
Super-FOFO + Dipole Simulations: Status Report II

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Emittance Exchange with Super FOFO Lattice

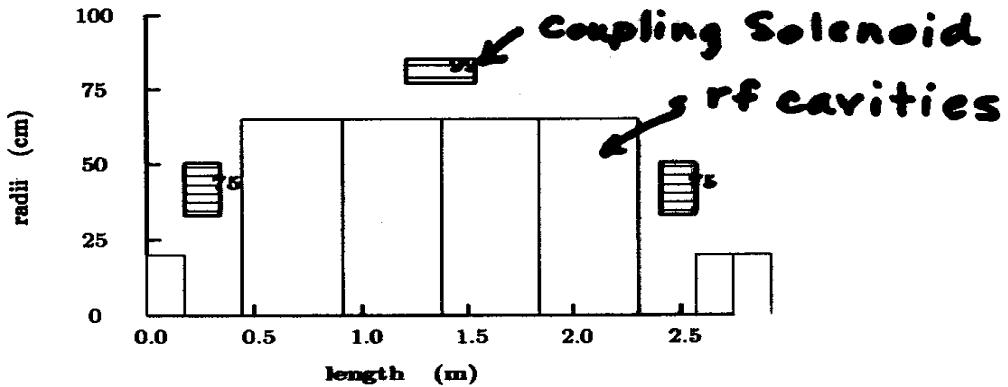
DRAFT

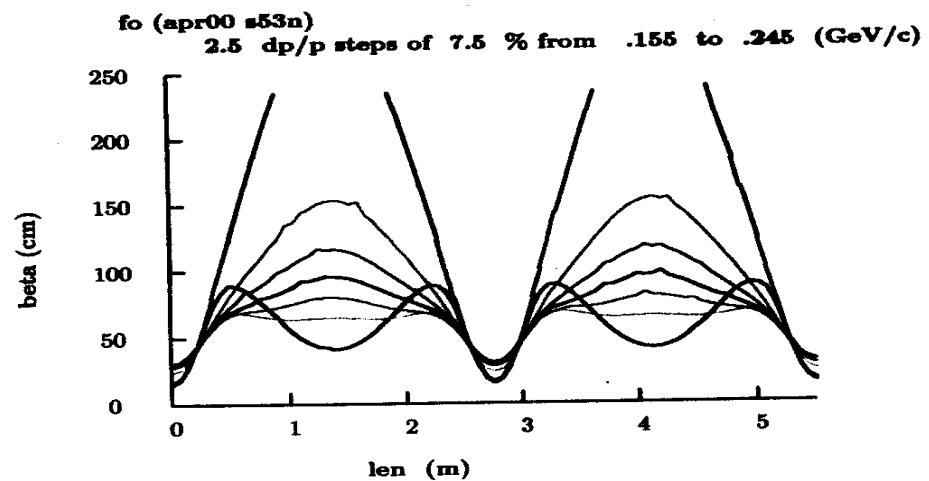
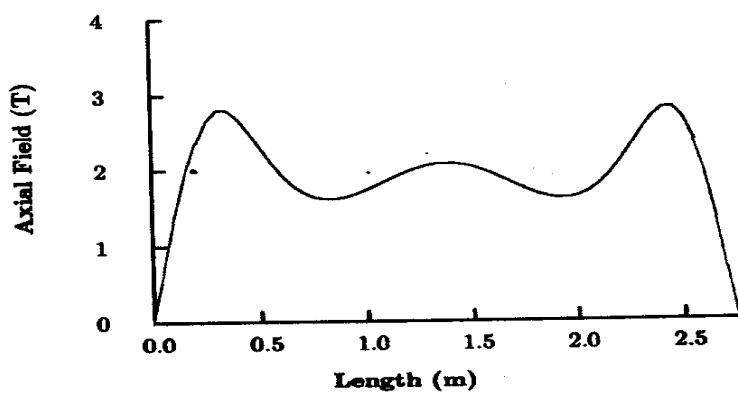
R.B. Palmer 8/17/00

0.1 Super FOFO

Consider modifying the initial super FOFO lattice as used now in the Study 2 simulations. The cell length is 2.75 m, the β at the absorber is $\approx .3$ m and it accepts an emittance ($\epsilon_n \approx 10$ mm rad), and a momentum spread of $+/- 22\%$.

Figs (a) and (b) show the coils and axial field. Fig (c) shows the beta functions for a range of momenta, as a function of position down the cell. Fig (d) shows the phase advance per cell vs momentum (it is relatively linear between limits at π and 2π where there are resonances), and the maximum and minimum β 's. It is seen that the β at the absorber (red) is relatively flat, but the β at the center is strongly momentum dependent, rising from about .4 m at one limit and 3 m at the other (this will turn out to be important).





- The cooling is done in six sections with steadily decreasing β 's. This is done to maximize the cooling rate. Too small a β at a given emittance results in too large divergence angles and particle loss. Too large a β gives small divergence angles and a greater relative emittance growth from coulomb scattering. The best β scales down with the emittance and is always such as to give a certain constant divergence angle ($\sigma_x' = \sigma_y' \approx 0.1$).

2.11 Cooling Latices

2.11.1 Introduction

The cooling is done in six sections with steadily decreasing β 's. This is done to maximize the cooling rate. Too small a β at a given emittance results in too large divergence angles and particle loss. Too large a β gives small divergence angles and a greater relative emittance growth from coulomb scattering. The best β scales down with the emittance and is always such as to give a certain constant divergence angle ($\sigma_x' = \sigma_y' \approx 0.1$).

The six sections are made from two different physical latices [(1) and (2)], with three different current setting in each: 1,1 1,2 1,3 in the first lattice, and 2,1 2,2 2,3, in the second. The final cooling section (2,3) is further broken into 2 parts (2,3a) and (2,3b) that differ only in their window sizes and thicknesses.

The lengths of the sections are:

	length m	from target m
cool 1,1	$6 \times 2.75 = 16.5$	444.9
cool 1,2	$6 \times 2.75 = 16.5$	461.4
cool 1,3	$6 \times 2.75 = 16.5$	477.9
cool 2,1	$14 \times 1.65 = 23.1$	501
cool 2,2	$10 \times 1.65 = 16.5$	517.5
cool 2,3a	$16 \times 1.65 = 26.4$	543.9
cool 2,3b	$16 \times 1.65 = 26.4$	570.3

The latices used have been named "Super FOFO". The "FOFO" refers to the basic sequence of alternating solenoids, that focus the beam and generate β , and thus beam size, minima between the solenoids. The "Super" part, proposed by A. Sessler, is the replacement of the simple alternating solenoids with alternating, but more complex, solenoid systems. In this case the systems consist of strong short "focusing" solenoids at either end, and a weaker "coupling" fields between them.

An example of the fields (for the last part) is given below

BEFORE COOLING

+0.1

P_{x0}

-0.1

-0.2 -0.1 x → 0.1 0.2

+0.1 -

P_Y

-0.1

-0.2 -0.1 y → 0.1 0.2

```

+
+
+ + + 2 +
+ 2 + 2++ + + +
+ + 2+2+2 3 + 2
+++ 223 22+ 2+
+ + 323322244322+ + +
2 5+ 3433825 62323+++22
+ 22+3+5656598664253+2
2 25+3543b877b62+++ 2+2 +
+++ + +2 3 237854769447+++2+
+ + +32+5+56+4++22++ +
22+ +3 42+32 +
+ + +++ 2 3+ 2 + + +
+ + + + +2 + +
+ + + + +
+ 2 +

```

```

+
+
+
+ + + + + +
+ +55 + + + +2 + +
2 23+ 222+3+ 2 +2
+ + 3 43+4++5++2 + +
422+++442674763335325+ +
2 334333b7685254+2 + 2
++242446858a76633+++3+++ + +
+ + 2366756a6438342 ++ +
+ + 32332+542324++ +22+ +
+ + 2+ + + 2+ 2+
+ + 22 + + + 2+ ++
+ + + + + +
+ + + + +
+
```

ICOOL	2.11
plot	= 1
var(x)	= x
dest(x)	= 1
var(y)	= Px
dest(y)	= 1
lo x	= -0.200E+00
hi x	= 0.200E+00
step x	= 0.100E-01
lo y	= -0.100E+00
hi y	= 0.100E+00
step y	= 0.100E-01
contents	= 500
under x	= 0
over x	= 0
under y	= 0
over y	= 0

ICOOL	2.11
plot	= 2
var(x)	= Y
dest(x)	= 1
var(y)	= Py
dest(y)	= 1
lo x	= -0.200E+00
hi x	= 0.200E+00
step x	= 0.100E-01
lo y	= -0.100E+00
hi y	= 0.100E+00
step y	= 0.100E-01
contents	= 500
under x	= 0
over x	= 0
under y	= 0
over y	= 0

AFTER COOLING

+0.1

\vec{P}_x

-0.1

-0.2 -0.1 0 1 0.2

$x \rightarrow$

+0.1

\vec{P}_y

-0.1

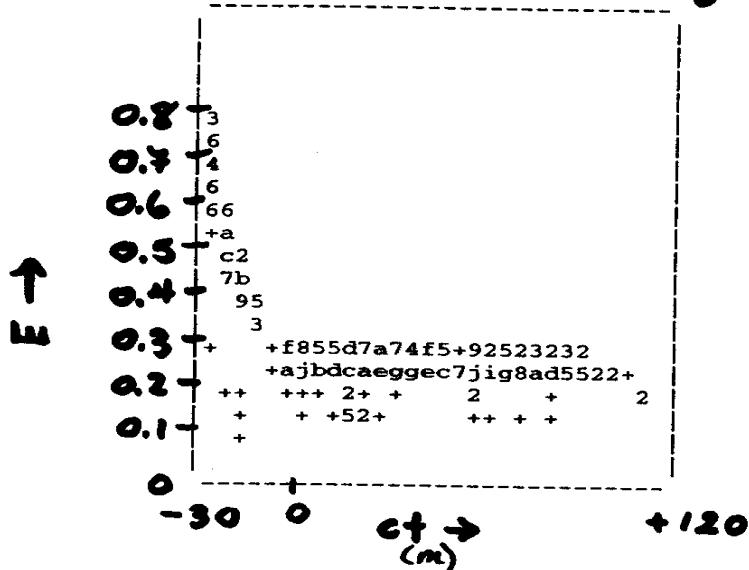
-0.2 -0.1 0 1 0.2

$y \rightarrow$

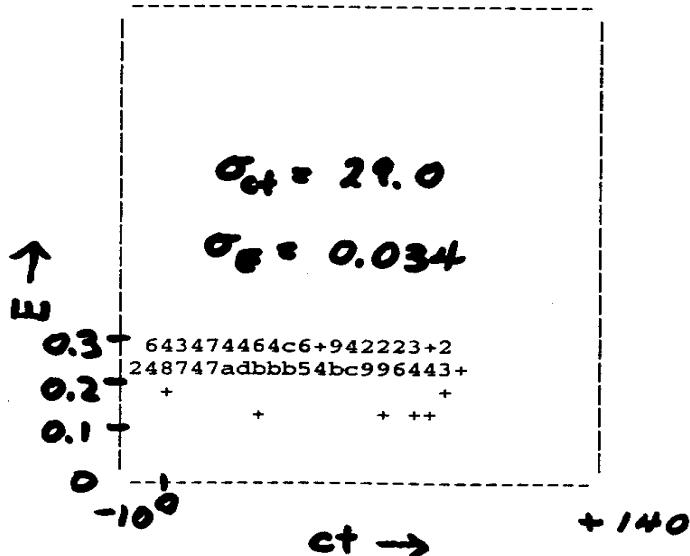
ICOOL	2.11
plot	= 18
var(x)	= X
dest(x)	= 1042
var(y)	= Px
dest(y)	= 1042
lo x	= -0.200E+00
hi x	= 0.200E+00
step x	= 0.100E-01
lo y	= -0.100E+00
hi y	= 0.100E+00
step y	= 0.100E-01
contents	= 200
under x	= 0
over x	= 0
under y	= 0
over y	= 0

ICOOL	2.11
plot	= 19
var(x)	= Y
dest(x)	= 1042
var(y)	= Py
dest(y)	= 1042
lo x	= -0.200E+00
hi x	= 0.200E+00
step x	= 0.100E-01
lo y	= -0.100E+00
hi y	= 0.100E+00
step y	= 0.100E-01
contents	= 200
under x	= 0
over x	= 0
under y	= 0
over y	= 0

Before Cooling



After 2,1



```

ICOOL          2.11
-----
plot          =      3
var(x)        =      ct
dest(x)       =      1
var(y)        =      E
dest(y)       =      1

lo x          = -0.300E+02
hi x          = 0.120E+03
step x        = 0.500E+01
lo y          = 0.000E+00
hi y          = 0.100E+01
step y        = 0.500E-01

contents      =      500
under x       =      0
over x        =      0
under y       =      0
over y        =      0

```

```

ICOOL          2.11

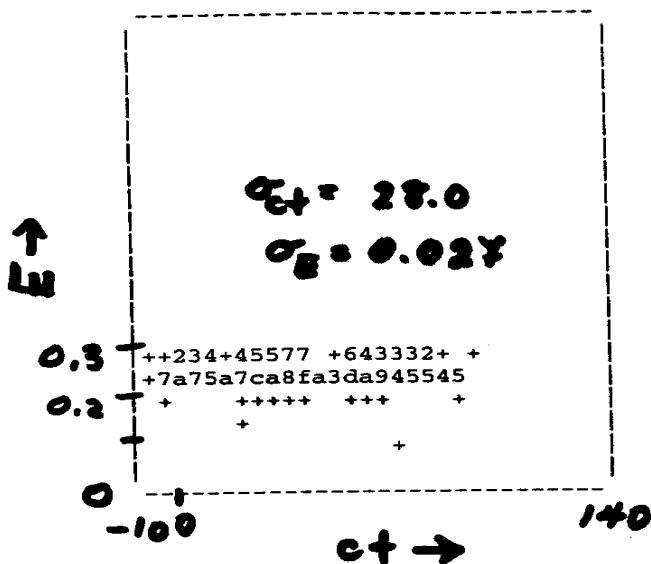
plot      =    14
var(x)   =   ct
dest(x)  =   496
var(y)   =   E
dest(y)  =   496

lo x    = -0.100E+02
hi x    = 0.140E+03
step x  = 0.500E+01
lo y    = 0.000E+00
hi y    = 0.100E+01
step y = 0.500E-01

contents = 248
under x = 9
over x  = 0
under y = 0
over y  = 0

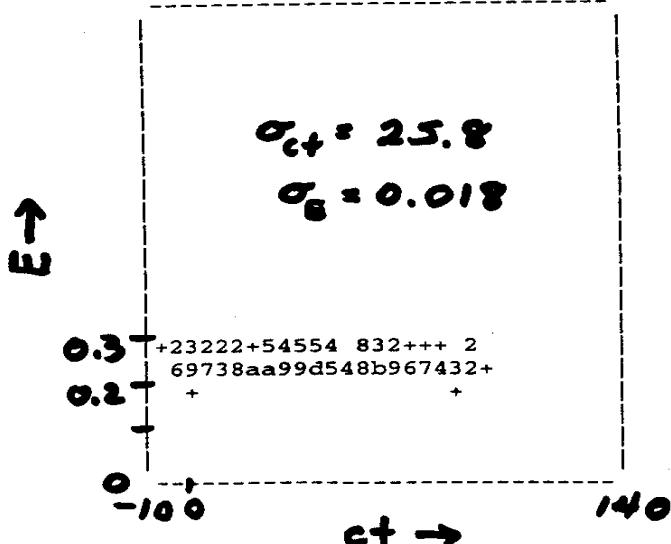
```

After 2,2

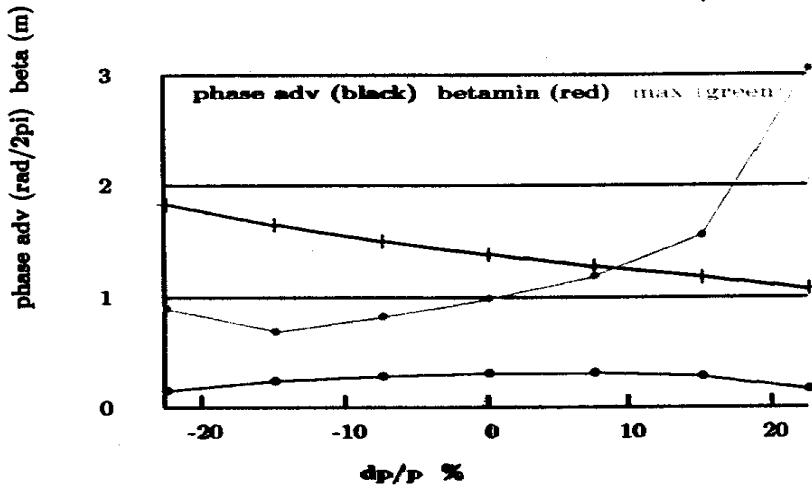


```
ICOOL          2.11
plot           =    17
var(x)        =    ct
dest(x)       =   626
var(y)        =    E
dest(y)       =   626
lo x          = -0.100E+02
hi x          =  0.140E+03
step x        =  0.500E+01
lo y          =  0.000E+00
hi y          =  0.100E+01
step y        =  0.500E-01
contents      =    236
under x       =      6
over x        =      0
under y       =      0
over y        =      0
```

After Cooling

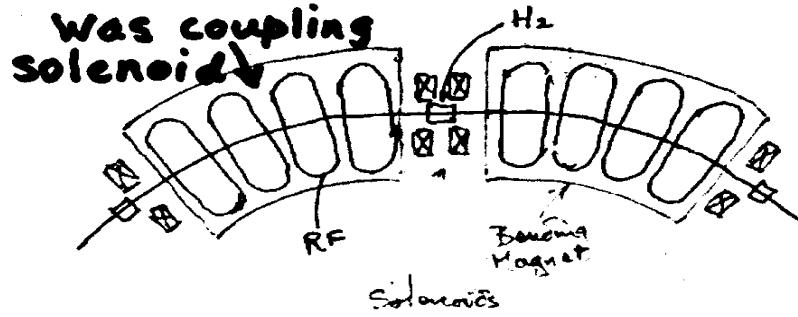


```
ICOOL          2.11
plot           =    20
var(x)        =    ct
dest(x)       =   1042
var(y)        =    E
dest(y)       =   1042
lo x          = -0.100E+02
hi x          =  0.140E+03
step x        =  0.500E+01
lo y          =  0.000E+00
hi y          =  0.100E+01
step y        =  0.500E-01
contents      =    200
under x       =      0
over x        =      0
under y       =      0
over y        =      0
```



0.2 Add a Bend

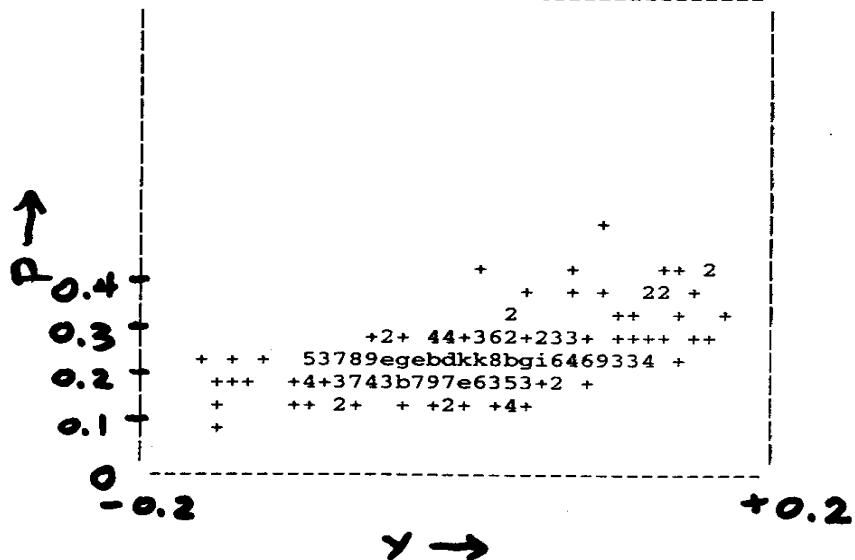
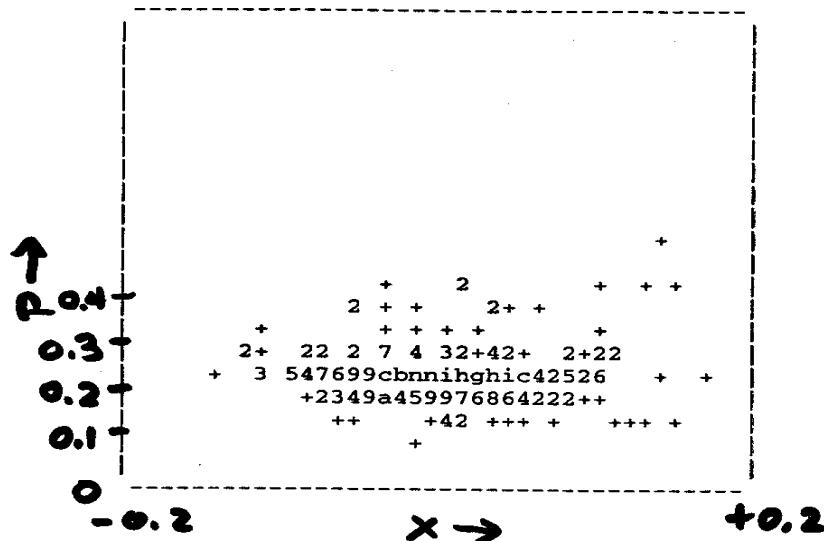
In order to use this lattice for emittance exchange we must add some bending to generate dispersion at the absorber. The absorber can then be made wedge like and the required exchange obtained. Consider putting the bend in the center of the cell. It now falls on top of the coupling solenoid, which is uncomfortable. But bending magnets have focussing: perhaps it can replace this solenoid.



$B_y = 0.12T$ Same
After 1st Absorber

| ICOOL

2.11



```

plot      =    7
var(x)   =    X
dest(x)  =    17
var(y)   =    Pm
dest(y)  =    17

lo x    = -0.200E+00
hi x    =  0.200E+00
step x  =  0.100E-01
lo y    =  0.000E+00
hi y    =  0.100E+01
step y  =  0.500E-01

contents =  406
under x =  0
over x  =  0
under y =  0
over y  =  0

```

| ICOOL 2.11

```

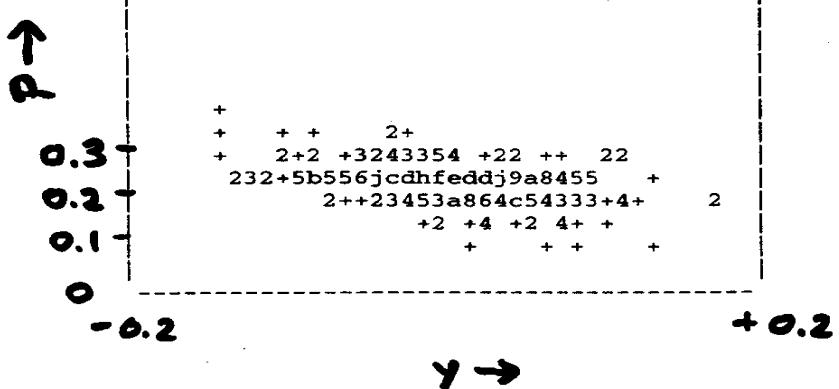
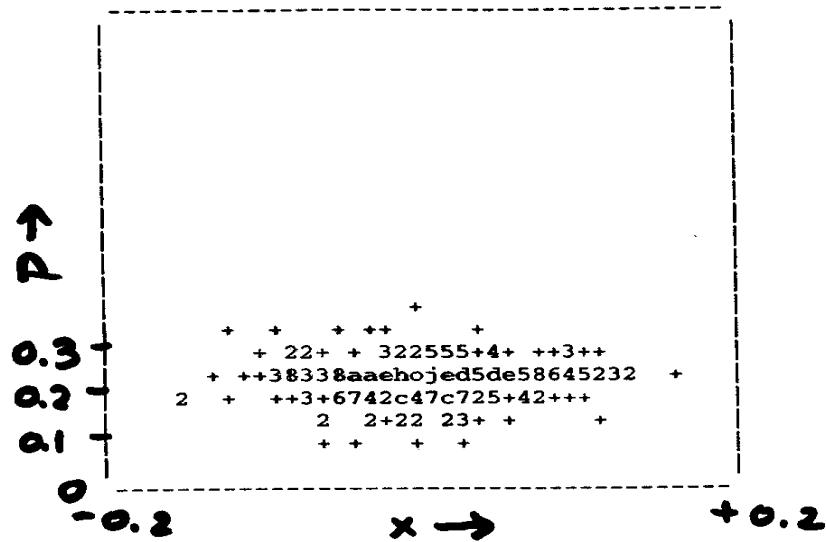
plot      =    8
var(x)   =    Y
dest(x)  =    17
var(y)   =    Pm
dest(y)  =    17

lo x    = -0.200E+00
hi x    =  0.200E+00
step x  =  0.100E-01
lo y    =  0.000E+00
hi y    =  0.100E+01
step y  =  0.500E-01

contents =  406
under x =  0
over x  =  0
under y =  0
over y  =  0

```

$B_y = 0.12 T$ Same
After 2nd Absorber



```
ICOOL      2.11
plot       =    9
var(x)    =   X
dest(x)   =   35
var(y)    =   Pm
dest(y)   =   35

lo x     = -0.200E+00
hi x     =  0.200E+00
step x  =  0.100E-01
lo y     =  0.000E+00
hi y     =  0.100E+01
step y  =  0.500E-01

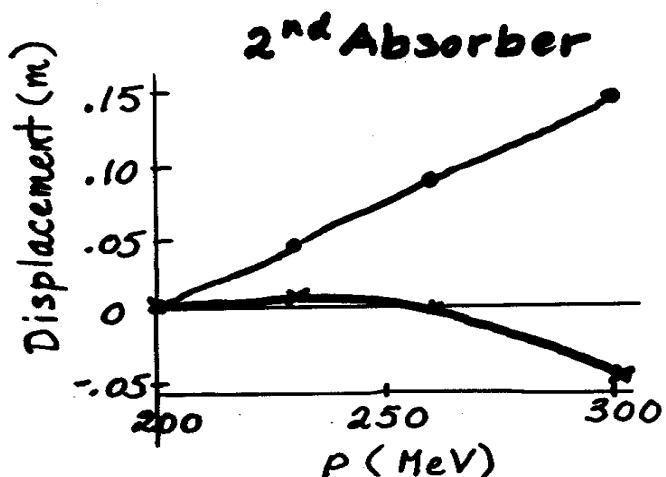
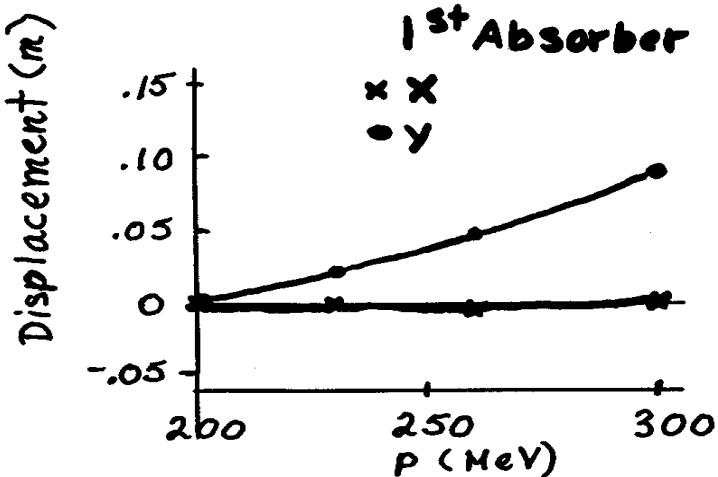
contents  =  374
under x  =   0
over x   =   0
under y  =   0
over y   =   0
```

```
ICOOL      2.11
plot       =    10
var(x)   =   Y
dest(x)  =   35
var(y)   =   Pm
dest(y)  =   35

lo x     = -0.200E+00
hi x     =  0.200E+00
step x  =  0.100E-01
lo y     =  0.000E+00
hi y     =  0.100E+01
step y  =  0.500E-01

contents  =  374
under x  =   0
over x   =   0
under y  =   0
over y   =   0
```

CONSTANT $B_y = 0.12 T$ FLIPPED



CONSTANT $B_y = 0.12 T$ SAME

