

Lecture 9

Lattice design with MAD-X

Dr. Suzie Sheehy
suzie.sheehy@physics.ox.ac.uk
@suziesheehy

Royal Society University Research Fellow

Lecture 9 – MADX introduction and examples

- This lecture is based on one by Ted Wilson (CERN) which was in turn based on a lecture by V. Ziemann (Uppsala University)
 - Installing/running MADX (on mac/windows)
 - Input of elements and beamlines
 - Beta functions, tunes, dispersion
 - Matching
 - Examples

All credit to V. Ziemann for example input files
More examples are on the MAD-X website!

What is MAD-X?

“A program for accelerator design and simulation with a long history”
 Developed from previous versions (MAD, MAD-8, then finally MAD-X in 2002).
 User guide: <http://mad.web.cern.ch/mad/uguide.html>

- Uses a sequence of elements placed sequentially along a *reference orbit*
- *Reference orbit* is path of a charged particle having the central design momentum of the accelerator through idealised magnets (no fringe fields)
- The reference orbit consists of a series of straight line segments and circular arcs
- local curvilinear right handed coordinate system (x, y, s)

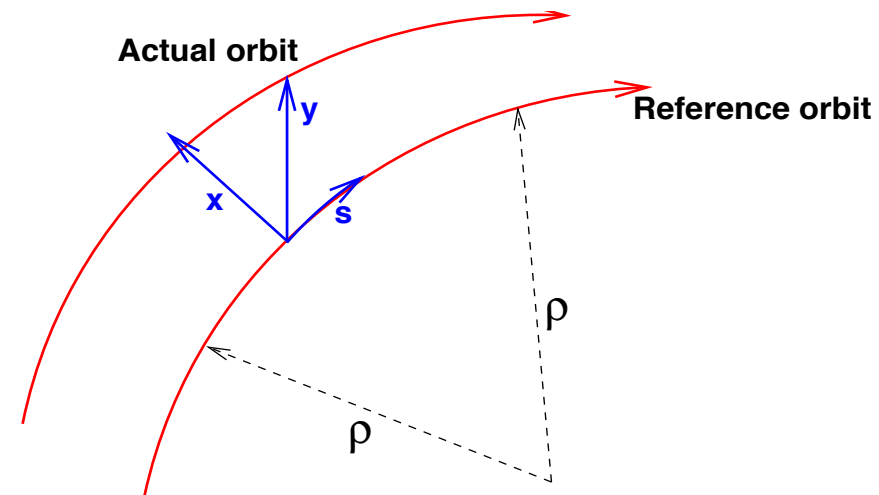


Figure 1: *Local coordinate system as used by MAD-X.*

What can it do?

- You can input magnets (& electrostatic elements) according to the manual
 - User guide: <http://madx.web.cern.ch/madx/>
- Calculate beta functions, tune, dispersion, chromaticity, momentum compaction numerically.
- Generates tables and plots (.ps)
- (tip: you might need to install ghostscript to view plots)

Why choose MAD-X?

- There are any number of tracking and beam optics codes, but MAD-X is *widely used* (especially at CERN), *well maintained* and *well documented*.
 - The more lattice design you do – the more you will appreciate this about MAD-X!
- What it can't do:
 - Acceleration and tracking simultaneously
 - Not so accurate at large excursions from closed orbit (as in an FFAG)
 - Complicated magnet geometries
 - Field maps

Installing MADX

<http://madx.web.cern.ch/madx/>

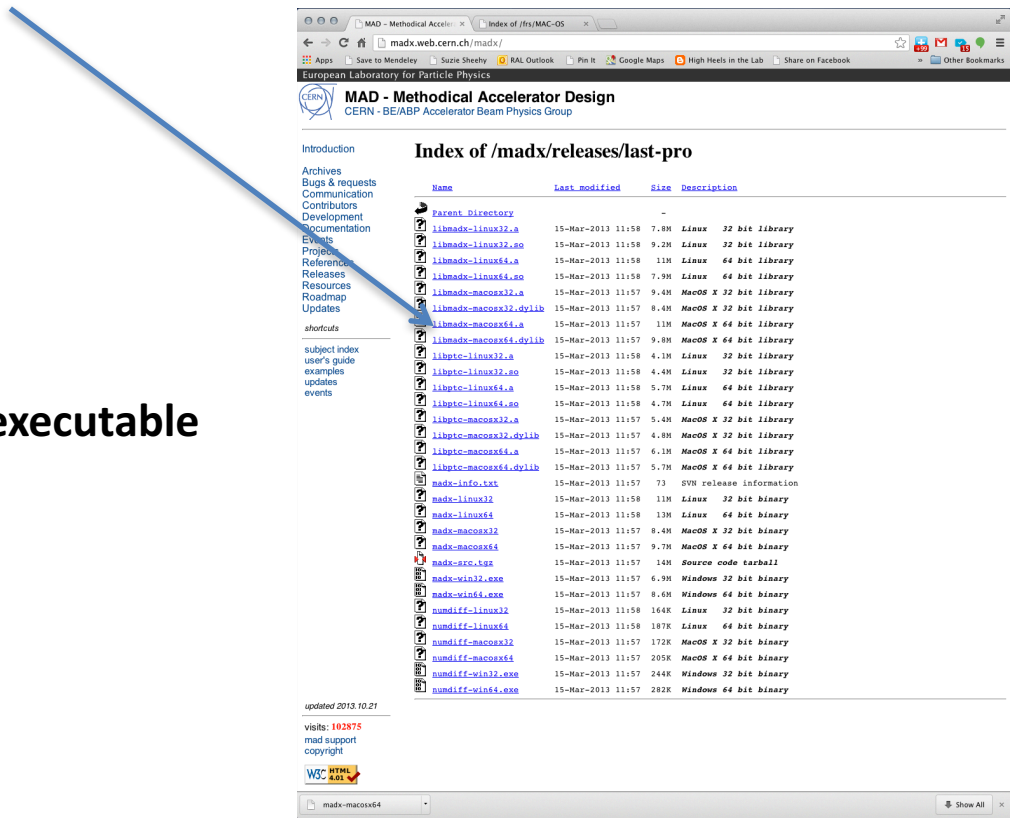
Go to 'Releases'

Get latest production version

Choose your system & download:

NOTE: You may need to make it executable

```
mv madx-macosx64 madx
chmod u+x madx
./madx
```



How to run MADX

- In command prompt:
- Go to directory (with madx.exe and input files)
>> `madx.exe < inputfile > outputfile`
- Or on Mac OSX in terminal:
>> `./madx < inputfile < outputfile`
- Or you can add the madx.exe location to your path (if you know how...)

An input file – Basic FODO

```

//                                     // comments out a line
// MADX Example 1: FODO cell
// Author: V. Ziemann, Uppsala University
// Date: 060910
// UPDATED SUZIE SHEEHY 04/11/2013 FOR MADX VERSION 5.01.00
TITLE, 'Example 1: FODO.MADX';
BEAM, PARTICLE=ELECTRON, PC=3.0;
D: DRIFT, L=1.0;
QF: QUADRUPOLE, L=0.5, K1=0.2;
QD: QUADRUPOLE, L=0.5, K1=-0.2;
FODO: LINE=(QF, 5*(D), QD, QD, 5*(D), QF);
SETPLOT, POST=2, FONT=-1;
USE, PERIOD=FODO;
TWISS, SAVE, BETX=15.0, BETY=5.0;
PLOT, HAXIS=S, VAXIS=BETX, BETY, NOVERSION=TRUE, TITLE='unmatched beta functions';
//Here MATCH is used as a single command this finds periodic solution
USE, PERIOD=FODO;
MATCH, SEQUENCE=FODO;
TWISS, SAVE;
PLOT, HAXIS=S, VAXIS=BETX, BETY, NOVERSION=TRUE, TITLE='matched beta functions';
Value, TABLE(SUMM, Q1);
Value, TABLE(SUMM, Q2);
WRITE, TABLE=SUMM, FILE=print.dat;

```

← TITLE at top of output

← Define particle type and momentum (pc) GeV/c
Or can use ENERGY in GeV.

← Elements are given in the manual including definitions of L, K1 etc...

← Define a 'line' –can be a cell or a whole beamline that you will USE.

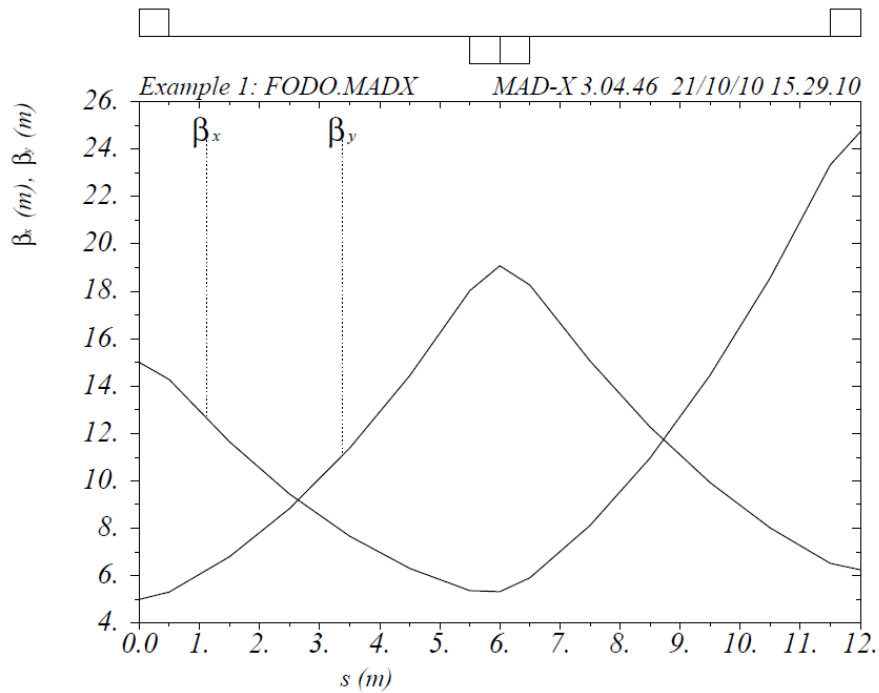
← Calculate beta functions from starting values

← Plot beta values from TWISS (internal table)

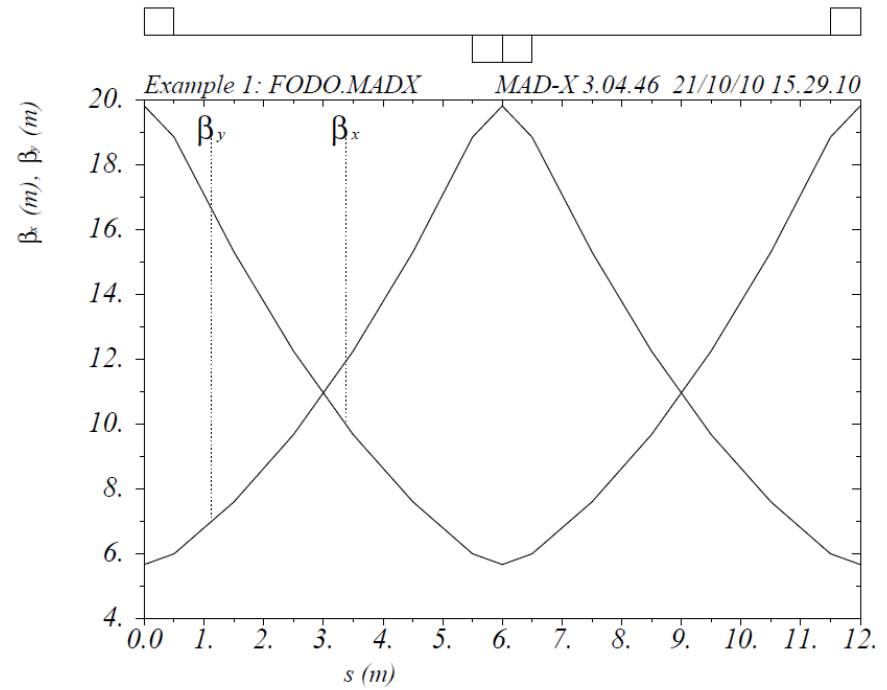
← Match the periodic solution (+Plot that)

← Output to tables

Result of a MADX run



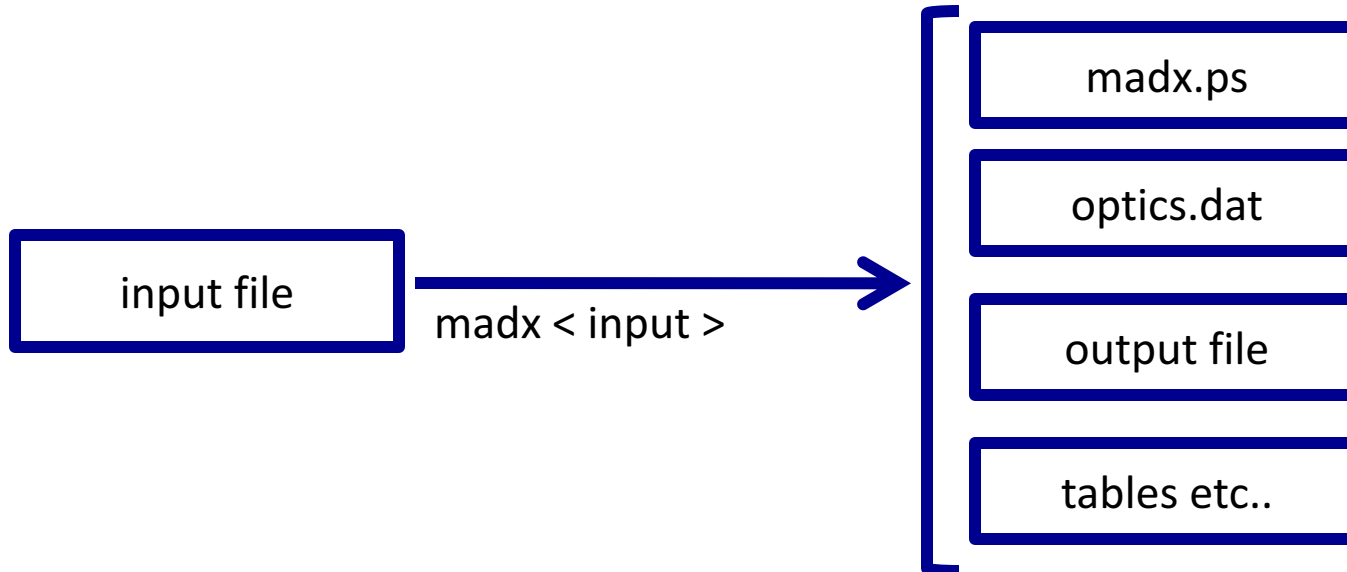
With β_x starting at 15m (β_y at 5m)



If the line is matched (periodic)
 $\beta_{\text{start}} = \beta_{\text{end}}$

Output files

- Optics.dat
- Your specified output file `madx< input > output`
- Can specify tables



Add bending magnets

- Can introduce your own parameters (watch the :=)
- Can use alternative 'sequence' format
- Let's add dipoles
- Look at dispersion

```

TITLE,'Example 2: FODO2.MADX';
BEAM, PARTICLE=ELECTRON,PC=3.0;
DEGREE:=PI/180.0;           // for readability

QF: QUADRUPOLE,L=0.5,K1=0.2;  // still half-length
QD: QUADRUPOLE,L=1.0,K1=-0.2; // changed to full length
B: SBEND,L=1.0,ANGLE=15.0*DEGREE; // added dipole

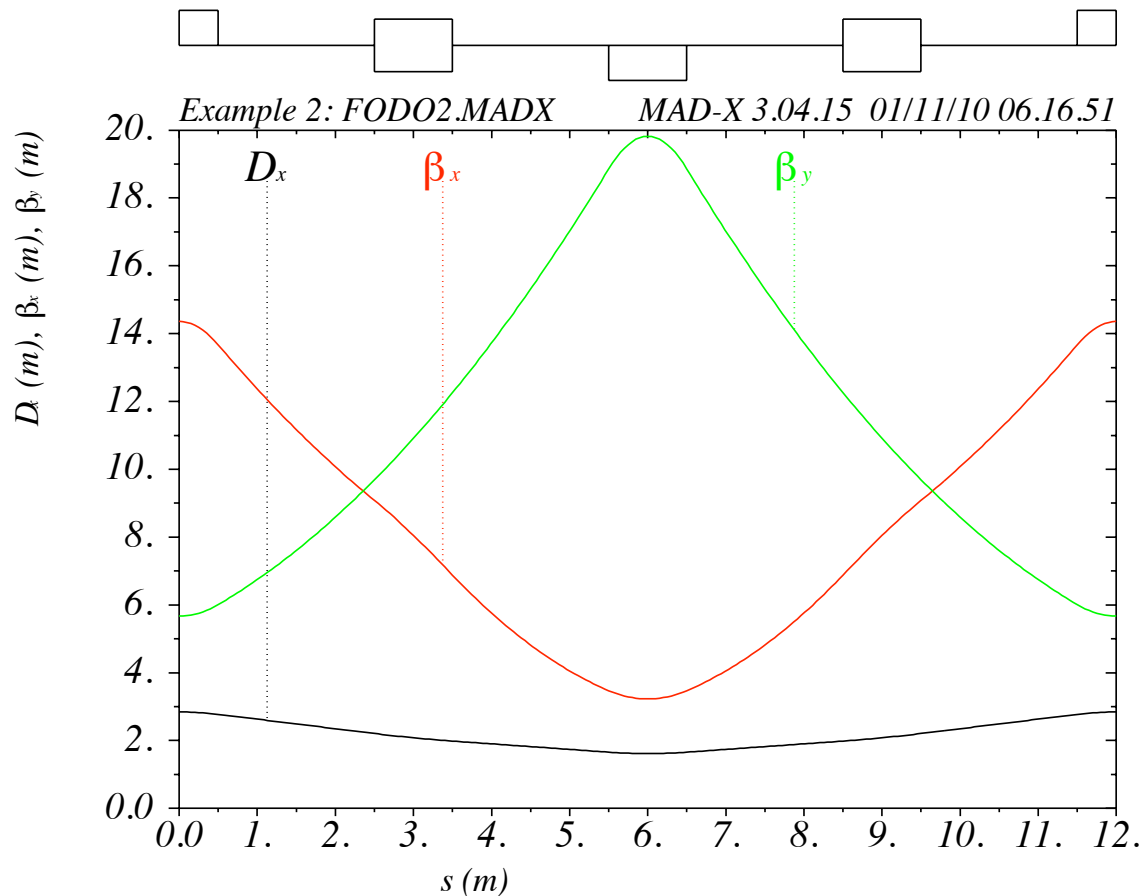
FODO: SEQUENCE,REFER=ENTRY,L=12.0;
  QF1: QF,  AT=0.0;
  B1:  B,  AT=2.5;
  QD1: QD,  AT=5.5;
  B2:  B,  AT=8.5;
  QF2: QF,  AT=11.5;
ENDSEQUENCE;

USE, PERIOD=FODO;
//MATCH, SEQUENCE=FODO; //Uncomment to match
SELECT,FLAG=SECTORMAP,clear;
SELECT,FLAG=TWISS,column=name,s,betx,bety;
TWISS, file=optics.dat,sectormap;

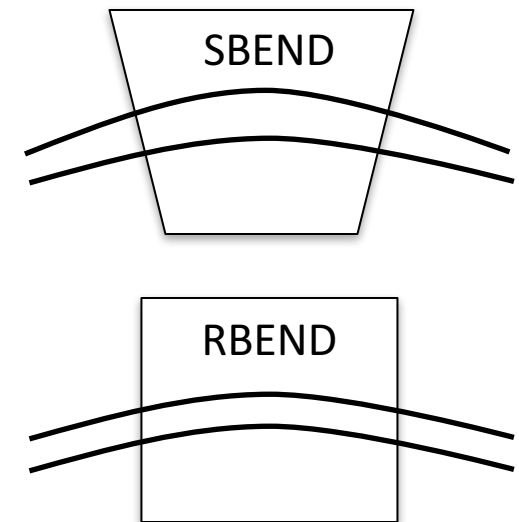
PLOT,HAXIS=S, COLOUR=100, VAXIS=DX, BETX, BETY,
INTERPOLATE=TRUE;

Value, TABLE(SUMM,Q1);
Value, TABLE(SUMM,Q2);
  
```

Output from FODO with dipole



Note the colours!
We now have dispersion
(from bending magnet)
Focusing changed as we
used SBEND



- If you add
SELECT, FLAG=SECTOR
SELECT, FLAG=TWISS,
TWISS, file=optics
- You will get
matrices
- and an out

You can customise
select, flag=my_sec
Or even select by comp
select, flag=my_sec

```

optics.dat
@ NAME           %05s "TWISS"
@ TYPE           %05s "TWISS"
@ SEQUENCE       %04s "FODO"
@ PARTICLE       %08s "ELECTRON"
@ MASS           %le      0.000510998902
@ CHARGE         %le      -1
@ ENERGY        %le      3.000000044
@ PC             %le      3
@ GAMMA          %le      5870.854187
@ KBUNCH         %le      1
@ BCURRENT       %le      0
@ SIGE           %le      0
@ SIGT           %le      0
@ NPART          %le      0
@ EX             %le      1
@ EY             %le      1
@ ET             %le      1
@ LENGTH         %le      12
@ ALFA           %le      0.0910331259
@ ORBITS         %le      -0
@ GAMMATR        %le      3.314364527
@ Q1             %le      0.2909501025
@ Q2             %le      0.1913459932
@ DQ1            %le      -0.2152485772
@ DQ2            %le      -0.2276020063
@ DXMAX          %le      2.84272104
@ DYNAMX         %le      0
@ XCOMAX         %le      0
@ YCOMAX         %le      0
@ BETXMAX        %le      14.36014477
@ BETYMAX        %le      18.85637615
@ XCORMS         %le      0
@ YCORMS         %le      0
@ DXRMS          %le      2.388159047
@ DYRMS          %le      0
@ DELTAP         %le      0
@ SYNCH_1        %le      0
@ SYNCH_2        %le      0
@ SYNCH_3        %le      0
@ SYNCH_4        %le      0
@ SYNCH_5        %le      0
@ TITLE          %21s "Example 2: FODO2.MADX"
@ ORIGIN         %20s "MAD-X 3.04.15 Darwin"
@ DATE           %08s "01/11/10"
@ TIME           %08s "08.52.11"
* NAME           S           BETX           BETY
$ %s             %le      %le      %le
"FODO$START"    0           14.36014477    5.674619595
"QF1"           0.5         13.67114589    6.007906055
"DRIFT_0"       2.5         9.081076899    9.690010744
"B1"            3.5         6.880079877    12.25970476
"DRIFT_0"       5.5         3.471640314    18.85637615
"QD1"           6.5         3.471640314    18.85637615
"DRIFT_0"       8.5         6.880079877    12.25970476
"B2"            9.5         9.081076899    9.690010744

```

Matching

- Matching lets MAD-X do the tedious work for you!
- Before MATCH select at least one sequence (USE)
- Initiated by the MATCH command
- Initiating:
 - MATCH, SEQUENCE='name1', 'name2', ..., 'nameX';
- Can define constraints & variables (magnets) to achieve aim

```
MATCH, SEQUENCE = FODO;
```

```
CONSTRAINT,SEQUENCE=FODO, RANGE=#E, MUX=0.1666666, MUY=0.25;
```

```
VARY, NAME=QF->K1, STEP=1E-6;
```

```
VARY, NAME=QD->K1, STEP=1E-6;
```

```
LMDIF,CALLS=500,TOLERANCE=1E-20;
```

```
ENDMATCH;
```

Matching input file

```
TITLE, 'Example 3: MATCH1.MADX';  
BEAM, PARTICLE=ELECTRON, PC=3.0;
```

```
D: DRIFT, L=1.0;  
QF: QUADRUPOLE, L=0.5, K1:=0.2;  
QD: QUADRUPOLE, L=0.5, K1:=-0.2;
```

```
FODO: LINE=(QF, 5*(D), QD, QD, 5*(D), QF);  
USE, PERIOD=FODO;
```

```
//....match phase advance at end of cell to 60 and 90 degrees  
MATCH, SEQUENCE=FODO;  
CONSTRAINT, SEQUENCE=FODO, RANGE=#E, MUX=0.16666666, MUY=0.25;  
VARY, NAME=QF->K1, STEP=1E-6;  
VARY, NAME=QD->K1, STEP=1E-6;  
LMDIF, CALLS=500, TOLERANCE=1E-20;  
ENDMATCH;
```

```
SELECT, FLAG=SECTORMAP, clear;  
SELECT, FLAG=TWISS, column=name, s, betx, alfx, bety, alfy;  
TWISS, file=optics.dat, sectormap;
```

```
PLOT, HAXIS=S, VAXIS=BETX, BETY;  
Value, TABLE(SUMM, Q1); // verify result  
Value, TABLE(SUMM, Q2);
```

Matching commands

Print out final values of matching

Matching example

- Demonstration MATCH1.MADX

Fitting beta functions

- Use MATCH2.MADX

```

Initial Penalty Function = 0.87329908E+02

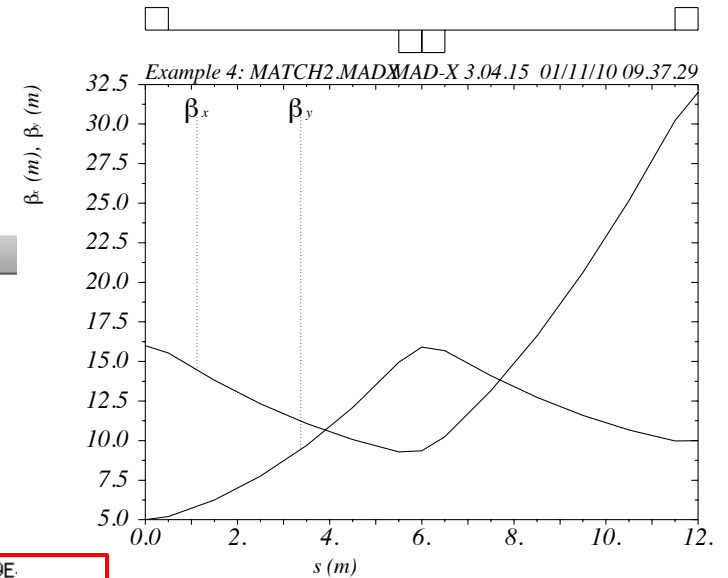
call:      5  Penalty function = 0.14687632E+02
call:      8  Penalty function = 0.52122712E+00
call:     11  Penalty function = 0.14867643E-03
call:     14  Penalty function = 0.17048760E-10
call:     17  Penalty function = 0.86820304E-19
call:     20  Penalty function = 0.66895405E-27
+++++++ LMDIF ended: converged successfully
call:     20  Penalty function = 0.66895405E-27
fodo$end:1      betx      4      3.20000000E+01      3.20000000E+01      5.04870979E-
fodo$end:1      bety      4      1.00000000E+01      1.00000000E+01      6.18466950E-20
qf->k1          1.21494427E-01      -1.00000000E+20      1.00000000E+20
qd->k1          -1.58047975E-01      -1.00000000E+20      1.00000000E+20

GXPLOT-X11 1.50 initialized

plot number = 1

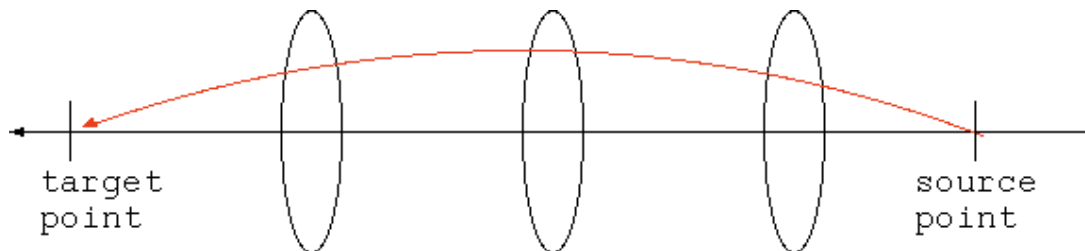
+++++++
+          MAD-X 3.04.15          +
+ Code Modification Date: 15.01.2008      +
+ Execution Time Stamp: 01.11.10 09.39.37 +
+++++++
//
// MADX Example 1: FODO matching final beta function

```



Transfer matrix matching

- Sometimes want to constrain transfer matrix elements to some value.
- For example $R_{16}=0$ and $R_{26}=0$ will make the horizontal position and angle independent of the momentum after a beamline.
- This is called an 'Achromat'.
- Other versions are imaginable
- point-to-point imaging $\rightarrow R_{12} = 0$.
 - This means $\sin(\mu)=0$ or a phase advance of a multiple of π .



Examples in MAD-X

- FODO arcs
- Dispersion suppressor
- ‘Telescopes’ for low- β
- Synchrotron radiation lattices + achromats

Is that it?

- ‘the not-so-ideal world’
- What happens to α, β, γ if we stop focusing for a distance?

$$\beta(s) = \beta_0 - 2\alpha_0 s + \gamma_0 s^2$$

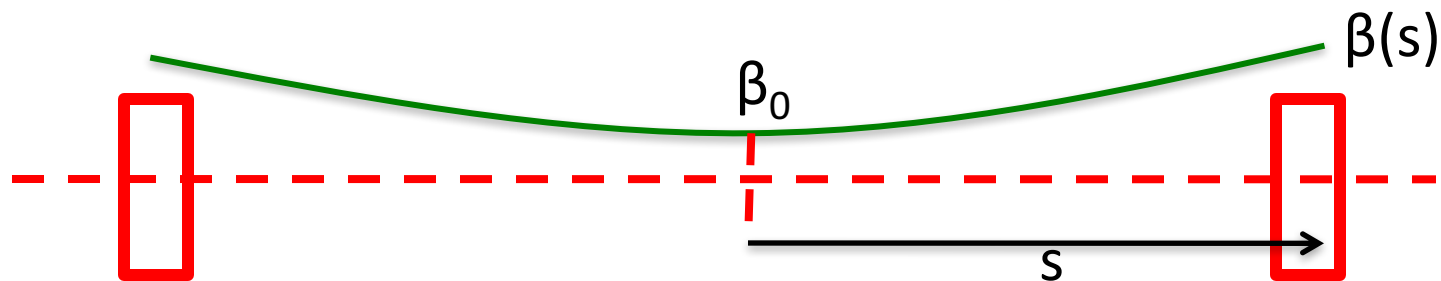
$$\alpha(s) = \alpha_0 - \gamma_0 s$$

$$\gamma(s) = \gamma_0$$

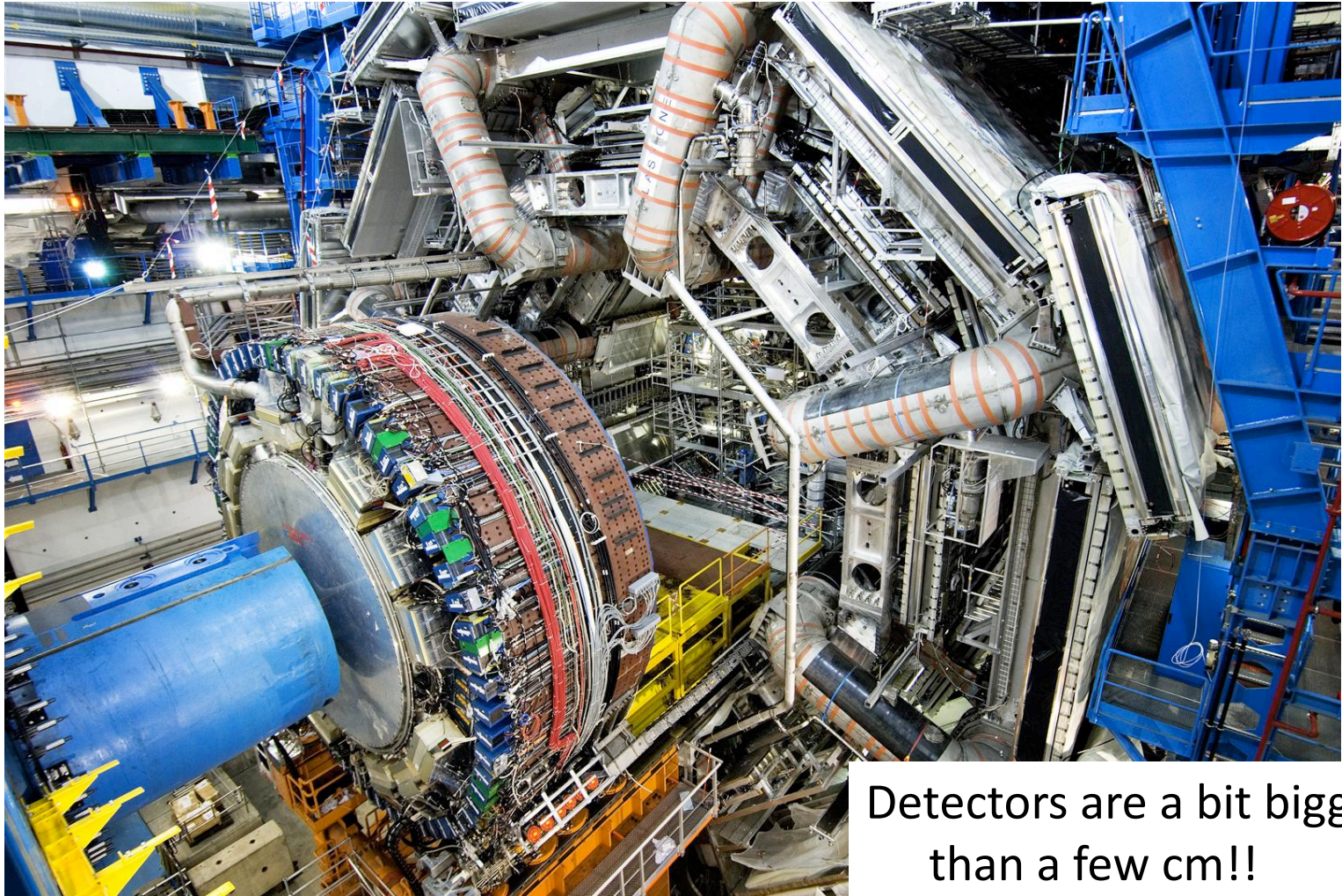
The drift length:

$$M = \begin{bmatrix} 1 & s \\ 0 & 1 \end{bmatrix}$$

- If we take the center of a drift ($\alpha_0=0$), we find $\beta(s) = \beta_0 + \frac{s^2}{\beta_0}$
- It doesn't matter what you do – β will grow!



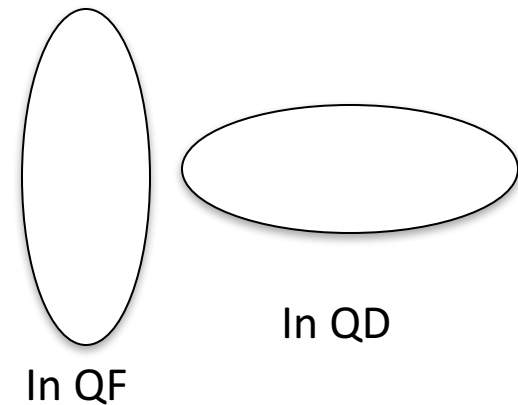
Seems fine, until...



Detectors are a bit bigger
than a few cm!!

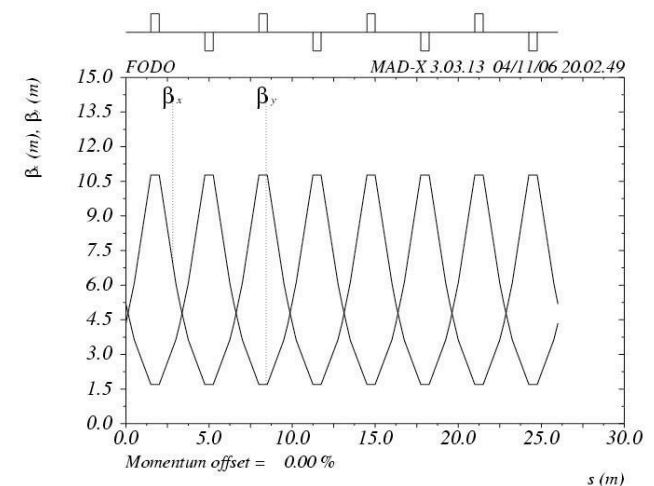
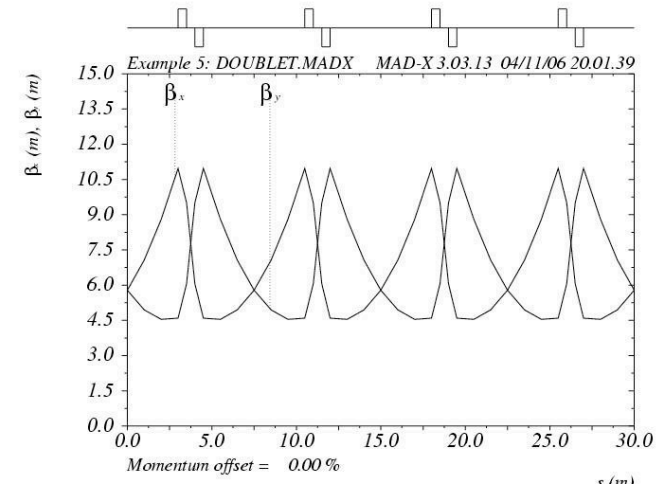
FODO Arcs

- Usually in colliders – take beam between interaction regions
- Simple and tunable (β_x large in QF, β_y at QD)
- Moderate quad strengths
- The beam is not round
- In arcs dipoles generate dispersion



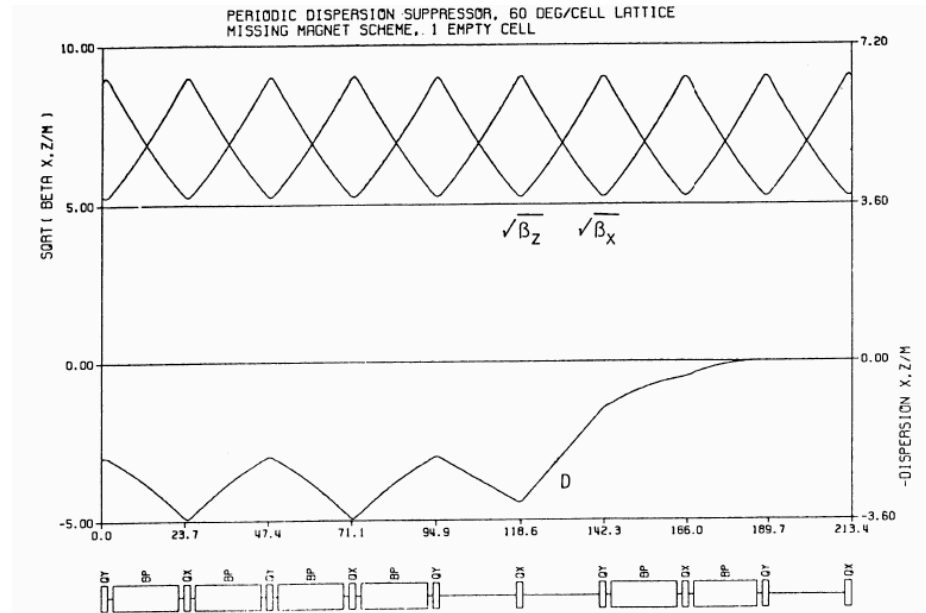
FD Doublet Lattice

- More space between quads
- Stronger quad strengths
- Round beams
- Used in CTF3 linac



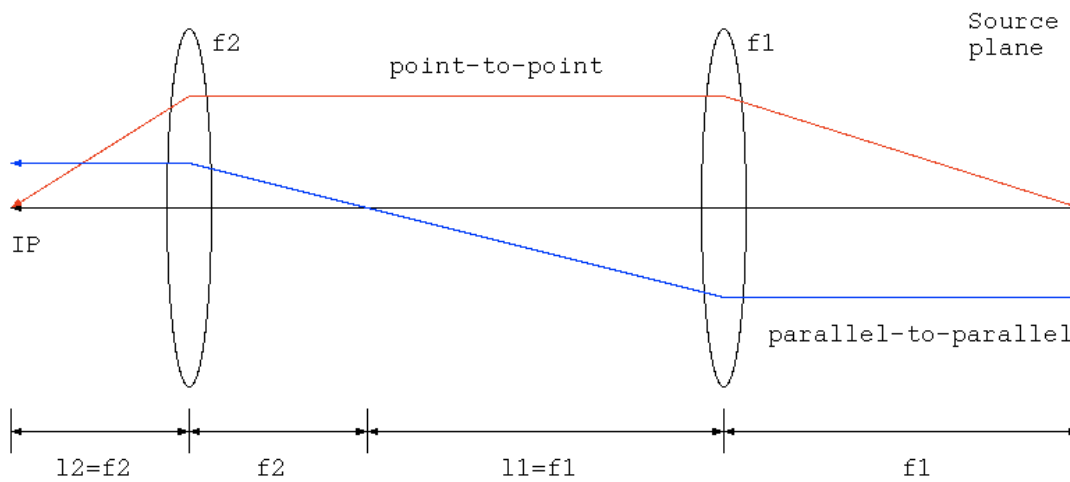
Dispersion suppressor

- Want small spot size at interaction point
- Spot size: $\sqrt{\epsilon\beta + (D\Delta p/p)^2}$
- Missing magnet dispersion suppression scheme
- Works with proper phase advance between elements



Bring D from FODO cell value to zero by a forced oscillation around $\sim\frac{1}{2}$ of that in the FODO cells

Telescope and low β



- Used in colliders to achieve small beam size at IP, assume $D=0$
- Doublet or Triplet
- Want:
 - Point-to-point $R_{12}=0$
 - Parallel to parallel $R_{21}=0$
 - R_{11} =demagnification
- Ratio of focal lengths
- Needs to work in both planes with doublets/triplets

$$\begin{pmatrix} 1 & l_1 \\ 0 & 1 \end{pmatrix} \begin{pmatrix} 1 & 0 \\ -1/f_1 & 1 \end{pmatrix} \begin{pmatrix} 1 & l_1 \\ 0 & 1 \end{pmatrix} = \begin{pmatrix} 1 - l_1/f_1 & 2l_1 - l_1^2/f_1 \\ -1/f_1 & 1 - l_1/f_1 \end{pmatrix}$$

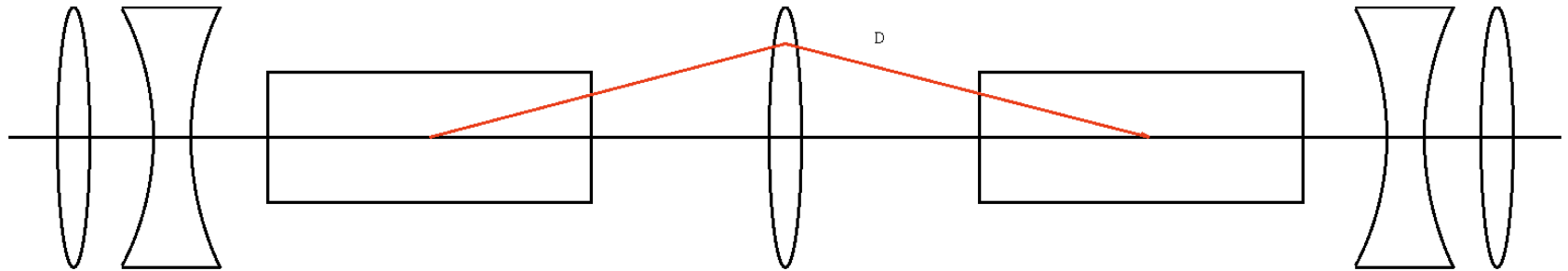
For one module with $l_1=f_1$

$$\begin{pmatrix} 0 & f_1 \\ -1/f_1 & 0 \end{pmatrix}$$

For both modules:

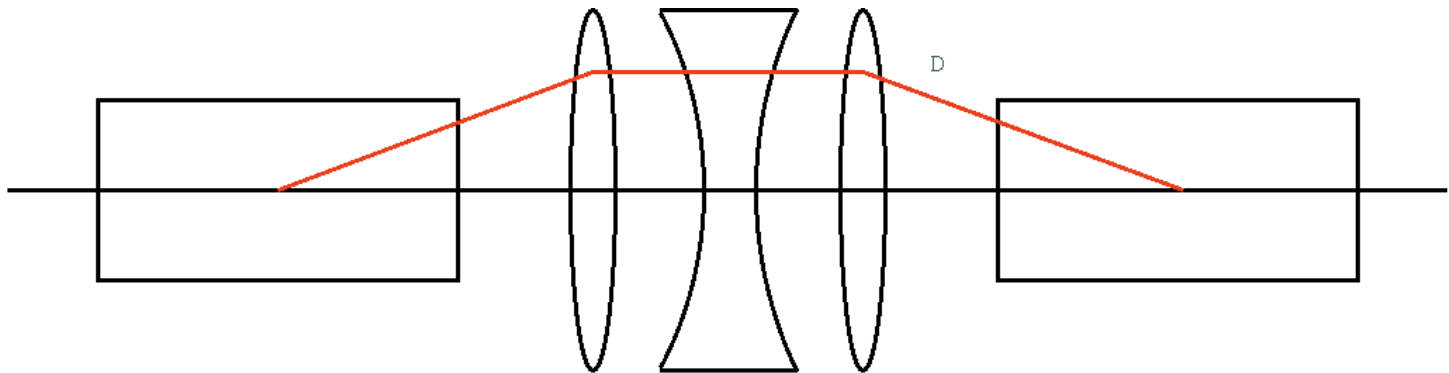
$$R = \begin{pmatrix} -f_2/f_1 & 0 \\ 0 & -f_1/f_2 \end{pmatrix}$$

Double bend achromat



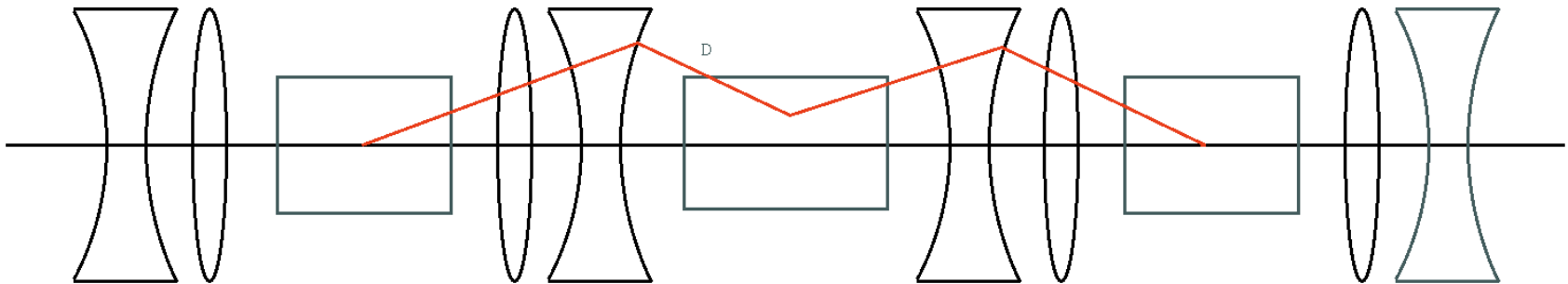
- One dipole generates dispersion and the next, which is 180 degrees apart will take it out again
- Remember: the dispersion is the orbit of a particle with slightly too high momentum w.r.t the reference particle
- Quadrupoles are used to make β_x in dipoles small

Triplet achromat



- Do the 180 degrees in the horizontal plane and the beta matching by quads between dipoles
- very compact, few magnets, but not flexible

Triple bend achromat



- Small emittance.
- Very flexible due to large number of quadrupoles.
- Adjacent drift space can be made long to accommodate undulators/wigglers.

Resources

- Many examples available at the MAD-X website
 - A helpful ‘primer’ by W. Herr:
http://madx.web.cern.ch/madx/doc/madx_primer.pdf

You can always ask me or another lecturer – though we can’t promise to know the answer!