

Fig. 3.39 A scan of the revolution time, from 0.02 to 1 MeV, and its dependence on the field index k. The right vertical axis only concerns the case k = 0 where the change in revolution time is weak and only due to the mass increase (in $T_{rev} = 2\pi \gamma m_0/qB$). The right graph shows, up to 5 MeV, the relatively important contribution of the focusing index, even a weak k=-0.03, compared to the effect of the mass increase (k=0 curve). Markers are from raytracing, solid lines are from theory

2587 **3.8 Ion Trajectories**

A zgoubi data file is set up for computation of particle trajectories, taking a field value on reference radius of $B_0(R_0) = 0.5$ T, and reference energy 200 keV (proton). These hypotheses determine the reference radius value. DIPOLE [16, *lookup* INDEX] is used (Tab. 3.21), for its greater flexibility in changing magnet parameters, field and radial field index amongst other, compared to using TOSCA and a field map.

(a) Transverse motion.

It first has to be checked that there is consistency between initial orbital radius Y_0 in OBJET at 200 keV proton energy and the value of the reference radius R_0 in DIPOLE (Eq. 3.35). Its theoretical value is $R_0 = BORO/5[kG] = 12.924889$ cm, a closed orbit finding using FIT can be performed, or it can be referred to the solutions of earlier exercises, to check agreement with raytracing outcomes.

(b) Wave numbers at 1 and 5 MeV.

These considerations result in the input data file given in Tab. 3.22, to compute multiturn trajectories. ; note that $R_0 = 12.924889$ cm therein, whereas a value of $R_0 = 50$ cm may be taken instead in other exercises. Field index derivatives k', k'', ... are taken null in the present exercise.

Three particles with paraxial radial and axial motions are raytraced over a few turns. Their starting radius is the closed orbit radius for the respective energies, while a 0.1 mrad take-off angle is imparted to each particle both vertically and horizontally.

The value of the focusing index k_E at an energy *E* can be expressed in terms of DIPOLE data which are, the index value *k* at R_0 (Eq. 3.11), reference radius R_0 , and field $B_0 = B_Z(R_0)$, namely,

$$k_E = \frac{R_E}{B_E} \frac{\partial B}{\partial R} = \frac{R_0 + \Delta R}{B_0 + \Delta B} \frac{\partial B}{\partial R} \approx k \frac{1 + \Delta R/R_0}{1 + k\Delta R/R_0} \approx k \left[1 + (1 - k) \frac{\Delta R}{R_0} \right]$$

3.4 Solutions of Exercises of Chapter 3: Classical Cyclotron

Table 3.21 Input data file 60DegSectorR200.inc: it defines DIPOLE as a sequence segment comprised between the "LABEL_1" type labels [16, Sect. 7.7] #S_60DegSectorR200 and #E_60DegSectorR200. DIPOLE here, has an index k = -0.03, reference radius $R_0 \equiv R_0(E_k = 200 \text{ keV}) = 12.924888 \text{ cm}$ and $B_0 = B(R_0) = 0.5 \text{ T}$. Note that (i) this file can be run on its own: it has been designed to provide the transport MATRIX of that DIPOLE; (ii) for the purpose of some of the exercises, IL=2 under DIPOLE, optional, causes the printout of particle data in zgoubi.plt, at each integration step (this is at the expense of CPU time, and memory volume)

```
60DegSectorR200.inc
       'OBIET
! 200keV proton.
0.01 0.001 0.01 0.001 0.0.0001
12.9248888074 0. 0. 0. 0. 1.
2.9248888074 0. 0. 0. 0. 1.
2.0124888
'Decode State CPU ime.
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       64.62444403717985

    ! Reference ;
    0.
    4.1455 2.2670 -.6395 1.1558 0.0.0.
    30.0.1.E6 -1.E6 1.E6 1.E6
    0.

                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            EFB 2.
                 .1455 2.2670 -.6395 1.1558 0.0.0.
       4
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    -30.
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0. 0.
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1.E6 -1.E6 1.E6 1.E6 0.
                         10.
    0 5
                                                                                                                                                                               ! Integration step size. Small enough for orbits to close accurately
                    0.0.
                                                                                                                                                                                                                                                                                                                                                ! Magnet positionning RE, TE, RS,
       2 0. 0. 0. 0.
'FAISCEAU' #E_60DegSectorR200
       'MATRIX'
     1 0
'END'
```

 Table 3.22
 Simulation input data file: raytrace a few turns around the cyclotron, three particles with different momenta, and 0.1 mrad horizontal and vertical take-off angles. The INCLUDE segment is taken from Tab. 3.21

```
'MARKER' ProbProjTraj_S
'OBJET'
64.62444403717985
                                         ! Reference Brho ("BORO" in the users' guide) -> 200keV proton.
31
12.924888 0.1 0. 0.1 0. 1. 'o'
30.107898 0.1 0. 0.1 0. 2.2365445 'm'
75.754671 0.1 0. 0.1 0. 5.0063900 'o'
                                            ! p[MeV/c]=969.934, Brho[kG.cm]=323.535, kin-E[MeV]=5.
'INCLUDE
6* 60DegSectorR200.inc[#S_60DegSectorR200:#E_60DegSectorR200]
                                                                        ! 6 sectors for an overall 360 deg.
'REBELOTE'
                                                                   ! There will be a total of 9+1=10 tunrs.
9 0.1 99
'SYSTEM'
1
gnuplot < ./gnuplot_Zplt_traj.gnu
'MARKER' ProbProjTraj_E</pre>
                                                               ! Plot the projected Y(s) and Z(s) motions.
'END'
```

A gnuplot script to obtain Fig. 3.41:

set ylabel 'Z [cm]'; plot for [sector=1:6] for [trj=1:3] 'zgoubi.plt' u (\$19==trj && \$42==sector1+2*(sector-1)? \$14*cm2m \
/(2.*pi*R[\$19]) :1/0):(\$12):(\$12):(\$19) w p ps .2 lc palette notit ; pause 1



with ΔR assumed small, $\partial B/\partial R = kB_E/R_E$ an energy independent quantity, and the index E denoting a quantity taken at the reference energy. The latter property is illustrated in Fig. 3.40, produced using the input data file of Tab. 3.23.

 Table 3.23
 Simulation input data file for a magnetic field scan. The INCLUDE segment is taken from Tab. 3.21

```
Field and derivative dB/dR, as a finction of R
'MARKER' ProbProjTrajB_S
'OBJET'
                                           ! Reference Brho ("BORO" in the users' guide) -> 200keV proton.
64.62444403717985
1 1
12.924888 0.1 0. 0.1 0. 1. 'o'
                                             ! Just one ion.
! A particle with kin-E=0.2 MeV and 0.1 mrad take-off angles.
'INCLUDE
                  ! IL=2 is necessary under DIPOLE, for step-by-step log of particle data in zgoubi.plt.
60DegSectorR200.inc[#S_60DegSectorR200:#E_60DegSectorR200]
                                                                                          One sector is enough
'FIT'
2 30 0 [12,80]
                                       ! Vary particle's Y0 at OBJET, to have it match its D (=Brho/BORO).
1 1e-20
3.1 1 2 #End 0. 1. 0
                                                                                         ! Consrain Y final=Y0.
'REBELOTE'
25 0.1 0 1
                                                        ! Scan parameter 35 (relative rigidity, D) in OBJET.
OBJET 35 1:5.00639
                                         ! Scan relative rigidity D from 1 (200 keV) to 5.0063900 (5 MeV).
'SYSTEM'
gnuplot < ./gnuplot_Zplt_field.gnu</pre>
                                                                        ! Plot B(R), as read fron zgoubi.plt.
'MARKER'
'END'
           ProbProjTrajB_E
```

A gnuplot script to obtain Fig. 3.40:

```
# gnuplot_Zplt_field.gnu
set xtics nomirror; set xZtics; set ytics; set xlabel 's /C_E '; set ylabel 'Y [cm]'
set palette defined (1 "red", 2 "blue", 3 "black"); unset colorbox
array R(3); R(1)=0.12924888; R[2]=0.3010789365; R[3]=0.75754671; pi = 4.*atan(1.); cm2m = 0.01
sector1=3 # number (NOEL) of 1st DIPULE in /zgoubi.rpl; pi = 4.*atan(1.); cm2m = 0.01
# in zgoubi.plt, col. 19; particle number; col. 42: keyword number; col. 14: distance; col. 10: Y; col. 12: YZ
plot for [i=1:6] for [trj=1:3]
'zgoubi.plt' u (S19==trj && $42==sector1 +2*(i-1) ? $14*cm2m /(2.*pi*R[$19]) :1/0) \
:(S10*cm2m_R[trj]):(S19) w p ps .2 lc palette notit ; pause 1
set ylabel 'Z [cm]' ;
plot for [i=1:6] for [trj=1:3]
'zgoubi.plt' u ($19==trj && $42==sector1 +2*(i-1) ? $14*cm2m \
/(2.*pi*R[$19]) :1/0):($12):($19) w p ps .2 lc palette notit ; pause 1
```



Fig. 3.41 Radial (left) and axial (right) paraxial motion around respectively the 200 keV (smallest amplitude), 1 MeV (intermediate) and 5 MeV (greatest amplitude) closed orbit (the latter is circular, in the median plane, with radius respectively $R_{200 \, keV} = 12.924888 \, \text{cm}, R_{1MeV} = 30.107898 \, \text{cm}$ and $R_{5MeV} = 75.754671 \, \text{cm}$). The horizontal axis in this graph is s/C_E : path length over closed orbit circumference at energy E, the vertical axis is the motion excursion

The resulting radial and axial motions over 10 turns are displayed in Fig. 3.41, which also illustrates, for paraxial motion at some reference energy, the energy dependence of the focusing strength (or wave number) and of the motion amplitude.

Table 3.24 Wave numbers, from numerical raytracing (columns denoted "ray-tr."), from theory, and from discrete Fourier transform ('DFT' cols.) from a multi-turn tracking

			$v_R =$			$v_Z =$	
E (MeV)	k_E	ray-tr.	$\sqrt{1+k}$	DFT	ray-tr.	$\sqrt{-k}$	DFT
0.2	-0.03	0.98520	0.9849	0.98513	0.17320	0.1732	0.17321
1	-0.07279	0.96187	0.96292	0.96291	0.26980	0.26979	0.26981
5	-0.20586	0.89083	0.89115	0.89115	0.45326	0.45371	0.45371

An estimate of the wave numbers can be obtained as the inverse of the number of turns per oscillation, namely,

$$v_R = \frac{C_E}{\Delta s_M}\Big|_E$$
 and $v_Z = \frac{C_E}{\Delta s_M}\Big|_E$

with Δs_M the measured distance between two consecutive maxima in the sinusoid of concern in Fig. 3.41, C_E the closed orbit length for the energy of concern. Both quantities are obtained from motion records in zgoubi.plt. This yields the values of Tab. 3.24, where they are compared with the theoretical expectations, namely (Eq. 3.18), $v_R = \sqrt{1 + k}$ and $v_Z = \sqrt{-k}$.

The maximum amplitude of the oscillation is obtained from zgoubi.plt records as well, this yields the results of Tab. 3.25. For comparison, the theoretical values are (Eqs. 3.16, 3.17 with respectively $x_0 = 0$, $x'_0 = T_0$ and $y_0 = 0$, $y'_0 = P_0$) $\hat{Y} = T_0 \frac{R_E}{\sqrt{1+k}}$ and $\hat{Z} = P_0 \frac{R_E}{\sqrt{-k}}$. wherein R_E denotes the closed orbit radius at energy E (for the record: $R_E \equiv R_0$ at energy E = 200 keV, in the foregoing).

Table 3.25 Maximum amplitude of the oscillation, from raytracing (columns denoted "ray-tr.") and from theory. R_E is the closed orbit radius for the energy of concern, $T_0 = P_0 = 0.1$ mrad is the trajectory angle at the origin, positions at the origin are zero

		Ŷ	Ź
E (MeV)	k	ray-tr. $T_0 \frac{R_E}{\sqrt{1+k}}$	ray-tr. $P_0 \frac{R_E}{\sqrt{-k}}$
(MeV)		$(\times 10^{-5})$	$(\times 10^{-5})$
0.2	-0.03	1.3123 1.3125	7.4622 7.4624
1	-0.072787	3.1270 3.1267	1.1160 1.1160
5	-0.20586	8.5010 8.5008	1.6697 1.6697

(c) Comparison with theory.

Figure 3.42 shows the difference between numerical and theoretical vertical motion excursion, using an *ad hoc* gnuplot script. An integration step size $\Delta s = 2$ cm is

used in the numerical integration.



(d) A scan of energy dependence of wave numbers.

A scan of the wave numbers over 200 keV-5 MeV energy range, computing tunes with MATRIX, is performed using the input data file given in Tab. 3.26 (essentially a copy of the input data file of Tab. 3.23, with an INCLUDE accounting for 6 DIPOLES [16, *lookup* INDEX]).

OBJET[KOBJ=5] generates 13 particles with paraxial horizontal, vertical and 2633 longitudinal sampling, proper to allow the computation of the first order transport 2634 coefficients and wave numbers by MATRIX. REBELOTE repeats MATRIX com-2635 putation for a series of different particle rigidities. It is preceded by FIT which finds 2636 the closed orbit. MATRIX includes a PRINT command, which causes the transport 2637 coefficients (and various other outcomes of MATRIX computation) to be logged 2638 in zgoubi.MATRIX.out. This allows producing the graphic in Fig. 3.43 - using the 2639 gnuplot script given at the bottom of Tab. 3.26. 2640

3.4 Solutions of Exercises of Chapter 3: Classical Cyclotron

 Table 3.26
 Simulation input data file: for this wave number scan, the INCLUDE segment is taken from Tab. 3.21

```
Field and derivative dB/dR, as a finction of R
'MARKER' ProbMATRIX_S
'OBJET'
                                                              ! Reference Brho ("BORO" in the users' guide) -> 200keV proton.
! Define 13 particles for MATRIX computation.
! Sampling of the initial coordinates.
! Reference: p[MeV/c]=193.739, Brho[K6.cm]=BORO, kin-E[MeV]=0.2.
64.62444403717985
.001 .01 .001 .01 .001 .00001
12.924888 0. 0. 0. 0. 1.
1 ! IL=2 is necessary under DIPOLE, for step-by-step log of particle data in zgoubi.plt.
6* 60PegSectorR200.inc[#S_60DegSectorR200:#E_60DegSectorR200] ! Six 60 degree sectors.
'FIT'
1
1

2 30 0 [12,80]

1 1e-10

3.1 1 2 #End 0. 1. 0

'MATRIX'
                                                          ! Vary particle's Y0 at OBJET, to have it match its D (=Brho/BORO).
                                                                                                                                   ! Consrain Y_final=Y0.
1 11 PRINT
'REBELOTE'
                                  ! PRINT: log computation outcome data to zgoubi.MATRIX.out, for further plotting.
                                                                                              ! Scan parameter 35 (particle 1's D) in OBJT.
25 0.1 0 1
1
OBJET 35 1:5.00639
'SYSTEM'
1
gnuplot < ./gnuplot_MATRIX_Qxy.gnu
'MARKER' ProbMATRIX_E
'END'
```

A gnuplot script to obtain Fig. 3.43:

 $+v_{y}^{2})^{1/2}$

v_x, (v_x²

Fig. 3.43 A scan of the energy dependence of the horizontal and vertical wave numbers. Markers are from raytracing, solid lines are from theory (Eq. 3.18). The figure also shows that the raytracing yields $v_R^2 + v_y^2 = 1$, $\forall E$, as expected

