
Example: Proton Treatment Plan with subsequent Isocenter shift

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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;
```

```
% dose calculation settings  
pln.propDoseCalc.doseGrid.resolution.x = 3; % [mm]  
pln.propDoseCalc.doseGrid.resolution.y = 3; % [mm]  
pln.propDoseCalc.doseGrid.resolution.z = 3; % [mm]
```

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 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.....:      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
alpha_pr  ls
    0 4.3797572e+002 0.00e+000 1.06e+000  0.0 0.00e+000 - 0.00e
+000 0.00e+000  0
```

Example: Proton Treatment Plan
with subsequent Isocenter shift

1	4.1078869e+002	0.00e+000	6.83e-002	-1.1	7.03e-002	-	
	9.93e-001	1.00e+000f	1				
2	7.5086992e+001	0.00e+000	1.81e-002	-1.7	1.38e+000	-	1.00e
	+000	1.00e+000f	1				
3	3.8671035e+001	0.00e+000	1.21e-002	-3.4	3.71e-001	-	
	9.79e-001	1.00e+000f	1				
4	3.1272594e+001	0.00e+000	1.01e-002	-3.9	2.84e-001	-	
	9.87e-001	1.00e+000f	1				
5	2.5003599e+001	0.00e+000	9.35e-003	-4.8	4.36e-001	-	
	9.99e-001	1.00e+000f	1				
6	2.0846199e+001	0.00e+000	1.23e-002	-5.5	7.04e-001	-	1.00e
	+000	1.00e+000f	1				
7	1.7591884e+001	0.00e+000	6.94e-003	-6.0	2.83e-001	-	1.00e
	+000	1.00e+000f	1				
8	1.6326338e+001	0.00e+000	5.67e-003	-7.2	2.37e-001	-	1.00e
	+000	1.00e+000f	1				
9	1.4806683e+001	0.00e+000	4.46e-003	-8.5	4.12e-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	1.2847415e+001	0.00e+000	3.64e-003	-9.4	6.27e-001	-	1.00e
	+000	1.00e+000f	1				
11	1.2215055e+001	0.00e+000	5.30e-003	-10.0	8.57e-001	-	1.00e
	+000	4.32e-001f	1				
12	1.2209742e+001	0.00e+000	5.29e-003	-11.0	3.82e-001	-	1.00e
	+000	3.47e-003f	1				
13	1.2205224e+001	0.00e+000	1.68e-002	-11.0	5.32e-001	-	1.00e
	+000	1.87e-003f	1				
14	1.1927952e+001	0.00e+000	4.79e-003	-8.7	6.61e-001	-	
	9.96e-001	9.68e-002f	1				
15	1.1917877e+001	0.00e+000	1.22e-002	-9.3	6.82e-001	-	1.00e
	+000	3.48e-003f	1				
16	1.1184059e+001	0.00e+000	5.15e-003	-10.3	8.20e-001	-	1.00e
	+000	2.53e-001f	1				
17	1.1175054e+001	0.00e+000	3.38e-003	-8.2	7.06e-001	-	
	9.28e-001	3.97e-003f	1				
18	1.1104923e+001	0.00e+000	9.96e-003	-6.7	8.32e-001	-	1.00e
	+000	2.76e-002f	1				
19	1.0978977e+001	0.00e+000	8.25e-003	-6.4	9.25e-001	-	1.00e
	+000	4.83e-002f	1				
	iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg) alpha_du
	alpha_pr	ls					
20	1.0861417e+001	0.00e+000	1.05e-002	-4.9	9.85e-001	-	
	9.98e-001	4.48e-002f	1				
21	1.0512857e+001	0.00e+000	1.39e-002	-5.3	1.10e+000	-	1.00e
	+000	1.36e-001f	1				
22	1.0232965e+001	0.00e+000	2.36e-002	-4.0	1.91e+000	-	
	7.07e-001	1.27e-001f	1				
23	1.0021448e+001	0.00e+000	9.38e-003	-6.3	1.09e+000	-	
	3.74e-001	1.07e-001f	1				
24	9.5832724e+000	0.00e+000	6.53e-003	-3.9	1.13e+000	-	
	9.00e-001	2.86e-001f	1				
25	9.4425217e+000	0.00e+000	8.83e-003	-4.9	8.39e-001	-	
	6.07e-001	1.33e-001f	1				

Example: Proton Treatment Plan
with subsequent Isocenter shift

```

26 9.2721003e+000 0.00e+000 1.31e-002 -4.6 6.98e-001 -
9.46e-001 2.44e-001f 1
27 9.0298465e+000 0.00e+000 3.82e-003 -4.3 8.57e-001 - 1.00e
+000 4.26e-001f 1
28 8.8806041e+000 0.00e+000 1.26e-002 -4.0 5.19e-001 -
8.42e-001 3.52e-001f 1
29 8.7654375e+000 0.00e+000 4.62e-003 -4.1 4.16e-001 -
5.65e-001 3.60e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr ls
30 8.6062147e+000 0.00e+000 6.24e-003 -4.1 5.72e-001 -
6.78e-001 3.74e-001f 1
31 8.5131785e+000 0.00e+000 5.41e-003 -4.1 6.41e-001 -
3.72e-001 1.70e-001f 1
32 8.3595389e+000 0.00e+000 4.83e-003 -4.3 7.22e-001 -
2.75e-001 2.77e-001f 1
33 8.1685741e+000 0.00e+000 4.60e-003 -4.1 7.79e-001 -
4.64e-001 5.53e-001f 1
34 8.0508215e+000 0.00e+000 1.17e-002 -3.4 4.84e-001 -
5.14e-001 1.00e+000f 1
35 7.9345906e+000 0.00e+000 1.37e-003 -4.3 2.32e-001 -
9.54e-001 1.00e+000f 1
36 7.8499631e+000 0.00e+000 9.59e-004 -5.0 1.94e-001 -
9.87e-001 1.00e+000f 1
37 7.7721251e+000 0.00e+000 1.05e-003 -4.8 1.93e-001 -
9.43e-001 8.12e-001f 1
38 7.6778951e+000 0.00e+000 3.94e-003 -5.0 3.20e-001 -
9.98e-001 5.09e-001f 1
39 7.6329474e+000 0.00e+000 6.16e-003 -5.3 5.41e-001 -
9.18e-001 1.30e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr ls
40 7.5158258e+000 0.00e+000 9.85e-003 -4.1 6.00e-001 -
6.09e-001 3.23e-001f 1
41 7.4610116e+000 0.00e+000 2.81e-003 -5.8 9.18e-001 -
7.10e-001 1.24e-001f 1
42 7.3606139e+000 0.00e+000 5.82e-003 -4.1 5.33e-001 -
3.63e-001 3.14e-001f 1
43 7.2971402e+000 0.00e+000 3.10e-003 -4.7 6.48e-001 -
4.13e-001 1.84e-001f 1
44 7.2251428e+000 0.00e+000 4.20e-003 -4.3 5.35e-001 -
4.94e-001 2.57e-001f 1
45 7.1573042e+000 0.00e+000 5.76e-003 -4.4 5.28e-001 -
9.00e-001 3.03e-001f 1
46 7.0772400e+000 0.00e+000 2.55e-003 -4.7 8.17e-001 -
6.37e-001 3.30e-001f 1
47 7.0552974e+000 0.00e+000 4.08e-003 -10.6 4.24e-001 -
3.78e-001 1.37e-001f 1
48 7.7538384e+000 0.00e+000 5.52e-003 -2.8 1.07e+001 -
3.80e-002 2.50e-001f 1
49 7.1955782e+000 0.00e+000 4.61e-003 -4.2 2.19e+000 - 1.00e
+000 5.36e-001f 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

50	6.9497764e+000	0.00e+000	9.67e-003	-4.2	6.84e-001	-	1.00e
+000 7.00e-001f 1							
51	6.8964778e+000	0.00e+000	5.85e-003	-4.7	5.27e-001	-	
8.88e-001 3.32e-001f 1							
52	6.8436481e+000	0.00e+000	4.56e-003	-5.5	6.25e-001	-	
8.31e-001 3.23e-001f 1							
53	6.8060607e+000	0.00e+000	5.08e-003	-4.9	5.26e-001	-	
6.07e-001 2.58e-001f 1							
54	6.7560011e+000	0.00e+000	1.18e-002	-4.3	4.92e-001	-	
5.02e-001 6.39e-001f 1							
55	6.7222675e+000	0.00e+000	2.60e-003	-4.8	6.01e-001	-	
4.96e-001 2.32e-001f 1							
56	6.6746449e+000	0.00e+000	3.10e-003	-4.2	5.90e-001	-	
4.52e-001 2.82e-001f 1							
57	6.6225337e+000	0.00e+000	2.46e-003	-4.4	6.42e-001	-	
4.69e-001 2.76e-001f 1							
58	8.2500213e+000	0.00e+000	5.30e-003	-2.4	1.83e+001	-	
1.29e-002 1.58e-001f 1							
59	6.6454438e+000	0.00e+000	2.99e-003	-4.1	3.07e+000	-	
2.02e-001 8.59e-001f 1							
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
alpha_pr ls							
60	6.5771759e+000	0.00e+000	5.09e-003	-4.1	4.99e-001	-	1.00e
+000 4.43e-001f 1							
61	6.5336113e+000	0.00e+000	1.52e-002	-4.8	6.41e-001	-	
9.88e-001 3.49e-001f 1							
62	6.4775403e+000	0.00e+000	3.52e-003	-5.8	8.64e-001	-	1.00e
+000 3.72e-001f 1							
63	6.4449004e+000	0.00e+000	1.16e-003	-4.1	4.05e-001	-	
4.98e-001 5.00e-001f 2							
64	6.4079727e+000	0.00e+000	2.04e-003	-4.1	9.41e-002	-	1.00e
+000 1.00e+000f 1							
65	6.3207261e+000	0.00e+000	8.75e-004	-5.0	9.01e-001	-	
2.89e-001 4.27e-001f 1							
66	6.3019422e+000	0.00e+000	3.06e-003	-4.8	6.05e-001	-	
4.83e-001 1.34e-001f 1							
67	6.2640220e+000	0.00e+000	5.83e-003	-4.9	7.61e-001	-	
8.39e-001 2.39e-001f 1							
68	6.2259820e+000	0.00e+000	4.14e-003	-4.9	7.61e-001	-	
8.14e-001 2.43e-001f 1							
69	6.2080493e+000	0.00e+000	3.86e-003	-10.8	1.05e+000	-	
3.96e-001 8.82e-002f 1							
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
alpha_pr ls							
70	6.1657456e+000	0.00e+000	5.84e-003	-5.5	1.08e+000	-	
3.68e-001 2.04e-001f 1							
71	6.1370482e+000	0.00e+000	7.76e-003	-5.5	1.04e+000	-	
8.72e-001 1.58e-001f 1							
72	6.0859933e+000	0.00e+000	3.35e-003	-4.4	8.28e-001	-	
4.81e-001 3.65e-001f 1							
73	6.0582691e+000	0.00e+000	4.07e-003	-4.4	4.89e-001	-	
6.27e-001 2.95e-001f 1							
74	6.0340044e+000	0.00e+000	2.01e-003	-4.5	7.03e-001	-	
3.79e-001 1.81e-001f 1							

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75	6.0067968e+000	0.00e+000	3.99e-003	-6.1	9.32e-001	-
	2.18e-001	1.66e-001f	1			
76	5.9761879e+000	0.00e+000	2.84e-003	-4.9	9.62e-001	-
	7.22e-001	1.94e-001f	1			
77	5.9514259e+000	0.00e+000	2.06e-003	-4.3	4.72e-001	-
	2.58e-001	2.51e-001f	1			
78	5.9301538e+000	0.00e+000	3.76e-003	-4.8	6.46e-001	-
	3.49e-001	1.92e-001f	1			
79	5.9019870e+000	0.00e+000	2.60e-003	-4.7	8.02e-001	-
	3.68e-001	2.34e-001f	1			
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg) alpha_du
	alpha_pr	ls				
80	5.8874722e+000	0.00e+000	8.73e-003	-5.0	6.37e-001	-
	7.07e-001	1.50e-001f	1			
81	5.8363212e+000	0.00e+000	3.03e-003	-4.9	8.38e-001	-
	7.90e-001	5.02e-001f	1			
82	5.8327784e+000	0.00e+000	6.00e-003	-6.8	7.40e-001	-
	6.14e-001	3.30e-002f	1			
83	5.8173565e+000	0.00e+000	4.21e-003	-4.1	3.19e-001	-
	5.30e-001	1.00e+000f	1			
84	5.7847139e+000	0.00e+000	1.97e-003	-4.7	5.54e-001	-
	5.71e-001	4.47e-001f	1			
85	5.7499361e+000	0.00e+000	2.99e-003	-4.7	6.04e-001	-
	6.67e-001	5.06e-001f	1			
86	5.7336923e+000	0.00e+000	1.61e-003	-4.8	7.17e-001	-
	4.68e-001	1.59e-001f	1			
87	5.7253575e+000	0.00e+000	7.80e-003	-6.1	5.68e-001	-
	5.37e-001	1.19e-001f	1			
88	5.6948046e+000	0.00e+000	2.88e-003	-5.7	1.12e+000	-
	7.93e-001	2.37e-001f	1			
89	5.6855603e+000	0.00e+000	3.69e-003	-5.2	8.07e-001	-
	2.18e-001	8.10e-002f	1			
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg) alpha_du
	alpha_pr	ls				
90	5.6929520e+000	0.00e+000	4.48e-003	-4.0	9.04e-001	-
	3.84e-001	1.00e+000f	1			
91	5.6654813e+000	0.00e+000	1.24e-003	-4.5	6.48e-001	-
	7.18e-001	2.63e-001f	1			
92	5.6386652e+000	0.00e+000	5.33e-003	-4.5	5.53e-001	-
	8.61e-001	3.76e-001f	1			
93	5.6056039e+000	0.00e+000	4.42e-003	-4.6	5.78e-001	-
	9.37e-001	5.01e-001f	1			
94	5.5906091e+000	0.00e+000	7.93e-003	-4.9	7.07e-001	-
	9.29e-001	2.00e-001f	1			
95	5.5715780e+000	0.00e+000	4.40e-003	-5.5	9.45e-001	-
	6.58e-001	1.92e-001f	1			
96	5.5526547e+000	0.00e+000	2.98e-003	-4.8	7.18e-001	-
	3.58e-001	2.36e-001f	1			
97	5.5319082e+000	0.00e+000	3.38e-003	-5.0	7.86e-001	-
	4.31e-001	2.42e-001f	1			
98	5.5109597e+000	0.00e+000	2.16e-003	-5.1	1.02e+000	-
	5.59e-001	2.08e-001f	1			
99	5.5074125e+000	0.00e+000	3.71e-003	-11.0	8.31e-001	-
	2.24e-001	3.91e-002f	1			

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with subsequent Isocenter shift

iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
alpha_pr	ls						
100	5.4795582e+000	0.00e+000	2.37e-003	-6.2	1.23e+000	-	
	2.30e-001	2.17e-001f	1				
101	5.4730004e+000	0.00e+000	2.73e-003	-5.5	1.25e+000	-	
	8.62e-001	4.46e-002f	1				
102	5.4536776e+000	0.00e+000	4.88e-003	-4.8	5.97e-001	-	
	3.83e-001	2.65e-001f	1				
103	5.6283167e+000	0.00e+000	4.17e-003	-3.2	1.15e+001	-	
	1.65e-002	1.12e-001f	1				
104	5.4248219e+000	0.00e+000	4.41e-003	-4.8	2.17e+000	-	
	1.36e-001	7.07e-001f	1				
105	5.4049961e+000	0.00e+000	2.04e-003	-4.8	6.80e-001	-	
	5.05e-001	1.43e-001f	1				
106	5.3995461e+000	0.00e+000	4.41e-003	-6.4	6.10e-001	-	
	6.68e-001	8.96e-002f	1				
107	5.3804253e+000	0.00e+000	4.01e-003	-5.2	7.00e-001	-	
	9.46e-001	2.51e-001f	1				
108	5.3657193e+000	0.00e+000	3.02e-003	-5.3	9.02e-001	-	
	4.51e-001	1.35e-001f	1				
109	5.3558920e+000	0.00e+000	3.91e-003	-11.0	7.10e-001	-	
	1.27e-001	1.21e-001f	1				
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
alpha_pr	ls						
110	5.3331654e+000	0.00e+000	2.63e-003	-5.6	1.08e+000	-	
	4.26e-001	1.91e-001f	1				
111	5.3164685e+000	0.00e+000	1.87e-003	-5.9	1.52e+000	-	
	3.16e-001	9.83e-002f	1				
112	5.3138576e+000	0.00e+000	4.91e-003	-11.0	7.43e-001	-	
	2.09e-001	3.23e-002f	1				
113	5.2866218e+000	0.00e+000	2.05e-003	-5.7	1.44e+000	-	
	6.32e-001	1.87e-001f	1				
114	5.3187769e+000	0.00e+000	1.15e-003	-4.3	3.22e-001	-	
	2.24e-001	1.00e+000f	1				
115	5.2958090e+000	0.00e+000	9.72e-004	-4.9	9.95e-001	-	
	3.45e-001	1.91e-001f	1				
116	5.2529006e+000	0.00e+000	1.87e-003	-4.9	1.06e+000	-	
	5.46e-001	3.35e-001f	1				
117	5.2305804e+000	0.00e+000	3.85e-003	-4.6	5.56e-001	-	
	6.24e-001	3.25e-001f	1				
118	5.2241578e+000	0.00e+000	5.17e-003	-5.1	4.32e-001	-	
	7.41e-001	1.41e-001f	1				
119	5.2041025e+000	0.00e+000	1.89e-003	-4.7	5.32e-001	-	
	6.69e-001	4.60e-001f	1				
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
alpha_pr	ls						
120	5.1971050e+000	0.00e+000	2.24e-003	-5.9	7.37e-001	-	
	5.66e-001	9.48e-002f	1				
121	5.1931953e+000	0.00e+000	7.09e-003	-7.2	6.80e-001	-	
	4.79e-001	5.76e-002f	1				
122	5.1662009e+000	0.00e+000	3.82e-003	-5.1	1.06e+000	-	
	4.43e-001	2.64e-001f	1				
123	5.1855544e+000	0.00e+000	2.08e-003	-4.3	2.93e-001	-	
	4.25e-001	1.00e+000f	1				

Example: Proton Treatment Plan
with subsequent Isocenter shift

124	5.1665936e+000	0.00e+000	1.79e-003	-4.7	5.79e-001	-
	4.57e-001	2.92e-001f	1			
125	5.1404920e+000	0.00e+000	2.37e-003	-4.7	5.12e-001	-
	5.74e-001	4.32e-001f	1			
126	5.1333176e+000	0.00e+000	2.13e-003	-4.8	3.93e-001	-
	4.80e-001	1.55e-001f	1			
127	5.1165762e+000	0.00e+000	8.45e-003	-5.1	5.01e-001	-
	7.10e-001	3.17e-001f	1			
128	5.1096427e+000	0.00e+000	7.93e-003	-5.7	5.42e-001	-
	5.92e-001	1.31e-001f	1			
129	5.0957870e+000	0.00e+000	5.99e-003	-5.9	7.84e-001	-
	8.70e-001	1.94e-001f	1			
iter objective inf_pr inf_du lg(mu) d lg(rg) alpha_du						
alpha_pr ls						
130	5.0774987e+000	0.00e+000	3.86e-003	-6.0	8.29e-001	-
	8.13e-001	2.64e-001f	1			
131	5.0675379e+000	0.00e+000	3.36e-003	-5.3	6.30e-001	-
	7.34e-001	1.87e-001f	1			
132	5.0563330e+000	0.00e+000	5.22e-003	-4.7	2.73e-001	-
	7.01e-001	5.58e-001f	1			
133	5.0460255e+000	0.00e+000	2.46e-003	-4.8	3.68e-001	-
	3.66e-001	3.04e-001f	1			
134	5.0359030e+000	0.00e+000	1.20e-003	-4.9	5.46e-001	-
	3.07e-001	2.03e-001f	1			
135	5.0255534e+000	0.00e+000	1.16e-003	-4.9	8.91e-001	-
	6.18e-001	1.51e-001f	1			
136	5.0175313e+000	0.00e+000	4.48e-003	-5.2	9.16e-001	-
	3.43e-001	1.20e-001f	1			
137	4.9935769e+000	0.00e+000	3.12e-003	-5.2	1.25e+000	-
	8.81e-001	2.90e-001f	1			
138	4.9855568e+000	0.00e+000	2.33e-003	-5.9	1.09e+000	-
	1.68e-001	1.12e-001f	1			
139	4.9694039e+000	0.00e+000	2.31e-003	-5.5	1.17e+000	-
	5.14e-001	2.23e-001f	1			
iter objective inf_pr inf_du lg(mu) d lg(rg) alpha_du						
alpha_pr ls						
140	4.9603623e+000	0.00e+000	1.77e-003	-11.0	1.46e+000	-
	3.97e-001	1.03e-001f	1			
141	4.9446081e+000	0.00e+000	1.04e-003	-4.6	2.53e-001	-
	3.41e-001	1.00e+000f	1			
142	4.9380850e+000	0.00e+000	1.71e-003	-5.5	9.58e-001	-
	4.63e-001	1.16e-001f	1			
143	4.9216076e+000	0.00e+000	3.37e-003	-4.9	8.19e-001	-
	6.27e-001	3.82e-001f	1			
144	4.9149058e+000	0.00e+000	5.37e-003	-5.2	5.81e-001	-
	4.35e-001	1.80e-001f	1			
145	4.9075415e+000	0.00e+000	2.79e-003	-5.4	8.23e-001	-
	6.80e-001	1.41e-001f	1			
146	4.8916279e+000	0.00e+000	3.41e-003	-5.2	8.31e-001	-
	3.70e-001	3.08e-001f	1			
147	4.8882254e+000	0.00e+000	3.68e-003	-11.0	7.99e-001	-
	4.47e-001	7.04e-002f	1			
148	4.8765019e+000	0.00e+000	2.35e-003	-7.3	9.92e-001	-
	5.03e-001	2.06e-001f	1			

Example: Proton Treatment Plan
with subsequent Isocenter shift

```

149 5.0009188e+000 0.00e+000 2.04e-003 -3.9 4.08e+000 -
2.11e-002 4.74e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr  ls
150 4.8935944e+000 0.00e+000 1.98e-003 -5.4 2.65e+000 -
4.78e-002 4.95e-001f 1
151 4.8912896e+000 0.00e+000 3.23e-003 -5.4 1.26e+000 - 1.00e
+000 2.40e-002f 1
152 4.8626442e+000 0.00e+000 7.21e-003 -5.4 1.10e+000 -
6.31e-001 3.85e-001f 1
153 4.8549814e+000 0.00e+000 7.76e-003 -5.6 7.11e-001 -
9.97e-001 1.78e-001f 1
154 4.8446526e+000 0.00e+000 2.93e-003 -5.2 4.96e-001 -
7.15e-001 3.39e-001f 1
155 4.8410230e+000 0.00e+000 6.75e-003 -5.9 4.81e-001 -
7.41e-001 1.30e-001f 1
156 4.8346397e+000 0.00e+000 5.65e-003 -6.9 6.05e-001 -
4.35e-001 1.86e-001f 1
157 4.8267163e+000 0.00e+000 5.43e-003 -6.3 8.08e-001 -
7.47e-001 1.74e-001f 1
158 4.8190755e+000 0.00e+000 3.28e-003 -6.2 8.35e-001 -
4.63e-001 1.63e-001f 1
159 4.8132600e+000 0.00e+000 2.30e-003 -5.3 7.23e-001 -
2.53e-001 1.31e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr  ls
160 4.8049667e+000 0.00e+000 9.16e-004 -4.8 1.95e-001 -
3.18e-001 1.00e+000f 1
161 4.7994937e+000 0.00e+000 8.13e-004 -4.9 5.44e-001 -
4.20e-001 2.36e-001f 1
162 4.7857813e+000 0.00e+000 7.53e-004 -5.0 1.11e+000 -
3.95e-001 2.88e-001f 1
163 4.7799209e+000 0.00e+000 2.31e-003 -5.3 1.08e+000 -
5.00e-001 1.41e-001f 1
164 4.7695435e+000 0.00e+000 1.52e-003 -5.7 1.22e+000 -
5.59e-001 2.02e-001f 1
165 4.7635091e+000 0.00e+000 2.62e-003 -6.6 1.40e+000 -
5.40e-001 1.02e-001f 1
166 4.7501624e+000 0.00e+000 1.66e-003 -4.7 6.98e-001 -
5.89e-001 5.39e-001f 1
167 4.7438780e+000 0.00e+000 1.53e-003 -4.9 6.37e-001 -
4.62e-001 2.49e-001f 1
168 4.7388600e+000 0.00e+000 2.27e-003 -4.9 6.10e-001 -
3.39e-001 2.44e-001f 1
169 4.7323564e+000 0.00e+000 1.30e-003 -4.9 6.79e-001 -
4.39e-001 2.88e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr  ls
170 4.7305070e+000 0.00e+000 1.82e-003 -5.4 9.51e-001 -
4.30e-001 5.42e-002f 1
171 4.7184328e+000 0.00e+000 6.89e-003 -6.1 1.12e+000 -
4.87e-001 3.12e-001f 1
172 4.7071066e+000 0.00e+000 5.39e-003 -6.6 1.55e+000 -
1.99e-001 2.02e-001f 1

```

Example: Proton Treatment Plan
with subsequent Isocenter shift

```

173 4.7058131e+000 0.00e+000 6.97e-003 -5.4 8.04e-001 -
6.30e-001 4.19e-002f 1
174 4.6935621e+000 0.00e+000 2.17e-003 -4.9 1.14e+000 -
6.91e-001 2.94e-001f 1
175 4.6875891e+000 0.00e+000 2.53e-003 -4.8 4.09e-001 -
3.98e-001 3.16e-001f 1
176 4.6799790e+000 0.00e+000 1.60e-003 -5.0 6.00e-001 -
2.59e-001 2.71e-001f 1
177 5.0853764e+000 0.00e+000 4.14e-003 -2.9 2.67e+001 -
3.98e-003 1.17e-001f 1
178 4.7143241e+000 0.00e+000 1.06e-003 -4.8 3.24e+000 -
5.77e-003 8.09e-001f 1
179 4.7145642e+000 0.00e+000 2.08e-002 -4.8 1.17e+001 -
3.48e-002 5.28e-003f 7
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr  ls
180 4.6959088e+000 0.00e+000 7.71e-004 -4.8 7.61e-001 -
9.57e-001 3.39e-001f 1
181 4.6727963e+000 0.00e+000 6.17e-003 -4.8 7.67e-001 - 1.00e
+000 8.67e-001f 1
182 4.6695099e+000 0.00e+000 5.98e-003 -5.3 3.91e-001 -
8.14e-001 2.00e-001f 1
183 4.6651104e+000 0.00e+000 3.99e-003 -5.7 5.42e-001 -
8.28e-001 1.81e-001f 1
184 4.6558348e+000 0.00e+000 2.99e-003 -5.9 7.23e-001 -
9.55e-001 2.98e-001f 1
185 4.6470223e+000 0.00e+000 1.83e-003 -6.1 8.62e-001 -
8.02e-001 2.81e-001f 1
186 4.6413442e+000 0.00e+000 1.90e-003 -5.3 3.96e-001 -
6.06e-001 4.20e-001f 1
187 4.6314216e+000 0.00e+000 2.51e-003 -5.1 3.95e-001 -
5.90e-001 1.00e+000f 1
188 4.6261166e+000 0.00e+000 3.88e-004 -5.2 7.67e-001 -
3.86e-001 2.84e-001f 1
189 4.6188570e+000 0.00e+000 2.72e-003 -5.4 1.35e+000 -
4.53e-001 2.37e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr  ls
190 4.6128095e+000 0.00e+000 1.20e-003 -5.6 1.99e+000 -
7.03e-001 1.39e-001f 1
191 4.6062801e+000 0.00e+000 2.69e-003 -5.5 1.08e+000 -
1.81e-001 2.96e-001f 1
192 4.6005286e+000 0.00e+000 1.40e-003 -5.5 1.83e+000 -
9.63e-001 1.45e-001f 1
193 4.5924764e+000 0.00e+000 1.79e-003 -5.5 2.13e+000 -
2.63e-001 1.70e-001f 1
194 4.5883923e+000 0.00e+000 2.40e-003 -5.5 2.18e+000 -
4.34e-001 8.11e-002f 1
195 4.5765015e+000 0.00e+000 1.80e-003 -5.5 2.48e+000 -
1.98e-001 2.04e-001f 1
196 4.5732161e+000 0.00e+000 1.47e-003 -6.1 2.46e+000 -
2.44e-001 5.55e-002f 1
197 4.5682429e+000 0.00e+000 1.93e-003 -6.3 2.56e+000 -
4.26e-001 7.89e-002f 1

```

Example: Proton Treatment Plan
with subsequent Isocenter shift

```

198 4.5604844e+000 0.00e+000 2.70e-003 -5.9 2.49e+000 -
3.70e-001 1.27e-001f 1
199 4.5513251e+000 0.00e+000 1.90e-003 -5.7 2.56e+000 -
1.51e-001 1.38e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr ls
200 4.5482728e+000 0.00e+000 4.44e-003 -5.2 1.19e+000 -
5.08e-001 1.02e-001f 1
201 4.6292471e+000 0.00e+000 4.05e-003 -3.4 6.24e+001 -
1.07e-002 3.64e-002f 1
202 4.6221226e+000 0.00e+000 3.90e-003 -5.2 3.15e+000 -
1.66e-002 5.25e-002f 1
203 4.5706463e+000 0.00e+000 5.69e-003 -5.2 2.89e+000 -
4.51e-001 4.42e-001f 1
204 4.5582694e+000 0.00e+000 2.64e-003 -5.2 1.84e+000 -
4.28e-001 2.09e-001f 1
205 4.5441603e+000 0.00e+000 3.61e-003 -5.2 1.45e+000 -
7.52e-001 3.35e-001f 1
206 4.5386418e+000 0.00e+000 3.57e-003 -5.1 7.98e-001 -
7.23e-001 2.44e-001f 1
207 4.5391802e+000 0.00e+000 8.09e-004 -4.9 5.12e-001 -
5.09e-001 1.00e+000f 1
208 4.5369142e+000 0.00e+000 1.84e-003 -5.1 3.96e-001 - 1.00e
+000 2.49e-001f 1
209 4.5287469e+000 0.00e+000 6.69e-003 -5.7 1.09e+000 -
9.01e-001 3.11e-001f 1
iter   objective   inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr ls
210 4.5224960e+000 0.00e+000 1.92e-003 -5.3 9.36e-001 -
6.42e-001 2.85e-001f 1
211 4.5204834e+000 0.00e+000 1.27e-003 -5.1 4.74e-001 -
4.18e-001 1.91e-001f 1
212 4.5102765e+000 0.00e+000 1.81e-003 -5.3 1.23e+000 -
3.55e-001 3.47e-001f 1
213 4.5059642e+000 0.00e+000 9.76e-004 -5.3 1.02e+000 -
3.67e-001 1.77e-001f 1
214 4.5051061e+000 0.00e+000 2.93e-003 -6.8 1.32e+000 -
2.74e-001 2.48e-002f 1
215 4.4981644e+000 0.00e+000 1.54e-003 -6.5 2.22e+000 -
3.03e-001 1.17e-001f 1
216 4.4925760e+000 0.00e+000 1.84e-003 -5.3 1.13e+000 -
2.23e-001 1.74e-001f 1
217 4.4903224e+000 0.00e+000 1.83e-003 -6.4 1.55e+000 -
1.83e-001 5.10e-002f 1
218 4.4871978e+000 0.00e+000 2.01e-003 -5.3 1.21e+000 -
3.11e-001 8.82e-002f 1
219 4.4846916e+000 0.00e+000 1.12e-003 -4.9 1.46e+000 -
2.15e-001 4.38e-001f 1

```

Number of Iterations.....: 219

```

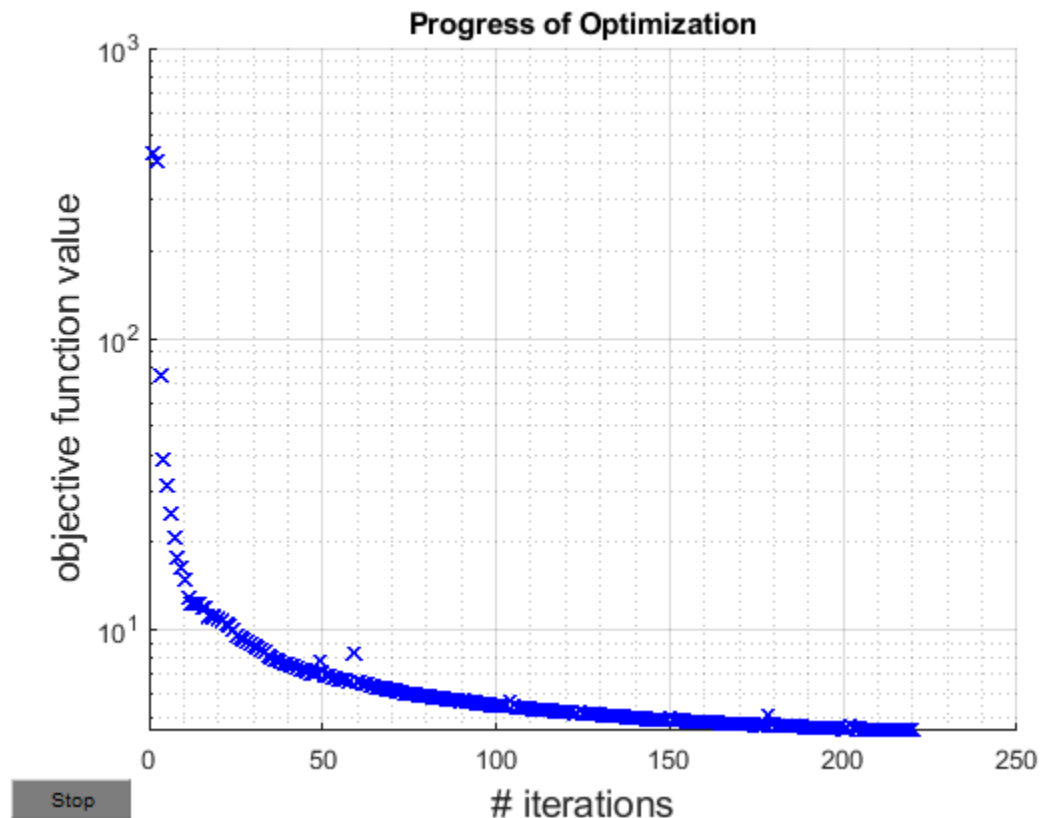
                                     (scaled)                               (unscaled)
Objective.....: 4.4846916045641736e+000
4.4846916045641736e+000

```

```
Dual infeasibility.....: 1.1246414861891072e-003
1.1246414861891072e-003
Constraint violation....: 0.0000000000000000e+000
0.0000000000000000e+000
Complementarity.....: 4.0607728791104100e-004
4.0607728791104100e-004
Overall NLP error.....: 1.1246414861891072e-003
1.1246414861891072e-003
```

```
Number of objective function evaluations      = 235
Number of objective gradient evaluations     = 220
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) = 69.821
Total CPU secs in NLP function evaluations    = 185.803
```

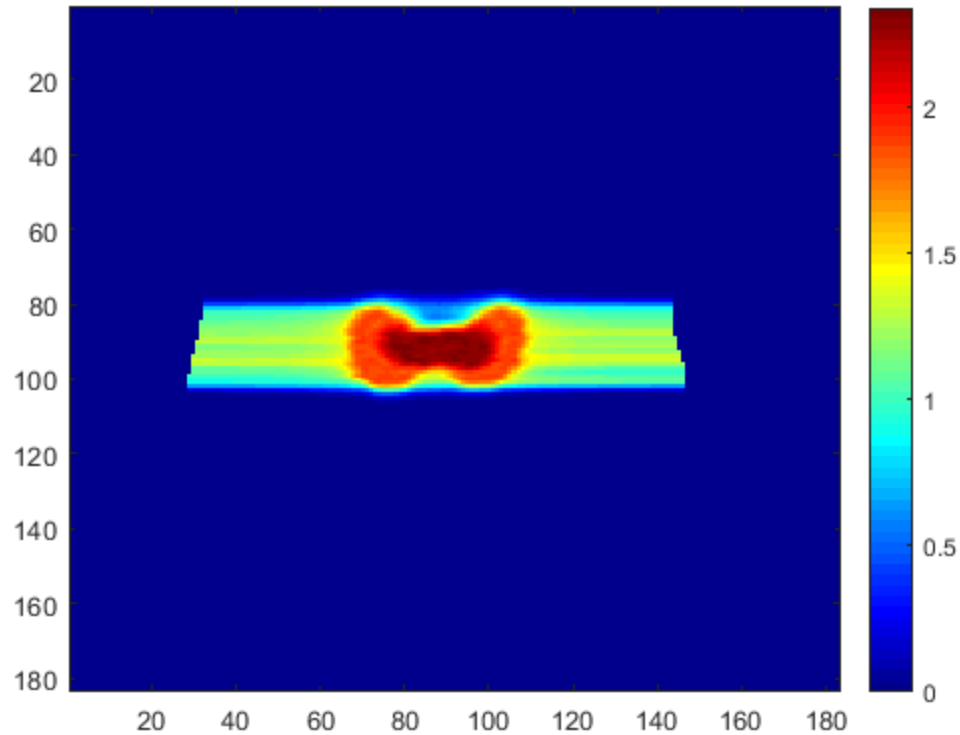
EXIT: Solved To Acceptable Level.
matRad: applying a constant RBE of 1.1



Plot the Resulting Dose Slice

Let's plot the transversal iso-center dose slice

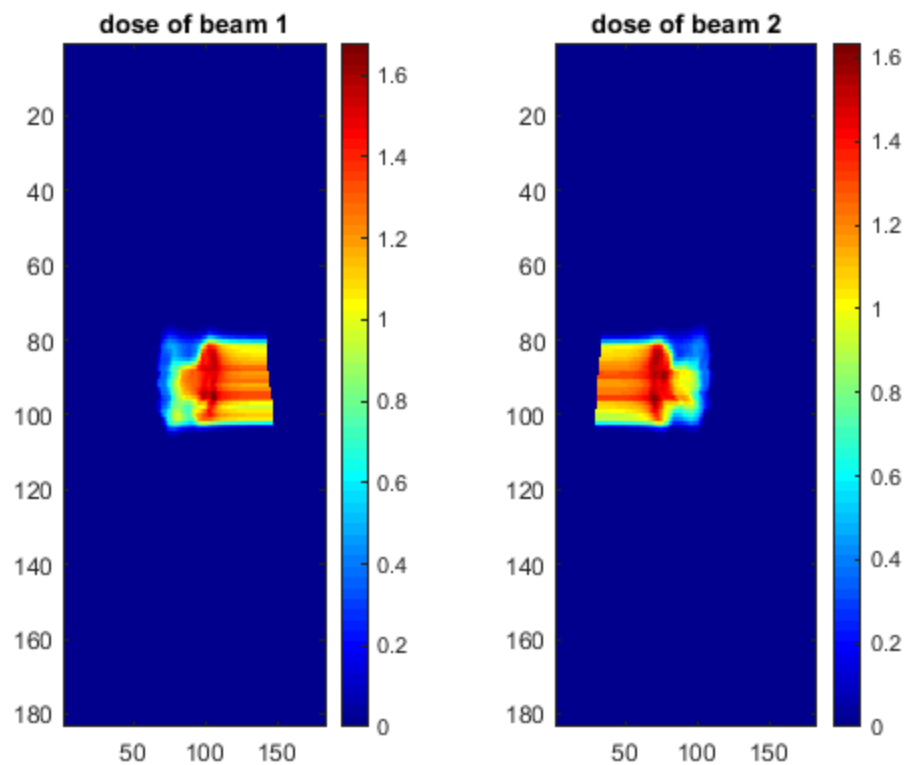
```
slice = round(pln.propStf.isoCenter(1,3)./ct.resolution.z);  
figure  
imagesc(resultGUI.RBExDose(:,:,slice)),colorbar,colormap(jet)
```



Plot the Resulting Beam Dose Slice

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

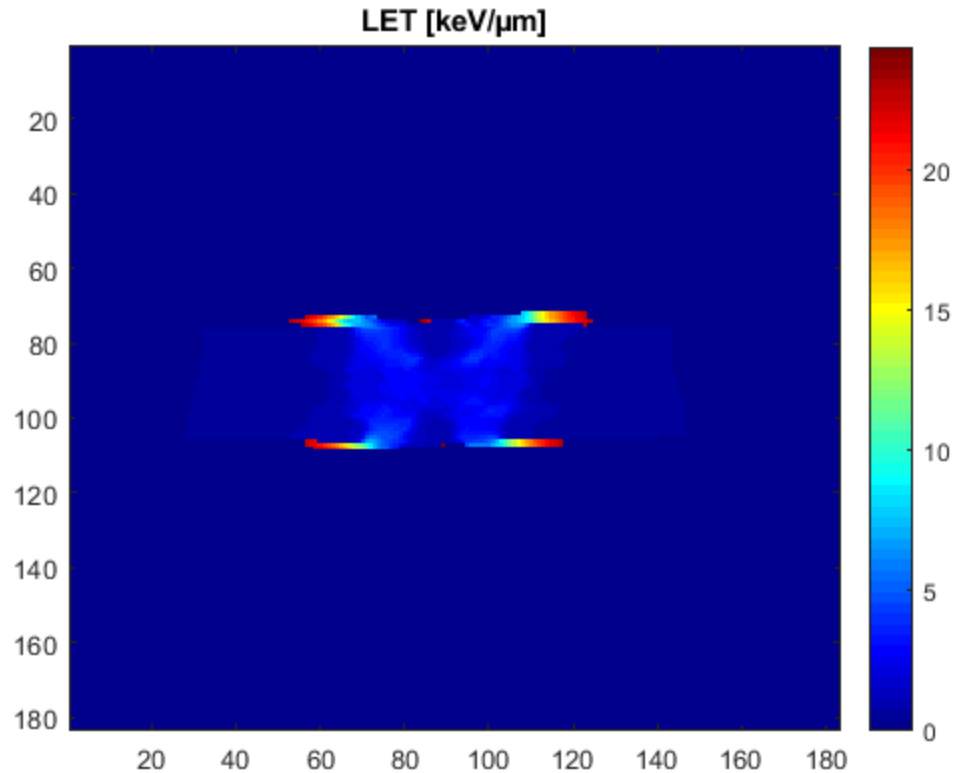
```
figure  
subplot(121),imagesc(resultGUI.RBExDose_beam1(:,:,slice)),colorbar,colormap(jet),t  
    of beam 1')  
subplot(122),imagesc(resultGUI.RBExDose_beam2(:,:,slice)),colorbar,colormap(jet),t  
    of beam 2')
```



and the corresponding LET distribution

Transversal iso-center slice

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



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.
```

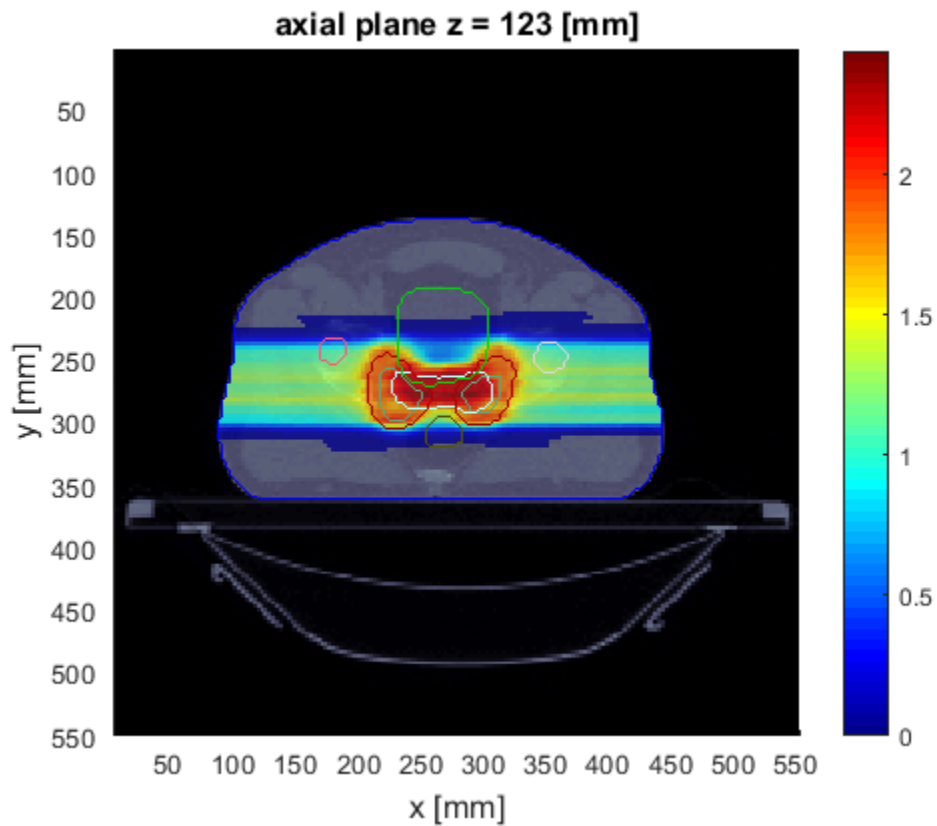
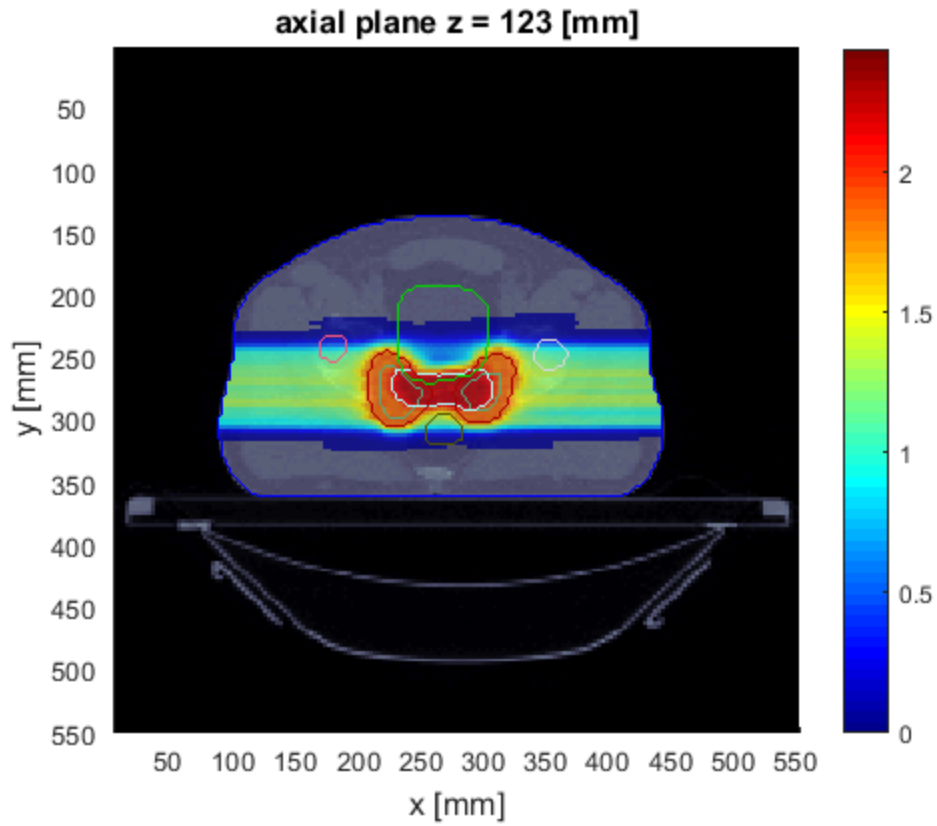


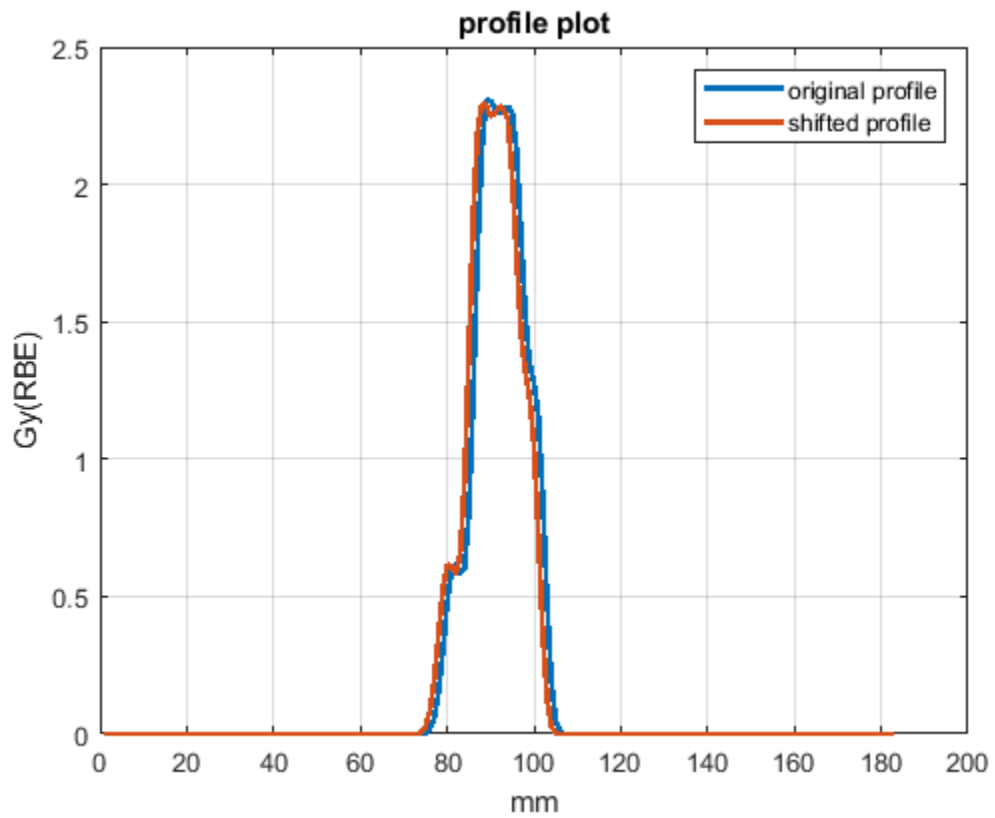
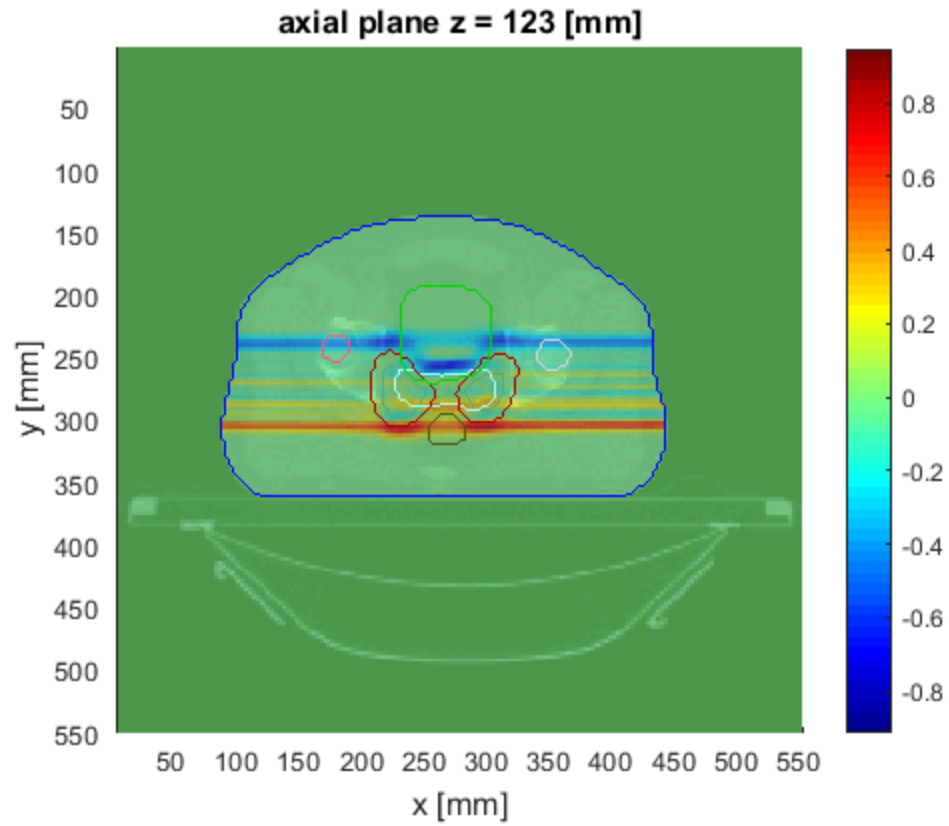
```
matRad: calculate lateral cutoff...done.  
Progress: 100.00 %  
matRad: applying a constant RBE of 1.1
```

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')
```

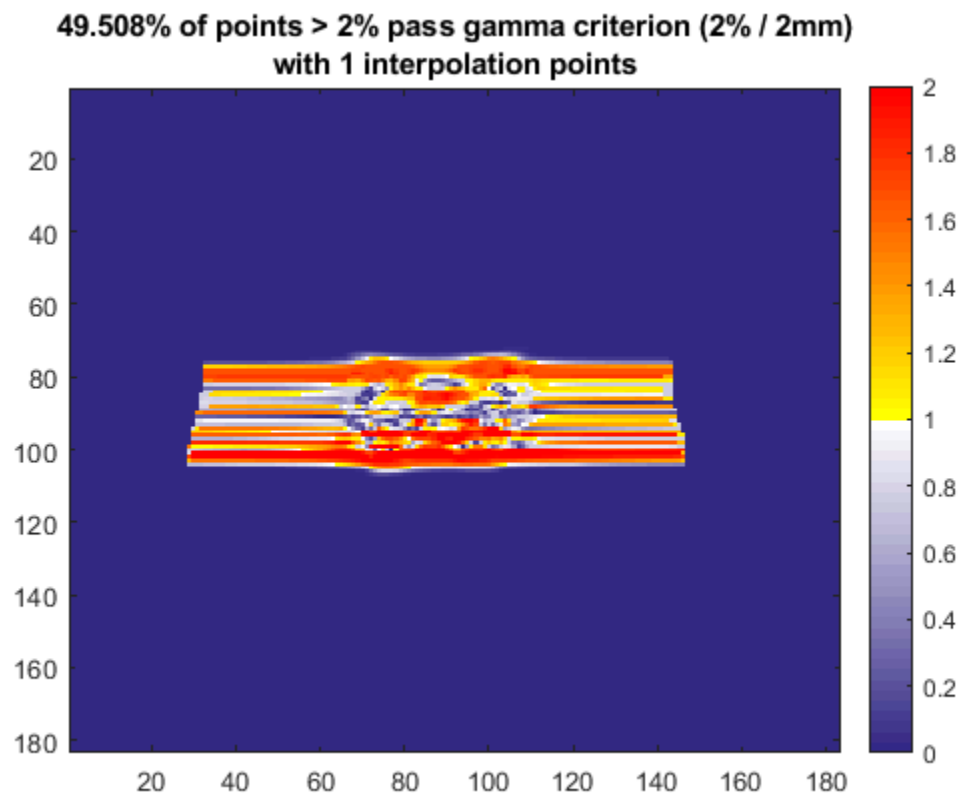


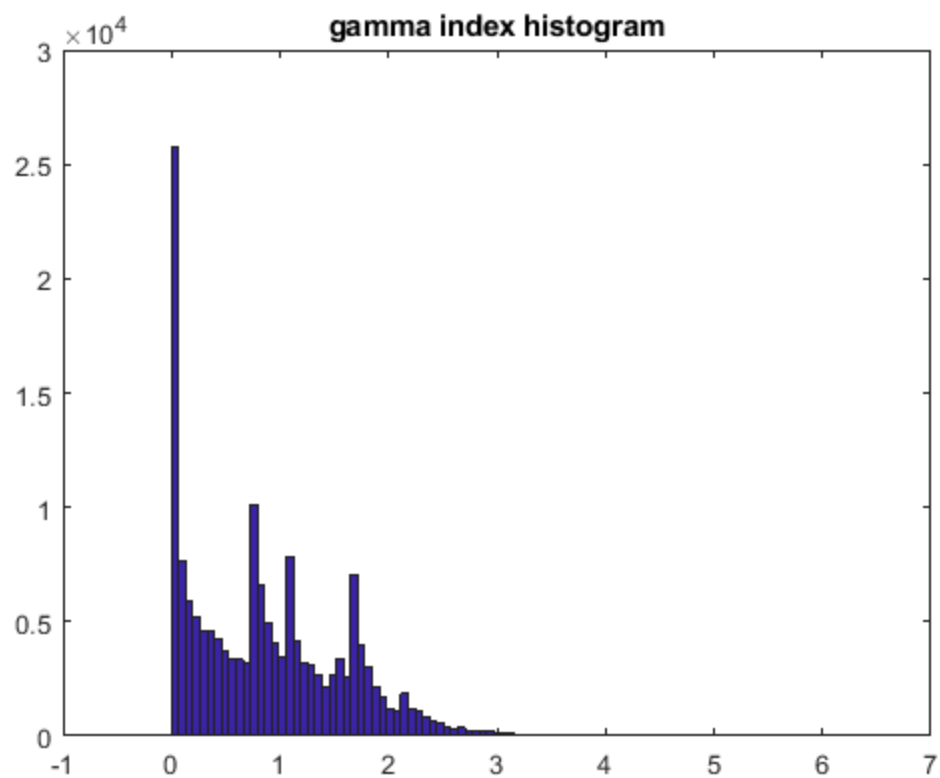


Quantitative Comparison of results

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.

```
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  
figure  
hist(gammaCube(gammaCube>0),100)  
title('gamma index histogram')
```





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