

---

# Example: Proton Treatment Plan with subsequent Isocenter shift

## Table of Contents

.....	1
Patient Data Import .....	1
Treatment Plan .....	1
Generate Beam Geometry STF .....	2
Dose Calculation .....	2
Inverse Optimization for IMPT .....	3
Plot the Resulting Dose Slice .....	14
Plot the Resulting Beam Dose Slice .....	14
and the corresponding LET distribution .....	15
Recalculate Plan .....	16
Visual Comparison of results .....	17
Quantitative Comparison of results .....	20

%%%

Copyright 2017 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 load patient data into matRad (ii) how to setup a proton dose calculation (iii) how to inversely optimize the pencil beam intensities directly from command window in MATLAB. (iv) how to simulate a lateral patient displacement by shifting the iso-center (v) how to recalculate the dose considering the shifted geometry and the previously optimized pencil beam intensities (vi) how to compare the two results

## Patient Data Import

Let's begin with a clear Matlab environment and import the prostate patient into your workspace

```
clc,clear,close all;  
load('PROSTATE.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.

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 protons 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 'proton\_Generic.mat'; consequently the machine has to be set accordingly

```
pln.radiationMode = 'protons';  
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 optimization; const\_RBExD: constant RBE of 1.1; LEMIV\_effect: effect-based optimization; LEMIV\_RBExD: optimization of RBE-weighted dose. As we use protons, we follow here the clinical standard and use a constant relative biological effectiveness of 1.1. Therefore we set bioOptimization to const\_RBExD

```
pln.propOpt.bioOptimization = 'const_RBExD';
```

for particles it is possible to also calculate the LET disutribution alongside the physical dose. Therefore you need to activate the corresponding option during dose calculation

```
pln.propDoseCalc.calcLET = 1;
```

Now we have to set the remaining plan parameters.

```
pln.numOfFractions      = 30;  
pln.propStf.gantryAngles = [90 270];  
pln.propStf.couchAngles  = [0 0];  
pln.propStf.bixelWidth   = 3;  
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  
Philips-AcQSimCT-ConvolutionKernel-000000_protons.hlut in hlutLibrary  
folder.  
matRad default HLUT loaded  
Progress: 100.00 %
```

## Dose Calculation

Lets generate dosimetric information by pre-computing dose influence matrices for unit beamlet intensities. Having dose influences available allows for subsequent inverse optimization.

```
dij = matRad_calcParticleDose(ct,stf,pln,cst);
```

```
Warning: Could not find HLUT  
Philips-AcQSimCT-ConvolutionKernel-000000_protons.hlut in hlutLibrary  
folder.  
matRad default HLUT loaded  
matRad: Using a constant RBE of 1.1  
matRad: Particle dose calculation...  
Beam 1 of 2:
```

```
matRad: calculate radiological depth cube...done.
matRad: calculate lateral cutoff...done.
Progress: 100.00 %
Beam 2 of 2:
matRad: calculate radiological depth cube...done.
matRad: calculate lateral cutoff...done.
Progress: 100.00 %
```

## Inverse Optimization for IMPT

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(dij,cst,pln);
```

```
Optimization initiating...
Press q to terminate the optimization...
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.....:      45367
      variables with only lower bounds:      45367
      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
0	4.3987636e+002	0.00e+000	1.07e+000	0.0	0.00e+000	-	0.00e
+000	0.00e+000	0					
1	4.0806688e+002	0.00e+000	7.31e-002	-1.1	7.92e-002	-	
9.91e-001	1.00e+000f	1					
2	7.2657697e+001	0.00e+000	2.00e-002	-1.8	1.33e+000	-	
9.92e-001	1.00e+000f	1					
3	3.8743259e+001	0.00e+000	1.34e-002	-3.4	3.85e-001	-	
9.73e-001	1.00e+000f	1					
4	3.1521547e+001	0.00e+000	1.15e-002	-3.9	2.83e-001	-	
9.87e-001	1.00e+000f	1					
5	2.5107747e+001	0.00e+000	1.03e-002	-4.7	4.47e-001	-	1.00e
+000	1.00e+000f	1					
6	2.1181842e+001	0.00e+000	1.45e-002	-5.4	7.07e-001	-	1.00e
+000	1.00e+000f	1					
7	1.7800111e+001	0.00e+000	7.17e-003	-5.9	2.63e-001	-	1.00e
+000	1.00e+000f	1					
8	1.6610129e+001	0.00e+000	6.24e-003	-7.2	2.05e-001	-	1.00e
+000	1.00e+000f	1					

Example: Proton Treatment Plan  
with subsequent Isocenter shift

---

```

  9 1.5046290e+001 0.00e+000 4.94e-003 -8.4 4.00e-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 1.3067048e+001 0.00e+000 3.87e-003 -9.3 5.97e-001 - 1.00e
+000 1.00e+000f 1
 11 1.2593017e+001 0.00e+000 8.31e-003 -9.6 7.94e-001 - 1.00e
+000 7.48e-001f 1
 12 1.2570649e+001 0.00e+000 8.25e-003 -10.9 2.37e-001 - 1.00e
+000 9.70e-003f 1
 13 1.2569264e+001 0.00e+000 1.03e-002 -11.0 3.04e-001 - 1.00e
+000 4.67e-004f 1
 14 1.2369437e+001 0.00e+000 7.78e-003 -11.0 3.93e-001 - 1.00e
+000 6.10e-002f 1
 15 1.2352346e+001 0.00e+000 1.47e-002 -11.0 4.45e-001 - 1.00e
+000 5.22e-003f 1
 16 1.1951431e+001 0.00e+000 6.80e-003 -11.0 5.12e-001 - 1.00e
+000 1.21e-001f 1
 17 1.1917942e+001 0.00e+000 1.30e-002 -11.0 5.08e-001 - 1.00e
+000 1.14e-002f 1
 18 1.1722141e+001 0.00e+000 1.10e-002 -7.6 5.35e-001 -
9.63e-001 6.57e-002f 1
 19 1.1434423e+001 0.00e+000 8.60e-003 -8.4 5.42e-001 - 1.00e
+000 1.05e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr  ls
 20 1.1392157e+001 0.00e+000 1.62e-002 -9.7 5.37e-001 - 1.00e
+000 1.71e-002f 1
 21 1.1174849e+001 0.00e+000 4.79e-003 -10.4 5.97e-001 - 1.00e
+000 8.66e-002f 1
 22 1.1012328e+001 0.00e+000 1.11e-002 -11.0 6.20e-001 - 1.00e
+000 6.97e-002f 1
 23 1.0914192e+001 0.00e+000 1.19e-002 -11.0 6.76e-001 - 1.00e
+000 4.29e-002f 1
 24 1.0761770e+001 0.00e+000 1.40e-002 -11.0 7.64e-001 - 1.00e
+000 6.53e-002f 1
 25 1.0553983e+001 0.00e+000 2.21e-002 -11.0 8.65e-001 - 1.00e
+000 8.99e-002f 1
 26 1.0389215e+001 0.00e+000 1.24e-002 -11.0 9.94e-001 - 1.00e
+000 6.82e-002f 1
 27 1.0330379e+001 0.00e+000 1.92e-002 -6.6 9.99e-001 -
8.00e-001 2.65e-002f 1
 28 1.0084019e+001 0.00e+000 9.26e-003 -6.0 1.03e+000 -
9.77e-001 1.19e-001f 1
 29 9.9262694e+000 0.00e+000 9.06e-003 -4.0 1.04e+000 -
6.20e-001 8.52e-002f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr  ls
 30 9.6642392e+000 0.00e+000 3.98e-003 -10.2 1.10e+000 -
4.66e-001 1.47e-001f 1
 31 9.4731493e+000 0.00e+000 9.53e-003 -5.0 1.20e+000 - 1.00e
+000 1.17e-001f 1
 32 9.0789575e+000 0.00e+000 8.79e-003 -4.1 1.33e+000 -
7.04e-001 2.82e-001f 1

```

---

Example: Proton Treatment Plan  
with subsequent Isocenter shift

---

33	8.9461172e+000	0.00e+000	1.52e-002	-4.0	7.48e-001	-	
	8.62e-001	1.78e-001f	1				
34	8.7941053e+000	0.00e+000	7.90e-003	-4.6	7.80e-001	-	
	6.48e-001	2.45e-001f	1				
35	8.5893624e+000	0.00e+000	4.11e-003	-4.0	8.33e-001	-	
	8.50e-001	3.72e-001f	1				
36	8.5143820e+000	0.00e+000	1.28e-002	-4.1	4.90e-001	-	
	8.04e-001	2.47e-001f	1				
37	8.3305167e+000	0.00e+000	5.35e-003	-4.5	6.33e-001	-	
	6.32e-001	5.10e-001f	1				
38	8.2320599e+000	0.00e+000	2.80e-003	-4.3	7.10e-001	-	
	4.92e-001	2.55e-001f	1				
39	8.1186421e+000	0.00e+000	1.18e-002	-3.7	3.12e-001	-	
	6.27e-001	5.35e-001f	1				
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
	alpha_pr	ls					
40	8.0377783e+000	0.00e+000	5.30e-003	-5.0	6.99e-001	-	
	4.41e-001	1.90e-001f	1				
41	7.8301566e+000	0.00e+000	2.02e-003	-4.2	6.67e-001	-	
	3.17e-001	5.30e-001f	1				
42	7.7279965e+000	0.00e+000	3.51e-003	-4.3	4.68e-001	-	
	9.81e-001	4.83e-001f	1				
43	7.6478581e+000	0.00e+000	3.92e-003	-4.4	2.70e-001	-	
	5.76e-001	6.99e-001f	1				
44	7.5645729e+000	0.00e+000	1.74e-003	-4.5	2.69e-001	-	
	7.20e-001	1.00e+000f	1				
45	7.4883816e+000	0.00e+000	4.54e-003	-3.9	1.86e-001	-	
	8.14e-001	1.00e+000f	1				
46	7.4251827e+000	0.00e+000	7.84e-004	-4.9	1.42e-001	-	
	6.56e-001	1.00e+000f	1				
47	7.3817034e+000	0.00e+000	1.94e-003	-4.8	3.06e-001	-	
	9.91e-001	3.39e-001f	1				
48	7.3246018e+000	0.00e+000	2.54e-003	-5.3	2.32e-001	-	1.00e
	+000 5.44e-001f	1					
49	7.2646499e+000	0.00e+000	1.92e-003	-4.8	2.85e-001	-	
	7.55e-001	4.00e-001f	1				
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
	alpha_pr	ls					
50	7.2188867e+000	0.00e+000	7.15e-003	-4.6	4.87e-001	-	
	9.02e-001	1.96e-001f	1				
51	7.1693624e+000	0.00e+000	3.30e-003	-4.4	6.45e-001	-	
	4.86e-001	1.74e-001f	1				
52	7.4995595e+000	0.00e+000	4.94e-003	-2.5	8.12e+000	-	
	2.03e-002	6.73e-002f	1				
53	7.1643888e+000	0.00e+000	2.95e-003	-4.1	1.17e+000	-	
	3.76e-001	4.96e-001f	1				
54	7.1487583e+000	0.00e+000	5.28e-003	-6.1	8.33e-001	-	
	7.73e-001	3.02e-002f	1				
55	7.0839869e+000	0.00e+000	2.63e-002	-5.4	6.48e-001	-	
	9.96e-001	2.17e-001f	1				
56	6.9769448e+000	0.00e+000	8.77e-003	-5.4	6.66e-001	-	
	9.47e-001	4.25e-001f	1				
57	6.9169707e+000	0.00e+000	5.90e-003	-6.3	6.83e-001	-	
	9.95e-001	2.88e-001f	1				

---

Example: Proton Treatment Plan  
with subsequent Isocenter shift

---

```

58 6.8645695e+000 0.00e+000 3.86e-003 -5.0 9.08e-001 -
9.31e-001 2.35e-001f 1
59 6.8418652e+000 0.00e+000 6.06e-003 -4.8 4.54e-001 -
3.54e-001 1.94e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr ls
60 6.7898912e+000 0.00e+000 2.69e-003 -5.0 7.29e-001 -
5.61e-001 2.88e-001f 1
61 6.7704529e+000 0.00e+000 5.39e-003 -5.3 1.05e+000 -
6.89e-001 7.15e-002f 1
62 6.7432305e+000 0.00e+000 4.70e-003 -4.6 6.16e-001 -
2.95e-001 1.77e-001f 1
63 6.7054188e+000 0.00e+000 2.99e-003 -4.6 8.15e-001 -
4.19e-001 1.99e-001f 1
64 6.6869537e+000 0.00e+000 5.89e-003 -5.2 6.70e-001 -
2.83e-001 1.23e-001f 1
65 6.6539826e+000 0.00e+000 3.88e-003 -5.0 8.80e-001 -
9.19e-001 1.68e-001f 1
66 6.6037914e+000 0.00e+000 2.74e-003 -4.3 1.07e+000 -
1.99e-001 2.28e-001f 1
67 6.5919268e+000 0.00e+000 5.58e-003 -4.4 4.80e-001 -
3.95e-001 1.10e-001f 1
68 7.4213451e+000 0.00e+000 9.22e-003 -3.3 4.10e+000 -
1.49e-001 5.29e-001f 1
69 6.9375644e+000 0.00e+000 6.33e-003 -4.1 1.55e+000 -
3.09e-002 3.76e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr ls
70 6.6330614e+000 0.00e+000 9.40e-003 -4.1 8.94e-001 -
4.57e-001 5.15e-001f 1
71 6.5703080e+000 0.00e+000 8.74e-003 -4.1 4.18e-001 -
5.93e-001 2.86e-001f 1
72 6.5166541e+000 0.00e+000 9.63e-003 -4.9 5.59e-001 -
5.80e-001 2.78e-001f 1
73 6.4815997e+000 0.00e+000 8.31e-003 -5.1 4.93e-001 - 1.00e
+000 2.61e-001f 1
74 6.4464187e+000 0.00e+000 6.27e-003 -5.4 4.39e-001 -
9.24e-001 3.58e-001f 1
75 6.4225014e+000 0.00e+000 6.98e-003 -6.2 3.90e-001 -
8.91e-001 3.35e-001f 1
76 6.3997716e+000 0.00e+000 8.38e-003 -6.8 4.47e-001 -
9.28e-001 3.25e-001f 1
77 6.3962007e+000 0.00e+000 7.90e-003 -7.7 4.77e-001 -
9.06e-001 4.78e-002f 1
78 6.3667565e+000 0.00e+000 6.57e-003 -7.2 8.04e-001 -
8.57e-001 2.31e-001f 1
79 6.3492239e+000 0.00e+000 4.80e-003 -5.9 8.41e-001 -
8.13e-001 1.25e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr ls
80 6.3296377e+000 0.00e+000 3.89e-003 -5.0 1.25e+000 -
6.24e-001 8.77e-002f 1
81 6.3026051e+000 0.00e+000 4.72e-003 -4.1 3.07e-001 -
2.61e-001 1.00e+000f 1

```

---

Example: Proton Treatment Plan  
with subsequent Isocenter shift

---

82	6.2531705e+000	0.00e+000	1.74e-003	-4.2	5.74e-001	-
5.12e-001	3.88e-001f	1				
83	6.2446393e+000	0.00e+000	6.71e-003	-5.2	5.46e-001	-
4.49e-001	7.87e-002f	1				
84	6.2001907e+000	0.00e+000	3.10e-003	-4.6	1.21e+000	-
5.62e-001	2.45e-001f	1				
85	6.1604640e+000	0.00e+000	5.67e-003	-4.6	1.17e+000	-
5.26e-001	2.27e-001f	1				
86	6.1406954e+000	0.00e+000	8.52e-003	-4.9	8.18e-001	-
7.87e-001	1.44e-001f	1				
87	6.1034531e+000	0.00e+000	4.65e-003	-4.9	1.17e+000	-
6.18e-001	2.08e-001f	1				
88	6.0805614e+000	0.00e+000	4.64e-003	-4.6	6.61e-001	-
2.62e-001	2.20e-001f	1				
89	6.0554092e+000	0.00e+000	1.39e-003	-4.1	4.69e-001	-
3.81e-001	4.75e-001f	1				
iter    objective    inf_pr    inf_du lg(mu)      d      lg(rg) alpha_du						
alpha_pr    ls						
90	6.0389098e+000	0.00e+000	5.88e-003	-4.7	4.70e-001	-
3.23e-001	2.29e-001f	1				
91	6.0281125e+000	0.00e+000	2.85e-003	-4.6	4.41e-001	-
6.18e-001	1.56e-001f	1				
92	6.8527112e+000	0.00e+000	4.04e-003	-2.5	2.73e+001	-
6.88e-003	1.34e-001f	1				
93	6.2434948e+000	0.00e+000	2.61e-003	-4.4	3.10e+000	-
2.13e-002	5.41e-001f	1				
94	6.0036749e+000	0.00e+000	5.51e-003	-4.4	1.34e+000	-
9.90e-001	8.26e-001f	1				
95	5.9946258e+000	0.00e+000	6.05e-003	-5.0	5.73e-001	-
9.05e-001	1.13e-001f	1				
96	5.9631774e+000	0.00e+000	7.35e-003	-5.3	7.10e-001	-
8.47e-001	3.25e-001f	1				
97	5.9490189e+000	0.00e+000	6.80e-003	-6.4	6.43e-001	-
8.43e-001	1.60e-001f	1				
98	5.9255408e+000	0.00e+000	4.22e-003	-4.8	5.65e-001	-
3.77e-001	3.04e-001f	1				
99	5.9047054e+000	0.00e+000	2.86e-003	-4.4	2.21e-001	-
4.71e-001	1.00e+000f	1				
iter    objective    inf_pr    inf_du lg(mu)      d      lg(rg) alpha_du						
alpha_pr    ls						
100	5.8941726e+000	0.00e+000	2.97e-003	-10.7	5.71e-001	-
4.40e-001	1.44e-001f	1				
101	5.8744665e+000	0.00e+000	3.07e-003	-4.9	7.10e-001	-
8.78e-001	2.15e-001f	1				
102	5.8606609e+000	0.00e+000	7.75e-003	-5.1	6.23e-001	-
7.73e-001	1.62e-001f	1				
103	5.8280724e+000	0.00e+000	3.26e-003	-5.3	9.69e-001	-
4.81e-001	3.25e-001f	1				
104	5.8024571e+000	0.00e+000	2.34e-003	-4.6	4.25e-001	-
2.23e-001	7.51e-001f	1				
105	5.7960339e+000	0.00e+000	2.71e-003	-4.9	4.76e-001	-
5.48e-001	1.66e-001f	1				
106	5.7812656e+000	0.00e+000	4.36e-003	-5.1	7.14e-001	-
7.90e-001	2.33e-001f	1				

---

Example: Proton Treatment Plan  
with subsequent Isocenter shift

---

```

107 5.7704164e+000 0.00e+000 3.21e-003 -11.0 8.52e-001 -
4.02e-001 1.24e-001f 1
108 5.7592528e+000 0.00e+000 4.66e-003 -6.5 8.26e-001 -
4.78e-001 1.18e-001f 1
109 5.7378901e+000 0.00e+000 4.98e-003 -6.2 9.97e-001 -
8.06e-001 1.86e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr ls
110 5.7133599e+000 0.00e+000 3.34e-003 -4.9 7.84e-001 -
2.54e-001 2.61e-001f 1
111 5.7048676e+000 0.00e+000 2.60e-003 -4.4 1.71e-001 -
5.37e-001 1.00e+000f 1
112 5.6911862e+000 0.00e+000 2.21e-003 -6.3 7.35e-001 -
4.84e-001 1.70e-001f 1
113 5.6695609e+000 0.00e+000 4.72e-003 -5.3 8.00e-001 -
5.94e-001 2.56e-001f 1
114 5.6585085e+000 0.00e+000 4.40e-003 -5.5 9.11e-001 -
5.94e-001 1.13e-001f 1
115 5.6378367e+000 0.00e+000 3.43e-003 -5.0 7.93e-001 -
3.11e-001 2.48e-001f 1
116 5.6222379e+000 0.00e+000 2.46e-003 -11.0 1.11e+000 -
2.46e-001 1.41e-001f 1
117 5.6147302e+000 0.00e+000 7.14e-003 -6.1 9.68e-001 -
5.28e-001 7.75e-002f 1
118 5.5919399e+000 0.00e+000 3.29e-003 -5.8 1.26e+000 -
4.63e-001 1.80e-001f 1
119 5.7414042e+000 0.00e+000 4.15e-003 -3.6 7.55e+000 -
1.50e-002 2.21e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr ls
120 5.6195013e+000 0.00e+000 4.15e-003 -5.2 3.11e+000 -
3.96e-003 3.44e-001f 1
121 5.5720778e+000 0.00e+000 4.37e-002 -5.2 2.12e+000 -
7.47e-001 2.06e-001f 1
122 5.7921276e+000 0.00e+000 4.39e-002 -3.3 9.60e+000 -
3.31e-003 1.84e-001f 1
123 5.7496429e+000 0.00e+000 4.20e-002 -5.1 3.60e+000 -
4.37e-002 7.80e-002f 1
124 5.6450420e+000 0.00e+000 7.46e-003 -5.1 2.92e+000 -
7.30e-001 2.59e-001f 1
125 5.5903529e+000 0.00e+000 3.66e-003 -5.1 2.66e+000 -
5.34e-001 3.21e-001f 1
126 5.5513051e+000 0.00e+000 2.18e-002 -5.1 7.83e-001 -
7.96e-001 2.69e-001f 1
127 5.5358297e+000 0.00e+000 1.52e-002 -5.6 6.43e-001 -
8.84e-001 1.49e-001f 1
128 5.5216565e+000 0.00e+000 2.10e-003 -4.3 3.63e-001 -
4.50e-001 1.00e+000f 1
129 5.5184830e+000 0.00e+000 9.68e-003 -5.4 6.33e-001 -
9.10e-001 5.12e-002f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr ls
130 5.4873388e+000 0.00e+000 1.00e-002 -6.4 8.35e-001 -
8.82e-001 3.96e-001f 1

```

---



Example: Proton Treatment Plan  
with subsequent Isocenter shift

---

131	5.4756909e+000	0.00e+000	2.39e-003	-4.8	4.34e-001	-
	7.55e-001	2.96e-001f	1			
132	5.4660276e+000	0.00e+000	2.99e-003	-4.7	3.93e-001	-
	3.40e-001	2.73e-001f	1			
133	5.4530848e+000	0.00e+000	3.30e-003	-4.6	4.31e-001	-
	3.00e-001	4.02e-001f	1			
134	5.4467846e+000	0.00e+000	4.41e-003	-6.6	1.04e+000	-
	3.05e-001	6.94e-002f	1			
135	5.4216824e+000	0.00e+000	2.28e-003	-4.9	1.19e+000	-
	5.59e-001	2.49e-001f	1			
136	5.4171900e+000	0.00e+000	4.65e-003	-10.9	7.89e-001	-
	3.19e-001	6.42e-002f	1			
137	5.3929768e+000	0.00e+000	2.40e-003	-5.3	1.44e+000	-
	5.05e-001	1.91e-001f	1			
138	5.3889651e+000	0.00e+000	5.98e-003	-5.2	8.33e-001	-
	3.02e-001	4.66e-002f	1			
139	5.3726731e+000	0.00e+000	4.80e-003	-5.9	1.31e+000	-
	1.52e-001	1.27e-001f	1			
iter    objective    inf_pr    inf_du lg(mu)      d      lg(rg) alpha_du						
alpha_pr    ls						
140	5.3633104e+000	0.00e+000	3.10e-003	-5.3	1.25e+000	-
	2.27e-001	7.79e-002f	1			
141	5.3562970e+000	0.00e+000	3.68e-003	-6.6	1.23e+000	-
	1.18e-001	5.95e-002f	1			
142	5.3344923e+000	0.00e+000	1.91e-003	-5.0	1.31e+000	-
	3.63e-001	1.83e-001f	1			
143	5.3267305e+000	0.00e+000	1.80e-003	-4.7	6.38e-001	-
	3.13e-001	1.25e-001f	1			
144	5.4286777e+000	0.00e+000	1.09e-003	-3.5	4.00e+000	-
	8.33e-003	2.51e-001f	1			
145	5.3869872e+000	0.00e+000	1.10e-003	-4.9	1.99e+000	-
	2.46e-002	1.90e-001f	1			
146	5.3220952e+000	0.00e+000	6.69e-003	-4.9	2.24e+000	-
	4.15e-001	4.01e-001f	1			
147	5.3033691e+000	0.00e+000	1.69e-002	-4.6	5.77e-001	-
	7.81e-001	2.93e-001f	1			
148	5.2944170e+000	0.00e+000	8.28e-003	-4.7	4.98e-001	-
	7.22e-001	1.70e-001f	1			
149	5.2778002e+000	0.00e+000	4.93e-003	-4.9	7.75e-001	-
	7.37e-001	2.71e-001f	1			
iter    objective    inf_pr    inf_du lg(mu)      d      lg(rg) alpha_du						
alpha_pr    ls						
150	5.2626170e+000	0.00e+000	5.50e-003	-5.3	9.81e-001	-
	8.08e-001	2.71e-001f	1			
151	5.2523915e+000	0.00e+000	6.22e-003	-6.0	1.08e+000	-
	6.95e-001	2.08e-001f	1			
152	5.2493030e+000	0.00e+000	9.96e-003	-6.9	1.09e+000	-
	6.62e-001	6.66e-002f	1			
153	5.2351958e+000	0.00e+000	3.73e-003	-7.3	1.79e+000	-
	7.49e-001	1.94e-001f	1			
154	5.2174130e+000	0.00e+000	3.73e-003	-4.9	1.53e+000	-
	1.20e-001	3.66e-001f	1			
155	5.2114748e+000	0.00e+000	2.26e-003	-4.6	3.60e-001	-
	5.37e-001	5.94e-001f	1			

---

Example: Proton Treatment Plan  
with subsequent Isocenter shift

---

```

156 5.2031256e+000 0.00e+000 3.69e-003 -4.8 1.55e+000 -
3.06e-001 1.79e-001f 1
157 5.1901584e+000 0.00e+000 1.54e-003 -5.0 2.64e+000 -
4.29e-001 1.67e-001f 1
158 5.1862996e+000 0.00e+000 3.22e-003 -10.9 1.81e+000 -
1.80e-001 7.38e-002f 1
159 5.1665320e+000 0.00e+000 2.34e-003 -5.8 2.97e+000 -
2.58e-001 2.40e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr ls
160 5.1616297e+000 0.00e+000 2.74e-003 -5.9 3.12e+000 -
4.05e-001 5.52e-002f 1
161 5.1472499e+000 0.00e+000 2.52e-003 -4.6 1.55e+000 -
1.64e-001 3.35e-001f 1
162 5.1424385e+000 0.00e+000 2.30e-003 -6.6 2.35e+000 -
2.97e-001 8.05e-002f 1
163 5.1275761e+000 0.00e+000 1.30e-003 -4.8 2.61e+000 -
4.60e-001 2.35e-001f 1
164 5.1256384e+000 0.00e+000 5.09e-003 -6.9 1.76e+000 -
3.75e-001 4.50e-002f 1
165 5.4700933e+000 0.00e+000 5.41e-003 -3.2 4.93e+001 -
7.91e-003 1.12e-001f 1
166 5.3429089e+000 0.00e+000 4.66e-003 -5.0 6.06e+000 -
1.00e-002 2.15e-001f 1
167 5.1843111e+000 0.00e+000 1.23e-002 -5.0 4.36e+000 -
5.24e-001 4.56e-001f 1
168 5.1578696e+000 0.00e+000 7.61e-003 -5.0 2.63e+000 -
4.73e-001 1.80e-001f 1
169 5.1352464e+000 0.00e+000 9.71e-003 -5.0 2.10e+000 -
8.42e-001 2.11e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr ls
170 5.1180486e+000 0.00e+000 7.40e-003 -5.0 1.36e+000 -
4.80e-001 2.53e-001f 1
171 5.1075336e+000 0.00e+000 1.75e-003 -4.4 3.73e-001 -
4.67e-001 1.00e+000f 1
172 5.1013251e+000 0.00e+000 6.39e-003 -5.0 7.14e-001 -
9.67e-001 3.03e-001f 1
173 5.0955675e+000 0.00e+000 1.13e-002 -5.8 1.13e+000 - 1.00e
+000 1.96e-001f 1
174 5.0875785e+000 0.00e+000 6.98e-003 -6.2 1.52e+000 - 1.00e
+000 2.18e-001f 1
175 5.0768727e+000 0.00e+000 6.78e-003 -6.9 1.68e+000 -
8.99e-001 2.76e-001f 1
176 5.0705059e+000 0.00e+000 5.58e-003 -6.2 1.56e+000 -
6.60e-001 1.78e-001f 1
177 5.0650544e+000 0.00e+000 2.59e-003 -5.5 1.74e+000 -
5.76e-001 1.34e-001f 1
178 5.0552089e+000 0.00e+000 3.10e-003 -5.2 1.98e+000 -
7.06e-001 2.00e-001f 1
179 5.0531568e+000 0.00e+000 2.47e-003 -5.8 1.53e+000 -
1.90e-001 5.29e-002f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr ls

```

---

Example: Proton Treatment Plan  
with subsequent Isocenter shift

---

```

180 5.0448146e+000 0.00e+000 2.11e-003 -5.5 2.60e+000 -
2.83e-001 1.19e-001f 1
181 5.0373794e+000 0.00e+000 9.72e-004 -4.5 7.64e-001 -
2.10e-001 7.16e-001f 1
182 5.0370542e+000 0.00e+000 7.63e-003 -11.0 1.50e+000 -
2.64e-001 7.27e-003f 1
183 5.0257458e+000 0.00e+000 3.43e-003 -4.9 1.55e+000 -
4.26e-001 2.16e-001f 1
184 5.0172499e+000 0.00e+000 1.94e-003 -4.8 1.56e+000 -
3.54e-001 2.75e-001f 1
185 5.0132233e+000 0.00e+000 5.19e-003 -5.0 9.11e-001 -
6.07e-001 1.47e-001f 1
186 5.0054286e+000 0.00e+000 4.90e-003 -5.0 9.00e-001 -
8.37e-001 3.00e-001f 1
187 4.9984315e+000 0.00e+000 1.54e-003 -5.0 8.88e-001 -
5.76e-001 2.83e-001f 1
188 4.9913142e+000 0.00e+000 3.42e-003 -11.0 1.03e+000 -
5.14e-001 2.54e-001f 1
189 4.9864613e+000 0.00e+000 2.63e-003 -5.3 9.05e-001 -
4.96e-001 1.92e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr  ls
190 4.9809391e+000 0.00e+000 2.38e-003 -5.1 1.13e+000 -
5.85e-001 1.53e-001f 1
191 4.9746679e+000 0.00e+000 3.07e-003 -6.5 1.25e+000 -
1.71e-001 1.56e-001f 1
192 5.0475730e+000 0.00e+000 3.00e-003 -3.2 4.88e+001 -
4.22e-003 3.95e-002f 1
193 5.0313487e+000 0.00e+000 2.72e-003 -5.1 3.51e+000 -
4.00e-002 1.04e-001f 1
194 4.9769608e+000 0.00e+000 8.58e-003 -5.1 3.00e+000 -
4.99e-001 4.37e-001f 1
195 4.9701926e+000 0.00e+000 3.97e-003 -5.1 1.70e+000 -
4.27e-001 1.06e-001f 1
196 4.9650930e+000 0.00e+000 6.61e-003 -5.1 1.16e+000 -
6.14e-001 1.14e-001f 1
197 4.9532675e+000 0.00e+000 6.07e-003 -5.2 1.16e+000 -
7.37e-001 2.72e-001f 1
198 4.9477936e+000 0.00e+000 4.43e-003 -6.1 1.04e+000 -
3.56e-001 1.46e-001f 1
199 4.9408636e+000 0.00e+000 2.25e-003 -5.7 1.21e+000 -
7.83e-001 1.70e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr  ls
200 4.9324313e+000 0.00e+000 3.37e-003 -6.0 9.93e-001 -
3.01e-001 2.55e-001f 1
201 4.9270468e+000 0.00e+000 4.61e-003 -6.2 1.17e+000 -
5.85e-001 1.44e-001f 1
202 4.9211038e+000 0.00e+000 2.69e-003 -5.5 1.11e+000 -
4.78e-001 1.68e-001f 1
203 4.9149753e+000 0.00e+000 2.96e-003 -5.1 8.40e-001 -
4.60e-001 2.28e-001f 1
204 4.9087203e+000 0.00e+000 1.45e-003 -4.8 1.68e-001 -
3.36e-001 1.00e+000f 1

```

Example: Proton Treatment Plan  
with subsequent Isocenter shift

```

205 4.9036409e+000 0.00e+000 9.91e-004 -5.1 8.32e-001 -
3.82e-001 1.71e-001f 1
206 4.8996718e+000 0.00e+000 4.65e-003 -5.4 9.18e-001 -
5.45e-001 1.23e-001f 1
207 4.8930047e+000 0.00e+000 2.54e-003 -5.8 1.29e+000 -
3.67e-001 1.45e-001f 1
208 4.8875684e+000 0.00e+000 2.27e-003 -7.0 1.33e+000 -
4.10e-001 1.15e-001f 1
209 4.8820797e+000 0.00e+000 2.04e-003 -11.0 1.29e+000 -
2.45e-001 1.17e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr ls
210 4.8759319e+000 0.00e+000 1.76e-003 -5.9 1.63e+000 -
2.04e-001 1.03e-001f 1
211 4.8732565e+000 0.00e+000 2.26e-003 -4.8 5.87e-001 -
2.61e-001 1.00e+000f 1
212 4.8684017e+000 0.00e+000 1.17e-003 -5.0 5.13e-001 -
6.46e-001 2.43e-001f 1
213 4.8592215e+000 0.00e+000 1.67e-003 -5.1 9.09e-001 -
5.14e-001 2.98e-001f 1
214 4.8545902e+000 0.00e+000 2.75e-003 -5.4 1.13e+000 -
4.23e-001 1.15e-001f 1
215 4.8491684e+000 0.00e+000 1.89e-003 -5.4 1.31e+000 -
1.52e-001 1.45e-001f 1
216 4.8445797e+000 0.00e+000 1.69e-003 -11.0 1.17e+000 -
1.66e-001 1.14e-001f 1
217 4.8402940e+000 0.00e+000 4.23e-003 -5.6 1.45e+000 -
4.32e-001 1.04e-001f 1
218 4.8344542e+000 0.00e+000 2.61e-003 -5.1 1.31e+000 -
2.45e-001 1.86e-001f 1
219 4.8289099e+000 0.00e+000 3.26e-003 -4.9 1.09e+000 -
4.32e-001 2.69e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr ls
220 4.9145895e+000 0.00e+000 2.04e-003 -3.1 1.32e+001 -
3.22e-003 1.07e-001f 1
221 4.8558000e+000 0.00e+000 2.06e-003 -5.0 2.74e+000 -
2.26e-002 4.85e-001f 1
222 4.8479361e+000 0.00e+000 3.10e-003 -5.0 1.74e+000 -
5.58e-001 1.30e-001f 1
223 4.8210721e+000 0.00e+000 7.79e-003 -5.0 1.79e+000 -
5.96e-001 5.64e-001f 1
224 4.8144870e+000 0.00e+000 3.01e-003 -5.0 9.77e-001 -
8.43e-001 4.04e-001f 1
225 4.8121123e+000 0.00e+000 2.35e-003 -5.1 7.21e-001 -
6.47e-001 2.22e-001f 1
226 4.8103997e+000 0.00e+000 6.61e-003 -5.7 8.31e-001 -
9.01e-001 1.25e-001f 1
227 4.8064874e+000 0.00e+000 5.83e-003 -6.2 1.20e+000 -
7.44e-001 1.87e-001f 1

```

Number of Iterations.....: 227

(scaled)

(unscaled)

Example: Proton Treatment Plan  
with subsequent Isocenter shift

---

```

Objective.....: 4.8064873895390905e+000
4.8064873895390905e+000
Dual infeasibility.....: 5.8305241778842656e-003
5.8305241778842656e-003
Constraint violation....: 0.0000000000000000e+000
0.0000000000000000e+000
Complementarity.....: 2.2980661315779770e-005
2.2980661315779770e-005
Overall NLP error.....: 5.8305241778842656e-003
5.8305241778842656e-003

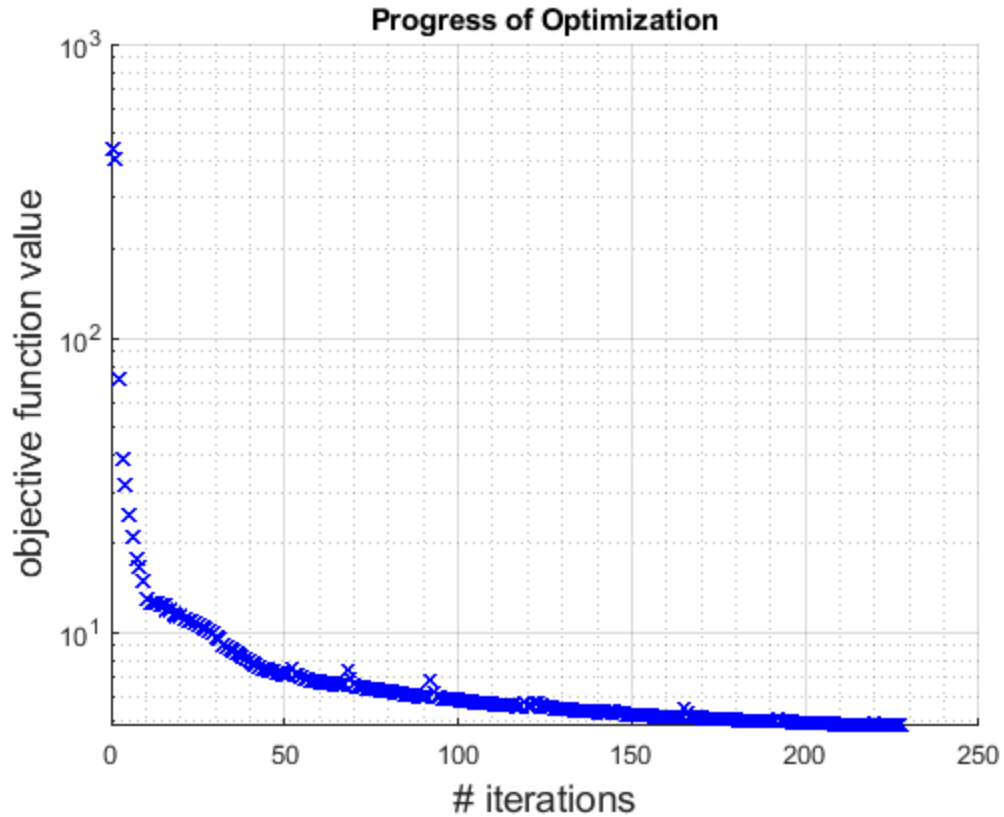
```

```

Number of objective function evaluations      = 228
Number of objective gradient evaluations     = 228
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) = 27.180
Total CPU secs in NLP function evaluations    = 190.721

```

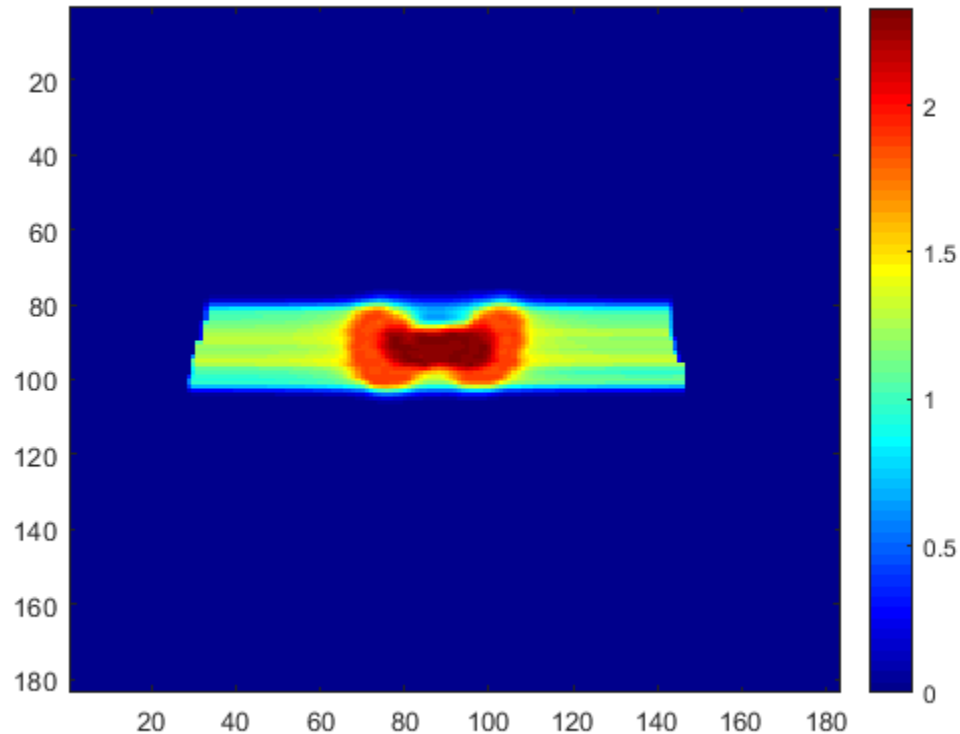
EXIT: Solved To Acceptable Level.  
Calculating final cubes...  
matRad: applying a constant RBE of 1.1



## Plot the Resulting Dose Slice

Let's plot the transversal iso-center dose slice

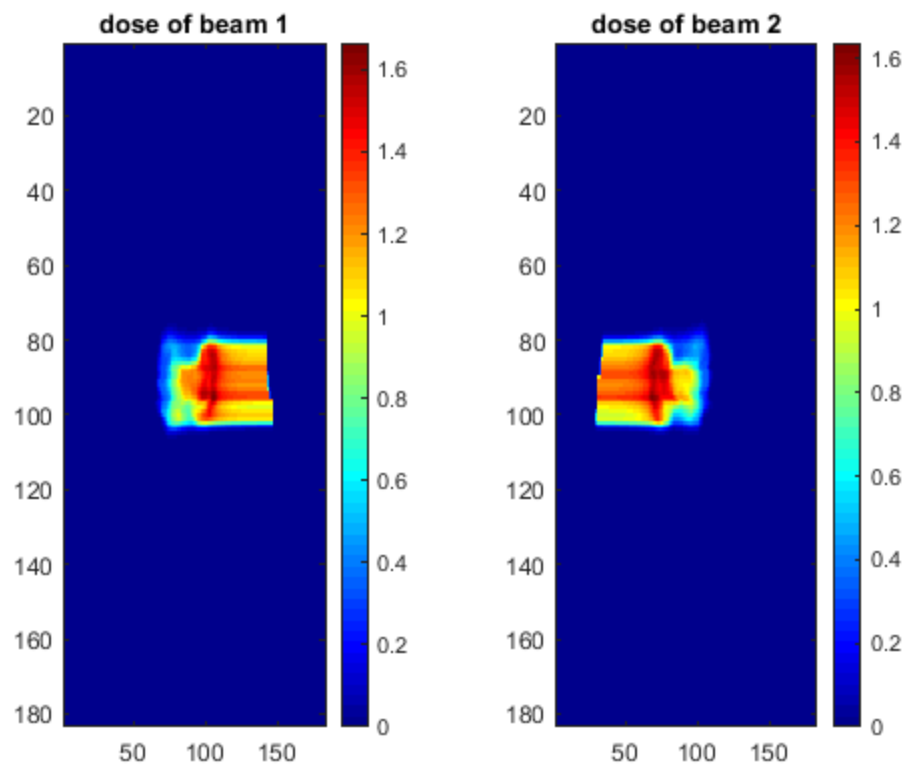
```
slice = round(pln.propStf.isoCenter(1,3)./ct.resolution.z);  
figure  
imagesc(resultGUI.RBExDose(:,:,slice)),colorbar,colormap(jet)
```



## Plot the Resulting Beam Dose Slice

Let's plot the transversal iso-center dose slice of beam 1 and beam 2 separately

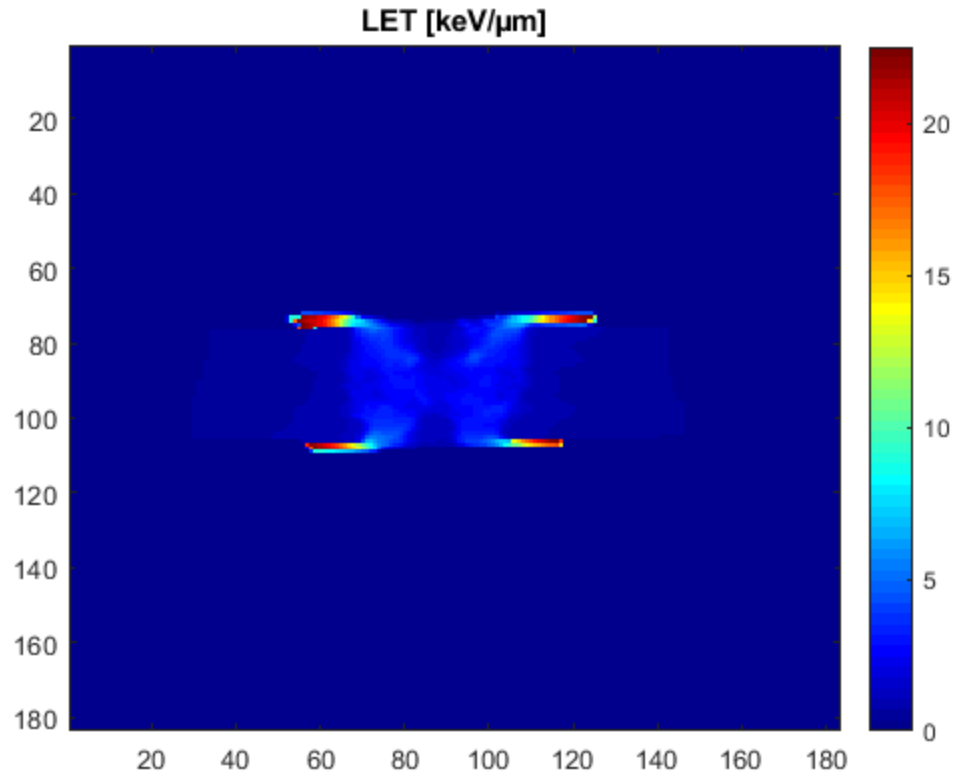
```
figure  
subplot(121),imagesc(resultGUI.RBExDose_beam1(:,:,slice)),colorbar,colormap(jet),t  
    'of beam 1')  
subplot(122),imagesc(resultGUI.RBExDose_beam2(:,:,slice)),colorbar,colormap(jet),t  
    'of beam 2')
```



## and the corresponding LET distribution

Transversal iso-center slice

```
figure
imagesc(resultGUI.LET(:,:,slice)),colormap(jet),colorbar,title('LET
[keV/μm]')
```



Now let's simulate a patient shift in y direction for both beams

```
stf(1).isoCenter(2) = stf(1).isoCenter(2) - 4;  
stf(2).isoCenter(2) = stf(2).isoCenter(2) - 4;  
pln.propStf.isoCenter      = reshape([stf.isoCenter],[3  
    pln.propStf.numOfBeams]));
```

## Recalculate Plan

Let's use the existing optimized pencil beam weights and recalculate the RBE weighted dose

```
resultGUI_isoShift =  
    matRad_calcDoseDirect(ct,stf,pln,cst,resultGUI.w);
```

```
Warning: Could not find HLUT  
Philips-AcQSimCT-ConvolutionKernel-000000_protons.hlut in hlutLibrary  
folder.  
matRad default HLUT loaded  
matRad: Using a constant RBE of 1.1  
matRad: Particle dose calculation...  
Beam 1 of 2:  
matRad: calculate radiological depth cube...done.  
matRad: calculate lateral cutoff...done.  
Progress: 100.00 %  
Beam 2 of 2:  
matRad: calculate radiological depth cube...done.
```

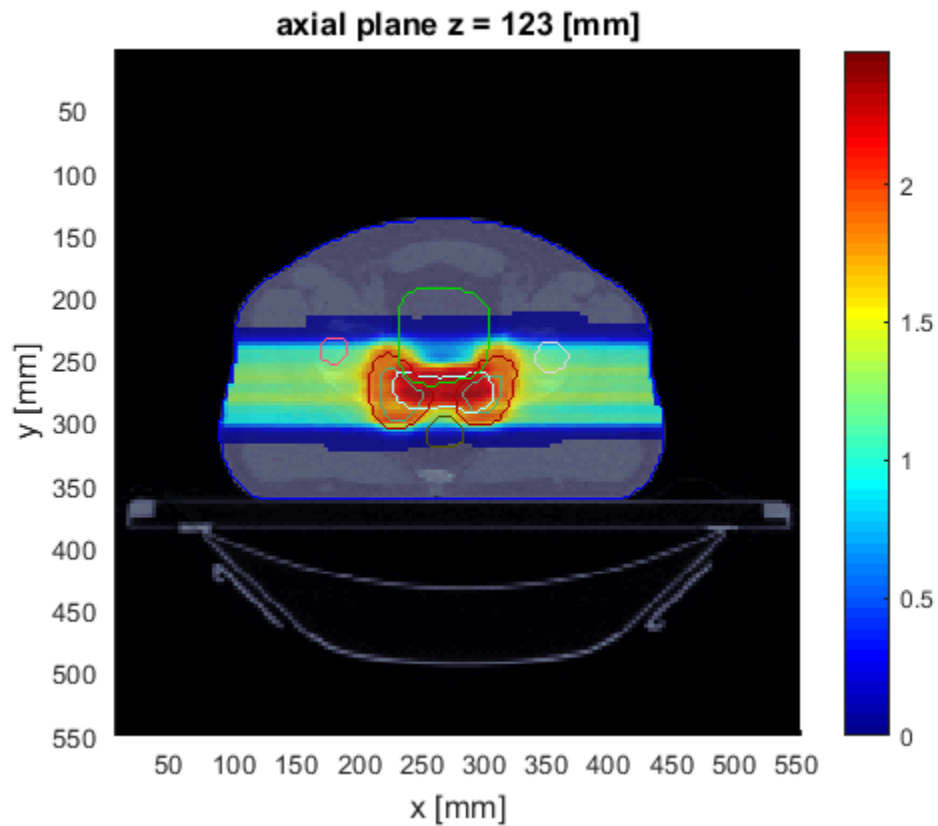
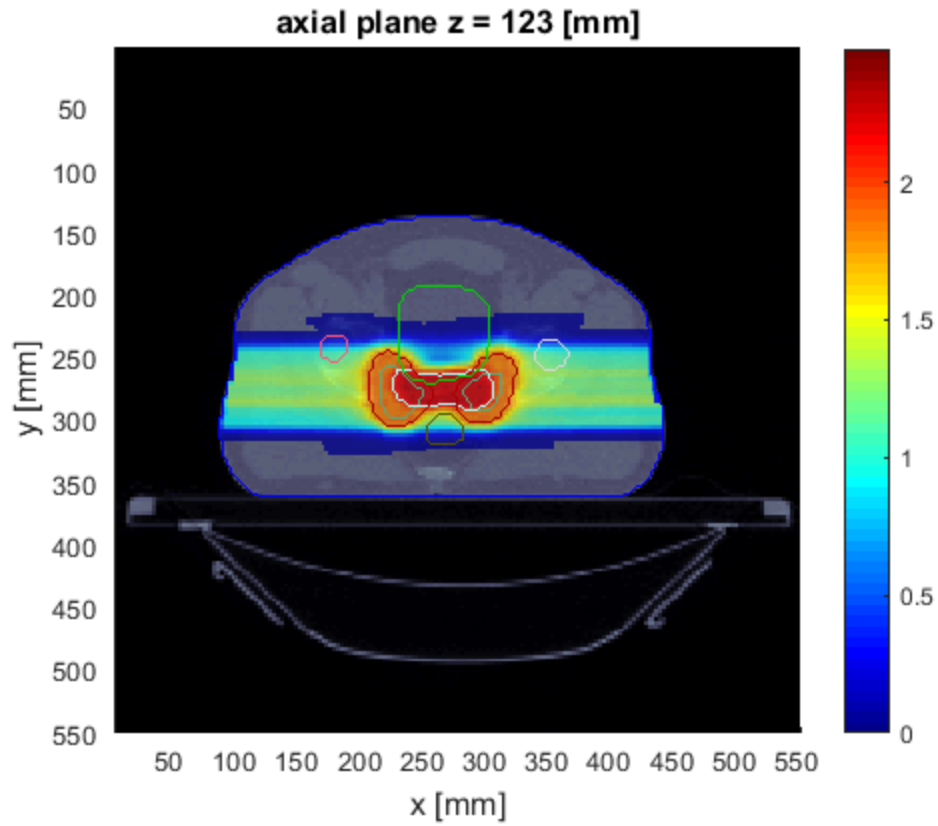


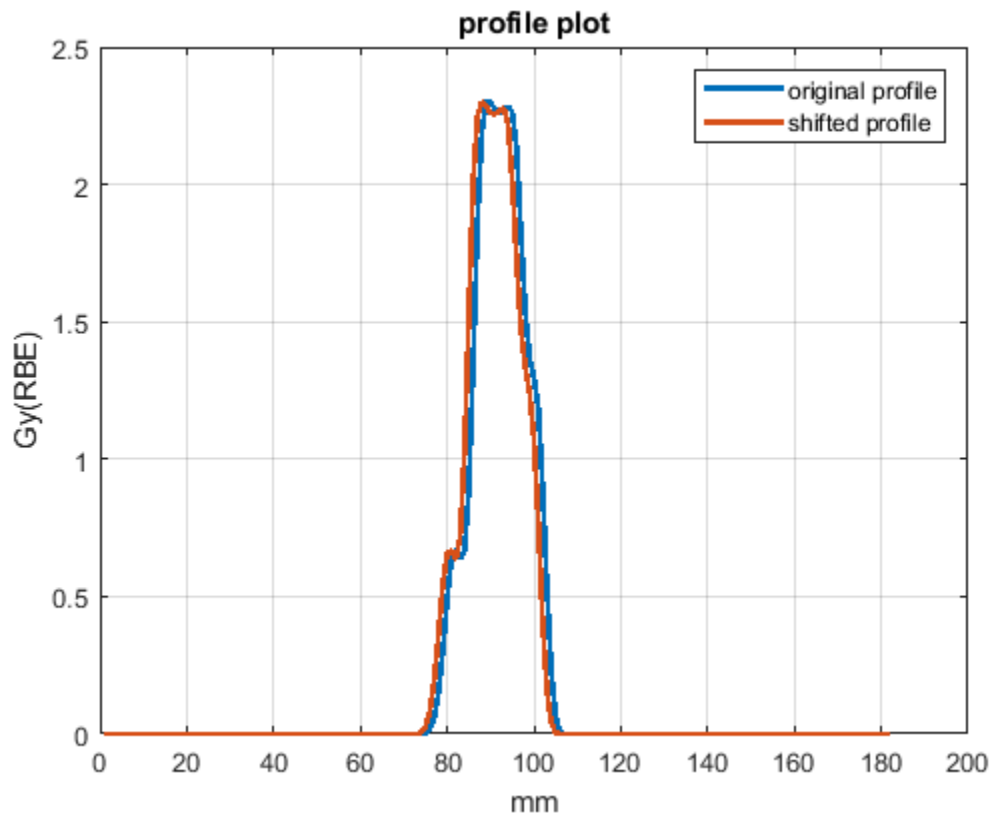
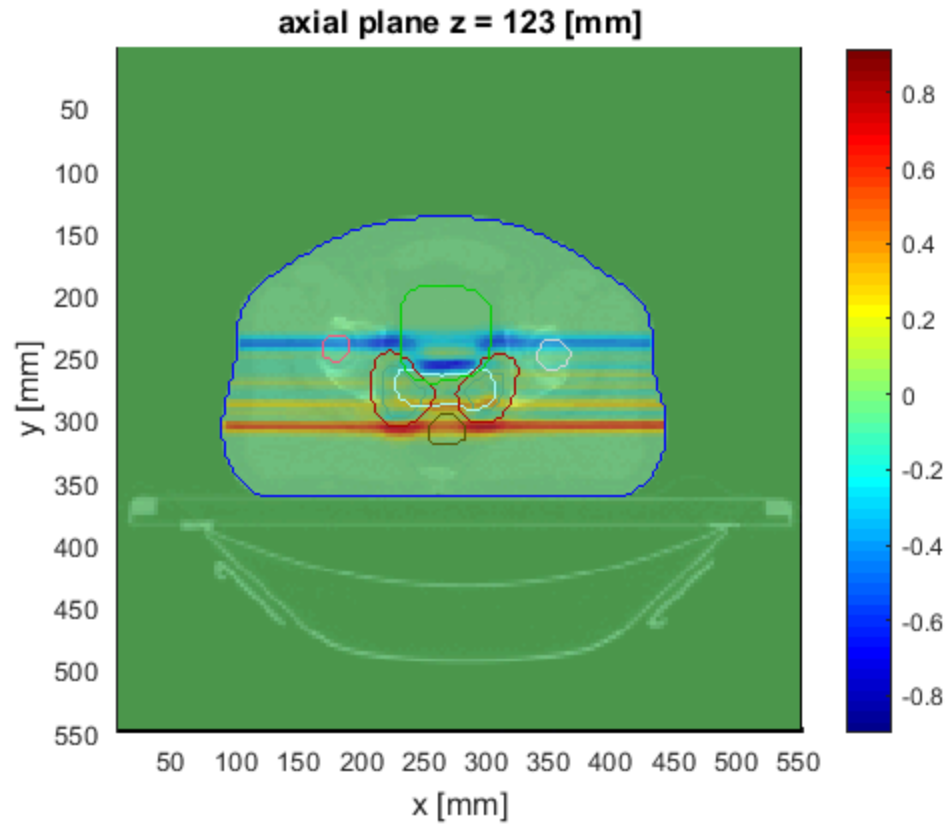
```
matRad: calculate lateral cutoff...done.  
Progress: 100.00 %  
matRad: applying a constant RBE of 1.1
```

## Visual Comparison of results

Let's compare the new recalculation against the optimization result.

```
plane = 3;  
doseWindow = [0 max([resultGUI.RBExDose(:);  
    resultGUI_isoShift.RBExDose(:)])];  
  
figure,title('original plan')  
matRad_plotSliceWrapper(gca,ct,cst,1,resultGUI.RBExDose,plane,slice,  
    [],0.75,colorcube,[],doseWindow,[]);  
figure,title('shifted plan')  
matRad_plotSliceWrapper(gca,ct,cst,1,resultGUI_isoShift.RBExDose,plane,slice,  
    [],0.75,colorcube,[],doseWindow,[]);  
  
absDiffCube = resultGUI.RBExDose-resultGUI_isoShift.RBExDose;  
figure,title('absolute difference')  
matRad_plotSliceWrapper(gca,ct,cst,1,absDiffCube,plane,slice,[],  
    [],colorcube);  
  
% Let's plot single profiles that are perpendicular to the beam  
% direction  
ixProfileY = round(pln.propStf.isoCenter(1,2)./ct.resolution.y);  
  
profileOriginal = resultGUI.RBExDose(:,ixProfileY,slice);  
profileShifted = resultGUI_isoShift.RBExDose(:,ixProfileY,slice);  
  
figure,plot(profileOriginal,'LineWidth',2),grid on,hold on,  
    plot(profileShifted,'LineWidth',2),legend({'original  
profile','shifted profile'}),  
    xlabel('mm'),ylabel('Gy(RBE)'),title('profile plot')
```





## Quantitative Comparison of results

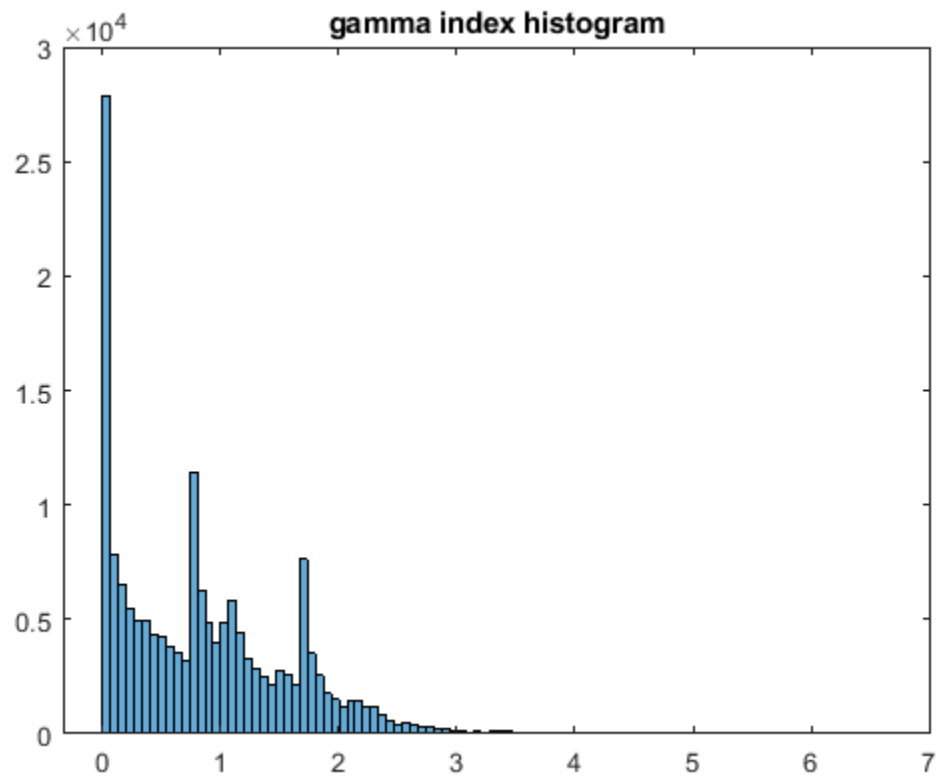
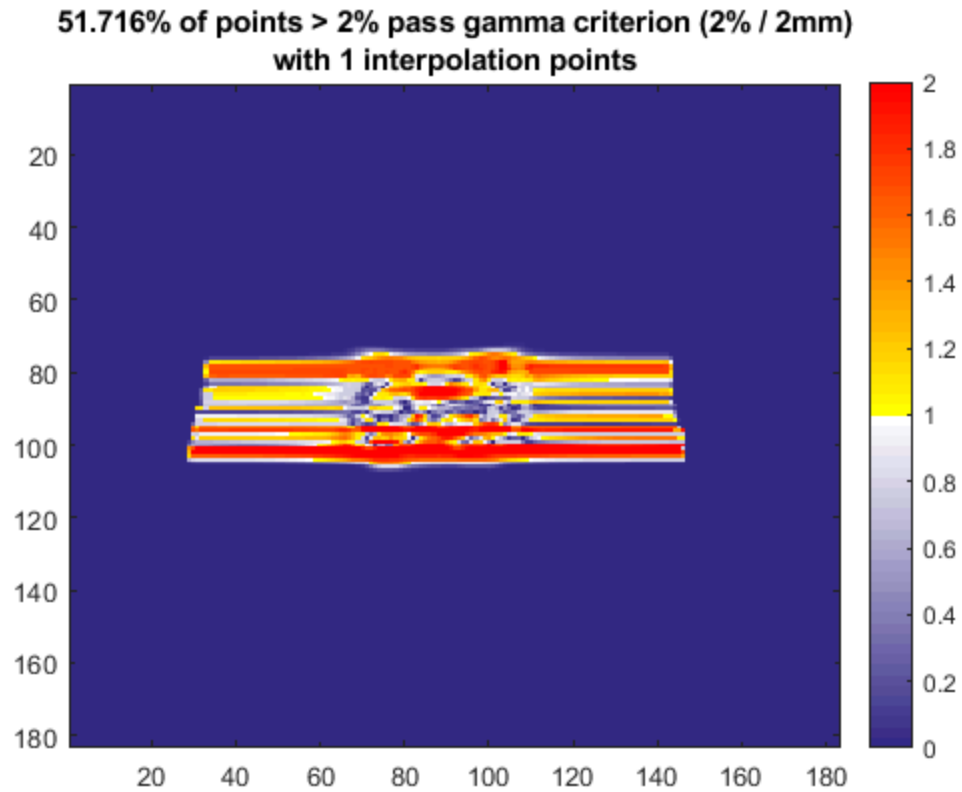
Compare the two dose cubes using a gamma-index analysis. The gamma index is a composite quality distribution equally taking into account a dose difference and a distance to agreement criterion in order to quantify differences between two dose cubes. A gamma-index value of smaller than 1 indicates a successful test and a value greater than 1 illustrates a failed test.

```
% add tools subdirectory
addpath([fileparts(fileparts(mfilename('fullpath'))
    filesep 'tools')]);

doseDifference = 2;
distToAgreement = 2;
n = 1;

[gammaCube,gammaPassRateCell] = matRad_gammaIndex(...
    resultGUI_isoShift.RBExDose,resultGUI.RBExDose,...
    [ct.resolution.x, ct.resolution.y, ct.resolution.z],...
    [doseDifference distToAgreement],slice,n,'global',cst);

[env, ~] = matRad_getEnvironment();
% Let's plot the gamma index histogram
switch env
    case 'MATLAB'
        figure,histogram(gammaCube(gammaCube>0),100),title('gamma
index histogram')
    case 'OCTAVE'
        figure,hist(gammaCube(gammaCube>0),100),title('gamma index
histogram')
end
```



*Published with MATLAB® R2018a*