factor
c  material 100, reaction type 107 (n,alpha)
c
c
c
c
c  MATERIAL CARDS
c

```

```

c   core mixture          atom density = 7.16E-02
m1  1001.60c 4.19E-01    8016.60c 4.20E-01    5010.60c 4.19E-05
    92235.60c 2.79E-03    92238.60c 9.78E-02    40000.60c 5.59E-02
    24050.60c 2.43E-05    24052.60c 4.68E-04    24053.60c 5.31E-05
    24054.60c 1.32E-05    26054.60c 3.30E-05    26056.60c 5.13E-04
    26057.60c 1.17E-04    26058.60c 1.57E-06    28058.60c 2.86E-03
    28060.60c 1.09E-03    28061.60c 4.73E-05    28062.60c 1.50E-04
    28064.60c 3.81E-05

c
c   core mixture          atom density = 7.16E-02
m2  1001.60c 4.19E-01    8016.60c 4.20E-01    5010.60c 4.19E-05
    92235.60c 2.79E-03    92238.60c 9.78E-02    40000.60c 5.59E-02
    24050.60c 2.43E-05    24052.60c 4.68E-04    24053.60c 5.31E-05
    24054.60c 1.32E-05    26054.60c 3.30E-05    26056.60c 5.13E-04
    26057.60c 1.17E-04    26058.60c 1.57E-06    28058.60c 2.86E-03
    28060.60c 1.09E-03    28061.60c 4.73E-05    28062.60c 1.50E-04
    28064.60c 3.81E-05

c
c   PLSA                  atom density = 7.16E-02
m3  1001.60c 4.19E-01    8016.60c 4.20E-01    5010.60c 4.19E-05
    92235.60c 2.79E-03    92238.60c 9.78E-02    40000.60c 5.59E-02
    24050.60c 2.43E-05    24052.60c 4.68E-04    24053.60c 5.31E-05
    24054.60c 1.32E-05    26054.60c 3.30E-05    26056.60c 5.13E-04
    26057.60c 1.17E-04    26058.60c 1.57E-06    28058.60c 2.86E-03
    28060.60c 1.09E-03    28061.60c 4.73E-05    28062.60c 1.50E-04
    28064.60c 3.81E-05

c
c   stainless steel       atom density = 8.69E-02
m4  6000.60c 2.73E-03    14000.60c 1.03E-02    25055.60c 1.75E-02
    24050.60c 8.70E-03    24052.60c 1.68E-01    24053.60c 1.90E-02
    24054.60c 4.73E-03    26054.60c 3.96E-02    26056.60c 6.15E-01
    26057.60c 1.41E-02    26058.60c 1.88E-03    28058.60c 6.72E-02
    28060.60c 2.57E-02    28061.60c 1.11E-03    28062.60c 3.53E-03
    28064.60c 8.95E-04

c
c   water                 atom density =
m5  1001.60c 6.67E-01    8016.60c 3.33E-01    5010.60c 8.00E-05

c
c   aluminum              atom density = 6.02E-02
m6  13027.60c 1.00E+00

c
c   response cross section for Cu-63(n,alpha) from CE ENOV/B-VI
m100 29063.60c 1.0

c
c   nonu      $ fission in all cells is treated as capture

c
c   neutron histories are terminated if their energy falls below 5 MeV
cut:n   j 5.00
c print 110 $ print table 110 - first 50 histories
print

c
c   void      $ void material regions
ctme 02      $ time cutoff card (minutes)
c   nps 100000 $ follow 1E+5 particle histories
c

```

file: venus.sdef

```
c
c SOURCE CARDS
sdef erg=d1 pos=d2
c
c SAILOR/BUGLE 47-group structure
sil h 1.000e-11 1.000e-7 4.140e-7 8.764e-7 1.855e-6 5.043e-6
1.068e-5 3.727e-5 1.013e-4 2.144e-4 4.540e-4 1.585e-3
3.355e-3 7.102e-3 1.503e-2 2.188e-2 2.418e-2 2.606e-2
3.183e-2 4.087e-2 6.738e-2 1.111e-1 1.832e-1 2.972e-1
3.688e-1 4.979e-1 6.081e-1 7.427e-1 8.208e-1 1.003e+0
1.353e+0 1.653e+0 1.920e+0 2.231e+0 2.346e+0 2.365e+0
2.466e+0 2.725e+0 3.012e+0 3.679e+0 4.966e+0 6.065e+0
7.408e+0 8.607e+0 1.000e+1 1.221e+1 1.419e+1 1.733e+1
c typical PWR fission spectrum
spl d 0 1.3464e-11 1.0080e-10 2.3789e-10 7.3247e-10 3.7774e-09 1.0123e-08
8.2702e-08 3.4003e-07 9.1003e-07 2.8049e-06 2.2923e-05 5.6253e-05
1.7296e-04 5.3045e-04 5.9161e-04 2.2158e-04 1.8892e-04 6.2068e-04
1.0843e-03 3.8297e-03 7.9326e-03 1.6195e-02 3.0840e-02 2.1559e-02
4.1541e-02 3.7103e-02 4.6232e-02 2.6923e-02 6.2179e-02 1.1339e-01
8.8005e-02 7.0140e-02 7.1859e-02 2.3970e-02 3.8290e-03 1.9742e-02
4.6115e-02 4.4058e-02 7.8188e-02 8.3999e-02 3.2243e-02 1.6918e-02
5.9238e-03 2.6238e-03 1.1094e-03 1.8428e-04 4.2592e-05
c
c source point positions
si2 1
6.93 .63 106.75 8.19 .63 106.75 9.45 .63 106.75
10.71 .63 106.75 11.97 .63 106.75 13.23 .63 106.75
14.49 .63 106.75 15.75 .63 106.75 17.01 .63 106.75
18.27 .63 106.75 19.53 .63 106.75 20.79 .63 106.75
22.05 .63 106.75 23.31 .63 106.75 24.57 .63 106.75
25.83 .63 106.75 27.09 .63 106.75 28.35 .63 106.75
29.61 .63 106.75 30.87 .63 106.75 6.93 1.89 106.75
...
...
c
c source point probabilities
sp2 d
.7865 .8719 .8842 .8914 .8886 .9096 .9168 .9309 .9331
.8850 .8751 .8599 .8356 .7946 .7575 .7409 .5792 .5665
.5511 .5350 .7770 .8658 .8880 .8842 .8985 .9163 .9320
.9331 .8822 .8698 .8552 .8314 .7940 .7528 .7441 .5806
.5657 .5451 .5321 .7675 .8658 .8866 .8946 .8955 .8998
.9125 .9344 .9248 .8774 .8632 .8446 .8157 .7840 .7372
.7424 .5821 .5630 .5392 .5292 .7491 .8708 .8857 .8946
...
...
```

file: sailor96_zaid.in

```
142      = mtp
1001    3  This file corresponds to the SAILOR96 library
5010    7
40000   27
92235   43
92238   47
 8016   51
24000   63
24050   63
24052   63
24053   63
24054   63
25055   67
26000   71
26054   71
26056   71
26057   71
26058   71
28000   75
28058   75
28060   75
28061   75
28062   75
28064   75
 6000   79
11023   95
12000   99
13027  103
14000  107
19000  111
20000  115
```

```
t
13
 1  2  3  4  5  6  7  8  9 10 11 12 13 14 15
16 17 18 18 20 21 22 23 24 25 26 27 28 29 30
31 32 33 34 35 36 37 38 39 40 41 42 43 44 45
46 47 48 49 50 51 52 53 54 55 56 57 58 59 60
61 62 63 64 65 66 67 68 69 70 71 72 73 74 75
76 77 78 79 80 81 82 83 84 85 86 87 88 89 90
91 92 93 94 95 96 97 98 99 100 101 102 103 104 105
106 107 108 109 110 111 112 113 114 115 116 117 118 119 120
121 122 123 124 125 126 127 128 129 130 131 132 133 134 135
136 137 138 139 140 141 142
```


file: venus.a31

```
c
c ----- A3MCNP parameters (step 1, Sn mesh generation)
c
c      isn isrcc igm nsctm iht ihm iups neut
sngrp   2  12  47  3  3  70  0  47
c
c      Sn spatial mesh input
c      x1  x2  y1  y2      z1      z2      xt  yt  zt pzlev
snmsh   0.   64.  0.   64.   95.    165.0  3.2  3.2  10.  -7
c
c      back-thinning parameters for different materials
snthn 3.2 3.2 3.2 3.2 3.2 3.2 3.2
c
c      Sn boundary condition - snbc()=(ibl,ibr,ibi,ibo,ibb,ibt)
c      ibl ibr  ibi  ibo  ibb  ibt
snbc    1  0   1  0   0  0
c
c      Sn energy group structure
c      SAILOR 47-group neutron structure
snsi    1.0e-7 4.14e-7 8.764e-7 1.855e-6 5.043e-6 1.068e-5
        3.727e-5 1.013e-4 2.144e-4 4.454e-4 1.585e-3 3.355e-3
        7.102e-3 1.503e-2 2.188e-2 2.418e-2 2.606e-2 3.183e-2
        4.087e-2 6.738e-2 1.111e-1 1.832e-1 2.972e-1 3.688e-1
        4.979e-1 6.081e-1 7.427e-1 8.208e-1 1.003 1.353 1.653
        1.920 2.231 2.346 2.365 2.466 2.725 3.012 3.679 4.966
        6.065 7.408 8.607 1.000e+1 1.221e+1 1.419e+1 1.733e+1
snsp    0.0 39r
        .003 .01 .012 .018 .02 .04 .03
snacc 2 2
        6.4 32.0 $coarse-mesh interfaces along x
        6.4 32.0 $coarse-mesh interfaces along y
```

file: venus.a32

```
c
c ----- A3MCNP parameters (step 2, variance reduction parameters generation)
c
c      wwp:n 5 3 5
c      wwa:n 47 0 1
```

file: venus.a33

```
c
c ----- A3MCNP parameters (step 3, accelerated simulation)
c
c      wwp:n 5 3 5
c      wwa:n 47 1 1
```

A³MCNP COMPUTER EXERCISE #2

Reduce the size of the detector (adjoint source)

1. Edit "venus" to divide Cell 12 into 2 cells: 12 and 121

vi venus

Before:

```
11 5 -1.00 2 -4 10 20 34 -19 -29 #12 $ outside vessel
12 6 -2.70 2 -4 -41 $ dosimeter
```

After:

```
11 5 -1.00 2 -4 10 20 34 -19 -29 #12 #121 $ outside vessel
12 6 -2.70 45 -4 -41 $ dosimeter
121 6 -2.70 2 -45 -41 $ dosimeter 2
```

2. Modify cell importance values

Before:

```
imp:n 1 13r 0 $ set all cell importances to unity
```

After:

```
imp:n 1 14r 0 $ set all cell importances to unity
```

3. Run MCNP4C and A³MCNP

4. Examine output files

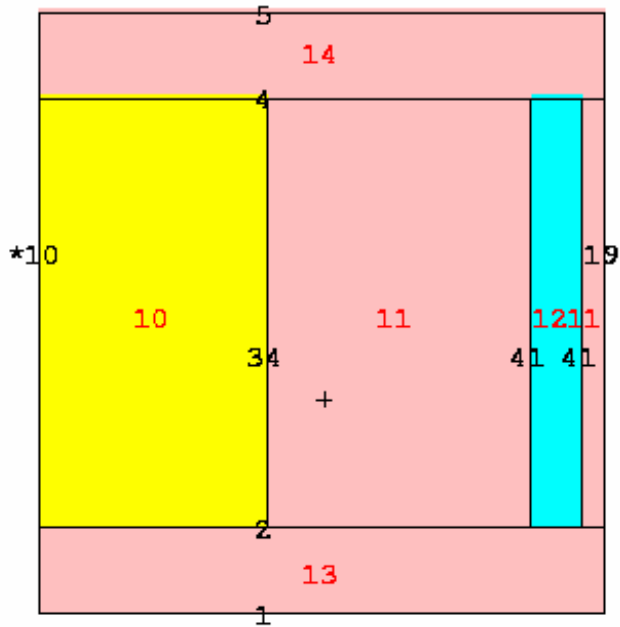
vi venus_0o

vi venus_3o

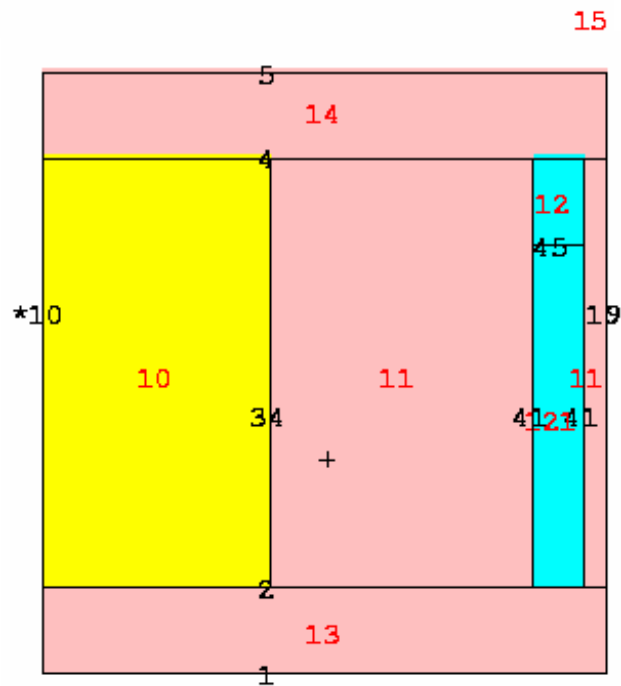
SIMPLIFIED VENUS-3 MODEL

Exercise #2

Before



After



15

A³MCNP COMPUTER EXERCISE #3

Coarsen fine meshes for S_N adjoint calculation using the back-thinning approach

1. Edit “venus.a31” to apply the back-thinning option

```
vi venus.a31
```

Before:

```
c    isn isrcc igm nsctm iht ihm iups neut  
sngp  2 12 47  3    3 70 0  47
```

After:

```
c    isn isrcc igm nsctm iht ihm iups neut  
sngp  1 12 47  3    3 70 0  47
```

2. Run A³MCNP
Observe the reduction of number of fine meshes

```
runa3_1 venus sailor96
```

```
runa3_2 venus
```

```
runa3_3 venus
```

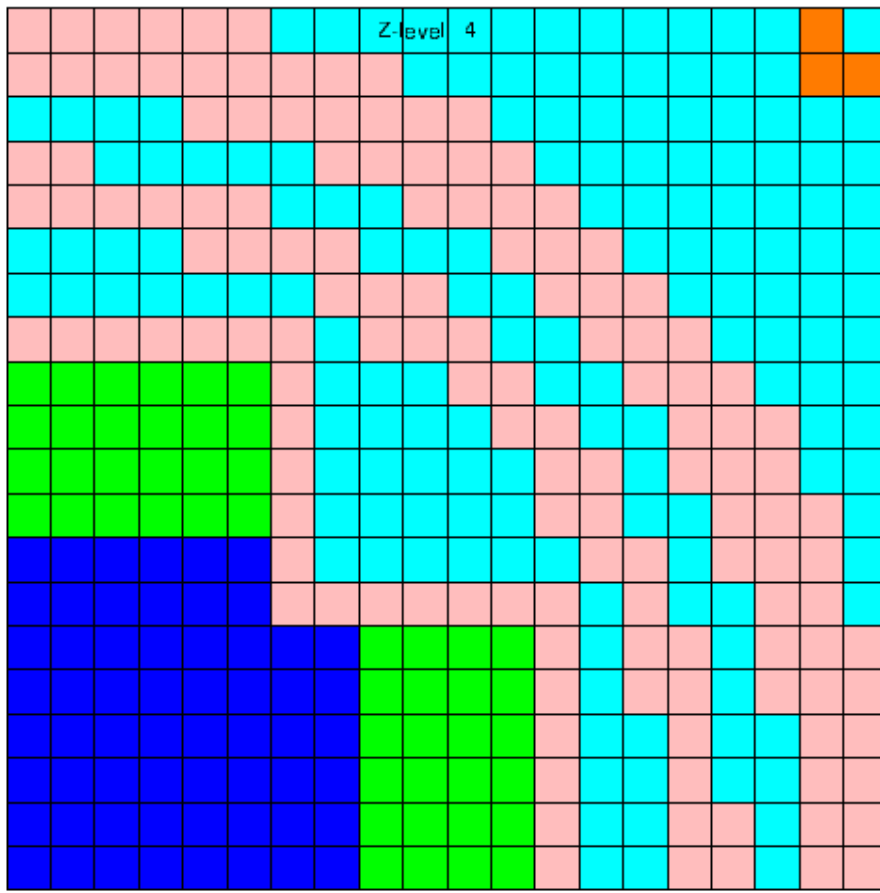
4. Examine output files
Look for tally fluctuation charts at the end of file
Compare results with and without back-thinning

```
vi venus_3o
```

SIMPLIFIED VENUS-3 MODEL

Exercise #3, uniform mesh distribution

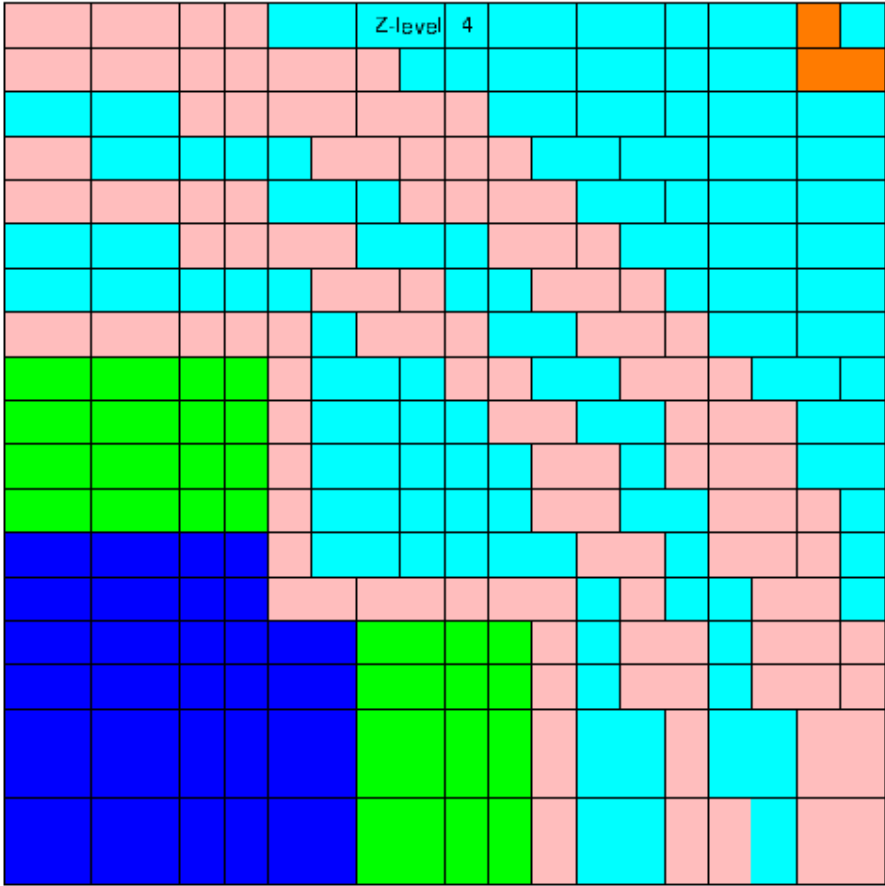
TORT Mesh Generation by A3MCNP



SIMPLIFIED VENUS-3 MODEL

Exercise #3, with back-thinning

TORT Mesh Generation by A3MCNP



original # of meshes = 2800
of meshes (after back-thinning) = 1488
reduction of 46.9%