
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	12
Plot the Resulting Beam Dose Slice	13
and the corresponding LET distribution	14
Recalculate Plan	15
Visual Comparison of results	16
Quantitative Comparison of results	19

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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 calculcation

```
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... 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);
```

```
matRad: Using a constant RBE of 1.1
```

```
Warning: Surface for SSD calculation starts directly in first voxel of  
CT
```

```
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);
```

This is Ipopt version 3.12.4, 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.....:      45333
      variables with only lower bounds:      45333
      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  4.3873631e+02  0.00e+00  1.07e+00   0.0  0.00e+00   -  0.00e+00
0.00e+00   0
  1  4.0759581e+02  0.00e+00  7.38e-02  -1.1  7.87e-02   -  9.91e-01
1.00e+00f  1
  2  7.3211108e+01  0.00e+00  2.02e-02  -1.7  1.37e+00   -  9.95e-01
1.00e+00f  1
  3  3.8669378e+01  0.00e+00  1.33e-02  -3.4  3.92e-01   -  9.76e-01
1.00e+00f  1
  4  3.1369069e+01  0.00e+00  1.09e-02  -3.9  2.89e-01   -  9.91e-01
1.00e+00f  1
  5  2.4979899e+01  0.00e+00  1.05e-02  -4.8  4.52e-01   -  9.98e-01
1.00e+00f  1
  6  2.0983319e+01  0.00e+00  1.42e-02  -5.5  7.01e-01   -  1.00e+00
1.00e+00f  1
  7  1.7675867e+01  0.00e+00  7.63e-03  -6.0  2.78e-01   -  1.00e+00
1.00e+00f  1
  8  1.6447624e+01  0.00e+00  6.12e-03  -7.2  2.32e-01   -  1.00e+00
1.00e+00f  1
  9  1.4931143e+01  0.00e+00  5.02e-03  -8.5  4.21e-01   -  1.00e+00
1.00e+00f  1
iter   objective    inf_pr   inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr  ls
 10  1.2970620e+01  0.00e+00  4.15e-03  -9.5  6.30e-01   -  1.00e+00
1.00e+00f  1
 11  1.2308371e+01  0.00e+00  4.93e-03 -10.1  9.11e-01   -  1.00e+00
3.28e-01f  1
 12  1.2304433e+01  0.00e+00  4.92e-03 -11.0  5.39e-01   -  1.00e+00
2.27e-03f  1
```

Example: Proton Treatment Plan
with subsequent Isocenter shift

13	1.2290515e+01	0.00e+00	1.57e-02	-11.0	7.62e-01	-	1.00e+00
5.00e-03f 1							
14	1.1995743e+01	0.00e+00	4.54e-03	-8.5	9.93e-01	-	9.33e-01
8.31e-02f 1							
15	1.1946252e+01	0.00e+00	4.47e-03	-6.5	1.12e+00	-	1.45e-01
1.25e-02f 1							
16	1.1292055e+01	0.00e+00	3.46e-03	-7.7	1.20e+00	-	1.00e+00
1.78e-01f 1							
17	1.1288120e+01	0.00e+00	6.81e-03	-8.6	1.13e+00	-	1.00e+00
1.28e-03f 1							
18	1.1154663e+01	0.00e+00	8.08e-03	-6.5	1.14e+00	-	1.53e-01
4.40e-02f 1							
19	1.0832457e+01	0.00e+00	1.88e-02	-5.3	1.19e+00	-	9.44e-01
1.11e-01f 1							
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
alpha_pr ls							
20	1.0772410e+01	0.00e+00	2.37e-02	-4.5	1.24e+00	-	1.00e+00
2.20e-02f 1							
21	1.0358877e+01	0.00e+00	8.77e-03	-4.8	1.29e+00	-	5.88e-01
1.61e-01f 1							
22	1.0083883e+01	0.00e+00	1.05e-02	-4.4	1.26e+00	-	1.00e+00
1.29e-01f 1							
23	9.7434140e+00	0.00e+00	1.02e-02	-4.4	1.23e+00	-	8.87e-01
1.99e-01f 1							
24	9.3764057e+00	0.00e+00	9.25e-03	-4.1	1.15e+00	-	9.30e-01
3.02e-01f 1							
25	9.2184245e+00	0.00e+00	6.87e-03	-5.3	8.83e-01	-	6.14e-01
2.07e-01f 1							
26	9.0714610e+00	0.00e+00	1.80e-02	-4.5	8.16e-01	-	8.50e-01
2.62e-01f 1							
27	8.9006009e+00	0.00e+00	5.45e-03	-4.1	7.78e-01	-	7.68e-01
3.86e-01f 1							
28	8.8003450e+00	0.00e+00	4.41e-03	-5.5	6.98e-01	-	4.63e-01
2.42e-01f 1							
29	8.7485973e+00	0.00e+00	4.80e-03	-4.2	7.40e-01	-	6.36e-01
1.10e-01f 1							
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
alpha_pr ls							
30	8.5225360e+00	0.00e+00	5.26e-03	-4.0	1.06e+00	-	3.89e-01
3.34e-01f 1							
31	8.3978542e+00	0.00e+00	3.97e-03	-3.9	7.74e-01	-	3.83e-01
2.09e-01f 1							
32	8.2488685e+00	0.00e+00	5.18e-03	-4.1	8.01e-01	-	4.06e-01
2.55e-01f 1							
33	8.1164631e+00	0.00e+00	6.27e-03	-4.1	7.81e-01	-	4.11e-01
2.25e-01f 1							
34	8.0065792e+00	0.00e+00	3.29e-03	-6.2	6.67e-01	-	2.72e-01
2.32e-01f 1							
35	7.9047812e+00	0.00e+00	9.28e-03	-4.6	6.79e-01	-	9.92e-01
2.47e-01f 1							
36	7.7590106e+00	0.00e+00	5.24e-03	-4.2	5.46e-01	-	4.06e-01
6.48e-01f 1							
37	7.6711185e+00	0.00e+00	2.45e-03	-4.7	3.43e-01	-	8.58e-01
8.37e-01f 1							

Example: Proton Treatment Plan
with subsequent Isocenter shift

```

38  7.5829603e+00  0.00e+00  5.90e-03  -4.3  1.21e-01  -  8.64e-01
1.00e+00f  1
39  7.4625863e+00  0.00e+00  1.59e-03  -4.4  2.17e-01  -  4.11e-01
1.00e+00f  1
iter  objective    inf_pr  inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr  ls
40  7.3769763e+00  0.00e+00  9.57e-04  -4.3  1.98e-01  -  8.20e-01
1.00e+00f  1
41  7.3356227e+00  0.00e+00  1.77e-03  -4.7  3.56e-01  -  9.86e-01
3.17e-01f  2
42  7.2741518e+00  0.00e+00  9.50e-04  -5.0  1.23e-01  -  1.00e+00
8.91e-01f  1
43  7.2330398e+00  0.00e+00  2.53e-03  -4.7  1.97e-01  -  1.00e+00
3.32e-01f  1
44  7.1402542e+00  0.00e+00  2.66e-03  -4.5  5.06e-01  -  5.84e-01
4.74e-01f  1
45  7.1103231e+00  0.00e+00  5.33e-03  -4.8  4.63e-01  -  8.85e-01
1.61e-01f  1
46  7.0866714e+00  0.00e+00  5.94e-03  -4.2  3.47e-01  -  9.11e-01
1.45e-01f  1
47  6.9915681e+00  0.00e+00  6.19e-03  -4.7  6.15e-01  -  6.14e-01
4.07e-01f  1
48  9.0032491e+00  0.00e+00  8.47e-03  -2.5  7.04e+00  -  1.05e-01
2.69e-01f  1
49  7.0843202e+00  0.00e+00  5.41e-03  -3.5  1.72e+00  -  8.10e-01
1.00e+00f  1
iter  objective    inf_pr  inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr  ls
50  6.9619232e+00  0.00e+00  5.40e-03  -3.5  2.39e-01  -  1.00e+00
1.00e+00f  1
51  6.9116610e+00  0.00e+00  5.27e-03  -4.3  4.37e-01  -  8.47e-01
3.72e-01f  1
52  6.8143634e+00  0.00e+00  3.23e-03  -4.6  6.70e-01  -  9.99e-01
7.55e-01f  1
53  6.7565957e+00  0.00e+00  6.17e-03  -5.0  5.38e-01  -  9.99e-01
4.45e-01f  1
54  6.7217306e+00  0.00e+00  3.08e-03  -4.5  2.46e-01  -  5.81e-01
5.49e-01f  1
55  7.6316222e+00  0.00e+00  5.94e-03  -2.6  1.07e+01  -  1.82e-02
2.18e-01f  1
56  7.0683086e+00  0.00e+00  7.95e-03  -4.3  2.93e+00  -  1.82e-01
7.50e-01f  1
57  6.8430036e+00  0.00e+00  1.60e-02  -4.3  1.04e+00  -  7.08e-01
2.31e-01f  1
58  6.7124481e+00  0.00e+00  1.24e-02  -4.3  6.09e-01  -  5.36e-01
2.49e-01f  1
59  6.5984020e+00  0.00e+00  1.87e-02  -4.6  6.57e-01  -  1.00e+00
3.19e-01f  1
iter  objective    inf_pr  inf_du lg(mu)  ||d||  lg(rg) alpha_du
alpha_pr  ls
60  6.5470229e+00  0.00e+00  1.12e-02  -4.9  5.53e-01  -  1.00e+00
2.34e-01f  1
61  6.5077251e+00  0.00e+00  1.12e-02  -5.4  5.26e-01  -  9.53e-01
2.33e-01f  1

```

Example: Proton Treatment Plan
with subsequent Isocenter shift

62	6.4672534e+00	0.00e+00	1.09e-02	-6.3	5.52e-01	-	1.00e+00
	2.99e-01f	1					
63	6.4422879e+00	0.00e+00	8.66e-03	-4.7	3.93e-01	-	7.01e-01
	2.90e-01f	1					
64	6.4229276e+00	0.00e+00	1.16e-02	-4.9	4.01e-01	-	7.92e-01
	2.47e-01f	1					
65	6.4026381e+00	0.00e+00	8.15e-03	-6.1	4.79e-01	-	4.37e-01
	2.42e-01f	1					
66	6.3920783e+00	0.00e+00	1.07e-02	-6.1	5.31e-01	-	8.45e-01
	1.16e-01f	1					
67	6.3669945e+00	0.00e+00	1.05e-02	-6.0	7.50e-01	-	8.48e-01
	1.96e-01f	1					
68	6.3458893e+00	0.00e+00	8.79e-03	-5.9	8.28e-01	-	6.61e-01
	1.46e-01f	1					
69	6.3302564e+00	0.00e+00	9.58e-03	-5.9	1.24e+00	-	8.73e-01
	6.83e-02f	1					
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
	alpha_pr	ls					
70	6.2672384e+00	0.00e+00	3.65e-03	-5.5	2.10e+00	-	1.00e+00
	1.58e-01f	1					
71	6.2290540e+00	0.00e+00	5.49e-03	-4.3	9.72e-01	-	3.61e-01
	1.87e-01f	1					
72	1.0157874e+01	0.00e+00	9.07e-03	-2.7	2.03e+01	-	7.57e-03
	3.39e-01f	1					
73	6.4117859e+00	0.00e+00	4.46e-03	-4.4	8.62e+00	-	6.27e-02
	7.49e-01f	1					
74	6.2934081e+00	0.00e+00	2.90e-03	-4.4	1.32e+00	-	7.46e-01
	2.03e-01f	1					
75	6.2579655e+00	0.00e+00	2.20e-02	-4.4	1.12e+00	-	9.56e-01
	9.95e-02f	1					
76	6.2086129e+00	0.00e+00	1.32e-02	-4.4	7.86e-01	-	6.65e-01
	1.85e-01f	1					
77	6.1189136e+00	0.00e+00	1.97e-02	-4.7	8.04e-01	-	9.63e-01
	4.18e-01f	1					
78	6.0968479e+00	0.00e+00	7.97e-03	-4.7	4.73e-01	-	8.50e-01
	2.06e-01f	1					
79	6.0729888e+00	0.00e+00	9.55e-03	-10.7	6.53e-01	-	4.96e-01
	1.90e-01f	1					
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
	alpha_pr	ls					
80	6.0455396e+00	0.00e+00	8.63e-03	-5.7	7.06e-01	-	9.41e-01
	2.30e-01f	1					
81	6.0183843e+00	0.00e+00	7.93e-03	-6.4	7.22e-01	-	9.30e-01
	2.47e-01f	1					
82	6.0038361e+00	0.00e+00	9.42e-03	-6.7	7.73e-01	-	7.93e-01
	1.33e-01f	1					
83	5.9894075e+00	0.00e+00	1.16e-02	-11.0	1.04e+00	-	6.92e-01
	1.00e-01f	1					
84	5.9769568e+00	0.00e+00	5.46e-03	-6.2	1.37e+00	-	9.25e-01
	6.66e-02f	1					
85	5.9416078e+00	0.00e+00	4.13e-03	-7.2	1.63e+00	-	5.26e-01
	1.67e-01f	1					
86	5.9180552e+00	0.00e+00	8.44e-03	-6.6	1.61e+00	-	7.26e-01
	1.16e-01f	1					

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with subsequent Isocenter shift

87	5.9046801e+00	0.00e+00	4.89e-03	-5.2	8.32e-01	-	3.93e-01
	1.31e-01f	1					
88	6.2383777e+00	0.00e+00	8.96e-03	-4.0	5.82e+00	-	2.31e-01
	1.00e+00f	1					
89	6.1256617e+00	0.00e+00	7.69e-03	-4.6	3.73e+00	-	3.93e-01
	2.09e-01f	1					
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
	alpha_pr	ls					
90	6.0929699e+00	0.00e+00	5.68e-03	-4.6	2.17e+00	-	4.21e-01
	7.37e-02f	1					
91	6.0213922e+00	0.00e+00	6.86e-03	-4.6	1.57e+00	-	2.30e-01
	1.84e-01f	1					
92	5.9421673e+00	0.00e+00	1.30e-02	-4.6	1.38e+00	-	5.93e-01
	2.71e-01f	1					
93	5.8985886e+00	0.00e+00	9.54e-03	-4.6	9.04e-01	-	6.42e-01
	2.52e-01f	1					
94	5.8653670e+00	0.00e+00	7.82e-03	-4.9	7.74e-01	-	8.19e-01
	2.58e-01f	1					
95	5.8445779e+00	0.00e+00	9.78e-03	-5.7	6.93e-01	-	5.89e-01
	2.06e-01f	1					
96	5.8269158e+00	0.00e+00	6.59e-03	-6.0	7.15e-01	-	8.88e-01
	1.89e-01f	1					
97	5.8078218e+00	0.00e+00	8.00e-03	-5.7	6.72e-01	-	6.22e-01
	2.36e-01f	1					
98	5.7930219e+00	0.00e+00	8.76e-03	-5.7	6.39e-01	-	7.29e-01
	2.05e-01f	1					
99	5.7768952e+00	0.00e+00	5.72e-03	-5.1	6.86e-01	-	6.84e-01
	2.16e-01f	1					
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
	alpha_pr	ls					
100	5.7561293e+00	0.00e+00	2.72e-03	-4.5	3.04e-01	-	5.18e-01
	6.18e-01f	1					
101	5.7374218e+00	0.00e+00	1.76e-03	-4.4	2.25e-01	-	3.92e-01
	6.50e-01f	1					
102	5.6981688e+00	0.00e+00	1.23e-03	-4.3	6.54e-01	-	4.54e-01
	4.60e-01f	1					
103	5.6777599e+00	0.00e+00	8.11e-03	-4.7	1.07e+00	-	4.64e-01
	1.47e-01f	1					
104	5.6720387e+00	0.00e+00	8.85e-03	-6.6	1.06e+00	-	4.02e-01
	4.33e-02f	1					
105	5.6376567e+00	0.00e+00	5.63e-03	-6.6	1.71e+00	-	3.23e-01
	1.69e-01f	1					
106	5.6292255e+00	0.00e+00	6.52e-03	-7.0	1.59e+00	-	3.14e-01
	4.33e-02f	1					
107	5.5967590e+00	0.00e+00	5.53e-03	-5.2	1.78e+00	-	3.41e-01
	1.59e-01f	1					
108	6.1655582e+00	0.00e+00	6.17e-03	-3.1	2.66e+01	-	3.70e-02
	1.55e-01f	1					
109	5.8513275e+00	0.00e+00	5.83e-03	-4.5	3.65e+00	-	9.30e-03
	3.49e-01f	1					
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
	alpha_pr	ls					
110	5.7363732e+00	0.00e+00	4.97e-03	-4.5	2.09e+00	-	5.53e-01
	2.32e-01f	1					

Example: Proton Treatment Plan
with subsequent Isocenter shift

111	5.6734649e+00	0.00e+00	8.09e-03	-4.5	1.47e+00	-	7.07e-01
	1.89e-01f	1					
112	5.5894769e+00	0.00e+00	1.24e-02	-4.5	1.04e+00	-	6.08e-01
	3.88e-01f	1					
113	5.5502883e+00	0.00e+00	6.74e-03	-4.1	2.41e-01	-	6.40e-01
	1.00e+00f	1					
114	5.5392051e+00	0.00e+00	6.02e-03	-5.0	4.35e-01	-	8.82e-01
	2.24e-01f	1					
115	5.5203424e+00	0.00e+00	1.26e-02	-5.5	7.24e-01	-	1.00e+00
	2.59e-01f	1					
116	5.5022283e+00	0.00e+00	7.64e-03	-6.2	8.39e-01	-	9.98e-01
	2.34e-01f	1					
117	5.4754477e+00	0.00e+00	4.87e-03	-6.0	1.21e+00	-	8.92e-01
	2.64e-01f	1					
118	5.4603671e+00	0.00e+00	4.89e-03	-5.5	9.19e-01	-	4.93e-01
	1.94e-01f	1					
119	5.5439585e+00	0.00e+00	1.08e-02	-4.1	3.14e+00	-	1.75e-01
	1.00e+00f	1					
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
	alpha_pr	ls					
120	5.5117901e+00	0.00e+00	6.41e-03	-4.9	1.57e+00	-	5.81e-01
	2.10e-01f	1					
121	5.4829909e+00	0.00e+00	1.71e-03	-4.9	2.00e+00	-	4.69e-01
	1.35e-01f	1					
122	5.4793397e+00	0.00e+00	9.14e-03	-4.9	1.03e+00	-	4.24e-01
	3.47e-02f	1					
123	5.4450212e+00	0.00e+00	6.94e-03	-4.9	1.49e+00	-	2.49e-01
	2.24e-01f	1					
124	5.4082859e+00	0.00e+00	4.04e-03	-4.6	3.80e+00	-	4.79e-01
	2.88e-01f	1					
125	5.4004431e+00	0.00e+00	5.71e-03	-5.0	4.19e+00	-	4.66e-01
	6.46e-02f	1					
126	5.3814757e+00	0.00e+00	9.82e-03	-5.3	5.10e+00	-	8.08e-01
	1.49e-01f	1					
127	5.7462797e+00	0.00e+00	8.83e-03	-3.1	1.66e+01	-	1.76e-02
	1.71e-01f	1					
128	5.6514869e+00	0.00e+00	8.50e-03	-4.7	4.69e+00	-	2.04e-02
	1.47e-01f	1					
129	5.3772949e+00	0.00e+00	5.93e-03	-4.7	4.27e+00	-	7.47e-01
	6.67e-01f	1					
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
	alpha_pr	ls					
130	5.3578723e+00	0.00e+00	1.13e-02	-4.4	1.57e+00	-	1.00e+00
	3.51e-01f	1					
131	5.6282087e+00	0.00e+00	1.03e-02	-2.4	5.81e+01	-	1.55e-03
	4.18e-02f	1					
132	5.4192280e+00	0.00e+00	2.97e-03	-4.4	4.69e+00	-	5.03e-01
	1.00e+00f	1					
133	5.3279444e+00	0.00e+00	1.39e-03	-4.4	1.36e+00	-	1.00e+00
	1.00e+00f	1					
134	5.3223453e+00	0.00e+00	7.44e-03	-5.2	6.65e-01	-	9.96e-01
	1.24e-01f	1					
135	5.3002813e+00	0.00e+00	5.86e-03	-6.4	1.01e+00	-	1.00e+00
	3.44e-01f	1					

Example: Proton Treatment Plan
with subsequent Isocenter shift

136	5.2797877e+00	0.00e+00	1.80e-03	-4.9	1.42e+00	-	7.73e-01
	4.76e-01f	1					
137	5.2765792e+00	0.00e+00	4.43e-03	-5.2	1.00e+00	-	3.47e-01
	1.04e-01f	1					
138	5.8966043e+00	0.00e+00	3.99e-03	-3.1	3.22e+01	-	4.07e-03
	1.82e-01f	1					
139	5.3153153e+00	0.00e+00	3.86e-03	-4.9	1.08e+01	-	2.95e-02
	6.93e-01f	1					
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
	alpha_pr	ls					
140	5.2811037e+00	0.00e+00	3.24e-03	-4.9	7.27e+00	-	5.02e-01
	5.29e-02f	2					
141	5.2741099e+00	0.00e+00	4.09e-03	-4.9	1.39e+00	-	7.07e-01
	8.37e-02f	1					
142	5.2518247e+00	0.00e+00	3.91e-03	-4.9	1.41e+00	-	9.98e-01
	2.79e-01f	1					
143	5.2422959e+00	0.00e+00	1.18e-02	-10.9	1.46e+00	-	7.03e-01
	1.43e-01f	1					
144	5.2396163e+00	0.00e+00	1.53e-02	-7.5	1.45e+00	-	7.97e-01
	4.51e-02f	1					
145	5.2161771e+00	0.00e+00	6.44e-03	-5.4	1.72e+00	-	5.50e-01
	3.86e-01f	1					
146	5.2070040e+00	0.00e+00	4.22e-03	-4.9	9.63e-01	-	6.17e-01
	3.31e-01f	1					
147	5.1971643e+00	0.00e+00	2.13e-03	-4.7	4.63e-01	-	3.97e-01
	1.00e+00f	1					
148	5.1939683e+00	0.00e+00	4.65e-03	-5.4	1.10e+00	-	5.88e-01
	1.32e-01f	1					
149	5.1872551e+00	0.00e+00	4.49e-03	-5.4	1.75e+00	-	9.97e-01
	1.69e-01f	1					
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
	alpha_pr	ls					
150	5.1764179e+00	0.00e+00	3.39e-03	-5.0	1.65e+00	-	6.78e-01
	2.86e-01f	1					
151	5.1608540e+00	0.00e+00	1.86e-03	-4.7	1.33e+00	-	3.21e-01
	5.08e-01f	1					
152	5.1517156e+00	0.00e+00	2.06e-03	-4.7	1.40e+00	-	2.74e-01
	2.48e-01f	1					
153	5.1168228e+00	0.00e+00	9.34e-04	-4.5	2.41e+00	-	3.86e-01
	5.96e-01f	1					
154	5.1107575e+00	0.00e+00	9.51e-03	-5.0	2.04e+00	-	6.42e-01
	1.10e-01f	1					
155	5.0984779e+00	0.00e+00	7.45e-03	-5.4	2.55e+00	-	8.22e-01
	1.98e-01f	1					
156	5.0943759e+00	0.00e+00	6.09e-03	-5.5	2.24e+00	-	4.97e-01
	7.61e-02f	1					
157	5.0792497e+00	0.00e+00	4.36e-03	-5.8	3.13e+00	-	7.71e-01
	2.13e-01f	1					
158	5.0732716e+00	0.00e+00	3.79e-03	-5.3	2.40e+00	-	2.99e-01
	1.13e-01f	1					
159	5.0651329e+00	0.00e+00	3.85e-03	-5.1	2.86e+00	-	6.67e-01
	1.31e-01f	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

160	5.0537938e+00	0.00e+00	5.00e-03	-5.3	2.82e+00	-	3.06e-01
	1.87e-01f	1					
161	5.0401234e+00	0.00e+00	3.72e-03	-11.0	3.73e+00	-	2.35e-01
	1.84e-01f	1					
162	5.0327249e+00	0.00e+00	4.53e-03	-5.9	3.63e+00	-	6.85e-01
	1.01e-01f	1					
163	5.0219279e+00	0.00e+00	3.07e-03	-6.1	3.28e+00	-	2.35e-01
	1.64e-01f	1					
164	5.1861227e+00	0.00e+00	3.19e-03	-3.7	2.17e+01	-	2.25e-02
	2.04e-01f	1					
165	5.1157534e+00	0.00e+00	3.45e-03	-5.2	7.92e+00	-	1.50e-02
	2.26e-01f	1					
166	5.1003448e+00	0.00e+00	5.27e-03	-5.2	6.20e+00	-	6.96e-01
	6.69e-02f	1					
167	5.0580055e+00	0.00e+00	7.72e-03	-5.2	5.29e+00	-	4.88e-01
	2.19e-01f	1					
168	5.0249617e+00	0.00e+00	6.12e-03	-5.2	3.94e+00	-	6.15e-01
	2.55e-01f	1					
169	5.0140981e+00	0.00e+00	3.45e-03	-5.2	2.49e+00	-	5.93e-01
	1.37e-01f	1					
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
	alpha_pr	ls					
170	4.9978747e+00	0.00e+00	8.21e-03	-5.3	2.07e+00	-	7.10e-01
	2.57e-01f	1					
171	5.1680528e+00	0.00e+00	8.31e-03	-3.4	2.12e+01	-	2.22e-02
	2.90e-01f	1					
172	4.9864678e+00	0.00e+00	7.83e-03	-4.9	7.80e+00	-	4.28e-01
	9.09e-01f	1					
173	4.9816180e+00	0.00e+00	8.17e-03	-5.6	1.82e+00	-	1.00e+00
	7.42e-02f	1					
174	4.9740463e+00	0.00e+00	8.97e-03	-7.0	1.53e+00	-	8.58e-01
	1.46e-01f	1					
175	4.9642860e+00	0.00e+00	6.45e-03	-6.3	1.52e+00	-	8.14e-01
	2.05e-01f	1					
176	4.9558072e+00	0.00e+00	1.13e-02	-5.9	1.35e+00	-	7.58e-01
	2.22e-01f	1					
177	4.9481910e+00	0.00e+00	9.49e-03	-6.0	1.47e+00	-	9.87e-01
	2.15e-01f	1					
178	4.9406065e+00	0.00e+00	5.13e-03	-5.6	1.29e+00	-	8.05e-01
	2.77e-01f	1					
179	4.9350252e+00	0.00e+00	2.65e-03	-5.0	4.71e-01	-	3.17e-01
	5.73e-01f	1					
iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du
	alpha_pr	ls					
180	4.9318309e+00	0.00e+00	1.91e-03	-5.0	2.36e-01	-	4.72e-01
	7.12e-01f	1					
181	4.9290659e+00	0.00e+00	5.47e-03	-5.5	7.47e-01	-	5.58e-01
	2.18e-01f	1					
182	4.9273414e+00	0.00e+00	4.15e-03	-5.6	1.24e+00	-	7.83e-01
	7.91e-02f	1					

Number of Iterations.....: 182

(scaled)

(unscaled)

Example: Proton Treatment Plan
with subsequent Isocenter shift

Objective.....: 4.9273413692900743e+00
4.9273413692900743e+00
Dual infeasibility.....: 4.1477500774598679e-03
4.1477500774598679e-03
Constraint violation....: 0.0000000000000000e+00
0.0000000000000000e+00
Complementarity.....: 1.4640735011706293e-05
1.4640735011706293e-05
Overall NLP error.....: 4.1477500774598679e-03
4.1477500774598679e-03

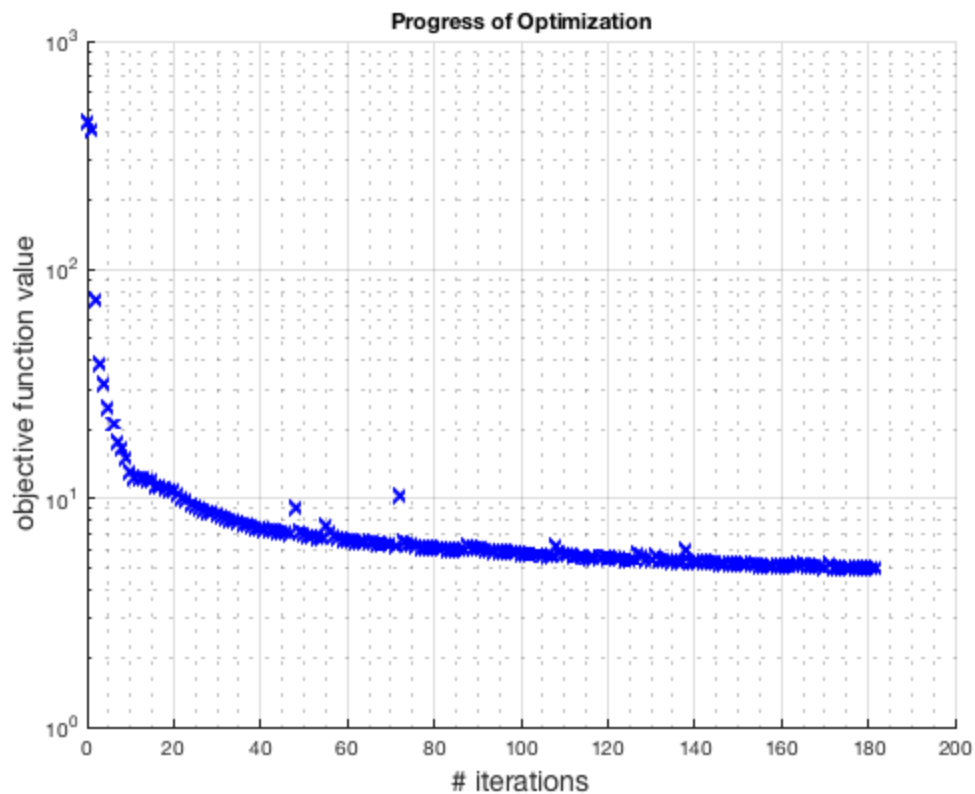
Number of objective function evaluations = 193
Number of objective gradient evaluations = 183
Number of equality constraint evaluations = 0
Number of inequality constraint evaluations = 0
Number of equality constraint Jacobian evaluations = 0
Number of inequality constraint Jacobian evaluations = 0
Number of Lagrangian Hessian evaluations = 0
Total CPU secs in IPOPT (w/o function evaluations) = 11.812
Total CPU secs in NLP function evaluations = 107.475

EXIT: Solved To Acceptable Level.

*** IPOPT DONE ***

Calculating final cubes...

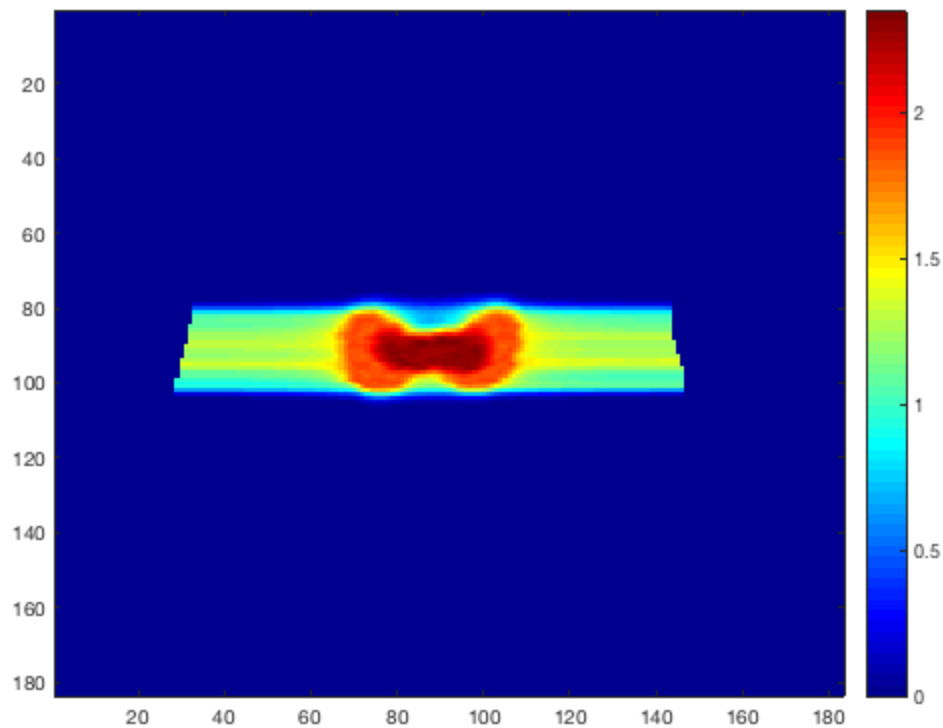
matRad: applying a constant RBE of 1.1



Plot the Resulting Dose Slice

Let's plot the transversal iso-center dose slice

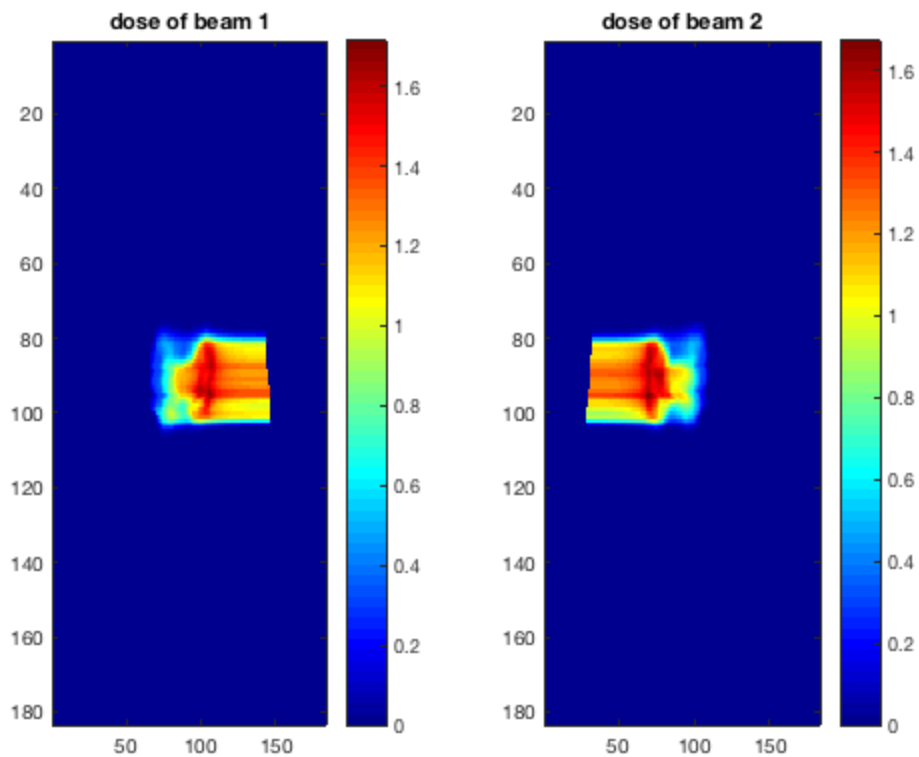
```
slice = round(pln.propStf.isoCenter(1,3)./ct.resolution.z);  
figure  
imagesc(resultGUI.RBExDose(:, :, slice)), colorbar, colormap(jet)
```



Plot the Resulting Beam Dose Slice

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

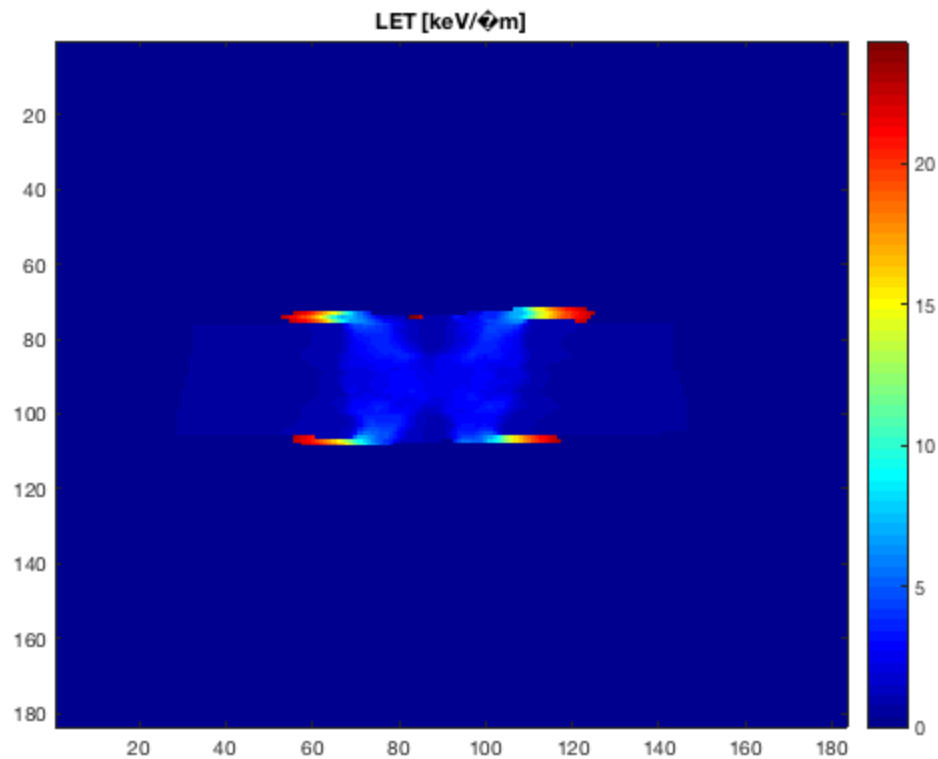
```
figure
subplot(121),imagesc(resultGUI.RBExDose_beam1(:,:,slice)),colorbar,colormap(jet),t
    'of beam 1')
subplot(122),imagesc(resultGUI.RBExDose_beam2(:,:,slice)),colorbar,colormap(jet),t
    'of beam 2')
```



and the corresponding LET distribution

Transversal iso-center slice

```
figure
imagesc(resultGUI.LET(:,:,slice)),colormap(jet),colorbar,title('LET
[keV/#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);  
  
matRad: Using a constant RBE of 1.1  
Warning: Surface for SSD calculation starts directly in first voxel of  
        CT  
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(:)])];

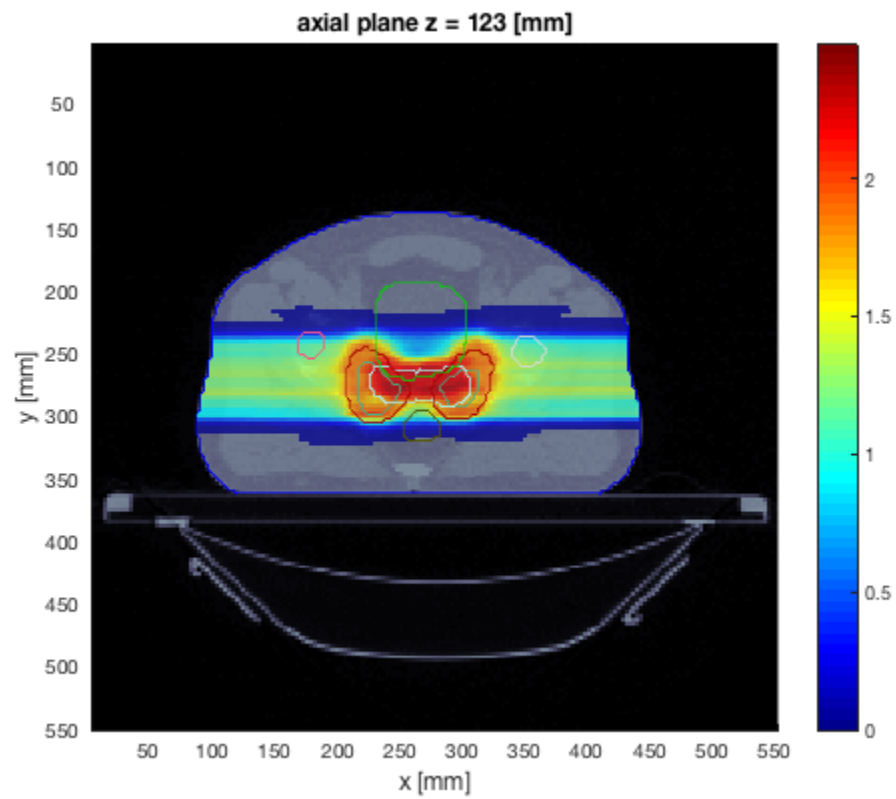
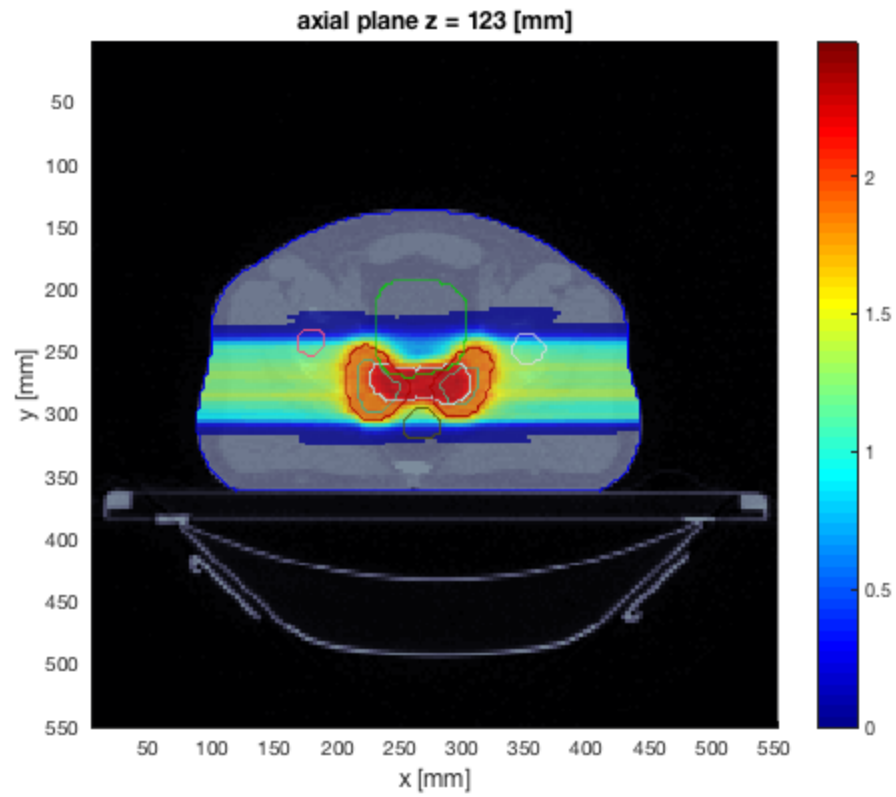
figure,title('original plan')
matRad_plotSliceWrapper(gca,ct,cst,1,resultGUI.RBExDose,plane,slice,
    [],0.75,colorcube,[],doseWindow,[]);
figure,title('shifted plan')
matRad_plotSliceWrapper(gca,ct,cst,1,resultGUI_isoShift.RBExDose,plane,slice,
    [],0.75,colorcube,[],doseWindow,[]);

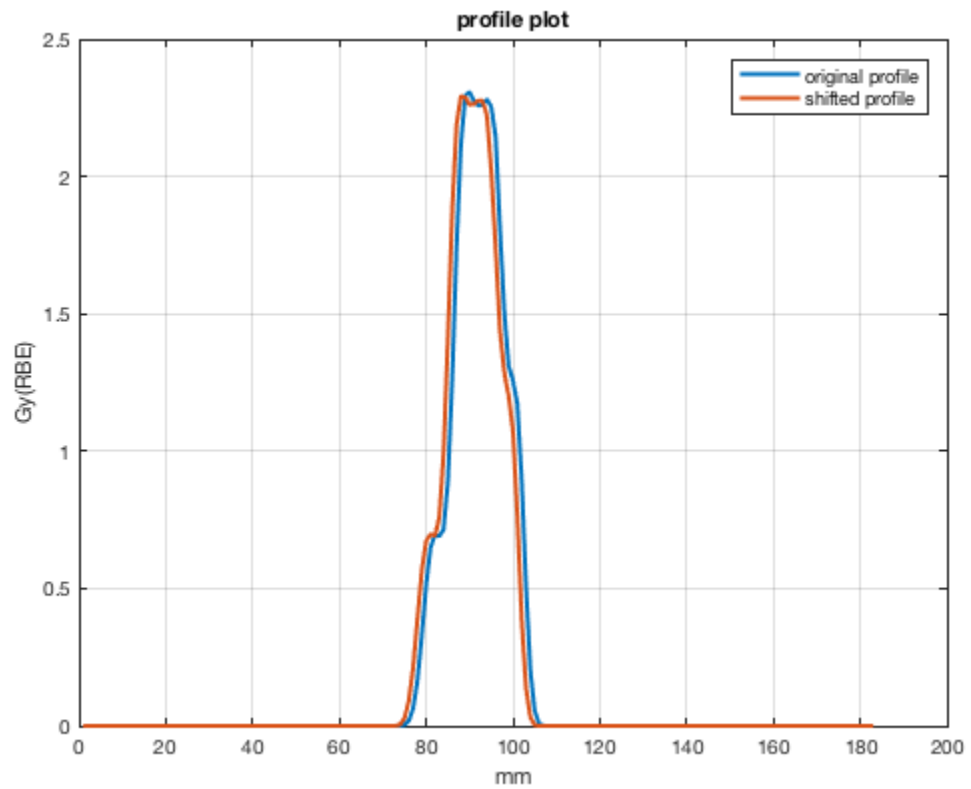
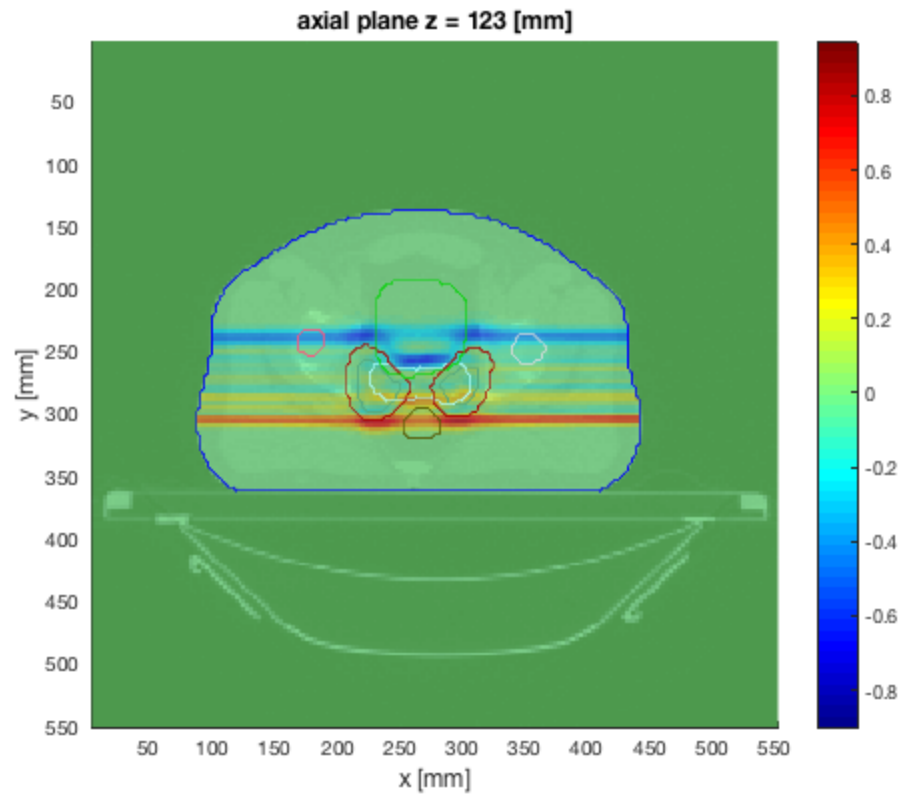
absDiffCube = resultGUI.RBExDose-resultGUI_isoShift.RBExDose;
figure,title('absolute difference')
matRad_plotSliceWrapper(gca,ct,cst,1,absDiffCube,plane,slice,[],
    [],colorcube);

% 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

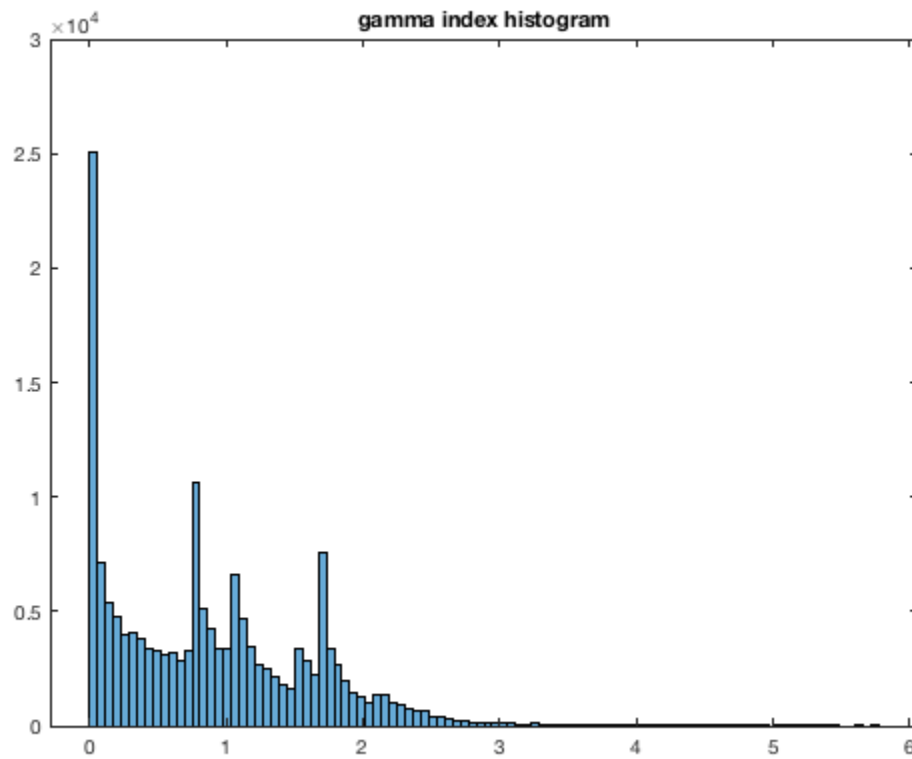
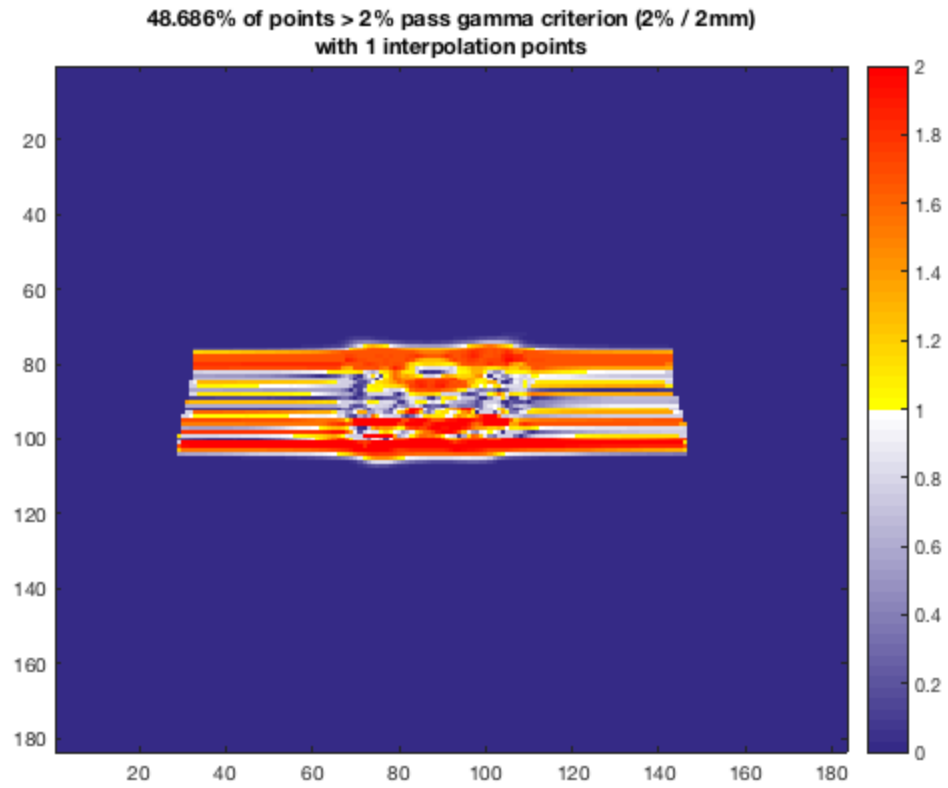
Compare the two dose cubes using a gamma-index analysis. The gamma index is a composite quality distribution equally taking into account a dose difference and a distance to agreement criterion in order to quantify differences between two dose cubes. A gamma-index value of smaller than 1 indicates a successful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



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