20.4 Solutions of Exercises of Chapter 9: Weak Focusing Synchrotron

9.1

Construct SATURNE I. Spin Resonances

A photo of SATURNE I synchrotron can be found in Fig. 9.1. A schematic layout of the ring and 90 deg cell is given in Fig. 9.20. This figure as well as Tab. 9.1 which lists the parameters of the synchrotron, will be referred to in building the SATURNE I ring in the following.

(a) A model of SATURNE I synchrotron.

DIPOLE is used to simulate the 90^o cell dipole, data are set for a hard-edge model in this exercise (for a DIPOLE model including fringe field, refer to the ZGS case, Exercise 9.2).

It is necessary to have Fig. 20.55 at hand (in addition to the users' guide), when filling up the data list under DIPOLE. Some guidance regarding these data:

- DIPOLE is defined in a cylindrical coordinate system.
- AT is given the value of the bending sector extent: AT=90 degrees. The dipole EFBs coincide with DIPOLE entrance and exit boundaries.
- RM is given the curvature radius value, $RM = B\rho/B = 0.274426548\,[\mathrm{T\,m}]/0.03259493\,[\mathrm{T}] = 8.4193\,\mathrm{m}$, as it fits the geometry of the optical axis around the ring. The field value matches the reference rigidity under OBJET, these are the injection energy values, 3.6 MeV, proton.
- ACENT=45 deg is the reference azimuth, for the positioning of the entrance and exit EFBs. It is taken half-way of the AT range, an arbitrary choice.

KPOS=2 allows cancelling the coordinates of particle 1 (considered here as the reference trajectory, coinciding with the optical axis around the ring) at entrance and exit of DIPOLE:

- The entrance and exit radii in and out of the AT sector for a particle on the closed orbit (*i.e.*, a particle travelling along the design optical axis) are $RE = RS = RM$.
- The angle TE identifies with the closed orbit angle at the entrance boundary: TE=0, the closed orbit is normal to the EFB. TS identifies with the closed orbit angle at the exit boundary: TE=0, the closed orbit is normal to the EFB.

A 90 deg sector in the hard edge model is given in Tab. 20.46; note that the sector has been split in two 45 deg halves, this is in order to allow a possible insertion of a beam monitor, so requiring $AT = 45\,\mathrm{deg}$, $\omega^+ = -\omega^- = 22.5\,\mathrm{deg}$. FAISCEAU located next to DIPOLE indicates that a trajectory entering DIPOLE at radius $R = RM$, normally to the EFB (thus, $Y_0 = 0$ and $T_0 = 0$ in OBJET) exits with $Y = 0$ and $T = 0$. Data validation at this stage can be performed by comparing DIPOLE's transport matrix computed with MATRIX (Tab. 20.47), and theoretical expectations (Sect. 18.2, Eq. 18.6):

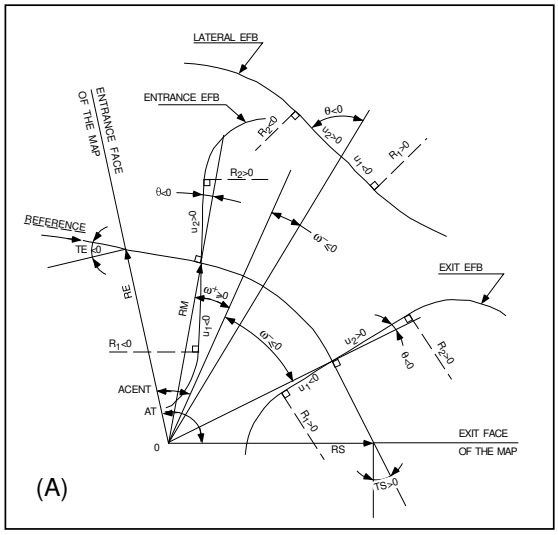

Fig. 20.55 A representation of the data that define a dipole magnet, using DIPOLE [1]

$$[T_{ij}]\ \begin{matrix}\alpha=\pi/2,\\ \rho=8.4193\\ n=0.6\\ =\\ (Eq.\ 18.6)\end{matrix}\ \begin{pmatrix} 0.545794 & 11.15444 & 0 & 0 & 0 & 9.560222\\ 0.062944 & 0.545794 & 0 & 0 & 0 & 1.324865\\ 0 & 0 & 0.346711 & 10.19506 & 0 & 0\\ 0 & 0 & -0.086295 & 0.346711 & 0 & 0\\ 1.324865 & 9.560222 & 0 & 0 & 1 & 5.17640\\ 0 & 0 & 0 & 0 & 0 & 1 \end{pmatrix} \tag{20.13}$$

Introducing fringe fields

The SATURNE ring simulations which follow use the hard edge model. However, it is leisurable, at this point, to choose to add fringe fields in the model; here are the changes which would be needed if so desired:

- The bending sector is 90 degrees, however the field region extent AT has to encompass the fringe fields, at both ends of the 90 deg sector. A 5 deg extension is taken (namely, $ACENT - \omega^+ = AT - ACENT + \omega^- = 5\,\text{deg}$), for a total AT=100 deg which allows $RM \times \tan(ACENT - \omega^+) \approx 74$ cm; this large extension ensures absence of truncation of the fringe fields at the AT sector boundaries, over the all radial excursion of the beam.
- ACENT=50 deg is the reference azimuth (an arbitrary value; taken half-way of the AT range for convenience), for the positioning of the entrance and exit EFBs.
- The entrance radius in the AT sector is $RE = RM/\cos(AT - \omega^+) = RM/\cos(5^o)$, with $\omega^+ = 45\,\text{deg}$ the positioning of the entrance EFB with respect to ACENT. And similarly for the positioning of the exit reference frame, $RS = RM/\cos(AT -$

Table 20.46 Simulation input data file: a pair of adjacent 45 degree sectors in the hard edge model. The magnet is split in order to allow insertion of FAISTORE or (here) FAISCEAU for beam monitoring. The reference optical axis has equal entrance (RE) and exit (RS) positions, and null angles (TE and TS), it coincides with the arc of radius $R = RM$ inside the sector. This input data file is named SatI_DIP.inc and defines the SATURNE I cell sequence segment S_SatI_DIP to E_SatI_DIP, for INCLUDE statements in subsequent exercises

```
File name: SatI_DIP.inc
! SATURNE I. Hard edge dipole model. Transport matrix.
'MARKER'   SatI_DIP.inc_S                                              ! Just for edition purposes.
'OBJET'
0.274426548e3                                                  ! Reference Brho: 3.6 MeV proton.
5                                          ! Create a 13 particle set, proper for MATRIX computation.
.001 .01 .001 .01 .001 .0001                                           ! Coordinate sampling.
0. 0. 0. 0. 0. 1.          ! Reference trajectory: all initial coordinates nul, relative rigidity D=1.
1

'MARKER' S_SatI_DIP     ! Cell dipole begins here. A marker used for INCLUDEs in subsequent exercises.

'DIPOLE'  upstream_half                                  ! Analytical modeling of a dipole magnet.
2                      ! IL=2 here, to log trajectory coordinates in zgoubi.plt, at integration steps.
45. 841.93                  ! Field region angle=90; reference radius set to curvature radius value.
22.5 0.3259493638 -0.6 0. 0.   ! Reference angle ACENT set to AT/2; Bo field at RM; indices, all zero.
.0  0.                                                                 ! EFB 1, hard-edged.
4  .1455   2.2670  -.6395  1.1558  0. 0.  0.                          ! Enge coefficients.
22.5 0.  1.E6  -1.E6  1.E6  1.E6                    ! Angle to ACENT; face angle; face is straight.
.0  0.                                                                 ! EFB 2, hard-edged.
4  .1455   2.2670  -.6395  1.1558  0. 0.  0.
-22.5 0.  1.E6  -1.E6  1.E6  1.E6
0. 0.                                                                      ! EFB 3. Unused.
0  0.      0.      0.      0.      0. 0.  0.
0. 0.  1.E6  -1.E6  1.E6  1.E6 0.
2 1                                       ! Degree of interpolation polynomial; flying grid sizing.
2.                                       ! Integration step size. It can be large in uniform field.
2 841.93 0. 841.93 0.                                ! Positioning of entrance and exit frames.
'MARKER' half-dipole !.plt              ! Uncomment LABEL_2='.plt' (may go with IL=2 under DIPOLE) to
                                                                ! log particle data in zgoubi.plt.
'FAISCEAU'        ! Provides local coordinates, and ellipse parameters, at center of SATURNE I dipole.
'DIPOLE'  downstream_half                                ! Analytical modeling of a dipole magnet.
2                      ! IL=2 here, to log trajectory coordinates in zgoubi.plt, at integration steps.
45. 841.93                  ! Field region angle=90; reference radius set to curvature radius value.
22.5 0.3259493638 -0.6 0. 0.   ! Reference angle ACENT set to AT/2; Bo field at RM; indices, all zero.
.0  0.                                                                 ! EFB 1, hard-edged.
4  .1455   2.2670  -.6395  1.1558  0. 0.  0.                          ! Enge coefficients.
22.5 0.  1.E6  -1.E6  1.E6  1.E6                    ! Angle to ACENT; face angle; face is straight.
.0  0.                                                                 ! EFB 2, hard-edged.
4  .1455   2.2670  -.6395  1.1558  0. 0.  0.
-22.5 0.  1.E6  -1.E6  1.E6  1.E6
0. 0.                                                                      ! EFB 3. Unused.
0  0.      0.      0.      0.      0. 0.  0.
0. 0.  1.E6  -1.E6  1.E6  1.E6 0.
2 1                                       ! Degree of interpolation polynomial; flying grid sizing.
2.                                       ! Integration step size. It can be large in uniform field.
2 841.93 0. 841.93 0.                                ! Positioning of entrance and exit frames.

'MARKER' E_SatI_DIP      ! Cell dipole ends here. A marker used for INCLUDEs in subsequent exercises.

'FAISCEAU'                                                         ! Local particle coordinates.
'MATRIX'                                    ! Compute transport matrix, from trajectory coordinates.
1 0

'MARKER'   SatI_DIP.inc_E                                              ! Just for edition purposes.
'END'
```

$(ACENT - \omega^-)) = RM/\cos(5^o)$ with $\omega^- = -45$ deg the positioning of the exit EFB. Note that $\omega^+ - \omega^- = 90^o$, the value of the bend angle.

- The entrance angle TE identifies with the angular increase of the sector: TE=5 deg. And similarly for the positioning of exit frame, 5 deg downstream of the exit EFB, thus TS=5 deg.
- Negative drifts with length $RM \times \tan(ACENT - \omega^+) = 0.7366545469$ cm need to be added upstream and downstream of DIPOLE, to account for the optical axis additional length over the 5 deg angular extent.

Table 20.47 Outcomes of the simulation file of Tab. 20.46

An excerpt from zgoubi.res. Coordinates of the first particle (*considered here as the reference trajectory*) *and its path length under FAISCEAU, at OBJET on the left hand side below, locally on the right hand side*:

```
         3  Keyword, label(s) :  FAISCEAU
                                          TRACE DU FAISCEAU
                                       (follows element #      2)
                                          13 TRAJECTOIRES
                                 OBJET                                          FAISCEAU
              D       Y(cm)     T(mr)     Z(cm)     P(mr)      S(cm)      D-1     Y(cm)     T(mr)     Z(cm)     P(mr)      S(cm)
   0  1   1.0000     0.000     0.000     0.000     0.000     0.0000    0.0000    0.000     0.000     0.000     0.000   1.322501E+03
```

Transport matrix of SATURNE I 90 *degree sector bend, in the hard edge model, two difference cases of integration step size, namely,* 4 *cm and* 1 *m* (*an excerpt of MATRIX computation, from zgoubi.res*). *It can be checked against matrix transport expectations. The "first order symplectic conditions" are very small in the* 4 *cm step size case, which is an indication of accurate numerical integration of the trajectories across DIPOLE*; *the reference trajectory* (*first one*) *exits better aligned* (*reference coordinates, before change of frame for MATRIX computation, are closer to zero*):

- Case of 4 *cm step size*:

```
         4  Keyword, label(s) :  MATRIX

      Reference, before change of frame (particle #  1  - D-1,Y,T,Z,s,time)  :
       0.00000000E+00   4.53054326E-07   6.27843350E-07   0.00000000E+00   0.00000000E+00   1.32250055E+03   4.41138700E-02

                     TRANSFER  MATRIX  ORDRE  1  (MKSA units)
               0.545795         11.1544         0.00000         0.00000         0.00000         9.56022
              -6.294423E-02    0.545795         0.00000         0.00000         0.00000         1.32487
                0.00000         0.00000        0.346711         10.1951         0.00000         0.00000
                0.00000         0.00000       -8.629576E-02    0.346711        0.00000         0.00000
                1.32487         9.56022         0.00000         0.00000         1.00000         5.17640
                0.00000         0.00000         0.00000         0.00000         0.00000         1.00000

               DetY-1 =       0.0000000278,    DetZ-1 =       0.0000000045
```

- Case of 1 *m step size*:

```
         4  Keyword, label(s) :  MATRIX

      Reference, before change of frame (particle #  1  - D-1,Y,T,Z,s,time)  :
       0.00000000E+00  -7.54923113E-03  -1.08904867E-02   0.00000000E+00   0.00000000E+00   1.32249873E+03   4.41138091E-02

                     TRANSFER  MATRIX  ORDRE  1  (MKSA units)
               0.545757         11.1567         0.00000         0.00000         0.00000         9.56154
              -6.295274E-02    0.546125         0.00000         0.00000         0.00000         1.32517
                0.00000         0.00000        0.346697         10.1954         0.00000         0.00000
                0.00000         0.00000       -8.629900E-02    0.346750        0.00000         0.00000
                1.32486         9.56148         0.00000         0.00000         1.00000         5.17692
                0.00000         0.00000         0.00000         0.00000         0.00000         1.00000

               DetY-1 =       0.0003978566,    DetZ-1 =       0.0000685588
```

(b) SATURNE I cell.

A cell with origin in the middle of the drift is given Tab. 20.48, it is comprised of the split dipole and a pair of 2 m half-drifts at each ends (Fig. 9.20).

Closed orbit; *chromatic closed orbit*

The on-momentum closed orbit has been set to zero along the drifts ($Y_{\text{c.o.}} \equiv 0$), above, by a proper choice of RE, RS radii and TE, TS incidence angles.

Table 20.48 Simulation input data file: SATURNE I cell, assembled by INCLUDE-ing DIPOLE taken from Tab. 20.46 together with two half-drifts. This input data file is named SatI_cell.inc and defines the SATURNE I cell sequence segment S_SatI_cell to E_SatI_cell, for INCLUDE statements in subsequent exercises

```
File name: SatI_cell.inc.
! SATURNE I, one cell of the 4-period ring.
'MARKER'   SatICellMATRIX_S                                              ! Just for edition purposes.
'OBJET'
0.274426548e3                                                        ! Reference Brho: 3.6 MeV proton.
5                                              ! Create a 13 particle set, proper for MATRIX computation.
.001 .01 .001 .01 .001 .0001                                                 ! Coordinate sampling.
0. 0. 0. 0. 0. 1.          ! Reference trajectory: all initial coordinates nul, relative rigidity D=1.

'MARKER'  S_SatI_cell
'DRIFT' half_drift
200.
'INCLUDE'
1
./SatI_DIP.inc[S_SatI_DIP:E_SatI_DIP]
'DRIFT' half_drift
200.
'MARKER'  E_SatI_cell
'FAISCEAU'                                                               ! Local particle coordinates.
'TWISS'    ! Produce transport matrix, beam matrix, and periodic optical functions along the sequence.
2  1. 1.
'MARKER'   SatICellMATRIX_E                                              ! Just for edition purposes.
'END'
```

The radial coordinate of an off-momentum chromatic orbit can be estimated from the dispersion, Eq. 9.26, namely,

$$Y_\delta = \frac{\rho_0}{1-n}\frac{\delta p}{p} = 841.93\frac{10^{-4}}{1-(-0.6))} \approx 0.21048\,\text{cm}$$

whereas the orbit angle is zero, around the ring (on- and off-momentum closed orbits are parallel to the optical axis).

Besides,

- computation of an accurate value of Y_δ is performed adding FIT at the end of the cell;

- in order to raytrace three particles, respectively on-momentum and at $\delta p/p = \pm 10^{-4}$, OBJET[KOBJ=2] is used;

- in order to raytrace around the ring, for the purpose of plotting the closed orbit coordinates, a 4-cell sequence follows the FIT procedure.

This results in the input data file given in Tab. 20.49. Running this input simulation file produces the following coordinates as per the FIT procedure (an excerpt from zgoubi.res):

```
 STATUS OF VARIABLES  (Iteration #     4 /    999 max.)
LMNT VAR PARAM  MINIMUM    INITIAL        FINAL        MAXIMUM     STEP        NAME   LBL1                LBL2
   2   1    30   0.168       0.211      0.21056000      0.253      1.040E-05  OBJET      -                   -
   2   2    40    0.00        0.00       0.0000000       0.00       0.00      OBJET      -                   -
   2   3    50  -0.253      -0.210     -0.21040403     -0.168      1.040E-05  OBJET      -                   -
 STATUS OF CONSTRAINTS (Target penalty =   1.0000E-10)
TYPE  I   J LMNT#     DESIRED          WEIGHT         REACHED         KI2     NAME   LBL1                LBL2      Nb param. [value]
  3   1   2    12    0.000000E+00    1.000E+00    1.466978E-06    6.70E-01 MARKER     E_SatI_cell         -                   0
  3   2   2    12    0.000000E+00    1.000E+00    6.028957E-07    1.13E-01 MARKER     E_SatI_cell         -                   0
  3   3   2    12    0.000000E+00    1.000E+00    8.357183E-07    2.17E-01 MARKER     E_SatI_cell         -                   0
 Fit reached penalty value   3.2139E-12
```

The local coordinates Y, T and initial coordinates Y_0, T_0 (as defined under OBJET) are identical to better than 5 μm, 0.5 μrad accuracy, respectively, confirming the

6724 periodicity of these chromatic trajectories. Orbit coordinates around the ring are displayed in Fig. 20.56.

Table 20.49 Simulation input data file: first find the periodic orbit through a cell, then complete a 4-cell turn

```
SatI_Orbits.INC.dat: SATURNE I, on-momentum and chromatic orbits.
'MARKER'   SatI_Orbits_S                                          ! Just for edition purposes.
'OBJET'
0.274426548e3                                                   ! Reference Brho: 3.6 MeV proton.
2                                                                ! Create particles individually.
3 1                                                                        ! Three particles.
+.210560 0. 0. 0. 0. 1.0001 'p' ! Chromatic orbit coordinates Y0 and T0 for D=1.001 relative rigidity.
0. 0. 0. 0. 0. 1.  'o'                                                     ! On-momentum orbit.
-.210404 0. 0. 0. 0. 0.9999 'm' ! Chromatic orbit coordinates Y0 and T0 for D=0.999 relative rigidity.
1 1 1

'INCLUDE'
1
./SatI_cell.inc[S_SatI_cell:E_SatI_cell]

'FIT'
2
2 30 0 .2                                                       ! Vary Y_0(particle 1) under OBJET.
2 50 0 .2                                                       ! Vary Y_0(particle 3) under OBJET.
2
3.1 1 2 #End 0. 1. 0                                  ! Constrain Y(particle 1)=Y_0(particle 1).
3.1 3 2 #End 0. 1. 0                                  ! Constrain Y(particle 1)=Y_0(particle 1).

!       When FIT is done converging on the constraints, execution quietly carries on with the periodic
!                          coordinates , raytracing through 4 cells to complete a turn around the ring.
'INCLUDE'
1
4 * ./SatI_cell.inc[S_SatI_cell:E_SatI_cell]

'SYSTEM'
1
gnuplot < gnuplot_Zplt_traj.gnu                                 ! Plot the orbit radial coordiante.
'MARKER'   SatI_Orbits_E                                              ! Just for edition purposes.
'END'
```

A gnuplot script (*excerpt*) *to obtain a graph of particle coordinates, from zgoubi.plt* (*as in Fig.* 20.56):

```
# gnuplot_Zplt_traj.gnu
traj1 = 1 ; traj2 = 3
plot \
for [i=traj1:traj2]  'zgoubi.plt' u ($19== i ? $14 *cm2m : 1/0):($10 *cm2m):($19) w p ps .4 lc palette
```

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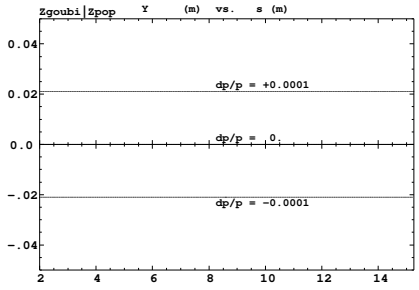

Fig. 20.56 Radial coordinate of the orbits around the ring, on-momentum, and for $dp/p = \pm 10^{-3}$. This graph is obtained using zpop, data read from zgoubi.plt: menu 7; 1/1 to open zgoubi.plt; 2/[6,2] for Y versus distance s; 7 to plot. A gnuplot script for a similar graph given is given in Tab. 20.49

Lattice parameters

The TWISS command down the sequence (Tab. 20.48) produces the periodic beam matrix results shown in Tab. 20.50; MATRIX[IFOC=11] would, as well. It also

Table 20.50 Results obtained running the simulation input data file of Tab. 20.48, SATURNE I cell - an excerpt from zgoubi.res

```
   14  Keyword, label(s) :  TWISS

Reference, before change of frame (particle #  1  - D-1,Y,T,Z,s,time)  :
 0.00000000E+00   6.02895730E-07   6.54169939E-07   0.00000000E+00   0.00000000E+00   1.72250055E+03   6.57784696E-01

     Beam  matrix  (beta/-alpha/-alpha/gamma) and  periodic  dispersion  (MKSA units)
        14.418595     0.000000     0.000000     0.000000     0.000000    21.048250
         0.000000     0.069355     0.000000     0.000000     0.000000     0.000000
         0.000000     0.000000    11.411041     0.000000     0.000000    -0.000000
         0.000000     0.000000     0.000000     0.087634     0.000000     0.000000
         0.000000     0.000000     0.000000     0.000000     0.000000     0.000000
         0.000000     0.000000     0.000000     0.000000     0.000000     0.000000

                          Betatron  tunes  (Q1 Q2 modes)
                 NU_Y =  0.18103144         NU_Z =  0.22214599

                          dL/L / dp/p =   1.9194487
(dp =  0.000000E+00      L(0)   =   1.72250E+03 cm,  L(0)-L(-dp) =    3.30606E-01 cm,  L(0)-L(+dp) =   -3.30645E-01 cm)

                               Transition gamma  =  7.21791469E-01

                               Chromaticities :
                 dNu_y / dp/p = -0.60221729                 dNu_z / dp/p =  0.38005442
```

produces a zgoubi.TWISS.out file which details the optical functions along the sequence (at the downstream end of the optical elements. The header of that file details the optical parameters of the structure (Tab. 20.51).

Table 20.51 An excerpt of zgoubi.TWISS.out file resulting from the execution of the SATURNE I cell simulation input data file of Tab. 20.48. Note that the ring (4-period) wave numbers are 4 times the cell values Q1, Q2 displayed here. Optical functions (betatron function and derivative, orbit, phase advance, etc.) along the optical sequence are listed as part of zgoubi.TWISS.out following th header. The top part and last line of that listing are given below

```
@ LENGTH          %le     17.22500552
@ ALFA            %le     1.919448707
@ ORBIT5          %le                -0
@ GAMMATR         %le    0.7217914685
@ Q1              %le    0.1810314404      [fractional]
@ Q2              %le    0.2221459901      [fractional]
@ DQ1             %le   -0.6022172911
@ DQ2             %le    0.3800544183
@ DXMAX           %le     2.10586311E+01    @ DXMIN           %le     2.10482503E+01
@ DYMAX           %le     0.00000000E+00    @ DYMIN           %le     0.00000000E+00
@ XCOMAX          %le     2.10528899E-01    @ XCOMIN          %le     0.00000000E+00
@ YCOMAX          %le     0.00000000E+00    @ YCOMIN          %le     0.00000000E+00
@ BETXMAX         %le     1.57006971E+01    @ BETXMIN         %le     1.44132839E+01
@ BETYMAX         %le     1.30884296E+01    @ BETYMIN         %le     1.14110171E+01
@ XCORMS          %le     6.05227342E-04
@ YCORMS          %le    0.    not computed
@ DXRMS           %le     2.98427468E-03
@ DYRMS           %le     0.00000000E+00
```

Optical functions listing zgoubi.TWISS.out (*there is more*: $D_{x,y}$, *etc.*: *lines are truncated, here*), *including the periodic* $\alpha_{x,y,l}$, $\beta_{x,y,l}$, $D_{x,y}$, *etc.*

```
# alfx          btx           alfy          bty           alfl          btl           Dx            Dxp
# 1             2             3             4             5             6             7             8
 1.3683565E-08  1.4426805E+01 -6.6336606E-09  1.1411067E+01  0.0000000E+00  0.0000000E+00  2.1058631E+01  1.1261490E-03
 2.3958789E-08  1.4426805E+01 -2.0952612E-10  1.1411067E+01  0.0000000E+00  0.0000000E+00  2.1048250E+01  3.4685148E-09
 2.3958789E-08  1.4426805E+01 -2.0952612E-10  1.1411067E+01  0.0000000E+00  0.0000000E+00  2.1048250E+01  3.4685148E-09
-1.3863081E-01  1.4704066E+01 -1.7526845E-01  1.1761604E+01  0.0000000E+00  0.0000000E+00  2.1048250E+01  3.4685148E-09
-1.3863081E-01  1.4704066E+01 -1.7526845E-01  1.1761604E+01  0.0000000E+00  0.0000000E+00  2.1048250E+01  3.4685148E-09
 5.1661104E-04  1.5700697E+01  2.2204071E-06  1.3088430E+01  0.0000000E+00  0.0000000E+00  2.1048250E+01  1.4621225E-09
 5.1661104E-04  1.5700697E+01  2.2204071E-06  1.3088430E+01  0.0000000E+00  0.0000000E+00  2.1048250E+01  1.4621225E-09
 1.3919474E-01  1.4692541E+01  1.7526999E-01  1.1761559E+01  0.0000000E+00  0.0000000E+00  2.1048250E+01  3.4685146E-09
 1.3919474E-01  1.4692541E+01  1.7526999E-01  1.1761559E+01  0.0000000E+00  0.0000000E+00  2.1048250E+01  3.4685146E-09
 4.3383067E-04  1.4413284E+01  7.7310157E-07  1.1411017E+01  0.0000000E+00  0.0000000E+00  2.1048250E+01  3.4685146E-09
 4.3383067E-04  1.4413284E+01  7.7310157E-07  1.1411017E+01  0.0000000E+00  0.0000000E+00  2.1048250E+01  3.4685146E-09
```

Moving the origin of the cell

The origin of the sequence can be moved by placing both drifts on one side of DIPOLE. It can also be taken in the middle of DIPOLE, as the latter has been split. A fully deployed input data sequence (INCLUDEs accounted for) is provided at the top of the execution listing zgoubi.res, it can be used to copy-paste pieces around. It can then be checked that betatron tunes, chromaticities, momentum compaction (Tab. 20.50) do not change, and that the beam matrix does.

Optical functions along the cell

They are computed by transporting the beam matrix, from the origin. A Fortran program available in zgoubi sourceforge package toolbox, betaFromPlt [1], performs this computation in the following way: OBJET[KOBJ=5.1] provides the initial beta function values (determined in the previous question); IL=2 under DIPOLE logs stepwise particle data in zgoubi.plt; 'split 10 2' added under DRIFT does it, too. The program betaFromPlt computes the transport matrix T_{step} from a step to the next by reading particle data from zgoubi.plt, and performing a similar computation to what

MATRIX does. The beam matrix σ is then transported, from step 1 to step 2, using (Eq. 19.11)

$$\sigma_2 = T_{\text{step}}\ \sigma_1\ \tilde{T}_{\text{step}}$$

The result is displayed in Fig. 20.57.

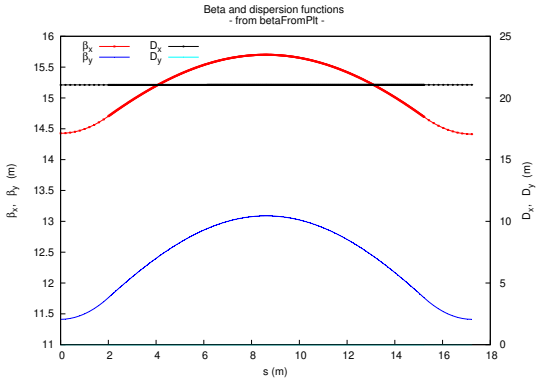

Fig. 20.57 Optical functions along SATURNE I cell. They are obtained from the transport of the beta functions, from the origin (at OBJET), using transport matrices computed from step-by-step particle coordinates stored in zgoubi.plt

Tune scan

A simulation is given in Tab. 20.52, derived from Tab. 20.48: MATRIX[IFOC=11] has been substituted to TWISS, a REBELOTE do loop repeatedly changes n. A graph of the scan is given in Fig. 20.58, a few values are detailed in Tab. 20.53.

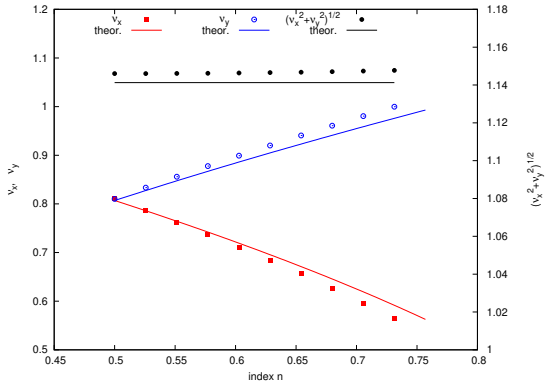

Fig. 20.58 A scan of the wave numbers, and of $\sqrt{\nu_Y^2 + \nu_Z^2} \approx \sqrt{R/\rho_0} = 1.141$, in SATURNE I for $0.5 \leq n \leq 0.757$

Table 20.52 Simulation input data file: tune scan, using REBELOTE to repeatedly change n. Beam matrix and wave numbers are computed by MATRIX, from the coordinates of the 13 particle sample generated by OBJET[KOBJ=5]

```
SATURNE I, tune scan.
'MARKER'   SatI_Qscan_S                                          ! Just for edition purposes.
'OBJET'
0.274426548e3                                                  ! Reference Brho: 3.6 MeV proton.
5                                        ! Create a 13 particle set, proper for MATRIX computation.
.001 .01 .001 .01 .001 .0001                                     ! Coordinate sampling.
0. 0. 0. 0. 0. 1.          ! Reference trajectory: all initial coordinates nul, relative rigidity D=1.
1

'MARKER'  S_SatI_cell
'DRIFT' half_drift
200.
'INCLUDE'
1
./SatI_DIP.inc[S_SatI_DIP:E_SatI_DIP]
'DRIFT' half_drift
200.
'MARKER'  E_SatI_cell
'FAISCEAU'                                                     ! Local particle coordinates.
'MATRIX'
1 11  PRINT   ! Comoute a 10+4 period transport matrix, and tunes. Save outcomes to zgoubi.MATRIX.out.

'REBELOTE'                    ! A do loop: repeat the section above commencing at the top of the file,
10  1.1  0 1                                                                    ! 10 times.
1
DIPOLE 6  -0.757:-0.5   ! Change the value of parameter 30 (namely, n) in DIPOLE (prior to repeating).
                                                                  ! in any DIPOLE in the sequence.
'SYSTEM'
1
gnuplot <./gnuplot_MATRIX_Qxy.gnu                                    ! Plot tunes vs index.
'MARKER'   SatI_Qscan_E                                        ! Just for edition purposes.
'END'
```

gnuplot script to obtain Fig. 20.58:

```
# ./gnuplot_MATRIX_Qxy.gnu
set xlabel "index n";set ylabel "{/Symbol n}_x,     ({/Symbol n}_x^2+{/Symbol n}_y^2)^{1/2}"
set y2label "{/Symbol n}_y"; set xtics; set ytics nomirror; set y2tics nomirror; ncell=4
set key t l; set key maxrow 2; set yrange [:1.3]; set y2range [:1.06]
n1 = -0.757; dn=(.757-.5)/10.; R=10.9658; rho=8.4193
plot \
"zgoubi.MATRIX.out" u (n1+($61-1)*dn): \
($61>1? $56 *ncell :1/0) w p pt 5 lt 1 lw .5 lc rgb "red"               tit "{/Symbol n}_x " ,\
"zgoubi.MATRIX.out" u (n1+($61-1)*dn):($61>1? sqrt((1+(n1+($61-1)*dn))*R/rho): \
1/0) w l lt 1  lc rgb "red" tit "theor. " ,\
"zgoubi.MATRIX.out" u (n1+($61-1)*dn): \
($61>1? $57 *ncell :1/0) axes x1y2 w p pt 6 lt 3 lw .5 lc rgb "blue"   tit "{/Symbol n}_y " ,\
"zgoubi.MATRIX.out" u (n1+($61-1)*dn): \
($61>1? sqrt((-(n1+($61-1)*dn))*R/rho):1/0) axes x2y2 w l lt 3  lc rgb "blue" tit "theor. " ,\
"zgoubi.MATRIX.out" u (n1+($61-1)*dn): \
($61>1? sqrt($56**2+$57**2) *ncell :1/0) w p pt 7 lt 1 lc rgb "black" tit "({/Symbol n}_x^2+{/Symbol n}_y^2)^{1/2}" ,\
"zgoubi.MATRIX.out" u (n1+($61-1)*dn):($61>1? sqrt(R/rho):1/0)  w l lt 1  lc rgb "black" tit "theor. "
pause 1
```

Table 20.53 Dependence of wave numbers on index n, from numerical raytracing (columns denoted "ray-tr.") and from theory

	ν_Y		ν_Z	
n	ray-tr.	$\sqrt{(1-n)\frac{R}{\rho_0}}$	ray-tr.	$\sqrt{n\frac{R}{\rho_0}}$
0.5	0.810353	0.806987	0.810353	0.806987
0.6	0.724125	0.721791	0.888583	0.884010
0.7	0.626561	0.625089	0.960806	0.954840
0.757	0.563635	0.562580	0.999804	0.992955