
Example: Generate your own phantom geometry

Table of Contents

.....	1
Create a CT image series	1
Create the VOI data for the phantom	2
Lets create either a cubic or a spheric phantom	2
Assign relative electron densities	5
Treatment Plan	5
Generate Beam Geometry STF	5
Dose Calculation	6
Inverse Optimization for intensity-modulated photon therapy	6
Plot the resulting dose slice	8

%%

Copyright 2018 the matRad development team.

This file is part of the matRad project. It is subject to the license terms in the LICENSE file found in the top-level directory of this distribution and at <https://github.com/e0404/matRad/LICENSES.txt>. No part of the matRad project, including this file, may be copied, modified, propagated, or distributed except according to the terms contained in the LICENSE file.

%%

In this example we will show (i) how to create arbitrary ct data (resolution, ct numbers) (ii) how to create a cst structure containing the volume of interests of the phantom (iii) generate a treatment plan for this phantom

```
clc, clear, close all
```

Create a CT image series

```
xDim = 200;
yDim = 200;
zDim = 50;

ct.cubeDim      = [xDim yDim zDim];
ct.resolution.x = 2;
ct.resolution.y = 2;
ct.resolution.z = 3;
ct.numOfCtScen  = 1;

% create an ct image series with zeros - it will be filled later
ct.cubeHU{1} = ones(ct.cubeDim) * -1000; % assign HU of Air
```

Create the VOI data for the phantom

Now we define structures a contour for the phantom and a target

```
ixOAR = 1;
ixPTV = 2;

% define general VOI properties
cst{ixOAR,1} = 0;
cst{ixOAR,2} = 'contour';
cst{ixOAR,3} = 'OAR';

cst{ixPTV,1} = 1;
cst{ixPTV,2} = 'target';
cst{ixPTV,3} = 'TARGET';

% define optimization parameter for both VOIs
cst{ixOAR,5}.TissueClass = 1;
cst{ixOAR,5}.alphaX      = 0.1000;
cst{ixOAR,5}.betaX       = 0.0500;
cst{ixOAR,5}.Priority     = 2;
cst{ixOAR,5}.Visible     = 1;
cst{ixOAR,6}.type        = 'square overdosing';
cst{ixOAR,6}.dose        = 30;
cst{ixOAR,6}.penalty     = 10;
cst{ixOAR,6}.EUD         = NaN;
cst{ixOAR,6}.volume      = NaN;
cst{ixOAR,6}.coverage    = NaN;
cst{ixOAR,6}.robustness  = 'none';

cst{ixPTV,5}.TissueClass = 1;
cst{ixPTV,5}.alphaX      = 0.1000;
cst{ixPTV,5}.betaX       = 0.0500;
cst{ixPTV,5}.Priority     = 1;
cst{ixPTV,5}.Visible     = 1;
cst{ixPTV,6}.type        = 'square deviation';
cst{ixPTV,6}.dose        = 60;
cst{ixPTV,6}.penalty     = 50;
cst{ixPTV,6}.EUD         = NaN;
cst{ixPTV,6}.volume      = NaN;
cst{ixPTV,6}.coverage    = NaN;
cst{ixPTV,6}.robustness  = 'none';
```

Lets create either a cubic or a spheric phantom

```
TYPE = 'spheric'; % either 'cubic' or 'spheric'

% first the OAR
cubeHelper = zeros(ct.cubeDim);

switch TYPE

    case {'cubic'}
```

```
xLowOAR = round(xDim/2 - xDim/4);
xHighOAR = round(xDim/2 + xDim/4);
yLowOAR = round(yDim/2 - yDim/4);
yHighOAR = round(yDim/2 + yDim/4);
zLowOAR = round(zDim/2 - zDim/4);
zHighOAR = round(zDim/2 + zDim/4);

for x = xLowOAR:1:xHighOAR
    for y = yLowOAR:1:yHighOAR
        for z = zLowOAR:1:zHighOAR
            cubeHelper(x,y,z) = 1;
        end
    end
end

case {'spheric'}

    radiusOAR = xDim/4;

    for x = 1:xDim
        for y = 1:yDim
            for z = 1:zDim
                currPost = [x y z] - round([ct.cubeDim./2]);
                if sqrt(sum(currPost.^2)) < radiusOAR
                    cubeHelper(x,y,z) = 1;
                end
            end
        end
    end

end

% extract the voxel indices and save it in the cst
cst{ixOAR,4}{1} = find(cubeHelper);

% second the PTV
cubeHelper = zeros(ct.cubeDim);

switch TYPE

    case {'cubic'}

        xLowPTV = round(xDim/2 - xDim/8);
        xHighPTV = round(xDim/2 + xDim/8);
        yLowPTV = round(yDim/2 - yDim/8);
        yHighPTV = round(yDim/2 + yDim/8);
        zLowPTV = round(zDim/2 - zDim/8);
        zHighPTV = round(zDim/2 + zDim/8);

        cubeHelper = zeros(ct.cubeDim);

        for x = xLowPTV:1:xHighPTV
```

```
        for y = yLowPTV:1:yHighPTV
            for z = zLowPTV:1:zHighPTV
                cubeHelper(x,y,z) = 1;
            end
        end
    end
end

case {'spheric'}

    radiusPTV = xDim/12;

    for x = 1:xDim
        for y = 1:yDim
            for z = 1:zDim
                currPost = [x y z] - round([ct.cubeDim./2]);
                if sqrt(sum(currPost.^2)) < radiusPTV
                    cubeHelper(x,y,z) = 1;
                end
            end
        end
    end
end

% extract the voxel indices and save it in the cst
cst{ixPTV,4}{1} = find(cubeHelper);

% now we have ct data and cst data for a new phantom
display(ct);
display(cst);

ct =

    struct with fields:

        cubeDim: [200 200 50]
        resolution: [1x1 struct]
        numOfCtScen: 1
        cubeHU: {[200x200x50 double]}

cst =

    2x6 cell array

    Columns 1 through 5

        {[0]}        {'contour'}        {'OAR'}        {1x1 cell}        {1x1 struct}
        {[1]}        {'target'}        {'TARGET'}        {1x1 cell}        {1x1 struct}
```

Column 6

```
{1x1 struct}  
{1x1 struct}
```

Assign relative electron densities

```
vIxoAR = cst{ixOAR,4}{1};  
vIxpTV = cst{ixPTV,4}{1};  
  
ct.cubeHU{1}(vIxoAR) = 1;  
ct.cubeHU{1}(vIxpTV) = 1;
```

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.

First of all, we need to define what kind of radiation modality we would like to use. Possible values are photons, protons or carbon. In this example we would like to use photons for treatment planning. Next, we need to define a treatment machine to correctly load the corresponding base data. matRad features generic base data in the file 'photons_Generic.mat'; consequently the machine has to be set to 'Generic'

```
pln.radiationMode = 'photons';  
pln.machine       = 'Generic';
```

Define the flavor of biological optimization for treatment planning along with the quantity that should be used for optimization. Possible values are (none: physical dose based optimization; const_RBExD: constant RBE of 1.1; LEMIV_effect: effect-based optimization; LEMIV_RBExD: optimization of RBE-weighted dose. As we use photons, we select 'none' as we want to optimize the physical dose.

```
pln.propOpt.bioOptimization = 'none';
```

The remaining plan parameters are set like in the previous example files

```
pln.numOfFractions      = 30;  
pln.propStf.gantryAngles = [0 45];  
pln.propStf.couchAngles  = [0 0];  
pln.propStf.bixelWidth   = 5;  
pln.propStf.numOfBeams   = numel(pln.propStf.gantryAngles);  
pln.propStf.isoCenter    = ones(pln.propStf.numOfBeams,1) *  
    matRad_getIsoCenter(cst,ct,0);  
pln.propOpt.runDAO       = 0;  
pln.propOpt.runSequencing = 0;
```

Generate Beam Geometry STF

```
stf = matRad_generateStf(ct,cst,pln);  
  
matRad: Generating stf struct... Warning: Could not find HLUT in  
    hlutLibrary folder. matRad default HLUT loaded  
Progress: 100.00 %
```

Dose Calculation

```
di = matRad_calcPhotonDose(ct,stf,pln,cst);

Warning: Could not find HLUT in hlutLibrary folder. matRad default
HLUT loaded
matRad: Photon dose calculation...
Beam 1 of 2:
matRad: calculate radiological depth cube...done
SSD = 901mm
matRad: Uniform primary photon fluence -> pre-compute kernel
convolution for SSD = 901 mm ...
Progress: 100.00 %
Beam 2 of 2:
matRad: calculate radiological depth cube...done
SSD = 900mm
matRad: Uniform primary photon fluence -> pre-compute kernel
convolution for SSD = 900 mm ...
Progress: 100.00 %
```

Inverse Optimization for intensity-modulated photon therapy

The goal of the fluence optimization is to find a set of bixel/spot weights which yield the best possible dose distribution according to the clinical objectives and constraints underlying the radiation treatment.

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

```
Optimization initiating...
Press q to terminate the optimization...
```

```
*****
This program contains Ipopt, a library for large-scale nonlinear
optimization.
Ipopt is released as open source code under the Eclipse Public
License (EPL).
For more information visit http://projects.coin-or.org/Ipopt
*****
```

```
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

Total number of variables.....:      534
      variables with only lower bounds:      534
      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
```

Example: Generate your
own phantom geometry

```

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 2.5998464e+000 0.00e+000 1.02e+000  0.0 0.00e+000 - 0.00e
+000 0.00e+000 0
  1 2.5357843e+000 0.00e+000 2.83e-002 -1.4 3.24e-002 -
9.99e-001 1.00e+000f 1
  2 1.9092313e+000 0.00e+000 4.52e-002 -2.7 2.07e-001 -
9.99e-001 1.00e+000f 1
  3 1.6142827e+000 0.00e+000 2.81e-002 -3.5 1.08e-001 -
9.94e-001 1.00e+000f 1
  4 1.2545654e+000 0.00e+000 1.96e-002 -4.3 3.69e-001 - 1.00e
+000 1.00e+000f 1
  5 1.1976709e+000 0.00e+000 2.20e-002 -4.6 1.50e-001 -
9.99e-001 1.00e+000f 1
  6 1.1517353e+000 0.00e+000 8.72e-003 -5.7 4.48e-002 - 1.00e
+000 1.00e+000f 1
  7 1.1162275e+000 0.00e+000 6.55e-003 -6.8 1.01e-001 - 1.00e
+000 1.00e+000f 1
  8 1.0932738e+000 0.00e+000 8.62e-003 -8.1 1.05e-001 - 1.00e
+000 1.00e+000f 1
  9 1.0687354e+000 0.00e+000 2.02e-002 -8.9 2.81e-001 - 1.00e
+000 1.00e+000f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr  ls
 10 1.0407284e+000 0.00e+000 1.04e-002 -9.8 1.50e-001 - 1.00e
+000 1.00e+000f 1
 11 1.0295842e+000 0.00e+000 7.99e-003 -9.7 3.43e-002 - 1.00e
+000 1.00e+000f 1
 12 1.0108424e+000 0.00e+000 7.89e-003 -10.5 1.00e-001 - 1.00e
+000 1.00e+000f 1
 13 9.9503865e-001 0.00e+000 1.50e-002 -11.0 2.23e-001 - 1.00e
+000 1.00e+000f 1
 14 9.9074286e-001 0.00e+000 1.38e-002 -11.0 1.85e-001 - 1.00e
+000 1.89e-001f 1
 15 9.9069925e-001 0.00e+000 1.38e-002 -11.0 3.40e-001 - 1.00e
+000 1.02e-003f 1
 16 9.9069723e-001 0.00e+000 3.16e-002 -11.0 2.53e-001 - 1.00e
+000 4.42e-005f 1
 17 9.9017677e-001 0.00e+000 2.48e-002 -11.0 3.73e-001 - 1.00e
+000 8.15e-003f 1

```

Number of Iterations.....: 17

```

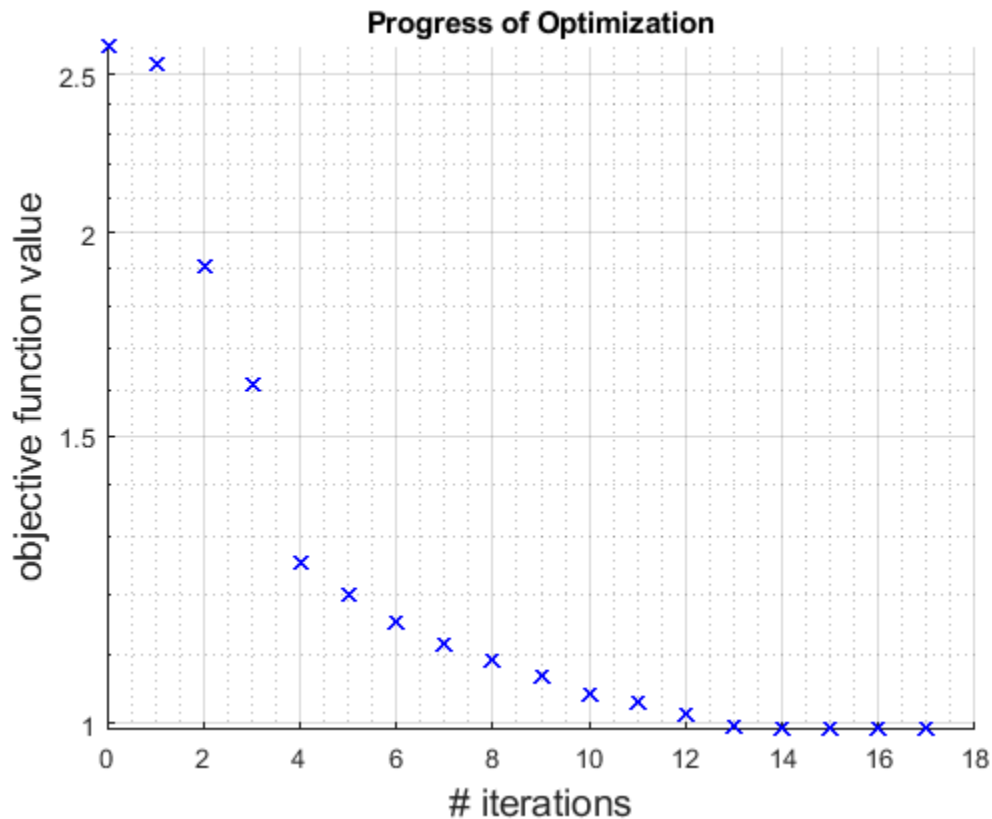
                                     (scaled)                (unscaled)
Objective.....: 9.9017677048753372e-001
9.9017677048753372e-001
Dual infeasibility.....: 2.4847495822923300e-002
2.4847495822923300e-002
Constraint violation.....: 0.0000000000000000e+000
0.0000000000000000e+000

```

```
Complementarity.....: 2.0831495922667587e-007
2.0831495922667587e-007
Overall NLP error.....: 2.4847495822923300e-002
2.4847495822923300e-002
```

```
Number of objective function evaluations      = 18
Number of objective gradient evaluations     = 18
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) = 1.273
Total CPU secs in NLP function evaluations   = 2.456
```

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



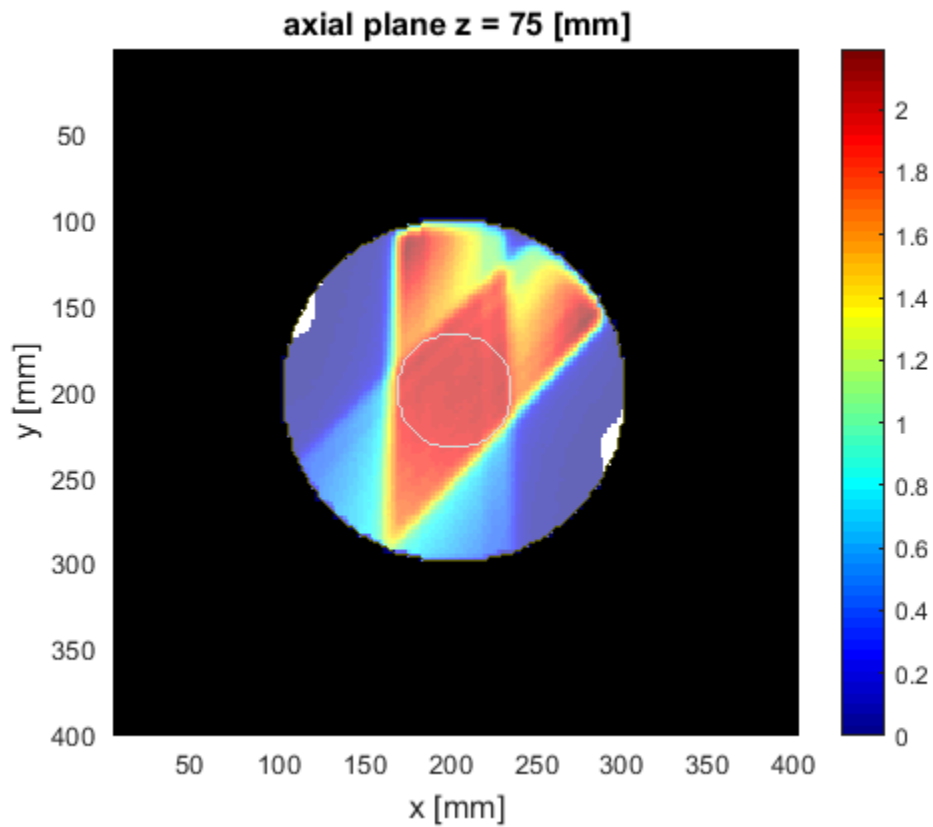
Plot the resulting dose slice

```
plane      = 3;
slice      = round(pln.propStf.isoCenter(1,3)./ct.resolution.z);
doseWindow = [0 max([resultGUI.physicalDose(:)])];

figure,title('phantom plan')
```



```
matRad_plotSliceWrapper(gca,ct,cst,1,resultGUI.physicalDose,plane,slice,  
[],[],colorcube,[],doseWindow,[]);
```



Published with MATLAB® R2018a