
Example: Photon Treatment Plan using VMC++ dose calculation

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%%%

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%%%

In this example we will show (i) how to load patient data into matRad (ii) how to setup a photon dose calculation based on the VMC++ Monte Carlo algorithm (iii) how to inversely optimize the beamlet intensities directly from command window in MATLAB. (iv) how to visualize the result

Patient Data Import

Let's begin with a clear Matlab environment and import the boxphantom into your workspace.

```
clc,clear,close all;  
load('BOXPHANTOM.mat');
```

Treatment Plan

The next step is to define your treatment plan labeled as 'pln'. This structure requires input from the treatment planner and defines the most important cornerstones of your treatment plan.

```
pln.radiationMode = 'photons';  
pln.machine       = 'Generic';  
pln.bioOptimization = 'none';  
pln.gantryAngles  = [0];  
pln.couchAngles   = [0];  
pln.bixelWidth    = 10;
```

```
pln.numOfFractions = 30;  
pln.numOfBeams     = numel(pln.gantryAngles);  
pln.numOfVoxels    = prod(ct.cubeDim);  
pln.voxelDimensions = ct.cubeDim;  
pln.isoCenter      = ones(pln.numOfBeams,1) *  
    matRad_getIsoCenter(cst,ct,0);  
pln.runSequencing  = 0;  
pln.runDAO         = 0;
```

Generate Beam Geometry STF

```
stf = matRad_generateStf(ct,cst,pln);  
  
matRad: Generating stf struct... Progress: 100.00 %
```

Dose Calculation

Calculate dose influence matrix for unit pencil beam intensities using the VMC++ monte carlo algorithm. We define the number of photons simulated per beamlet to be 700. You can find compatible VMC++ files at <http://www.cerr.info/download.php> which have to be located in matRadrootDirectory\vmc++.

```
dij = matRad_calcPhotonDoseVmc(ct,stf,pln,cst);  
  
Warning: Number of photons simulated per bixel (nCasePerBixel) and  
        number  
of parallel MC simulations (numOfParallelMCSimulations) not specified  
by  
user. Use default settings with nCasePerBixel = 5000 and  
numOfParallelMCSimulations = 4 in vmc++ calculations.  
matRad: VMC++ photon dose calculation...
```

```
Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/  
MCpencilbeam_temp_1.vmc
```

```
Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/  
MCpencilbeam_temp_2.vmc
```

```
Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/  
MCpencilbeam_temp_3.vmc
```

```
Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/  
MCpencilbeam_temp_4.vmc  
Reading MS data ... OK  
Parsing input file ... construct_mmc_geometry: urs = 1  
Reading MS data ... OK  
Parsing input file ... construct_mmc_geometry: urs = 1
```

```

Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
=====
                          Beamlet Source
=====
      charge = 0
virtual source position = (-76,24,24)
      Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,22.5,23.5)
=====
XYZ Geometry
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
                          Monte Carlo Parameter
=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition   : 0
Cut-off energy for KERMA approx.    : 1.10239 MeV
Bremsstrahlung transport mode       : 1
CSDA approximation                  : 0

```

Example: Photon Treatment Plan
using VMC++ dose calculation

Fractional energy loss/step at E_p : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 18971 2927

=====

Variance Reduction

=====

f_{repeat} = 0.251
split photons = 1
photon split factor = -40

=====

Quasi Random Numbers

=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

Scoring and output options

=====

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.179 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK

=====

Beamlet Source

=====

charge = 0

Example: Photon Treatment Plan
using VMC++ dose calculation

```
virtual source position = (-76,24,24)
    Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,22.5,22.5)
=====
XYZ Geometry
=====
    name: CT id: 0
    global smax: 1e+030
    Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
    Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
    Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
Monte Carlo Parameter
=====
Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition   : 0
Cut-off energy for KERMA approx.    : 1.10239 MeV
Bremsstrahlung transport mode       : 1
CSDA approximation                  : 0
Fractional energy loss/step at Ep   : 10%
Max. 1st elastic moment per step    : 0.5
Max. acceptable energy loss/step    : 0.6 MeV
alpha and beta                      : 0.0298764 0.420741
Fano calculation                     : 0
Exact Compton                       : 0
Electron transport mode              : VMC++
=====
MC_Control
=====
will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 24442 27174
=====
Variance Reduction
=====
f_repeat = 0.251
split photons = 1
photon split factor = -40
=====
Quasi Random Numbers
=====
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```

number of generators: 1
      1st:  base = 2 dimensions = 60 warm-up = 1
=====
Scoring and output options
=====
number of dose scoring objects: 1
dose output options for geometry CT
  dump dose: 2
  dose scans:

CPU time so far: 2.24 seconds

Will run approximately 5000 particle sets
  with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK
=====
Beamlet Source
=====
  charge = 0
  virtual source position = (-76,24,24)
  Energy = 6 MeV
  number of beamlets = 1
  beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,22.5,24)
=====
XYZ Geometry
=====
  name: CT  id: 0
  global smax: 1e+030
  Number of x-planes: 160  uniform with Xo = 0.15 Dx = 0.3
  Number of y-planes: 160  uniform with Yo = 0.15 Dx = 0.3
  Number of z-planes: 160  uniform with Zo = 0.15 Dx = 0.3
=====
Monte Carlo Parameter
=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition    : 0
Cut-off energy for KERMA approx.     : 1.10239 MeV
Bremsstrahlung transport mode        : 1
CSDA approximation                  : 0
Fractional energy loss/step at Ep    : 10%
Max. 1st elastic moment per step     : 0.5
Max. acceptable energy loss/step     : 0.6 MeV
alpha and beta                       : 0.0298764 0.420741

```

Example: Photon Treatment Plan
using VMC++ dose calculation

```
Fano calculation           : 0
Exact Compton             : 0
Electron transport mode   : VMC++
```

```
=====
=====
```

MC_Control

```
=====
```

```
will use fixed number of histories
number of batches      : 10
histories per batch    : 500
total histories        : 5000
initial rng seeds      : 8355 16407
```

```
=====
```

```
=====
```

Variance Reduction

```
=====
```

```
f_repeat          = 0.251
split photons     = 1
photon split factor = -40
```

```
=====
```

```
=====
```

Quasi Random Numbers

```
=====
```

```
number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1
```

```
=====
```

```
=====
```

Scoring and output options

```
=====
```

```
number of dose scoring objects: 1
dose output options for geometry CT
  dump dose: 2
  dose scans:
```

CPU time so far: 2.338 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation

OK

```
=====
```

Beamlet Source

```
=====
```

```
charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,22.5,23)
```

Example: Photon Treatment Plan
using VMC++ dose calculation

=====
=====
XYZ Geometry
=====

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====

Monte Carlo Parameter
=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++
=====

=====
=====
MC_Control
=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 3810 27402
=====

Variance Reduction
=====

f_repeat = 0.251
split photons = 1
photon split factor = -40
=====

Quasi Random Numbers
=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1
=====

Scoring and output options

```
=====
number of dose scoring objects: 1
dose output options for geometry CT
  dump dose: 2
  dose scans:
```

CPU time so far: 2.458 seconds

Will run approximately 5000 particle sets
 with 1 particles per set on average
will use 2 quasi numbers to sample the source

```
Starting MC simulation
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.043  500  500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.045  500  500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.045  500  500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.045  500  500
+++finished batch 2 cpu time: 0.307  500  500
+++finished batch 2 cpu time: 0.364  500  500
+++finished batch 2 cpu time: 0.204  500  500
+++finished batch 2 cpu time: 0.144  500  500
+++finished batch 3 cpu time: 0.383  500  500
+++finished batch 3 cpu time: 0.44  500  500
+++finished batch 3 cpu time: 0.311  500  500
+++finished batch 3 cpu time: 0.215  500  500
+++finished batch 4 cpu time: 0.504  500  500
+++finished batch 4 cpu time: 0.449  500  500
+++finished batch 4 cpu time: 0.371  500  500
+++finished batch 4 cpu time: 0.285  500  500
+++finished batch 5 cpu time: 0.58  500  500
+++finished batch 5 cpu time: 0.533  500  500
+++finished batch 5 cpu time: 0.46  500  500
+++finished batch 6 cpu time: 0.655  500  500
+++finished batch 5 cpu time: 0.398  500  500
+++finished batch 6 cpu time: 0.619  500  500
+++finished batch 6 cpu time: 0.528  500  500
+++finished batch 7 cpu time: 0.718  500  500
+++finished batch 7 cpu time: 0.68  500  500
+++finished batch 6 cpu time: 0.473  500  500
+++finished batch 7 cpu time: 0.589  500  500
+++finished batch 8 cpu time: 0.784  500  500
+++finished batch 8 cpu time: 0.767  500  500
+++finished batch 7 cpu time: 0.569  500  500
+++finished batch 8 cpu time: 0.686  500  500
+++finished batch 9 cpu time: 0.877  500  500
+++finished batch 8 cpu time: 0.641  500  500
+++finished batch 9 cpu time: 0.861  500  500
```

```
+++finished batch 9 cpu time: 0.763 500 500
+++finished batch 10 cpu time: 0.955 500 500
finished simulation, cpu time = 0.982 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.982
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.939 500 500
finished simulation, cpu time = 0.952 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.952
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.856 500 500
+++finished batch 9 cpu time: 0.746 500 500
finished simulation, cpu time = 0.871 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.871
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.843 500 500
finished simulation, cpu time = 0.857 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.857
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
  max dose is 0.0342552 in region 1956847
  max dose is 0.0343312 in region 2033647
  max dose is 0.0343228 in region 1803246
  max dose is 0.0345088 in region 1880046
```

```
ICRU-2K efficiency: 849.507 1/s
+++ total
max dose is 0.0342552 in region 1956847
ICRU-2K efficiency: 849.507 1/s
=====
ICRU-2K efficiency: 858.861 1/s
+++ total
max dose is 0.0343228 in region 1803246
ICRU-2K efficiency: 858.861 1/s
=====
ICRU-2K efficiency: 927.296 1/s
+++ total
max dose is 0.0343312 in region 2033647
ICRU-2K efficiency: 987.342 1/s
+++ total
max dose is 0.0345088 in region 1880046
ICRU-2K efficiency: 927.296 1/s
=====
ICRU-2K efficiency: 987.342 1/s
=====
Completed 4 of 49 beamlets...
```

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_2.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_3.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_1.vmc

```
Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/  
MCpencilbeam_temp_4.vmc  
Reading MS data ... OK  
Parsing input file ... construct_mmc_geometry: urs = 1  
Reading MS data ... OK  
Parsing input file ... construct_mmc_geometry: urs = 1  
Reading MS data ... OK  
Parsing input file ... construct_mmc_geometry: urs = 1  
Reading MS data ... OK  
Parsing input file ... construct_mmc_geometry: urs = 1  
DE_GeometryFactory::get_region_size():  
get_region_size: urs = 1 ignore = 1  
region size: 0.3  
geometry 1 region size: 0.3  
DE_GeometryFactory::get_region_size():
```

```

get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
=====
                        Beamlet Source
=====
      charge = 0
virtual source position = (-76,24,24)
      Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,22.5,25)
=====
=====
                        XYZ Geometry
=====
      name: CT id: 0
      global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
                        Monte Carlo Parameter
=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition   : 0
Cut-off energy for KERMA approx.    : 1.10239 MeV
Bremsstrahlung transport mode       : 1
CSDA approximation                  : 0
Fractional energy loss/step at Ep   : 10%
Max. 1st elastic moment per step    : 0.5
Max. acceptable energy loss/step    : 0.6 MeV
alpha and beta                      : 0.0298764 0.420741
Fano calculation                    : 0
Exact Compton                      : 0
Electron transport mode              : VMC++
=====

```

Example: Photon Treatment Plan
using VMC++ dose calculation

```
=====
MC_Control
=====
will use fixed number of histories
number of batches      : 10
histories per batch    : 500
total histories        : 5000
initial rng seeds      : 4729 29118
=====
=====
Variance Reduction
=====

f_repeat              = 0.251
split photons         = 1
photon split factor   = -40
=====
=====
Quasi Random Numbers
=====
number of generators: 1
                    1st: base = 2 dimensions = 60 warm-up = 1
=====
=====
Scoring and output options
=====
number of dose scoring objects: 1
dose output options for geometry CT
    dump dose: 2
    dose scans:

CPU time so far: 2.371 seconds

Will run approximately 5000 particle sets
    with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.052 500 500
OK
=====
Beamlet Source
=====
charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,23,22.5)
=====
=====
XYZ Geometry
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====

Monte Carlo Parameter
=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

MC_Control
=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 24009 4257

=====

Variance Reduction
=====

f_repeat = 0.251
split photons = 1
photon split factor = -40

=====

Quasi Random Numbers
=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

Scoring and output options
=====

number of dose scoring objects: 1
```

Example: Photon Treatment Plan
using VMC++ dose calculation

dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.45 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK

=====

Beamlet Source

=====

charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,22.5,24.5)

=====

=====

XYZ Geometry

=====

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

=====

MC_Control

=====

Example: Photon Treatment Plan
using VMC++ dose calculation

```
=====
will use fixed number of histories
number of batches      : 10
histories per batch    : 500
total histories        : 5000
initial rng seeds      : 28726  28947
=====
=====
Variance Reduction
=====
=====

f_repeat              = 0.251
split photons         = 1
photon split factor   = -40
=====
=====
Quasi Random Numbers
=====
=====
number of generators: 1
                    1st: base = 2 dimensions = 60 warm-up = 1
=====
=====
Scoring and output options
=====
=====
number of dose scoring objects: 1
dose output options for geometry CT
    dump dose: 2
    dose scans:

CPU time so far: 2.571 seconds

Will run approximately 5000 particle sets
    with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK
=====
Beamlet Source
=====
=====
charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,22.5,25.5)
=====
=====
XYZ Geometry
=====
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
```


Example: Photon Treatment Plan
using VMC++ dose calculation

Number of y-planes: 160 uniform with $Y_0 = 0.15$ $Dx = 0.3$
Number of z-planes: 160 uniform with $Z_0 = 0.15$ $Dx = 0.3$

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at E_p : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 28716 14562

=====

Variance Reduction

=====

$f_{repeat} = 0.251$
split photons = 1
photon split factor = -40

=====

Quasi Random Numbers

=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

Scoring and output options

=====

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.584 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation

```
+++finished batch 2 cpu time: 0.13 500 500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.042 500 500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.042 500 500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.069 500 500
+++finished batch 3 cpu time: 0.305 500 500
+++finished batch 2 cpu time: 0.109 500 500
+++finished batch 2 cpu time: 0.226 500 500
+++finished batch 2 cpu time: 0.137 500 500
+++finished batch 4 cpu time: 0.375 500 500
+++finished batch 3 cpu time: 0.175 500 500
+++finished batch 3 cpu time: 0.305 500 500
+++finished batch 3 cpu time: 0.203 500 500
+++finished batch 5 cpu time: 0.444 500 500
+++finished batch 4 cpu time: 0.238 500 500
+++finished batch 4 cpu time: 0.374 500 500
+++finished batch 4 cpu time: 0.269 500 500
+++finished batch 5 cpu time: 0.301 500 500
+++finished batch 6 cpu time: 0.512 500 500
+++finished batch 5 cpu time: 0.439 500 500
+++finished batch 5 cpu time: 0.353 500 500
+++finished batch 7 cpu time: 0.578 500 500
+++finished batch 6 cpu time: 0.384 500 500
+++finished batch 6 cpu time: 0.514 500 500
+++finished batch 6 cpu time: 0.419 500 500
+++finished batch 8 cpu time: 0.642 500 500
+++finished batch 7 cpu time: 0.448 500 500
+++finished batch 7 cpu time: 0.595 500 500
+++finished batch 7 cpu time: 0.482 500 500
+++finished batch 9 cpu time: 0.704 500 500
+++finished batch 8 cpu time: 0.51 500 500
+++finished batch 8 cpu time: 0.665 500 500
+++finished batch 8 cpu time: 0.552 500 500
+++finished batch 10 cpu time: 0.769 500 500
finished simulation, cpu time = 0.787 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
geometry: CT
cpu time: 0.787
number of histories: 5000
```

```

number of batches: 10
+++ beamlet 0
+++finished batch 9 cpu time: 0.586 500 500
+++finished batch 9 cpu time: 0.615 500 500
+++finished batch 9 cpu time: 0.74 500 500
+++finished batch 10 cpu time: 0.648 500 500
finished simulation, cpu time = 0.662 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
geometry: CT
cpu time: 0.662
number of histories: 5000
number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.8 500 500
+++finished batch 10 cpu time: 0.684 500 500
finished simulation, cpu time = 0.697 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
geometry: CT
cpu time: 0.697
number of histories: 5000
number of batches: 10
+++ beamlet 0
finished simulation, cpu time = 0.834 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
geometry: CT
cpu time: 0.834
number of histories: 5000
number of batches: 10
+++ beamlet 0
max dose is 0.0341063 in region 2187246
max dose is 0.0346184 in region 2110447
max dose is 0.0345979 in region 2264047
max dose is 0.0338019 in region 1803726
ICRU-2K efficiency: 1039.73 1/s
+++ total
max dose is 0.0341063 in region 2187246
ICRU-2K efficiency: 1039.73 1/s
=====
ICRU-2K efficiency: 1279.31 1/s
+++ total
max dose is 0.0346184 in region 2110447
ICRU-2K efficiency: 1279.31 1/s

```

```
=====
ICRU-2K efficiency: 962.513 1/s
+++ total
max dose is 0.0338019 in region 1803726
ICRU-2K efficiency: 962.513 1/s
=====
ICRU-2K efficiency: 1206.82 1/s
+++ total
max dose is 0.0345979 in region 2264047
ICRU-2K efficiency: 1206.82 1/s
=====
Completed 8 of 49 beamlets...
```

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_1.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_2.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_4.vmc

```
Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_3.vmc
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```

region size: 0.3
geometry 1 region size: 0.3
OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
=====
                        Beamlet Source
=====
        charge = 0
    virtual source position = (-76,24,24)
        Energy = 6 MeV
    number of beamlets = 1
    beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,23,23.5)
=====
=====
                        XYZ Geometry
=====
        name: CT id: 0
    global smax: 1e+030
    Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
    Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
    Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
                        Monte Carlo Parameter
=====

Delta particle production threshold:    0.447479 MeV
Bremsstrahlung production threshold:    0.05 MeV
Min. electron transport energy          :    0.447479 MeV
Min. photon transport energy            :    0.05 MeV
Local track-end energy deposition        :    0
Cut-off energy for KERMA approx.         :    1.10239 MeV
Bremsstrahlung transport mode           :    1
CSDA approximation                     :    0
Fractional energy loss/step at Ep        :    10%
Max. 1st elastic moment per step         :    0.5
Max. acceptable energy loss/step         :    0.6 MeV
alpha and beta                          :    0.0298764  0.420741
Fano calculation                        :    0
Exact Compton                          :    0
Electron transport mode                  :    VMC++
=====
=====
                        MC_Control
=====
    will use fixed number of histories
    number of batches : 10
    histories per batch : 500
    total histories : 5000
    initial rng seeds : 23767 28785
=====

```

Example: Photon Treatment Plan
using VMC++ dose calculation

=====

Variance Reduction

=====

f_repeat = 0.251
split photons = 1
photon split factor = -40

=====

=====

Quasi Random Numbers

=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

=====

Scoring and output options

=====

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.277 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK

=====

Beamlet Source

=====

charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,23,24.5)

=====

=====

XYZ Geometry

=====

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV

Example: Photon Treatment Plan
using VMC++ dose calculation

```

Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition    : 0
Cut-off energy for KERMA approx.    : 1.10239 MeV
Bremsstrahlung transport mode       : 1
CSDA approximation                  : 0
Fractional energy loss/step at Ep   : 10%
Max. 1st elastic moment per step    : 0.5
Max. acceptable energy loss/step    : 0.6 MeV
alpha and beta                      : 0.0298764 0.420741
Fano calculation                     : 0
Exact Compton                       : 0
Electron transport mode              : VMC++

```

```

=====
=====
MC_Control
=====

```

```

will use fixed number of histories
number of batches      : 10
histories per batch    : 500
total histories        : 5000
initial rng seeds      : 25474 28020
=====

```

```

=====
Variance Reduction
=====

```

```

f_repeat      = 0.251
split photons  = 1
photon split factor = -40
=====

```

```

=====
Quasi Random Numbers
=====

```

```

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1
=====

```

```

=====
Scoring and output options
=====

```

```

number of dose scoring objects: 1
dose output options for geometry CT
  dump dose: 2
  dose scans:

```

CPU time so far: 2.353 seconds

Will run approximately 5000 particle sets
 with 1 particles per set on average
 will use 2 quasi numbers to sample the source

Example: Photon Treatment Plan
using VMC++ dose calculation

Starting MC simulation

OK

=====

Beamlet Source

=====

charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,23,23)

=====

=====

XYZ Geometry

=====

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 12653 27473

=====

Variance Reduction

=====

Example: Photon Treatment Plan
using VMC++ dose calculation

```
f_repeat          = 0.251
split photons     = 1
photon split factor = -40
=====
=====
                        Quasi Random Numbers
=====
number of generators: 1
                    1st: base = 2 dimensions = 60 warm-up = 1
=====
=====
                        Scoring and output options
=====
number of dose scoring objects: 1
dose output options for geometry CT
    dump dose: 2
    dose scans:

CPU time so far: 2.475 seconds

Will run approximately 5000 particle sets
    with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK
=====
                        Beamlet Source
=====
charge = 0
virtual source position = (-76,24,24)
    Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,23,24)
=====
=====
                        XYZ Geometry
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
=====
                        Monte Carlo Parameter
=====
Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition   : 0
```

Example: Photon Treatment Plan
using VMC++ dose calculation

Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at E_p : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 19673 1072

=====

Variance Reduction

=====

f_{repeat} = 0.251
split photons = 1
photon split factor = -40

=====

Quasi Random Numbers

=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

Scoring and output options

=====

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.548 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.059 500 500

Example: Photon Treatment Plan
using VMC++ dose calculation

```
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.043 500 500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.043 500 500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.045 500 500
+++finished batch 2 cpu time: 0.293 500 500
+++finished batch 2 cpu time: 0.195 500 500
+++finished batch 2 cpu time: 0.39 500 500
+++finished batch 2 cpu time: 0.135 500 500
+++finished batch 3 cpu time: 0.367 500 500
+++finished batch 3 cpu time: 0.264 500 500
+++finished batch 3 cpu time: 0.471 500 500
+++finished batch 3 cpu time: 0.197 500 500
+++finished batch 4 cpu time: 0.339 500 500
+++finished batch 4 cpu time: 0.452 500 500
+++finished batch 4 cpu time: 0.542 500 500
+++finished batch 4 cpu time: 0.273 500 500
+++finished batch 5 cpu time: 0.516 500 500
+++finished batch 5 cpu time: 0.41 500 500
+++finished batch 5 cpu time: 0.605 500 500
+++finished batch 5 cpu time: 0.343 500 500
+++finished batch 6 cpu time: 0.579 500 500
+++finished batch 6 cpu time: 0.673 500 500
+++finished batch 6 cpu time: 0.485 500 500
+++finished batch 6 cpu time: 0.415 500 500
+++finished batch 7 cpu time: 0.654 500 500
+++finished batch 7 cpu time: 0.565 500 500
+++finished batch 7 cpu time: 0.779 500 500
+++finished batch 7 cpu time: 0.486 500 500
+++finished batch 8 cpu time: 0.75 500 500
+++finished batch 8 cpu time: 0.649 500 500
+++finished batch 8 cpu time: 0.862 500 500
+++finished batch 8 cpu time: 0.577 500 500
+++finished batch 9 cpu time: 0.817 500 500
+++finished batch 9 cpu time: 0.743 500 500
+++finished batch 9 cpu time: 0.947 500 500
+++finished batch 9 cpu time: 0.677 500 500
+++finished batch 10 cpu time: 0.895 500 500
finished simulation, cpu time = 0.911 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
geometry: CT
cpu time: 0.911
number of histories: 5000
number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.822 500 500
finished simulation, cpu time = 0.836 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
```

```

total number of particles: 5000
average sampled energy: 6 +/- 0
+++finished batch 10 cpu time: 1.03 500 500
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.836
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
finished simulation, cpu time = 1.046 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 1.046
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.751 500 500
finished simulation, cpu time = 0.765 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.765
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
    max dose is 0.0342117 in region 2110926
    max dose is 0.0342145 in region 1880527
    max dose is 0.0340826 in region 1957327
    max dose is 0.0344097 in region 2034127
    ICRU-2K efficiency: 967.774 1/s
+++ total
    max dose is 0.0342145 in region 1880527
    ICRU-2K efficiency: 967.774 1/s
=====
    ICRU-2K efficiency: 819.593 1/s
+++ total
    max dose is 0.0340826 in region 1957327
    ICRU-2K efficiency: 819.593 1/s
=====
    ICRU-2K efficiency: 925.181 1/s
+++ total
    max dose is 0.0342117 in region 2110926
    ICRU-2K efficiency: 1070.74 1/s
+++ total
    max dose is 0.0344097 in region 2034127
    ICRU-2K efficiency: 925.181 1/s
=====

```

```
ICRU-2K efficiency: 1070.74 1/s
=====
Completed 12 of 49 beamlets...

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_1.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_4.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_2.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_3.vmc
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
=====
                        Beamlet Source
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```

=====
      charge = 0
    virtual source position = (-76,24,24)
      Energy = 6 MeV
    number of beamlets = 1
    beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,23.5,23)
=====
=====
                        XYZ Geometry
=====
      name: CT   id: 0
    global smax: 1e+030
    Number of x-planes: 160   uniform with Xo = 0.15 Dx = 0.3
    Number of y-planes: 160   uniform with Yo = 0.15 Dx = 0.3
    Number of z-planes: 160   uniform with Zo = 0.15 Dx = 0.3
=====
                        Monte Carlo Parameter
=====

Delta particle production threshold:    0.447479 MeV
Bremsstrahlung production threshold:    0.05 MeV
Min. electron transport energy         :    0.447479 MeV
Min. photon transport energy           :    0.05 MeV
Local track-end energy deposition       :    0
Cut-off energy for KERMA approx.        :    1.10239 MeV
Bremsstrahlung transport mode           :    1
CSDA approximation                     :    0
Fractional energy loss/step at Ep       :    10%
Max. 1st elastic moment per step        :    0.5
Max. acceptable energy loss/step        :    0.6 MeV
alpha and beta                          :    0.0298764  0.420741
Fano calculation                        :    0
Exact Compton                          :    0
Electron transport mode                  :    VMC++
=====
=====
                        MC_Control
=====
    will use fixed number of histories
    number of batches   : 10
    histories per batch : 500
    total histories     : 5000
    initial rng seeds   : 21182  955
=====
=====
                        Variance Reduction
=====

f_repeat          = 0.251
split photons     = 1
photon split factor = -40
=====
=====

```

Example: Photon Treatment Plan
using VMC++ dose calculation

Quasi Random Numbers

```
=====
number of generators: 1
      1st:  base = 2 dimensions = 60 warm-up = 1
=====
```

Scoring and output options

```
=====
number of dose scoring objects: 1
dose output options for geometry CT
  dump dose: 2
  dose scans:
=====
```

CPU time so far: 2.487 seconds

Will run approximately 5000 particle sets
 with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation

OK

Beamlet Source

```
=====
charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,23.5,22.5)
=====
```

XYZ Geometry

```
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
```

Monte Carlo Parameter

```
=====
Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition    : 0
Cut-off energy for KERMA approx.    : 1.10239 MeV
Bremsstrahlung transport mode       : 1
CSDA approximation                  : 0
Fractional energy loss/step at Ep   : 10%
Max. 1st elastic moment per step    : 0.5
=====
```

Example: Photon Treatment Plan
using VMC++ dose calculation

Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 19665 5136

=====

Variance Reduction

=====

f_repeat = 0.251
split photons = 1
photon split factor = -40

=====

Quasi Random Numbers

=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

Scoring and output options

=====

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.557 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK

=====

Beamlet Source

=====

charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV

Example: Photon Treatment Plan
using VMC++ dose calculation

```

number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,23,25)
=====
XYZ Geometry
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
Monte Carlo Parameter
=====
Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++
=====
MC_Control
=====
will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 20363 22733
=====
Variance Reduction
=====
f_repeat = 0.251
split photons = 1
photon split factor = -40
=====
Quasi Random Numbers
=====
number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

```

Example: Photon Treatment Plan
using VMC++ dose calculation

```
=====
=====
                        Scoring and output options
=====
```

```
number of dose scoring objects: 1
dose output options for geometry CT
    dump dose: 2
    dose scans:
```

CPU time so far: 2.583 seconds

Will run approximately 5000 particle sets
 with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.049 500 500
OK

```
=====
                        Beamlet Source
=====
```

```
charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,23,25.5)
=====
=====
```

```
=====
                        XYZ Geometry
=====
```

```
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
=====
```

```
=====
                        Monte Carlo Parameter
=====
```

```
Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition    : 0
Cut-off energy for KERMA approx.     : 1.10239 MeV
Bremsstrahlung transport mode        : 1
CSDA approximation                  : 0
Fractional energy loss/step at Ep    : 10%
Max. 1st elastic moment per step     : 0.5
Max. acceptable energy loss/step     : 0.6 MeV
alpha and beta                       : 0.0298764 0.420741
```

Example: Photon Treatment Plan
using VMC++ dose calculation

Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 22294 11767

=====

=====

Variance Reduction

=====

f_repeat = 0.251
split photons = 1
photon split factor = -40

=====

=====

Quasi Random Numbers

=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

=====

Scoring and output options

=====

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.669 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.041 500 500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.046 500 500
+++finished batch 2 cpu time: 0.145 500 500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.043 500 500
+++finished batch 2 cpu time: 0.179 500 500
+++finished batch 2 cpu time: 0.186 500 500

Example: Photon Treatment Plan
using VMC++ dose calculation

```

+++finished batch 3 cpu time: 0.278 500 500
+++finished batch 2 cpu time: 0.111 500 500
+++finished batch 3 cpu time: 0.249 500 500
+++finished batch 3 cpu time: 0.255 500 500
+++finished batch 4 cpu time: 0.35 500 500
+++finished batch 3 cpu time: 0.169 500 500
+++finished batch 4 cpu time: 0.315 500 500
+++finished batch 4 cpu time: 0.317 500 500
+++finished batch 5 cpu time: 0.408 500 500
+++finished batch 4 cpu time: 0.237 500 500
+++finished batch 5 cpu time: 0.378 500 500
+++finished batch 5 cpu time: 0.381 500 500
+++finished batch 6 cpu time: 0.472 500 500
+++finished batch 5 cpu time: 0.299 500 500
+++finished batch 6 cpu time: 0.441 500 500
+++finished batch 7 cpu time: 0.536 500 500
+++finished batch 6 cpu time: 0.451 500 500
+++finished batch 6 cpu time: 0.381 500 500
+++finished batch 7 cpu time: 0.511 500 500
+++finished batch 7 cpu time: 0.513 500 500
+++finished batch 8 cpu time: 0.606 500 500
+++finished batch 7 cpu time: 0.448 500 500
+++finished batch 8 cpu time: 0.572 500 500
+++finished batch 9 cpu time: 0.668 500 500
+++finished batch 8 cpu time: 0.583 500 500
+++finished batch 8 cpu time: 0.515 500 500
+++finished batch 9 cpu time: 0.634 500 500
+++finished batch 10 cpu time: 0.725 500 500
+++finished batch 9 cpu time: 0.642 500 500
finished simulation, cpu time = 0.742 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.742
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
+++finished batch 9 cpu time: 0.577 500 500
+++finished batch 10 cpu time: 0.699 500 500
+++finished batch 10 cpu time: 0.702 500 500
finished simulation, cpu time = 0.716 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.716
  number of histories: 5000
  number of batches: 10
+++ beamlet 0

```

```
finished simulation, cpu time = 0.719 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.719
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.65  500  500
finished simulation, cpu time = 0.663 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.663
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
    max dose is 0.0340248 in region 1881006
    max dose is 0.0343928 in region 2187726
    max dose is 0.03426 in region 1804207
    max dose is 0.0343579 in region 2264526
    ICRU-2K efficiency: 1123.43 1/s
+++ total
    max dose is 0.0340248 in region 1881006
    ICRU-2K efficiency: 1123.43 1/s
=====
    ICRU-2K efficiency: 1192.13 1/s
+++ total
    max dose is 0.03426 in region 1804207
    ICRU-2K efficiency: 1192.13 1/s
=====
    ICRU-2K efficiency: 1179.88 1/s
+++ total
    max dose is 0.0343928 in region 2187726
    ICRU-2K efficiency: 1179.88 1/s
=====
    ICRU-2K efficiency: 1254.3 1/s
+++ total
    max dose is 0.0343579 in region 2264526
    ICRU-2K efficiency: 1254.3 1/s
=====
Completed 16 of 49 beamlets...
```

```
Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_1.vmc
```

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_2.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_3.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_4.vmc

```
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
=====
                        Beamlet Source
=====
        charge = 0
        virtual source position = (-76,24,24)
        Energy = 6 MeV
        number of beamlets = 1
        beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,23.5,25)
=====
=====
                        XYZ Geometry
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====

Monte Carlo Parameter
=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

MC_Control
=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 28507 1034

=====

Variance Reduction
=====

f_repeat = 0.251
split photons = 1
photon split factor = -40

=====

Quasi Random Numbers
=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

Scoring and output options
=====

number of dose scoring objects: 1
```

Example: Photon Treatment Plan
using VMC++ dose calculation

dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.295 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK

=====

Beamlet Source

=====

charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,23.5,23.5)

=====

=====

XYZ Geometry

=====

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

MC_Control

Example: Photon Treatment Plan
using VMC++ dose calculation

```
=====
will use fixed number of histories
number of batches      : 10
histories per batch    : 500
total histories        : 5000
initial rng seeds      : 8308 1386
=====

=====
Variance Reduction
=====

f_repeat              = 0.251
split photons         = 1
photon split factor   = -40
=====

=====
Quasi Random Numbers
=====

number of generators: 1
                    1st: base = 2 dimensions = 60 warm-up = 1
=====

=====
Scoring and output options
=====

number of dose scoring objects: 1
dose output options for geometry CT
    dump dose: 2
    dose scans:

CPU time so far: 2.363 seconds

Will run approximately 5000 particle sets
    with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK
=====

Beamlet Source
=====

charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,23.5,24.5)
=====

XYZ Geometry
=====

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
```

Example: Photon Treatment Plan
using VMC++ dose calculation

Number of y-planes: 160 uniform with $Y_0 = 0.15$ $Dx = 0.3$
Number of z-planes: 160 uniform with $Z_0 = 0.15$ $Dx = 0.3$

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at E_p : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 20845 9513

=====

Variance Reduction

=====

$f_{repeat} = 0.251$
split photons = 1
photon split factor = -40

=====

Quasi Random Numbers

=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

Scoring and output options

=====

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.466 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation

OK

=====

Beamlet Source

=====

charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,23.5,24)

=====

=====

XYZ Geometry

=====

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500

Example: Photon Treatment Plan
using VMC++ dose calculation

```
total histories      : 5000
initial rng seeds   : 2914 24704
=====
=====
Variance Reduction
=====
=====

f_repeat            = 0.251
split photons       = 1
photon split factor = -40
=====
=====

Quasi Random Numbers
=====
=====

number of generators: 1
                    1st: base = 2 dimensions = 60 warm-up = 1
=====
=====

Scoring and output options
=====
=====

number of dose scoring objects: 1
dose output options for geometry CT
    dump dose: 2
    dose scans:

CPU time so far: 2.577 seconds

Will run approximately 5000 particle sets
    with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.058 500 500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.042 500 500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.042 500 500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.041 500 500
+++finished batch 2 cpu time: 0.33 500 500
+++finished batch 2 cpu time: 0.31 500 500
+++finished batch 2 cpu time: 0.203 500 500
+++finished batch 2 cpu time: 0.136 500 500
+++finished batch 3 cpu time: 0.405 500 500
+++finished batch 3 cpu time: 0.376 500 500
+++finished batch 3 cpu time: 0.29 500 500
+++finished batch 3 cpu time: 0.203 500 500
+++finished batch 4 cpu time: 0.472 500 500
+++finished batch 4 cpu time: 0.45 500 500
+++finished batch 4 cpu time: 0.347 500 500
+++finished batch 4 cpu time: 0.269 500 500
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```
+++finished batch 5 cpu time: 0.536 500 500
+++finished batch 5 cpu time: 0.511 500 500
+++finished batch 5 cpu time: 0.408 500 500
+++finished batch 5 cpu time: 0.327 500 500
+++finished batch 6 cpu time: 0.608 500 500
+++finished batch 6 cpu time: 0.584 500 500
+++finished batch 6 cpu time: 0.499 500 500
+++finished batch 6 cpu time: 0.392 500 500
+++finished batch 7 cpu time: 0.68 500 500
+++finished batch 7 cpu time: 0.652 500 500
+++finished batch 7 cpu time: 0.57 500 500
+++finished batch 7 cpu time: 0.47 500 500
+++finished batch 8 cpu time: 0.767 500 500
+++finished batch 8 cpu time: 0.739 500 500
+++finished batch 8 cpu time: 0.648 500 500
+++finished batch 8 cpu time: 0.545 500 500
+++finished batch 9 cpu time: 0.828 500 500
+++finished batch 9 cpu time: 0.811 500 500
+++finished batch 9 cpu time: 0.712 500 500
+++finished batch 9 cpu time: 0.615 500 500
+++finished batch 10 cpu time: 0.875 500 500
+++finished batch 10 cpu time: 0.915 500 500
+++finished batch 10 cpu time: 0.777 500 500
+++finished batch 10 cpu time: 0.677 500 500
finished simulation, cpu time = 0.899 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.899
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
finished simulation, cpu time = 0.796 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
finished simulation, cpu time = 0.937 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.796
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.937
```

```
number of histories: 5000
number of batches: 10
+++ beamlet 0
finished simulation, cpu time = 0.694 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
geometry: CT
cpu time: 0.694
number of histories: 5000
number of batches: 10
+++ beamlet 0
    max dose is 0.0347457 in region 1957807
    max dose is 0.0344827 in region 2188206
    max dose is 0.0349114 in region 2034607
    max dose is 0.0346036 in region 2111407
    ICRU-2K efficiency: 951.083 1/s
+++ total
    max dose is 0.0347457 in region 1957807
    ICRU-2K efficiency: 951.083 1/s
=====
    ICRU-2K efficiency: 934.841 1/s
+++ total
    max dose is 0.0344827 in region 2188206
    ICRU-2K efficiency: 1088.52 1/s
+++ total
    max dose is 0.0346036 in region 2111407
    ICRU-2K efficiency: 1246.27 1/s
+++ total
    max dose is 0.0349114 in region 2034607
    ICRU-2K efficiency: 934.841 1/s
=====
    ICRU-2K efficiency: 1088.52 1/s
=====
    ICRU-2K efficiency: 1246.27 1/s
=====
Completed 20 of 49 beamlets...
```

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_1.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_2.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_3.vmc

```

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_4.vmc
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
=====
                        Beamlet Source
=====
        charge = 0
        virtual source position = (-76,24,24)
        Energy = 6 MeV
        number of beamlets = 1
        beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,24,23.5)
=====
=====
                        XYZ Geometry
=====
        name: CT id: 0
        global smax: 1e+030
        Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
        Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
        Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
                        Monte Carlo Parameter
=====

```

Example: Photon Treatment Plan
using VMC++ dose calculation

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at E_p : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====
=====
MC_Control
=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 13368 19390
=====

=====
=====
Variance Reduction
=====

f_{repeat} = 0.251
split photons = 1
photon split factor = -40
=====

=====
=====
Quasi Random Numbers
=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1
=====

=====
=====
Scoring and output options
=====

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.273 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average

Example: Photon Treatment Plan
using VMC++ dose calculation

will use 2 quasi numbers to sample the source

Starting MC simulation

OK

=====

Beamlet Source

=====

charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
*beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,24,23)*

=====

=====

XYZ Geometry

=====

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 5607 14693

=====

Variance Reduction

=====

Example: Photon Treatment Plan
using VMC++ dose calculation

```
=====
f_repeat          = 0.251
split photons     = 1
photon split factor = -40
=====
=====
Quasi Random Numbers
=====
number of generators: 1
                    1st: base = 2 dimensions = 60 warm-up = 1
=====
=====
Scoring and output options
=====
number of dose scoring objects: 1
dose output options for geometry CT
    dump dose: 2
    dose scans:

CPU time so far: 2.36 seconds

Will run approximately 5000 particle sets
    with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK
=====
Beamlet Source
=====
charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,23.5,25.5)
=====
=====
XYZ Geometry
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
Monte Carlo Parameter
=====
Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
```

Example: Photon Treatment Plan
using VMC++ dose calculation

Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at E_p : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 13163 11447

=====

Variance Reduction

=====

f_repeat = 0.251
split photons = 1
photon split factor = -40

=====

Quasi Random Numbers

=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

Scoring and output options

=====

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.48 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation

Example: Photon Treatment Plan
using VMC++ dose calculation

OK

=====

Beamlet Source

=====

charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,24,22.5)

=====

=====

XYZ Geometry

=====

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 22966 23856

=====

=====

Variance Reduction

=====

f_repeat = 0.251
split photons = 1

Example: Photon Treatment Plan
using VMC++ dose calculation

```
photon split factor = -40
=====
=====
Quasi Random Numbers
=====
number of generators: 1
    1st: base = 2 dimensions = 60 warm-up = 1
=====
=====
Scoring and output options
=====
number of dose scoring objects: 1
dose output options for geometry CT
    dump dose: 2
    dose scans:

CPU time so far: 2.583 seconds

Will run approximately 5000 particle sets
    with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.043  500  500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.042  500  500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.045  500  500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.043  500  500
+++finished batch 2 cpu time: 0.289  500  500
+++finished batch 2 cpu time: 0.181  500  500
+++finished batch 2 cpu time: 0.368  500  500
+++finished batch 2 cpu time: 0.117  500  500
+++finished batch 3 cpu time: 0.356  500  500
+++finished batch 3 cpu time: 0.432  500  500
+++finished batch 3 cpu time: 0.249  500  500
+++finished batch 3 cpu time: 0.182  500  500
+++finished batch 4 cpu time: 0.496  500  500
+++finished batch 4 cpu time: 0.426  500  500
+++finished batch 4 cpu time: 0.336  500  500
+++finished batch 4 cpu time: 0.244  500  500
+++finished batch 5 cpu time: 0.56  500  500
+++finished batch 5 cpu time: 0.492  500  500
+++finished batch 5 cpu time: 0.401  500  500
+++finished batch 5 cpu time: 0.31  500  500
+++finished batch 6 cpu time: 0.628  500  500
+++finished batch 6 cpu time: 0.559  500  500
+++finished batch 6 cpu time: 0.465  500  500
+++finished batch 6 cpu time: 0.386  500  500
+++finished batch 7 cpu time: 0.704  500  500
```

```
+++finished batch 7 cpu time: 0.526 500 500
+++finished batch 7 cpu time: 0.644 500 500
+++finished batch 7 cpu time: 0.454 500 500
+++finished batch 8 cpu time: 0.785 500 500
+++finished batch 8 cpu time: 0.721 500 500
+++finished batch 8 cpu time: 0.61 500 500
+++finished batch 8 cpu time: 0.537 500 500
+++finished batch 9 cpu time: 0.853 500 500
+++finished batch 9 cpu time: 0.676 500 500
+++finished batch 9 cpu time: 0.795 500 500
+++finished batch 9 cpu time: 0.606 500 500
+++finished batch 10 cpu time: 0.928 500 500
+++finished batch 10 cpu time: 0.746 500 500
finished simulation, cpu time = 0.942 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
geometry: CT
cpu time: 0.942
number of histories: 5000
number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.871 500 500
finished simulation, cpu time = 0.772 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
geometry: CT
cpu time: 0.772
number of histories: 5000
number of batches: 10
+++ beamlet 0
finished simulation, cpu time = 0.885 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
geometry: CT
cpu time: 0.885
number of histories: 5000
number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.674 500 500
finished simulation, cpu time = 0.689 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
```

```
geometry: CT
cpu time: 0.689
number of histories: 5000
number of batches: 10
+++ beamlet 0
    max dose is 0.0343987 in region 1958286
    max dose is 0.0344111 in region 2265006
    max dose is 0.0339486 in region 1804686
    max dose is 0.0343559 in region 1881486
    ICRU-2K efficiency: 914.102 1/s
+++ total
    max dose is 0.0343987 in region 1958286
    ICRU-2K efficiency: 914.102 1/s
=====
    ICRU-2K efficiency: 1128.79 1/s
+++ total
    max dose is 0.0344111 in region 2265006
    ICRU-2K efficiency: 1251.46 1/s
+++ total
    max dose is 0.0339486 in region 1804686
    ICRU-2K efficiency: 977.628 1/s
+++ total
    max dose is 0.0343559 in region 1881486
    ICRU-2K efficiency: 1251.46 1/s
=====
    ICRU-2K efficiency: 1128.79 1/s
=====
    ICRU-2K efficiency: 977.628 1/s
=====
Completed 24 of 49 beamlets...
```

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_1.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_4.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_2.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_3.vmc
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1

```

Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK

```

=====

Beamlet Source

=====

```

charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,24,25.5)

```

=====

XYZ Geometry

=====

```

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3

```

=====

Monte Carlo Parameter

=====

```

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition    : 0
Cut-off energy for KERMA approx.     : 1.10239 MeV
Bremsstrahlung transport mode        : 1
CSDA approximation                   : 0

```


Example: Photon Treatment Plan
using VMC++ dose calculation

Fractional energy loss/step at E_p : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 3570 14951

=====

Variance Reduction

=====

f_repeat = 0.251
split photons = 1
photon split factor = -40

=====

Quasi Random Numbers

=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

Scoring and output options

=====

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.223 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK

=====

Beamlet Source

=====

charge = 0

Example: Photon Treatment Plan
using VMC++ dose calculation

```

virtual source position = (-76,24,24)
    Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,24,24.5)
=====
XYZ Geometry
=====
    name: CT id: 0
    global smax: 1e+030
    Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3    Number of
y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
    Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
Monte Carlo Parameter
=====

Delta particle production threshold:    0.447479 MeV
Bremsstrahlung production threshold:    0.05 MeV
Min. electron transport energy         :    0.447479 MeV
Min. photon transport energy           :    0.05 MeV
Local track-end energy deposition       :    0
Cut-off energy for KERMA approx.        :    1.10239 MeV
Bremsstrahlung transport mode           :    1
CSDA approximation                     :    0
Fractional energy loss/step at Ep       :    10%
Max. 1st elastic moment per step        :    0.5
Max. acceptable energy loss/step        :    0.6 MeV
alpha and beta                          :    0.0298764  0.420741
Fano calculation                        :    0
Exact Compton                           :    0
Electron transport mode                  :    VMC++
=====
MC_Control
=====
    will use fixed number of histories
    number of batches : 10
    histories per batch : 500
    total histories   : 5000
    initial rng seeds : 8281 20392
=====
Variance Reduction
=====

f_repeat          = 0.251
split photons     = 1
photon split factor = -40
=====
Quasi Random Numbers
=====

```

Example: Photon Treatment Plan
using VMC++ dose calculation

```

number of generators: 1
      1st:  base = 2 dimensions = 60 warm-up = 1
=====
Scoring and output options
=====
number of dose scoring objects: 1
dose output options for geometry CT
  dump dose: 2
  dose scans:

CPU time so far: 2.231 seconds

Will run approximately 5000 particle sets
  with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK
=====
      Beamlet Source
=====
      charge = 0
virtual source position = (-76,24,24)
      Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,24,24)
=====
      XYZ Geometry
=====
      name: CT  id: 0
      global smax: 1e+030
Number of x-planes: 160  uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160  uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160  uniform with Zo = 0.15 Dx = 0.3
=====
      Monte Carlo Parameter
=====

Delta particle production threshold:  0.447479 MeV
Bremsstrahlung production threshold:  0.05 MeV
Min. electron transport energy       :  0.447479 MeV
Min. photon transport energy         :  0.05 MeV
Local track-end energy deposition     :  0
Cut-off energy for KERMA approx.      :  1.10239 MeV
Bremsstrahlung transport mode         :  1
CSDA approximation                   :  0
Fractional energy loss/step at Ep     :  10%
Max. 1st elastic moment per step      :  0.5
Max. acceptable energy loss/step      :  0.6 MeV
alpha and beta                        :  0.0298764  0.420741

```

Example: Photon Treatment Plan
using VMC++ dose calculation

```
Fano calculation           : 0
Exact Compton             : 0
Electron transport mode   : VMC++
```

```
=====
MC_Control
=====
```

```
will use fixed number of histories
number of batches      : 10
histories per batch    : 500
total histories        : 5000
initial rng seeds      : 21281 22641
=====
```

```
=====
Variance Reduction
=====
```

```
f_repeat      = 0.251
split photons  = 1
photon split factor = -40
=====
```

```
=====
Quasi Random Numbers
=====
```

```
number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1
=====
```

```
=====
Scoring and output options
=====
```

```
number of dose scoring objects: 1
dose output options for geometry CT
  dump dose: 2
  dose scans:
```

CPU time so far: 2.359 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK

```
=====
Beamlet Source
=====
```

```
charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,24,25)
```

Example: Photon Treatment Plan
using VMC++ dose calculation

=====

XYZ Geometry

=====

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 19653 4879

=====

Variance Reduction

=====

f_repeat = 0.251
split photons = 1
photon split factor = -40

=====

Quasi Random Numbers

=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

Scoring and output options

```
=====
number of dose scoring objects: 1
dose output options for geometry CT
  dump dose: 2
  dose scans:
```

CPU time so far: 2.345 seconds

Will run approximately 5000 particle sets
 with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation

```
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.053  500  500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.04  500  500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.041  500  500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.052  500  500
+++finished batch 2 cpu time: 0.223  500  500
+++finished batch 2 cpu time: 0.21  500  500
+++finished batch 2 cpu time: 0.108  500  500
+++finished batch 2 cpu time: 0.147  500  500
+++finished batch 3 cpu time: 0.297  500  500
+++finished batch 3 cpu time: 0.189  500  500
+++finished batch 3 cpu time: 0.3  500  500
+++finished batch 3 cpu time: 0.21  500  500
+++finished batch 4 cpu time: 0.384  500  500
+++finished batch 4 cpu time: 0.263  500  500
+++finished batch 4 cpu time: 0.389  500  500
+++finished batch 4 cpu time: 0.296  500  500
+++finished batch 5 cpu time: 0.444  500  500
+++finished batch 5 cpu time: 0.332  500  500
+++finished batch 5 cpu time: 0.453  500  500
+++finished batch 6 cpu time: 0.509  500  500
+++finished batch 5 cpu time: 0.385  500  500
+++finished batch 6 cpu time: 0.413  500  500
+++finished batch 6 cpu time: 0.528  500  500
+++finished batch 7 cpu time: 0.586  500  500
+++finished batch 6 cpu time: 0.453  500  500
+++finished batch 7 cpu time: 0.487  500  500
+++finished batch 7 cpu time: 0.607  500  500
+++finished batch 7 cpu time: 0.521  500  500
+++finished batch 8 cpu time: 0.675  500  500
+++finished batch 8 cpu time: 0.567  500  500
+++finished batch 8 cpu time: 0.684  500  500
+++finished batch 9 cpu time: 0.734  500  500
+++finished batch 8 cpu time: 0.6  500  500
+++finished batch 9 cpu time: 0.631  500  500
```

```
+++finished batch 9 cpu time: 0.764 500 500
+++finished batch 9 cpu time: 0.666 500 500
+++finished batch 10 cpu time: 0.81 500 500
finished simulation, cpu time = 0.824 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.824
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.714 500 500
+++finished batch 10 cpu time: 0.835 500 500
finished simulation, cpu time = 0.727 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.727
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.725 500 500
finished simulation, cpu time = 0.85 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.85
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
finished simulation, cpu time = 0.748 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.748
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
  max dose is 0.0347752 in region 2265486
  max dose is 0.034533 in region 2035086
  max dose is 0.0344639 in region 2111887
  max dose is 0.0347516 in region 2188686
```

```
ICRU-2K efficiency: 1050.29 1/s
+++ total
max dose is 0.0347752 in region 2265486
ICRU-2K efficiency: 1050.29 1/s
=====
ICRU-2K efficiency: 1160.63 1/s
+++ total
max dose is 0.034533 in region 2035086
ICRU-2K efficiency: 1010.14 1/s
+++ total
max dose is 0.0344639 in region 2111887
ICRU-2K efficiency: 1160.63 1/s
=====
ICRU-2K efficiency: 1169.03 1/s
+++ total
max dose is 0.0347516 in region 2188686
ICRU-2K efficiency: 1010.14 1/s
=====
ICRU-2K efficiency: 1169.03 1/s
=====
Completed 28 of 49 beamlets...
```

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_1.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_2.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_4.vmc

```
Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/  
MCpencilbeam_temp_3.vmc  
Reading MS data ... OK  
Parsing input file ... construct_mmc_geometry: urs = 1  
Reading MS data ... OK  
Parsing input file ... construct_mmc_geometry: urs = 1  
Reading MS data ... OK  
Parsing input file ... construct_mmc_geometry: urs = 1  
Reading MS data ... OK  
Parsing input file ... construct_mmc_geometry: urs = 1  
DE_GeometryFactory::get_region_size():  
get_region_size: urs = 1 ignore = 1  
region size: 0.3  
geometry 1 region size: 0.3  
DE_GeometryFactory::get_region_size():
```



```

get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
=====
                        Beamlet Source
=====
      charge = 0
virtual source position = (-76,24,24)
      Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,24.5,24)
=====
=====
                        XYZ Geometry
=====
      name: CT id: 0
      global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
                        Monte Carlo Parameter
=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition   : 0
Cut-off energy for KERMA approx.    : 1.10239 MeV
Bremsstrahlung transport mode       : 1
CSDA approximation                  : 0
Fractional energy loss/step at Ep   : 10%
Max. 1st elastic moment per step    : 0.5
Max. acceptable energy loss/step    : 0.6 MeV
alpha and beta                      : 0.0298764 0.420741
Fano calculation                    : 0
Exact Compton                      : 0
Electron transport mode              : VMC++
=====

```

```
=====
MC_Control
=====
will use fixed number of histories
number of batches      : 10
histories per batch    : 500
total histories        : 5000
initial rng seeds      : 15179 20973
=====
Variance Reduction
=====

f_repeat              = 0.251
split photons         = 1
photon split factor   = -40
=====
Quasi Random Numbers
=====
number of generators: 1
                    1st: base = 2 dimensions = 60 warm-up = 1
=====
Scoring and output options
=====
number of dose scoring objects: 1
dose output options for geometry CT
    dump dose: 2
    dose scans:

CPU time so far: 2.38 seconds

Will run approximately 5000 particle sets
    with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK
=====
Beamlet Source
=====
charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,24.5,22.5)
=====
XYZ Geometry
=====
name: CT id: 0
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
```

```
=====
Monte Carlo Parameter
=====
```

```
Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition    : 0
Cut-off energy for KERMA approx.    : 1.10239 MeV
Bremsstrahlung transport mode       : 1
CSDA approximation                  : 0
Fractional energy loss/step at Ep   : 10%
Max. 1st elastic moment per step    : 0.5
Max. acceptable energy loss/step    : 0.6 MeV
alpha and beta                      : 0.0298764 0.420741
Fano calculation                    : 0
Exact Compton                      : 0
Electron transport mode              : VMC++
```

```
=====
MC_Control
=====
```

```
will use fixed number of histories
number of batches      : 10
histories per batch    : 500
total histories        : 5000
initial rng seeds      : 28793 10212
```

```
=====
Variance Reduction
=====
```

```
f_repeat      = 0.251
split photons  = 1
photon split factor = -40
```

```
=====
Quasi Random Numbers
=====
```

```
number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1
```

```
=====
Scoring and output options
=====
```

```
number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
```

Example: Photon Treatment Plan
using VMC++ dose calculation

dose scans:

CPU time so far: 2.475 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation

OK

=====

Beamlet Source

=====

charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,24.5,23)

=====

=====

XYZ Geometry

=====

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

=====

MC_Control

=====

will use fixed number of histories

Example: Photon Treatment Plan
using VMC++ dose calculation

```
number of batches      : 10
histories per batch    : 500
total histories        : 5000
initial rng seeds      : 17559  6715
=====
=====
Variance Reduction
=====
=====
f_repeat              = 0.251
split photons         = 1
photon split factor   = -40
=====
=====
Quasi Random Numbers
=====
=====
number of generators: 1
                    1st: base = 2 dimensions = 60 warm-up = 1
=====
=====
Scoring and output options
=====
=====
number of dose scoring objects: 1
dose output options for geometry CT
    dump dose: 2
    dose scans:

CPU time so far: 2.471 seconds

Will run approximately 5000 particle sets
    with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.045  500  500
OK
=====
Beamlet Source
=====
=====
charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,24.5,23.5)
=====
=====
XYZ Geometry
=====
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
```

Example: Photon Treatment Plan
using VMC++ dose calculation

Number of y-planes: 160 uniform with $Y_0 = 0.15$ $Dx = 0.3$
Number of z-planes: 160 uniform with $Z_0 = 0.15$ $Dx = 0.3$

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at E_p : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 22539 7653

=====

Variance Reduction

=====

$f_{repeat} = 0.251$
split photons = 1
photon split factor = -40

=====

Quasi Random Numbers

=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

Scoring and output options

=====

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.54 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

```
Starting MC simulation
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.041  500  500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.06  500  500
+++finished batch 2 cpu time: 0.161  500  500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.045  500  500
+++finished batch 2 cpu time: 0.179  500  500
+++finished batch 2 cpu time: 0.181  500  500
+++finished batch 3 cpu time: 0.293  500  500
+++finished batch 2 cpu time: 0.124  500  500
+++finished batch 3 cpu time: 0.249  500  500
+++finished batch 3 cpu time: 0.266  500  500
+++finished batch 4 cpu time: 0.362  500  500
+++finished batch 3 cpu time: 0.187  500  500
+++finished batch 4 cpu time: 0.343  500  500
+++finished batch 4 cpu time: 0.347  500  500
+++finished batch 5 cpu time: 0.424  500  500
+++finished batch 4 cpu time: 0.277  500  500
+++finished batch 5 cpu time: 0.424  500  500
+++finished batch 5 cpu time: 0.418  500  500
+++finished batch 6 cpu time: 0.518  500  500
+++finished batch 5 cpu time: 0.374  500  500
+++finished batch 6 cpu time: 0.498  500  500
+++finished batch 6 cpu time: 0.503  500  500
+++finished batch 7 cpu time: 0.588  500  500
+++finished batch 6 cpu time: 0.433  500  500
+++finished batch 7 cpu time: 0.586  500  500
+++finished batch 8 cpu time: 0.671  500  500
+++finished batch 7 cpu time: 0.609  500  500
+++finished batch 7 cpu time: 0.523  500  500
+++finished batch 9 cpu time: 0.757  500  500
+++finished batch 8 cpu time: 0.688  500  500
+++finished batch 8 cpu time: 0.71  500  500
+++finished batch 8 cpu time: 0.619  500  500
+++finished batch 10 cpu time: 0.838  500  500
+++finished batch 9 cpu time: 0.767  500  500
finished simulation, cpu time = 0.851 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
+++finished batch 9 cpu time: 0.683  500  500
===== DE_ScoreDose::analyze =====
geometry: CT
```

```

cpu time: 0.851
number of histories: 5000
number of batches: 10
+++ beamlet 0
+++finished batch 9 cpu time: 0.794 500 500
+++finished batch 10 cpu time: 0.833 500 500
finished simulation, cpu time = 0.848 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
+++finished batch 10 cpu time: 0.864 500 500
+++finished batch 10 cpu time: 0.77 500 500
===== DE_ScoreDose::analyze =====
geometry: CT
cpu time: 0.848
number of histories: 5000
number of batches: 10
+++ beamlet 0
finished simulation, cpu time = 0.88 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
geometry: CT
cpu time: 0.88
number of histories: 5000
number of batches: 10
+++ beamlet 0
finished simulation, cpu time = 0.8 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
geometry: CT
cpu time: 0.8
number of histories: 5000
number of batches: 10
+++ beamlet 0
max dose is 0.0343716 in region 2035566
max dose is 0.0341074 in region 1805166
max dose is 0.0341867 in region 1881967
max dose is 0.0346977 in region 1958767
ICRU-2K efficiency: 975.077 1/s
+++ total
max dose is 0.0343716 in region 2035566
ICRU-2K efficiency: 975.077 1/s
=====
ICRU-2K efficiency: 988.245 1/s
+++ total
max dose is 0.0341074 in region 1805166
ICRU-2K efficiency: 988.245 1/s

```



```
=====
    ICRU-2K efficiency: 991.515 1/s
+++ total
    max dose is 0.0341867 in region 1881967
    ICRU-2K efficiency: 1135.3 1/s
+++ total
    max dose is 0.0346977 in region 1958767
    ICRU-2K efficiency: 991.515 1/s
=====
    ICRU-2K efficiency: 1135.3 1/s
=====
Completed 32 of 49 beamlets...
```

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_1.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_3.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_4.vmc

```
Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_2.vmc
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
```

```

region size: 0.3
geometry 1 region size: 0.3
OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
=====
                        Beamlet Source
=====
        charge = 0
    virtual source position = (-76,24,24)
        Energy = 6 MeV
    number of beamlets = 1
    beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,24.5,24.5)
=====
=====
                        XYZ Geometry
=====
        name: CT id: 0
    global smax: 1e+030
    Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
    Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
    Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
                        Monte Carlo Parameter
=====

Delta particle production threshold:    0.447479 MeV
Bremsstrahlung production threshold:    0.05 MeV
Min. electron transport energy         :    0.447479 MeV
Min. photon transport energy           :    0.05 MeV
Local track-end energy deposition       :    0
Cut-off energy for KERMA approx.        :    1.10239 MeV
Bremsstrahlung transport mode           :    1
CSDA approximation                     :    0
Fractional energy loss/step at Ep       :    10%
Max. 1st elastic moment per step        :    0.5
Max. acceptable energy loss/step        :    0.6 MeV
alpha and beta                          :    0.0298764  0.420741
Fano calculation                        :    0
Exact Compton                          :    0
Electron transport mode                  :    VMC++
=====
=====
                        MC_Control
=====
    will use fixed number of histories
    number of batches : 10
    histories per batch : 500
    total histories : 5000
    initial rng seeds : 26728 28779
=====

```

Example: Photon Treatment Plan
using VMC++ dose calculation

=====

Variance Reduction

=====

f_repeat = 0.251
split photons = 1
photon split factor = -40

=====

=====

Quasi Random Numbers

=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

=====

Scoring and output options

=====

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.411 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK

=====

Beamlet Source

=====

charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,24.5,25.5)

=====

=====

XYZ Geometry

=====

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV

Example: Photon Treatment Plan
using VMC++ dose calculation

```

Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition    : 0
Cut-off energy for KERMA approx.    : 1.10239 MeV
Bremsstrahlung transport mode       : 1
CSDA approximation                  : 0
Fractional energy loss/step at Ep   : 10%
Max. 1st elastic moment per step    : 0.5
Max. acceptable energy loss/step    : 0.6 MeV
alpha and beta                      : 0.0298764 0.420741
Fano calculation                    : 0
Exact Compton                       : 0
Electron transport mode              : VMC++

```

```

=====
=====
MC_Control
=====

```

```

will use fixed number of histories
number of batches      : 10
histories per batch    : 500
total histories        : 5000
initial rng seeds      : 4479 7726
=====

```

```

=====
Variance Reduction
=====

```

```

f_repeat      = 0.251
split photons = 1
photon split factor = -40
=====

```

```

=====
Quasi Random Numbers
=====

```

```

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1
=====

```

```

=====
Scoring and output options
=====

```

```

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

```

CPU time so far: 2.47 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Example: Photon Treatment Plan
using VMC++ dose calculation

Starting MC simulation

Running part 1 of 1 ...

+++finished batch 1 cpu time: 0.065 500 500

OK

=====

Beamlet Source

=====

charge = 0

virtual source position = (-76,24,24)

Energy = 6 MeV

number of beamlets = 1

beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,25,22.5)

=====

=====

XYZ Geometry

=====

name: CT id: 0

global smax: 1e+030

Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3

Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3

Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV

Bremsstrahlung production threshold: 0.05 MeV

Min. electron transport energy : 0.447479 MeV

Min. photon transport energy : 0.05 MeV

Local track-end energy deposition : 0

Cut-off energy for KERMA approx. : 1.10239 MeV

Bremsstrahlung transport mode : 1

CSDA approximation : 0

Fractional energy loss/step at Ep : 10%

Max. 1st elastic moment per step : 0.5

Max. acceptable energy loss/step : 0.6 MeV

alpha and beta : 0.0298764 0.420741

Fano calculation : 0

Exact Compton : 0

Electron transport mode : VMC++

=====

=====

MC_Control

=====

will use fixed number of histories

number of batches : 10

histories per batch : 500

total histories : 5000

initial rng seeds : 25222 7629

=====

=====

Variance Reduction

=====

Example: Photon Treatment Plan
using VMC++ dose calculation

```
=====
f_repeat          = 0.251
split photons     = 1
photon split factor = -40
=====
=====
Quasi Random Numbers
=====
number of generators: 1
                    1st: base = 2 dimensions = 60 warm-up = 1
=====
=====
Scoring and output options
=====
number of dose scoring objects: 1
dose output options for geometry CT
    dump dose: 2
    dose scans:

CPU time so far: 2.571 seconds

Will run approximately 5000 particle sets
    with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK
=====
Beamlet Source
=====
charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,24.5,25)
=====
=====
XYZ Geometry
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
Monte Carlo Parameter
=====
Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```

Min. photon transport energy      : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx.  : 1.10239 MeV
Bremsstrahlung transport mode     : 1
CSDA approximation                : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step  : 0.5
Max. acceptable energy loss/step  : 0.6 MeV
alpha and beta                    : 0.0298764 0.420741
Fano calculation                   : 0
Exact Compton                     : 0
Electron transport mode           : VMC++

```

```

=====
=====
MC_Control
=====

```

```

will use fixed number of histories
number of batches      : 10
histories per batch    : 500
total histories        : 5000
initial rng seeds      : 16417 4159

```

```

=====
=====
Variance Reduction
=====

```

```

f_repeat      = 0.251
split photons  = 1
photon split factor = -40

```

```

=====
=====
Quasi Random Numbers
=====

```

```

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

```

```

=====
=====
Scoring and output options
=====

```

```

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

```

CPU time so far: 2.674 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation

Example: Photon Treatment Plan
using VMC++ dose calculation

```
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.042 500 500
+++finished batch 2 cpu time: 0.135 500 500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.039 500 500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.045 500 500
+++finished batch 2 cpu time: 0.287 500 500
+++finished batch 3 cpu time: 0.35 500 500
+++finished batch 2 cpu time: 0.181 500 500
+++finished batch 2 cpu time: 0.11 500 500
+++finished batch 4 cpu time: 0.413 500 500
+++finished batch 3 cpu time: 0.246 500 500
+++finished batch 3 cpu time: 0.374 500 500
+++finished batch 3 cpu time: 0.175 500 500
+++finished batch 5 cpu time: 0.474 500 500
+++finished batch 4 cpu time: 0.311 500 500
+++finished batch 4 cpu time: 0.443 500 500
+++finished batch 4 cpu time: 0.237 500 500
+++finished batch 6 cpu time: 0.534 500 500
+++finished batch 5 cpu time: 0.37 500 500
+++finished batch 5 cpu time: 0.507 500 500
+++finished batch 7 cpu time: 0.591 500 500
+++finished batch 6 cpu time: 0.429 500 500
+++finished batch 5 cpu time: 0.318 500 500
+++finished batch 6 cpu time: 0.573 500 500
+++finished batch 8 cpu time: 0.657 500 500
+++finished batch 7 cpu time: 0.492 500 500
+++finished batch 7 cpu time: 0.63 500 500
+++finished batch 6 cpu time: 0.411 500 500
+++finished batch 9 cpu time: 0.721 500 500
+++finished batch 8 cpu time: 0.574 500 500
+++finished batch 8 cpu time: 0.707 500 500
+++finished batch 7 cpu time: 0.486 500 500
+++finished batch 10 cpu time: 0.801 500 500
+++finished batch 9 cpu time: 0.64 500 500
finished simulation, cpu time = 0.815 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
geometry: CT
cpu time: 0.815
number of histories: 5000
number of batches: 10
+++ beamlet 0
+++finished batch 9 cpu time: 0.777 500 500
+++finished batch 8 cpu time: 0.557 500 500
+++finished batch 10 cpu time: 0.699 500 500
finished simulation, cpu time = 0.713 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
```



```
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.713
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.832  500  500
+++finished batch 9  cpu time: 0.618  500  500
finished simulation, cpu time = 0.85 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.85
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.683  500  500
      max dose is 0.0344799 in region 2112366
finished simulation, cpu time = 0.697 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.697
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
      max dose is 0.0337888 in region 1805646
      max dose is 0.0342375 in region 2265967
      max dose is 0.0347981 in region 2189167
      ICRU-2K efficiency: 1043.99 1/s
+++ total
      max dose is 0.0344799 in region 2112366
      ICRU-2K efficiency: 1043.99 1/s
=====
      ICRU-2K efficiency: 1111.75 1/s
+++ total
      max dose is 0.0337888 in region 1805646
      ICRU-2K efficiency: 994.682 1/s
+++ total
      max dose is 0.0342375 in region 2265967
      ICRU-2K efficiency: 994.682 1/s
=====
      ICRU-2K efficiency: 1111.75 1/s
=====
      ICRU-2K efficiency: 1243.91 1/s
+++ total
      max dose is 0.0347981 in region 2189167
```

```
ICRU-2K efficiency: 1243.91 1/s
=====
Completed 36 of 49 beamlets...

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_1.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_2.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_3.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_4.vmc
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
=====
                        Beamlet Source
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```
=====
      charge = 0
virtual source position = (-76,24,24)
      Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,25,23)
=====
XYZ Geometry
=====
      name: CT id: 0
      global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
Monte Carlo Parameter
=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition   : 0
Cut-off energy for KERMA approx.    : 1.10239 MeV
Bremsstrahlung transport mode       : 1
CSDA approximation                  : 0
Fractional energy loss/step at Ep   : 10%
Max. 1st elastic moment per step    : 0.5
Max. acceptable energy loss/step    : 0.6 MeV
alpha and beta                       : 0.0298764 0.420741
Fano calculation                     : 0
Exact Compton                       : 0
Electron transport mode              : VMC++
=====
MC_Control
=====
will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 24429 7306
=====
Variance Reduction
=====

f_repeat = 0.251
split photons = 1
photon split factor = -40
=====
```

Example: Photon Treatment Plan
using VMC++ dose calculation

Quasi Random Numbers

```
=====
number of generators: 1
      1st:  base = 2 dimensions = 60 warm-up = 1
=====
Scoring and output options
=====
number of dose scoring objects: 1
dose output options for geometry CT
  dump dose: 2
  dose scans:
```

CPU time so far: 2.186 seconds

Will run approximately 5000 particle sets
 with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.055 500 500
OK

Beamlet Source

```
=====
charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,25,24)
=====
```

XYZ Geometry

```
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
```

Monte Carlo Parameter

```
=====
Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition   : 0
Cut-off energy for KERMA approx.    : 1.10239 MeV
Bremsstrahlung transport mode       : 1
CSDA approximation                  : 0
=====
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```

Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

```

```

=====
=====
MC_Control
=====

```

```

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 5898 7533
=====

```

```

=====
Variance Reduction
=====

```

```

f_repeat = 0.251
split photons = 1
photon split factor = -40
=====

```

```

=====
Quasi Random Numbers
=====

```

```

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1
=====

```

```

=====
Scoring and output options
=====

```

```

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

```

CPU time so far: 2.234 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK

```

=====
Beamlet Source
=====

```

```

charge = 0

```

Example: Photon Treatment Plan
using VMC++ dose calculation

```

virtual source position = (-76,24,24)
    Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,25,23.5)
=====
XYZ Geometry
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
Monte Carlo Parameter
=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition    : 0
Cut-off energy for KERMA approx.    : 1.10239 MeV
Bremsstrahlung transport mode       : 1
CSDA approximation                  : 0
Fractional energy loss/step at Ep   : 10%
Max. 1st elastic moment per step    : 0.5
Max. acceptable energy loss/step    : 0.6 MeV
alpha and beta                       : 0.0298764 0.420741
Fano calculation                     : 0
Exact Compton                       : 0
Electron transport mode              : VMC++
=====
MC_Control
=====
will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 27878 10500
=====
Variance Reduction
=====

f_repeat = 0.251
split photons = 1
photon split factor = -40
=====
Quasi Random Numbers
=====

```

Example: Photon Treatment Plan
using VMC++ dose calculation

```
number of generators: 1
      1st: base = 2 dimensions = 60 warm-up = 1
=====
Scoring and output options
=====
number of dose scoring objects: 1
dose output options for geometry CT
  dump dose: 2
  dose scans:

CPU time so far: 2.351 seconds

Will run approximately 5000 particle sets
  with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK
=====
Beamlet Source
=====
  charge = 0
  virtual source position = (-76,24,24)
  Energy = 6 MeV
  number of beamlets = 1
  beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,25,24.5)
=====
XYZ Geometry
=====
  name: CT id: 0
  global smax: 1e+030
  Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
  Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
  Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
Monte Carlo Parameter
=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
Min. photon transport energy        : 0.05 MeV
Local track-end energy deposition    : 0
Cut-off energy for KERMA approx.    : 1.10239 MeV
Bremsstrahlung transport mode       : 1
CSDA approximation                  : 0
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.043 500 500
Fractional energy loss/step at Ep   : 10%
Max. 1st elastic moment per step    : 0.5
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```

Max. acceptable energy loss/step   :   0.6 MeV
alpha and beta                     :   0.0298764   0.420741
Fano calculation                   :   0
Exact Compton                     :   0
Electron transport mode            :   VMC++

```

```

=====
MC_Control
=====

```

```

will use fixed number of histories
number of batches   : 10
histories per batch : 500
total histories     : 5000
initial rng seeds   : 18482 14199

```

```

=====
Variance Reduction
=====

```

```

f_repeat           = 0.251
split photons      = 1
photon split factor = -40

```

```

=====
Quasi Random Numbers
=====

```

```

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

```

```

=====
Scoring and output options
=====

```

```

number of dose scoring objects: 1
dose output options for geometry CT
  dump dose: 2
  dose scans:

```

CPU time so far: 2.456 seconds

Will run approximately 5000 particle sets
 with 1 particles per set on average
 will use 2 quasi numbers to sample the source

Starting MC simulation

```

+++finished batch 2 cpu time: 0.147   500   500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.042   500   500
+++finished batch 2 cpu time: 0.181   500   500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.042   500   500
+++finished batch 3 cpu time: 0.354   500   500

```


Example: Photon Treatment Plan
using VMC++ dose calculation

```

+++finished batch 2 cpu time: 0.188 500 500
+++finished batch 3 cpu time: 0.326 500 500
+++finished batch 2 cpu time: 0.119 500 500
+++finished batch 4 cpu time: 0.44 500 500
+++finished batch 3 cpu time: 0.283 500 500
+++finished batch 4 cpu time: 0.413 500 500
+++finished batch 3 cpu time: 0.196 500 500
+++finished batch 4 cpu time: 0.354 500 500
+++finished batch 5 cpu time: 0.541 500 500
+++finished batch 5 cpu time: 0.493 500 500
+++finished batch 4 cpu time: 0.276 500 500
+++finished batch 5 cpu time: 0.422 500 500
+++finished batch 6 cpu time: 0.561 500 500
+++finished batch 6 cpu time: 0.627 500 500
+++finished batch 5 cpu time: 0.358 500 500
+++finished batch 6 cpu time: 0.517 500 500
+++finished batch 7 cpu time: 0.696 500 500
+++finished batch 7 cpu time: 0.656 500 500
+++finished batch 6 cpu time: 0.449 500 500
+++finished batch 7 cpu time: 0.582 500 500
+++finished batch 8 cpu time: 0.78 500 500
+++finished batch 7 cpu time: 0.515 500 500
+++finished batch 8 cpu time: 0.74 500 500
+++finished batch 8 cpu time: 0.659 500 500
+++finished batch 9 cpu time: 0.856 500 500
+++finished batch 8 cpu time: 0.589 500 500
+++finished batch 9 cpu time: 0.823 500 500
+++finished batch 9 cpu time: 0.733 500 500
+++finished batch 10 cpu time: 0.925 500 500
+++finished batch 10 cpu time: 0.891 500 500
finished simulation, cpu time = 0.954 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.954
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
+++finished batch 9 cpu time: 0.674 500 500
finished simulation, cpu time = 0.92 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
+++finished batch 10 cpu time: 0.815 500 500
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.92
  number of histories: 5000
  number of batches: 10
+++ beamlet 0

```

Example: Photon Treatment Plan
using VMC++ dose calculation

```
finished simulation, cpu time = 0.835 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.835
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.734 500 500
finished simulation, cpu time = 0.758 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.758
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
    max dose is 0.0340882 in region 1882447
    max dose is 0.0342205 in region 2036047
    max dose is 0.0343793 in region 1959246
    max dose is 0.0343678 in region 2112847
    ICRU-2K efficiency: 915.378 1/s
+++ total
    max dose is 0.0342205 in region 2036047
    ICRU-2K efficiency: 877.636 1/s
+++ total
    max dose is 0.0340882 in region 1882447
    ICRU-2K efficiency: 915.378 1/s
=====
    ICRU-2K efficiency: 877.636 1/s
=====
    ICRU-2K efficiency: 1092.44 1/s
+++ total
    max dose is 0.0343793 in region 1959246
    ICRU-2K efficiency: 1110.6 1/s
+++ total
    max dose is 0.0343678 in region 2112847
    ICRU-2K efficiency: 1092.44 1/s
=====
    ICRU-2K efficiency: 1110.6 1/s
=====
Completed 40 of 49 beamlets...
```

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_1.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_3.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_4.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_2.vmc

```
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
  geometry 1 region size: 0.3
OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
=====
                        Beamlet Source
=====
      charge = 0
virtual source position = (-76,24,24)
      Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,25.5,23)
=====
=====
                        XYZ Geometry
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====

Monte Carlo Parameter
=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====
=====

MC_Control
=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 22717 22612
=====
=====

Variance Reduction
=====

f_repeat = 0.251
split photons = 1
photon split factor = -40
=====
=====

Quasi Random Numbers
=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1
=====
=====

Scoring and output options
=====

number of dose scoring objects: 1
=====
```

Example: Photon Treatment Plan
using VMC++ dose calculation

dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.459 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK

=====

Beamlet Source

=====

charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,25,25)

=====

=====

XYZ Geometry

=====

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

=====

MC_Control

Example: Photon Treatment Plan
using VMC++ dose calculation

```
=====
will use fixed number of histories
number of batches      : 10
histories per batch    : 500
total histories        : 5000
initial rng seeds      : 10550  24925
=====
=====
Variance Reduction
=====
=====

f_repeat              = 0.251
split photons         = 1
photon split factor   = -40
=====
=====
Quasi Random Numbers
=====
=====
number of generators: 1
                    1st: base = 2 dimensions = 60 warm-up = 1
=====
=====
Scoring and output options
=====
=====
number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
OK
=====
Beamlet Source
=====
=====
charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,25.5,22.5)
=====
=====
XYZ Geometry
=====
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
dose scans:

CPU time so far: 2.566 seconds

Will run approximately 5000 particle sets
    with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
```

Example: Photon Treatment Plan
using VMC++ dose calculation

Number of y-planes: 160 uniform with $Y_0 = 0.15$ $Dx = 0.3$
Number of z-planes: 160 uniform with $Z_0 = 0.15$ $Dx = 0.3$

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at E_p : 10%
Max. 1st elastic moment per step : 0.5
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.048 500 500
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 27516 8576

=====

Variance Reduction

=====

$f_{\text{repeat}} = 0.251$
split photons = 1
photon split factor = -40

=====

Quasi Random Numbers

=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

Scoring and output options

=====

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2

Example: Photon Treatment Plan
using VMC++ dose calculation

dose scans:

CPU time so far: 2.671 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation

OK

=====

Beamlet Source

=====

charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,25,25.5)

=====

=====

XYZ Geometry

=====

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

=====

MC_Control

=====

will use fixed number of histories

Example: Photon Treatment Plan
using VMC++ dose calculation

```
number of batches      : 10
histories per batch    : 500
total histories        : 5000
initial rng seeds      : 17558 16492
=====
=====
Variance Reduction
=====
=====
f_repeat              = 0.251
split photons         = 1
photon split factor   = -40
=====
=====
Quasi Random Numbers
=====
=====
number of generators: 1
                    1st: base = 2 dimensions = 60 warm-up = 1
=====
=====
Scoring and output options
=====
=====
number of dose scoring objects: 1
dose output options for geometry CT
    dump dose: 2
    dose scans:

CPU time so far: 2.654 seconds

Will run approximately 5000 particle sets
    with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.042  500  500
+++finished batch 2 cpu time: 0.157  500  500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.04  500  500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.048  500  500
+++finished batch 2 cpu time: 0.165  500  500
+++finished batch 3 cpu time: 0.277  500  500
+++finished batch 2 cpu time: 0.105  500  500
+++finished batch 2 cpu time: 0.126  500  500
+++finished batch 3 cpu time: 0.238  500  500
+++finished batch 4 cpu time: 0.358  500  500
+++finished batch 3 cpu time: 0.173  500  500
+++finished batch 4 cpu time: 0.304  500  500
+++finished batch 3 cpu time: 0.2  500  500
+++finished batch 5 cpu time: 0.429  500  500
+++finished batch 4 cpu time: 0.242  500  500
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```
+++finished batch 5 cpu time: 0.364 500 500
+++finished batch 4 cpu time: 0.27 500 500
+++finished batch 6 cpu time: 0.487 500 500
+++finished batch 5 cpu time: 0.301 500 500
+++finished batch 6 cpu time: 0.423 500 500
+++finished batch 5 cpu time: 0.337 500 500
+++finished batch 7 cpu time: 0.547 500 500
+++finished batch 6 cpu time: 0.361 500 500
+++finished batch 7 cpu time: 0.511 500 500
+++finished batch 8 cpu time: 0.624 500 500
+++finished batch 6 cpu time: 0.419 500 500
+++finished batch 7 cpu time: 0.435 500 500
+++finished batch 8 cpu time: 0.59 500 500
+++finished batch 9 cpu time: 0.696 500 500
+++finished batch 7 cpu time: 0.493 500 500
+++finished batch 8 cpu time: 0.506 500 500
+++finished batch 9 cpu time: 0.652 500 500
+++finished batch 10 cpu time: 0.758 500 500
finished simulation, cpu time = 0.775 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
+++finished batch 8 cpu time: 0.565 500 500
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.775
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
+++finished batch 9 cpu time: 0.575 500 500
+++finished batch 10 cpu time: 0.728 500 500
+++finished batch 10 cpu time: 0.635 500 500
+++finished batch 9 cpu time: 0.635 500 500
finished simulation, cpu time = 0.746 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.746
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
finished simulation, cpu time = 0.653 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.653
  number of histories: 5000
```

```
number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.731 500 500
    max dose is 0.0342376 in region 1882926
finished simulation, cpu time = 0.744 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
geometry: CT
cpu time: 0.744
number of histories: 5000
number of batches: 10
+++ beamlet 0
    max dose is 0.0340721 in region 2189646
    max dose is 0.033951 in region 1806126
    max dose is 0.0350693 in region 2266446
    ICRU-2K efficiency: 1080.68 1/s
+++ total
    max dose is 0.0342376 in region 1882926
    ICRU-2K efficiency: 1080.68 1/s
=====
    ICRU-2K efficiency: 1074.01 1/s
+++ total
    max dose is 0.0340721 in region 2189646
    ICRU-2K efficiency: 1290.31 1/s
+++ total
    max dose is 0.033951 in region 1806126
    ICRU-2K efficiency: 1074.01 1/s
=====
    ICRU-2K efficiency: 1290.31 1/s
=====
    ICRU-2K efficiency: 1178.38 1/s
+++ total
    max dose is 0.0350693 in region 2266446
    ICRU-2K efficiency: 1178.38 1/s
=====
Completed 44 of 49 beamlets...
```

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_1.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_2.vmc

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_3.vmc

```

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_4.vmc
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
Initializing cross sections ... OK
=====
                        Beamlet Source
=====
        charge = 0
        virtual source position = (-76,24,24)
        Energy = 6 MeV
        number of beamlets = 1
        beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,25.5,25)
=====
=====
                        XYZ Geometry
=====
        name: CT id: 0
        global smax: 1e+030
        Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
        Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
        Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
                        Monte Carlo Parameter
=====

```

Example: Photon Treatment Plan
using VMC++ dose calculation

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at E_p : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 28021 3898

=====

Variance Reduction

=====

f_{repeat} = 0.251
split photons = 1
photon split factor = -40

=====

Quasi Random Numbers

=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

Scoring and output options

=====

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

CPU time so far: 2.206 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average

Example: Photon Treatment Plan
using VMC++ dose calculation

will use 2 quasi numbers to sample the source

Starting MC simulation

OK

=====

Beamlet Source

=====

charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,25.5,24.5)

=====

=====

XYZ Geometry

=====

name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3

=====

Monte Carlo Parameter

=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

=====

MC_Control

=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 15924 23376

=====

Variance Reduction

=====

Example: Photon Treatment Plan
using VMC++ dose calculation

```
=====
f_repeat          = 0.251
split photons     = 1
photon split factor = -40
=====
=====
Quasi Random Numbers
=====
number of generators: 1
                    1st: base = 2 dimensions = 60 warm-up = 1
=====
=====
Scoring and output options
=====
number of dose scoring objects: 1
dose output options for geometry CT
    dump dose: 2
    dose scans:

CPU time so far: 2.242 seconds

Will run approximately 5000 particle sets
    with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
OK
=====
Beamlet Source
=====
charge = 0
virtual source position = (-76,24,24)
Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,25.5,24)
=====
=====
XYZ Geometry
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====
Monte Carlo Parameter
=====
Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy      : 0.447479 MeV
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```

Min. photon transport energy      : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx.  : 1.10239 MeV
Bremsstrahlung transport mode     : 1
CSDA approximation                : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step  : 0.5
Max. acceptable energy loss/step  : 0.6 MeV
alpha and beta                    : 0.0298764 0.420741
Fano calculation                  : 0
Exact Compton                     : 0
Electron transport mode           : VMC++

```

```

=====
=====
MC_Control
=====

```

```

will use fixed number of histories
number of batches      : 10
histories per batch    : 500
total histories        : 5000
initial rng seeds      : 2276 1619
=====

```

```

=====
Variance Reduction
=====

```

```

f_repeat      = 0.251
split photons  = 1
photon split factor = -40
=====

```

```

=====
Quasi Random Numbers
=====

```

```

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1
=====

```

```

=====
Scoring and output options
=====

```

```

number of dose scoring objects: 1
dose output options for geometry CT
dump dose: 2
dose scans:

```

CPU time so far: 2.359 seconds

Will run approximately 5000 particle sets
with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation

Example: Photon Treatment Plan
using VMC++ dose calculation

```
OK
=====
                        Beamlet Source
=====
      charge = 0
virtual source position = (-76,24,24)
      Energy = 6 MeV
number of beamlets = 1
beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,25.5,23.5)
=====
=====
                        XYZ Geometry
=====
      name: CT   id: 0
      global smax: 1e+030
Number of x-planes: 160   uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160   uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160   uniform with Zo = 0.15 Dx = 0.3
=====
=====
                        Monte Carlo Parameter
=====

Delta particle production threshold:    0.447479 MeV
Bremsstrahlung production threshold:    0.05 MeV
Min. electron transport energy         :    0.447479 MeV
Min. photon transport energy           :    0.05 MeV
Local track-end energy deposition       :    0
Cut-off energy for KERMA approx.        :    1.10239 MeV
Bremsstrahlung transport mode           :    1
CSDA approximation                     :    0
Fractional energy loss/step at Ep       :    10%
Max. 1st elastic moment per step        :    0.5
Max. acceptable energy loss/step        :    0.6 MeV
alpha and beta                          :    0.0298764  0.420741
Fano calculation                        :    0
Exact Compton                          :    0
Electron transport mode                  :    VMC++

=====
=====
                        MC_Control
=====
      will use fixed number of histories
      number of batches   : 10
      histories per batch : 500
      total histories     : 5000
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.057  500  500
      initial rng seeds   : 11414  17035
=====
=====
                        Variance Reduction
=====
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```
f_repeat          = 0.251
split photons     = 1
photon split factor = -40
=====
=====
                        Quasi Random Numbers
=====
number of generators: 1
                    1st: base = 2 dimensions = 60 warm-up = 1
=====
=====
                        Scoring and output options
=====
number of dose scoring objects: 1
dose output options for geometry CT
    dump dose: 2
    dose scans:

CPU time so far: 2.364 seconds

Will run approximately 5000 particle sets
    with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.041  500  500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.041  500  500
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.058  500  500
+++finished batch 2 cpu time: 0.235  500  500
+++finished batch 2 cpu time: 0.224  500  500
+++finished batch 2 cpu time: 0.123  500  500
+++finished batch 2 cpu time: 0.155  500  500
+++finished batch 3 cpu time: 0.322  500  500
+++finished batch 3 cpu time: 0.195  500  500
+++finished batch 3 cpu time: 0.324  500  500
+++finished batch 3 cpu time: 0.223  500  500
+++finished batch 4 cpu time: 0.411  500  500
+++finished batch 4 cpu time: 0.281  500  500
+++finished batch 4 cpu time: 0.395  500  500
+++finished batch 4 cpu time: 0.308  500  500
+++finished batch 5 cpu time: 0.491  500  500
+++finished batch 5 cpu time: 0.355  500  500
+++finished batch 5 cpu time: 0.469  500  500
+++finished batch 5 cpu time: 0.397  500  500
+++finished batch 6 cpu time: 0.563  500  500
+++finished batch 6 cpu time: 0.422  500  500
+++finished batch 6 cpu time: 0.554  500  500
+++finished batch 6 cpu time: 0.471  500  500
+++finished batch 7 cpu time: 0.626  500  500
```

```

+++finished batch 7 cpu time: 0.623 500 500
+++finished batch 7 cpu time: 0.514 500 500
+++finished batch 7 cpu time: 0.55 500 500
+++finished batch 8 cpu time: 0.71 500 500
+++finished batch 8 cpu time: 0.588 500 500
+++finished batch 8 cpu time: 0.712 500 500
+++finished batch 8 cpu time: 0.625 500 500
+++finished batch 9 cpu time: 0.79 500 500
+++finished batch 9 cpu time: 0.783 500 500
+++finished batch 9 cpu time: 0.678 500 500
+++finished batch 9 cpu time: 0.716 500 500
+++finished batch 10 cpu time: 0.868 500 500
finished simulation, cpu time = 0.895 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
+++finished batch 10 cpu time: 0.863 500 500
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.895
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.756 500 500
finished simulation, cpu time = 0.783 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.783
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
finished simulation, cpu time = 0.897 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.897
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
+++finished batch 10 cpu time: 0.797 500 500
finished simulation, cpu time = 0.829 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====

```

```

geometry: CT
cpu time: 0.829
number of histories: 5000
number of batches: 10
+++ beamlet 0
    max dose is 0.0343218 in region 2190126
    max dose is 0.0341925 in region 2036526
    max dose is 0.0342907 in region 2113326
    max dose is 0.0339002 in region 1959726
    ICRU-2K efficiency: 931.074 1/s
+++ total
    max dose is 0.0343218 in region 2190126
    ICRU-2K efficiency: 931.074 1/s
=====
    ICRU-2K efficiency: 1078.04 1/s
+++ total
    max dose is 0.0341925 in region 2036526
    ICRU-2K efficiency: 1078.04 1/s
=====
    ICRU-2K efficiency: 1054.03 1/s
+++ total
    max dose is 0.0339002 in region 1959726
    ICRU-2K efficiency: 994.057 1/s
+++ total
    max dose is 0.0342907 in region 2113326
    ICRU-2K efficiency: 1054.03 1/s
=====
    ICRU-2K efficiency: 994.057 1/s
=====
Completed 48 of 49 beamlets...

Input file is: C:\Home\Bangertm\Git\matRad\vmc++\runs/
MCpencilbeam_temp_1.vmc
Reading MS data ... OK
Parsing input file ... construct_mmc_geometry: urs = 1
DE_GeometryFactory::get_region_size():
get_region_size: urs = 1 ignore = 1
region size: 0.3
    geometry 1 region size: 0.3
    OK
Initializing cross sections ... OK
=====
                        Beamlet Source
=====
    charge = 0
    virtual source position = (-76,24,24)
    Energy = 6 MeV
    number of beamlets = 1
    beamlet 1: size = 0.5x0.5 cm**2 midpoint = (-26,25.5,25.5)
=====
=====
                        XYZ Geometry

```

Example: Photon Treatment Plan
using VMC++ dose calculation

```
=====
name: CT id: 0
global smax: 1e+030
Number of x-planes: 160 uniform with Xo = 0.15 Dx = 0.3
Number of y-planes: 160 uniform with Yo = 0.15 Dx = 0.3
Number of z-planes: 160 uniform with Zo = 0.15 Dx = 0.3
=====

Monte Carlo Parameter
=====

Delta particle production threshold: 0.447479 MeV
Bremsstrahlung production threshold: 0.05 MeV
Min. electron transport energy : 0.447479 MeV
Min. photon transport energy : 0.05 MeV
Local track-end energy deposition : 0
Cut-off energy for KERMA approx. : 1.10239 MeV
Bremsstrahlung transport mode : 1
CSDA approximation : 0
Fractional energy loss/step at Ep : 10%
Max. 1st elastic moment per step : 0.5
Max. acceptable energy loss/step : 0.6 MeV
alpha and beta : 0.0298764 0.420741
Fano calculation : 0
Exact Compton : 0
Electron transport mode : VMC++

=====

MC_Control
=====

will use fixed number of histories
number of batches : 10
histories per batch : 500
total histories : 5000
initial rng seeds : 17065 14082

=====

Variance Reduction
=====

f_repeat = 0.251
split photons = 1
photon split factor = -40

=====

Quasi Random Numbers
=====

number of generators: 1
1st: base = 2 dimensions = 60 warm-up = 1

=====

Scoring and output options
=====

number of dose scoring objects: 1
```

```
dose output options for geometry CT
  dump dose: 2
  dose scans:

CPU time so far: 1.745 seconds

Will run approximately 5000 particle sets
  with 1 particles per set on average
will use 2 quasi numbers to sample the source

Starting MC simulation
Running part 1 of 1 ...
+++finished batch 1 cpu time: 0.036  500  500
+++finished batch 2 cpu time: 0.086  500  500
+++finished batch 3 cpu time: 0.139  500  500
+++finished batch 4 cpu time: 0.191  500  500
+++finished batch 5 cpu time: 0.245  500  500
+++finished batch 6 cpu time: 0.303  500  500
+++finished batch 7 cpu time: 0.364  500  500
+++finished batch 8 cpu time: 0.408  500  500
+++finished batch 9 cpu time: 0.457  500  500
+++finished batch 10 cpu time: 0.507  500  500
finished simulation, cpu time = 0.522 seconds
total particle fluence from the source: 20000
total number of particle sets: 5000
total number of particles: 5000
average sampled energy: 6 +/- 0
===== DE_ScoreDose::analyze =====
  geometry: CT
  cpu time: 0.522
  number of histories: 5000
  number of batches: 10
+++ beamlet 0
  max dose is 0.034158 in region 2266926
  ICRU-2K efficiency: 1586.02 1/s
+++ total
  max dose is 0.034158 in region 2266926
  ICRU-2K efficiency: 1586.02 1/s
=====
Completed 49 of 49 beamlets...
```

Inverse Optimization for IMRT

```
resultGUI = matRad_fluenceOptimization(dij,cst,pln);
```

This is Ipopt version 3.11.8, running with linear solver ma57.

```
Number of nonzeros in equality constraint Jacobian...:      0
Number of nonzeros in inequality constraint Jacobian.:      0
Number of nonzeros in Lagrangian Hessian.....:      0
```

Example: Photon Treatment Plan
using VMC++ dose calculation

```

Total number of variables.....: 49
      variables with only lower bounds: 49
      variables with lower and upper bounds: 0
      variables with only upper bounds: 0
Total number of equality constraints.....: 0
Total number of inequality constraints.....: 0
      inequality constraints with only lower bounds: 0
      inequality constraints with lower and upper bounds: 0
      inequality constraints with only upper bounds: 0

```

```

iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr  ls
  0  4.7141751e+001  0.00e+000  1.40e+001   0.0  0.00e+000   -  0.00e
+000  0.00e+000   0
  1  5.5399399e+001  0.00e+000  1.99e+001   0.7  6.17e+000   -
  4.91e-001  1.76e-002f  5
  2  4.2089461e+001  0.00e+000  1.05e+001   0.2  5.45e-002   -  1.00e
+000  1.00e+000f  1
  3  4.1347678e+001  0.00e+000  1.35e+000  -1.1  1.99e-002   -
  9.99e-001  1.00e+000f  1
  4  4.1316763e+001  0.00e+000  9.35e-001  -2.8  4.59e-003   -
  9.99e-001  1.00e+000f  1
  5  4.1284000e+001  0.00e+000  2.40e-001  -4.5  1.22e-002   -  1.00e
+000  1.00e+000f  1
  6  4.1281126e+001  0.00e+000  3.16e-001  -6.1  3.71e-003   -  1.00e
+000  1.00e+000f  1

```

Number of Iterations.....: 6

```

                                     (scaled)                               (unscaled)
Objective.....: 4.1281126121699344e+001
4.1281126121699344e+001
Dual infeasibility.....: 3.1588500984810963e-001
3.1588500984810963e-001
Constraint violation.....: 0.0000000000000000e+000
0.0000000000000000e+000
Complementarity.....: 8.5126569666667240e-007
8.5126569666667240e-007
Overall NLP error.....: 3.1588500984810963e-001
3.1588500984810963e-001

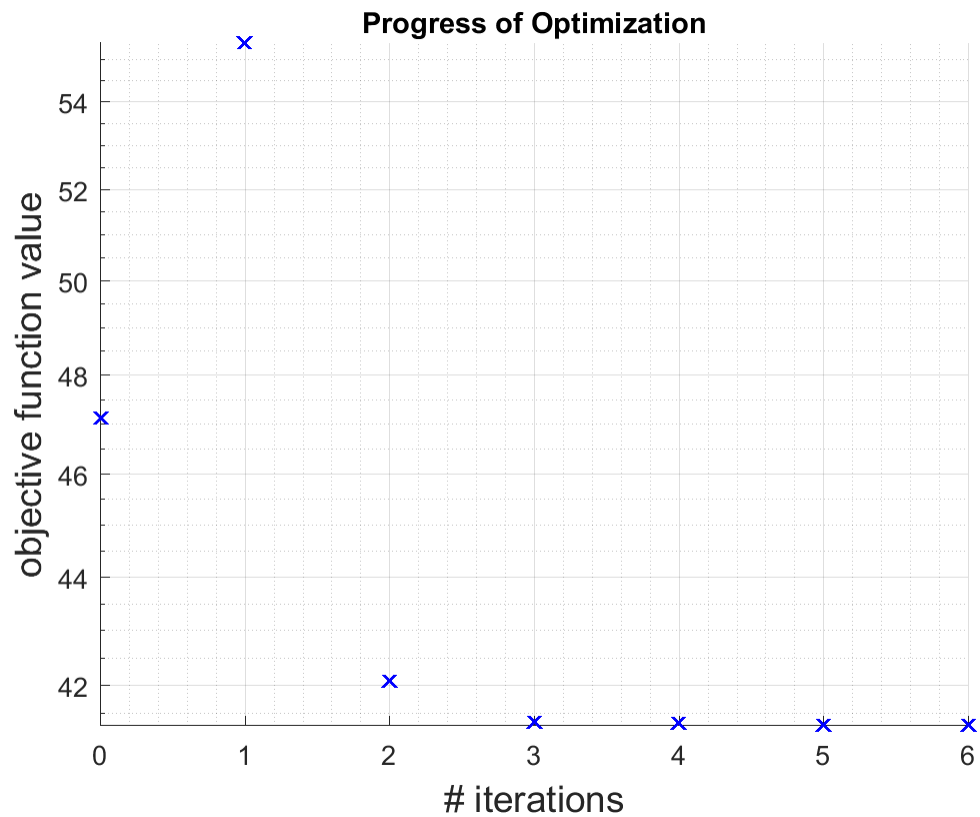
```

```

Number of objective function evaluations      = 15
Number of objective gradient evaluations      = 7
Number of equality constraint evaluations      = 0
Number of inequality constraint evaluations    = 0
Number of equality constraint Jacobian evaluations = 0
Number of inequality constraint Jacobian evaluations = 0
Number of Lagrangian Hessian evaluations     = 0
Total CPU secs in IPOPT (w/o function evaluations) = 0.463
Total CPU secs in NLP function evaluations    = 0.810

```

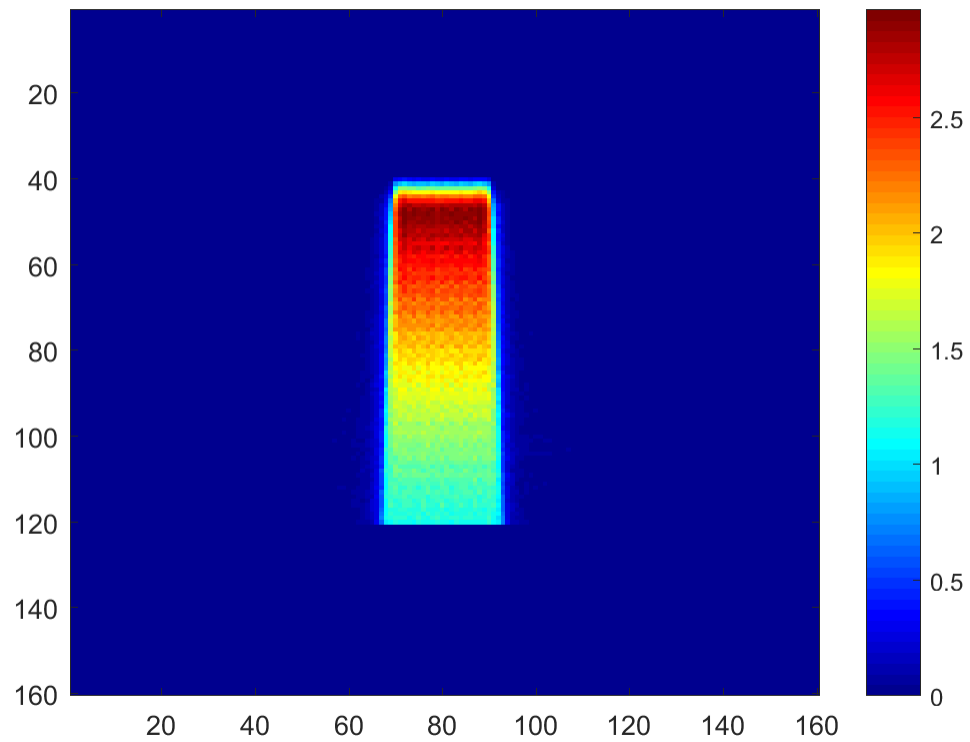
EXIT: Solved To Acceptable Level.
Calculating final cubes...



Plot the Resulting Dose Slice

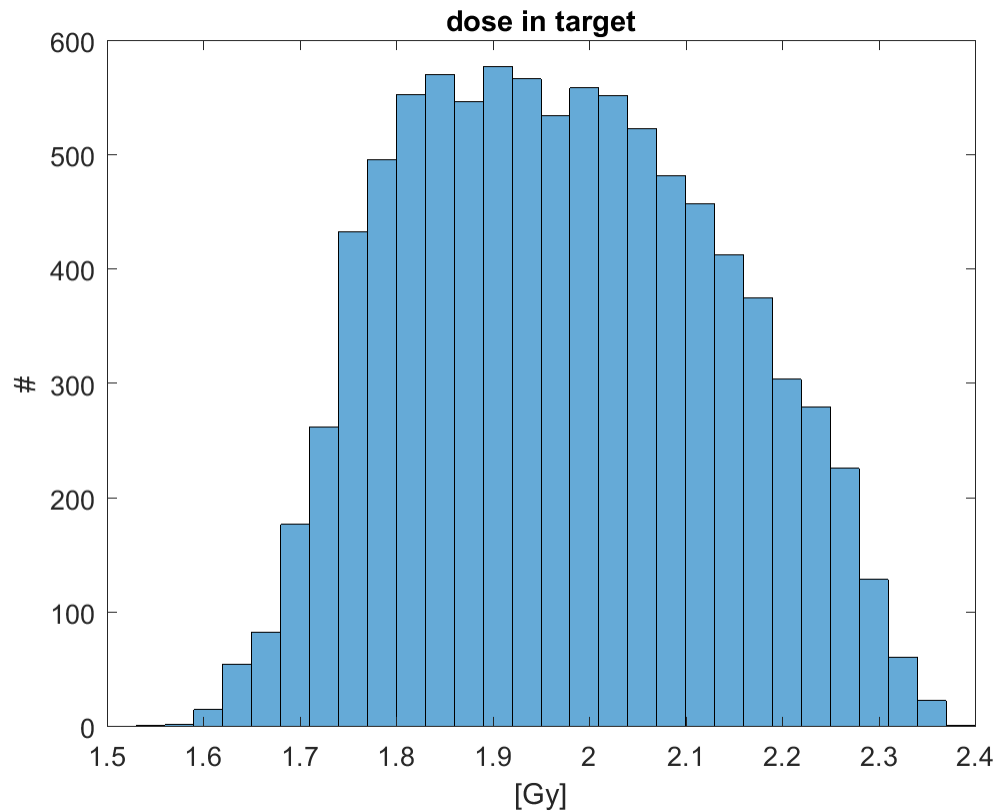
Just let's plot the transversal iso-center dose slice

```
slice = round(pln.isoCenter(1,3)./ct.resolution.z);  
figure,  
imagesc(resultGUI.physicalDose(:,:,slice)),colorbar, colormap(jet)
```

Exemplary, we show how to obtain the dose in the target and plot the histogram

```
ixTarget      = cst{2,4}{1};  
doseInTarget = resultGUI.physicalDose(ixTarget);  
figure  
histogram(doseInTarget);  
title('dose in target'),xlabel('[Gy]'),ylabel('#');
```



Start the GUI for Visualization

matRadGUI

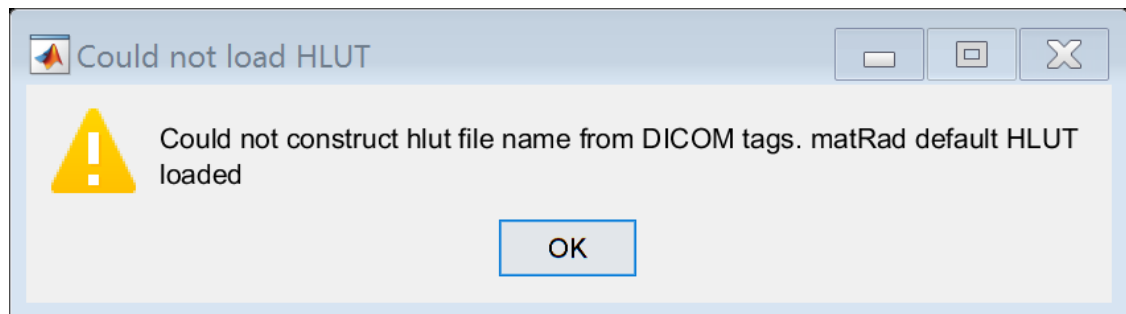
Warning: matRad default HLUT loaded

Reconversion of HU values could not be done because HLUT is not bijective.

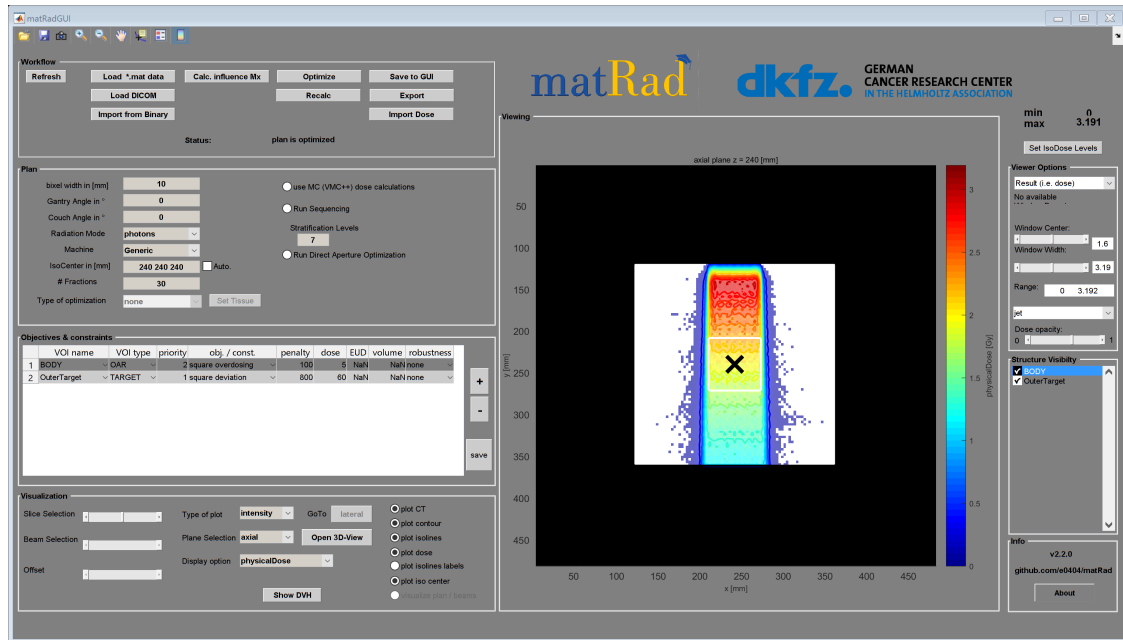
Warning: 'popupmenu' control requires that 'Value' be an integer within

String range

Control will not be rendered until all of its parameter values are valid



Example: Photon Treatment Plan using VMC++ dose calculation



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