

Definitions

- What is FPP?
- What is PTC?

Cornell-KEK support: Profs. Dave Rubin and Makoto Tobiyama

History

- Start FPP around 1998/Start PTC around 2000
- Put into MAD-X with help of Frank Schmidt and Eric Macintosh completed by 2002
- Spin added around 2007 in FPP and PTC (Barber pressure)
- Magnet modulation added in FPP and PTC (Schmidt pressure)
- Continuous collaboration with David Sagan since the beginning: FPP/PTC fully added in the early 2010 in BMAD
- PTC painfully added in ORBIT under the guidance Alexander Molodozhentsev (It is unfortunately a “hack” but it worked)
- Made SAD compatible under the guidance of David Sagan
- Analysis tools fully “complexified”, old Fortran 77 routines are obsolescent (Deniau pressure)

FPP : taylor series + analysis

- FPP overloads the “DA” or “TPSA” package of Berz
- It also overloads my Fortran 77 analysis tools (Normal Form etc...)
- It has a polymorphic type (REAL_8): real and Taylor
- Since 2011, it has a new analysis package based on a complex TPSA
- Normal form with magnet modulation, radiative maps, coasting beam, spin with $SO(3)$ and quaternions
- Maps of all sorts and kinds: Lie transforms, generating functions, partially inverted maps

PTC

- It is (generally symplectic) integrator: a tracking code
- Novel structures: physical beam lines are not identified to tracking structures: figure 8, colliders, recirculators
- Has the concept of a fibre to deal with the novel structures
- It uses a polymorphic type so anything can become a Taylor series: phase space, spin, a_n , b_n , etc..
- Radiation, beam envelopes, magnet modulation, spin, etc...
- It is inside MAD-X (Schmidt/Macintosh)
- It is inside BMAD (Sagan)
- Hacked job in ORBIT (Modolzhentsev)
- Pretty much supports SAD magnets (as does BMAD, Sagan pressure)

FPP and PTC: New?

- FPP is new technology but nothing scientifically new

Explained in a Fermilab Lecture in the late 1980s

- 1) Berz Taylor TPSA overloaded
- 2) Polymorph type created: can be real or Taylor at execution time (Bengtsson)
- 3) A complex TPSA with normal form for transverse, spin (quaternion), magnet modulation.

- PTC has a different beam structure: New idea

Realization of a fibre bundle structure (idea of Hirata and Forest 1992)

- 1) Trackable beam line is succession of containers called “fibres”
- 2) The fibre is discrete equivalent of the “s” variable
- 3) Each fibre points to a physical magnet
- 4) Magnets are not cloned as in the real world

Production of Taylor Maps (Sagan)

- It is possible to replace a magnet by a Taylor map **from PTC or elsewhere**
- It is possible to produce Taylor maps for tracking and incorporate it elsewhere
- Symplectic/Unitary algorithm are possible on the Taylor maps (Orbital-Spin)

FPP alone

- a_scratch_size.f90
- b_da_arrays_all.f90
- c_dabnew.f90
- cb_da_arrays_all.f90
- cc_dabnew.f90
- Ci_tpsa.f90
- d_lielib.f90
- h_definition.f90
- i_tpsa.f90
- j_tpsalie.f90
- k_tpsalie_analysis.f90
- l_complex_taylor.f90
- m_real_polymorph.f90
- n_complex_polymorph.f90
- o_tree_element.f90

Mini tracking code



y_multipole.f90

Main program



Z_FPP_normal_form.f90

One-turn Taylor map computation

```
program FPP
  use my_multipole
  implicit none
  integer order,nd
  real(dp) closed_orbit(6)
  ! polymorph
  type(real_8) x(6)
  longprint = .false.
  call set_multipole
  call find_closed_orbit(closed_orbit)
  write(6,*) " closed orbit ", closed_orbit

  order=8

  nd=1
  call c_init_all(order,nd,0,0)
  call alloc(x)

  x=closed_orbit

call print(x)

  call multipole(x)

  call print(x)

  x=closed_orbit

  x(1)=closed_orbit(1)+dz_8(1)
  x(2)=closed_orbit(2)+dz_8(2)
  call print(x)
  call multipole(x)

  call print(x)
end program FPP
```

```
program FPP_id
  use my_multipole
  implicit none
  integer order,nd
  real(dp) closed_orbit(6)
  ! polymorph
  type(real_8) x(6)
  ! TPSA Maps
  type(c_damap) id,one_turn
  longprint = .false.
  call set_multipole
  call find_closed_orbit(closed_orbit)
  write(6,*) " closed orbit ", closed_orbit

  order=8

  nd=1
  call c_init_all(order,nd,0,0)
  call alloc(x);call alloc(id,one_turn)

  x=closed_orbit

  call print(x)

  call multipole(x)

  call print(x)

  id=1
  x= closed_orbit + id

  call print(x)

  call multipole(x)

  one_turn=x
  call print(one_turn)
end program FPP_id
```

```
subroutine multipolep(x)
  implicit none
  type(real_8) x(2)
  x(1)=x(1)+Lpol*x(2)
  x(2)=x(2)-bnpol(1)-bnpol(2)*x(1) &
    -bnpol(3)*x(1)**2-bnpol(4)*x(1)**3
end subroutine multipolep
```


program FPP

```

closed orbit -1.127016653792583E-002
-7.531047606624180E-018
0
-1.127016653792583E-002
-7.531047606624180E-018
-1.127016653792583E-002
-7.535205098774256E-018

Properties, NO = 8, NV = 2, INA = 26
*****

0 -0.1127016653792583E-01 0 0
1 1.0000000000000000 1 0

Properties, NO = 8, NV = 2, INA = 27
*****

0 -0.7531047606624180E-17 0 0
1 1.0000000000000000 0 1

Properties, NO = 8, NV = 2, INA = 26
*****

0 -0.1127016653792583E-01 0 0
1 1.0000000000000000 1 0
1 0.1000000000000000 0 1

Properties, NO = 8, NV = 2, INA = 27
*****

0 -0.7535205098774256E-17 0 0
1 -0.7745966692414834E-01 1 0
1 0.9922540333075851 0 1
2 -1.0000000000000000 2 0
2 -0.2000000000000000 1 1
2 -0.1000000000000000E-01 0 2

```

program FPP_id

```

closed orbit -1.127016653792583E-002 -7.531047606624180E-018
0
-1.127016653792583E-002
-7.531047606624180E-018
-1.127016653792583E-002
-7.535205098774256E-018

Properties, NO = 8, NV = 2, INA = 25
*****

0 -0.1127016653792583E-01 0 0
1 1.0000000000000000 1 0

Properties, NO = 8, NV = 2, INA = 31
*****

0 -0.7531047606624180E-17 0 0
1 1.0000000000000000 0 1

2 Dimensional DA map (around chosen orbit in map%x0)

Properties, NO = 8, NV = 2, INA = 161
*****

0 -0.1127016653792583E-01 0.0000000000000000 0 0
1 1.0000000000000000 0.0000000000000000 1 0
1 0.1000000000000000 0.0000000000000000 0 1

Properties, NO = 8, NV = 2, INA = 160
*****

0 -0.7535205098774256E-17 0.0000000000000000 0 0
1 -0.7745966692414834E-01 0.0000000000000000 1 0
1 0.9922540333075851 0.0000000000000000 0 1
2 -1.0000000000000000 0.0000000000000000 2 0
2 -0.2000000000000000 0.0000000000000000 1 1
2 -0.1000000000000000E-01 0.0000000000000000 0 2

No Spin Matrix
c_quaternion is identity
No Stochastic Radiation

```

Normal form

```
program Z_FPP_normal_form
```

```
use my_multipole
implicit none
integer order,nd
real(dp) closed_orbit(6)
! polymorph
type(real_8) x(6)
! TPSA Maps
type(c_damap) id,one_turn,diagonal,rotation
type(c_normal_form) normal_form
longprint = .false.
call set_multipole
call find_closed_orbit(closed_orbit)
write(6,*) " closed orbit ", closed_orbit
```

```
order=3
```

```
nd=1
call c_init_all(order,nd,0,0)
call alloc(x);call alloc(id,one_turn);
call alloc(normal_form);
call alloc(diagonal,rotation)
```

```
id=1
x= closed_orbit + id
```

```
call multipole(x)
```

```
one_turn=x
```

```
call c_normal(one_turn,normal_form)
```

```
rotation=normal_form%atot**(-1)*one_turn*normal_form%atot
call clean(rotation,rotation,1.d-6)
call print(rotation)
diagonal=c_phasor()*(-1)*rotation*c_phasor()
call clean(diagonal,diagonal,1.d-6)
call print(diagonal)
```

```
end program Z_FPP_normal_form
```

```
closed orbit -1.127016653792583E-002 -7.531047606624180E-018
0
1 0.99612701665379 0.87925915931623E-01
2 0.99612701665379 -0.87925915931623E-01
The order of the planes has been guessed using the algorithm in c_locate_planes
Hopefully it is correct! Please check!
Order guessed -> 1
```

```
2 Dimensional DA map (around chosen orbit in map%x0)
```

```
Properties, NO = 3, NV = 2, INA = 258
*****
```

1	0.9961270166537924	0.0000000000000000	1	0
1	0.8792591593162297E-01	0.0000000000000000	0	1
3	0.6114646075503689	0.0000000000000000	3	0
3	-6.927382090420237	0.0000000000000000	2	1
3	0.6114646075504240	0.0000000000000000	1	2
3	-6.927382090420256	0.0000000000000000	0	3

```
Properties, NO = 3, NV = 2, INA = 259
*****
```

1	-0.8792591593162302E-01	0.0000000000000000	1	0
1	0.9961270166537927	0.0000000000000000	0	1
3	6.927382090420240	0.0000000000000000	3	0
3	0.6114646075503596	0.0000000000000000	2	1
3	6.927382090420270	0.0000000000000000	1	2
3	0.6114646075503209	0.0000000000000000	0	3

```
No Spin Matrix
c_quaternion is identity
No Stochastic Radiation
```

```
2 Dimensional DA map (around chosen orbit in map%x0)
```

```
Properties, NO = 3, NV = 2, INA = 243
*****
```

1	0.9961270166537926	-0.8792591593162299E-01	1	0
3	0.6114646075503567	6.927382090420251	2	1

```
Properties, NO = 3, NV = 2, INA = 244
*****
```

1	0.9961270166537926	0.8792591593162299E-01	0	1
3	0.6114646075503567	-6.927382090420251	1	2

```
No Spin Matrix
c_quaternion is identity
No Stochastic Radiation
```

FPP + PTC

- a_scratch_size.f90
- b_da_arrays_all.f90
- c_dabnew.f90
- cb_da_arrays_all.f90
- cc_dabnew.f90
- Ci_tpsa.f90
- d_lielib.f90
- h_definition.f90
- i_tpsa.f90
- j_tpsalie.f90
- k_tpsalie_analysis.f90
- l_complex_taylor.f90
- m_real_polymorph.f90
- n_complex_polymorph.f90
- o_tree_element.f90

← FPP

PTC →

- Sa_extend_poly.f90
- Sb_sagan_pol_arbitrary.f90
- Sc_euclidean.f90
- Sd_frame.f90
- Se_status.f90
- Sf_def_all_kinds.f90
- Sg_sagan_wiggler.f90
- Sh_def_kind.f90
- Si_def_element.f90
- Sk_link_list.f90
- Sl_family.f90
- Sm_tracking.f90
- Sma0_beam_beam_ptc.f90
- Sma_multiparticle.f90
- Sn_mad_like.f90
- So_fitting.f90
- Sp_keywords.f90
- Spb_fake_gino_sub.f90
- Sq_orbit_ptc.f90
- Sr_spin.f90
- Sra_fitting.f90
- Ss_fake_mad.f90
- St_pointers.f90

User defined module →

Main program →

- y_als_lattice.f90
- z_spin_phase_advance_isf_for_sad.f90

Courant-Snyder Loop

```

state=nocavity0+SPINO
CALL FIND_ORBIT(ALS,CLOSED,1,STATE,c_1d_5)
state=nocavity0
no=2
call init_all(STATE,no,1)

!!!! Polymorphic probe is created in the usual manner
XS0=CLOSED
ID_S=1
XS=XS0+ID_S

!!!! get spin polymorphic probe after one turn
CALL propagate(ALS,XS,+STATE,FIBRE1=1)

! Copy probe_8 into a complex damap
c_map=XS

! Normalize it c_map = C_n%atot^(-1) o R o C_n%atot
call c_normal(c_map,c_n,dospin=my_true,phase=phase,nu_spin=nu_spin)

U=c_n%Atot ! Non-descript U exiting normal form
call c_full_canonise(U,U_c,D,F,A,b,R,phase,nu_spin)

p => als%start

do i=1,als%n

CALL propagate(ALS,XS,+STATE,FIBRE1=i,fibre2=i+1)

xs0=xs ! Saving orbit ! (9b)
U=XS ! copying in map ! (9c)
! U = U_c o R = D o f o A o b o R
call c_full_canonise(U,U_c,D,F,A,b,R,phase,nu_spin)

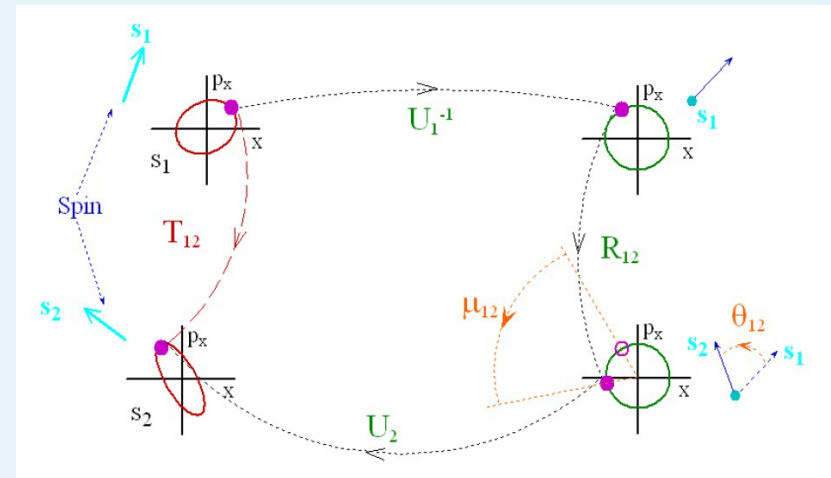
```

COMPUTE SOMETHING HERE

```

XS=XS0+U_c
p=>p%next
enddo

```



Compute something?

```
write(mft,*) "position, Element ", i, p%mag%name

betax_1=(U_c%v(1).sub.'1000')**2+(U_c%v(1).sub.'0100')**2
betax_2=(U_c%v(1).sub.'0010')**2+(U_c%v(1).sub.'0001')**2

if(use_quaternion) call makeso3(D)
ISF=2
ISF=D%s*ISF

write(mft,*) " 2< x^2 > "
call print(fonction_FLOQUET,mft)

write(mft,*) " Ripken Beta_x_1 Beta_x_2 ",betax_1,betax_2

Write(mft,*) " ISF vector n "
call print(ISF,mft,prec)

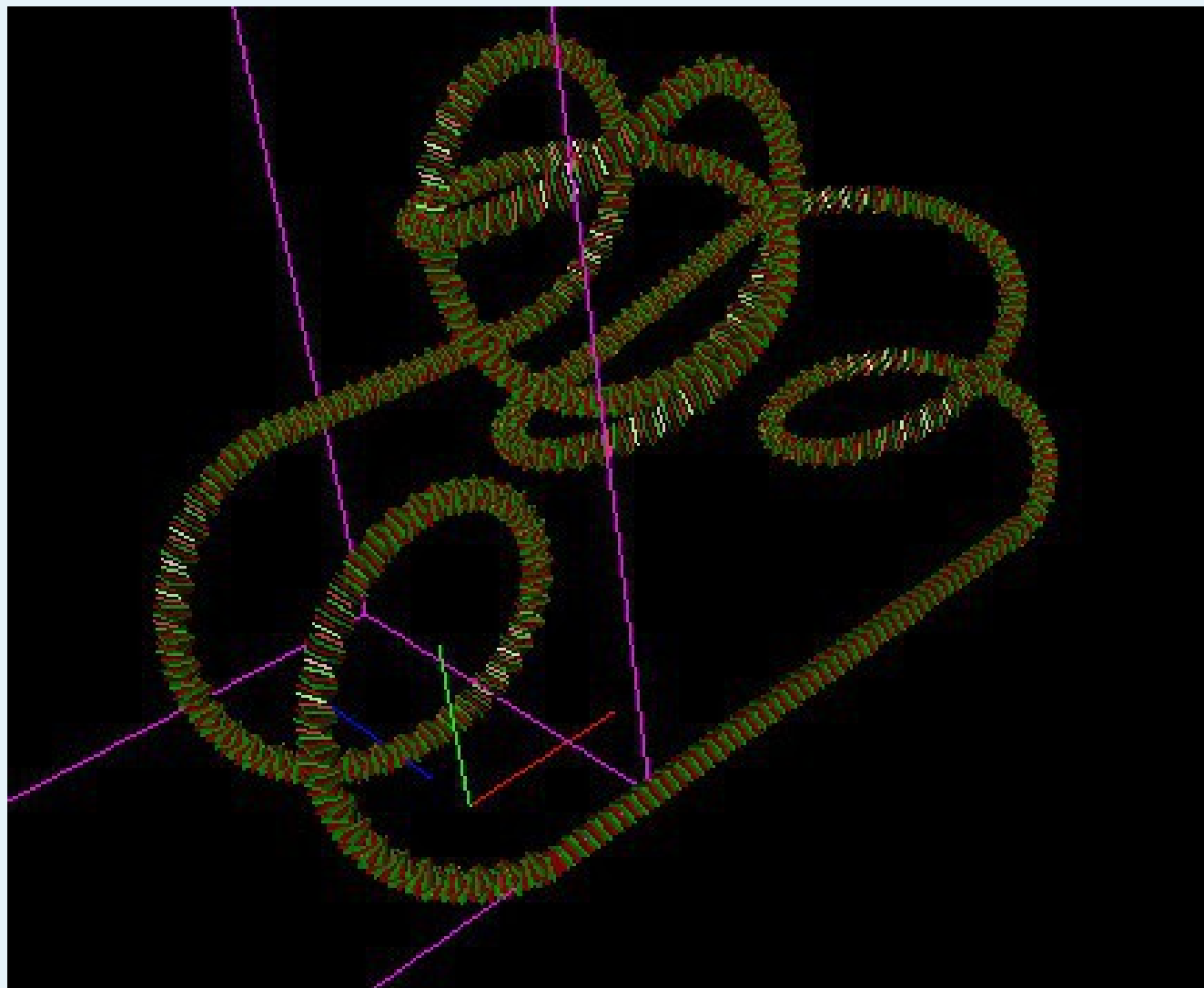
write(mft,*) " phase x"
call print(phase(1),mft,prec)

write(mft,*) " phase y"
call print(phase(2),mft,prec)

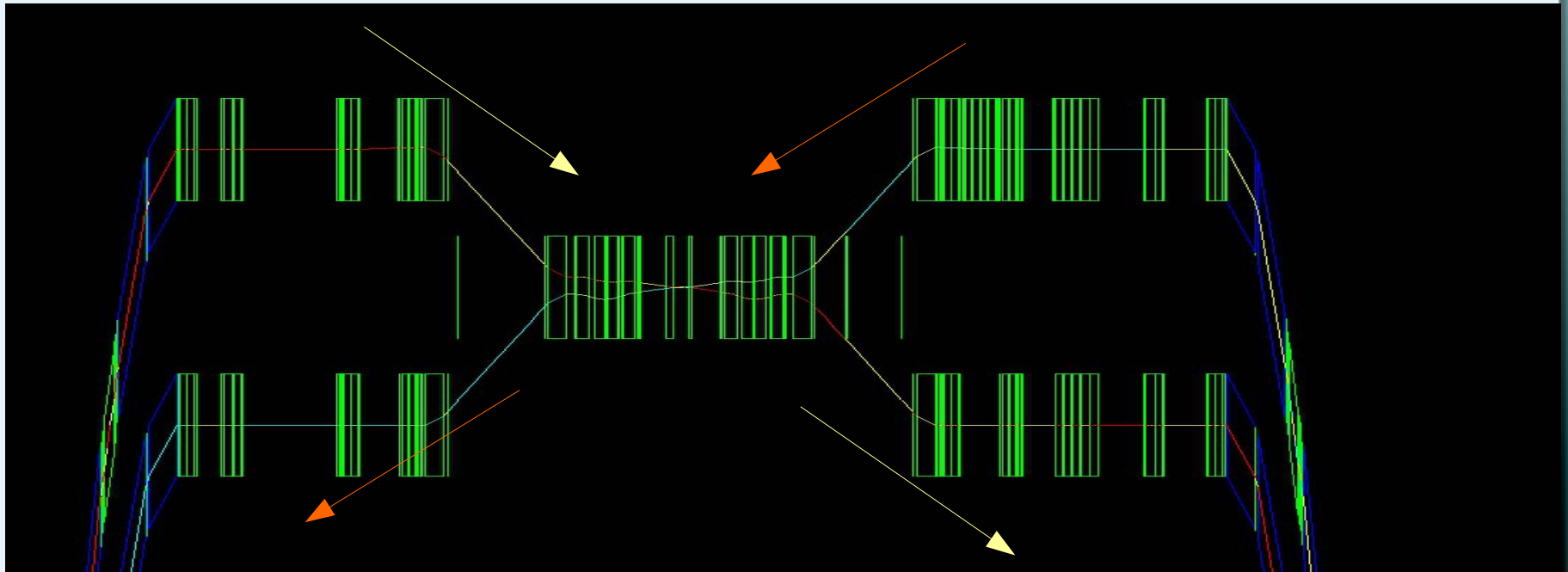
if(state%nocavity) then
write(mft,*) " Time "
call print(phase(3),mft,prec)
endif

write(mft,*) " phase spin"
call print(nu_spin,mft,prec)
```

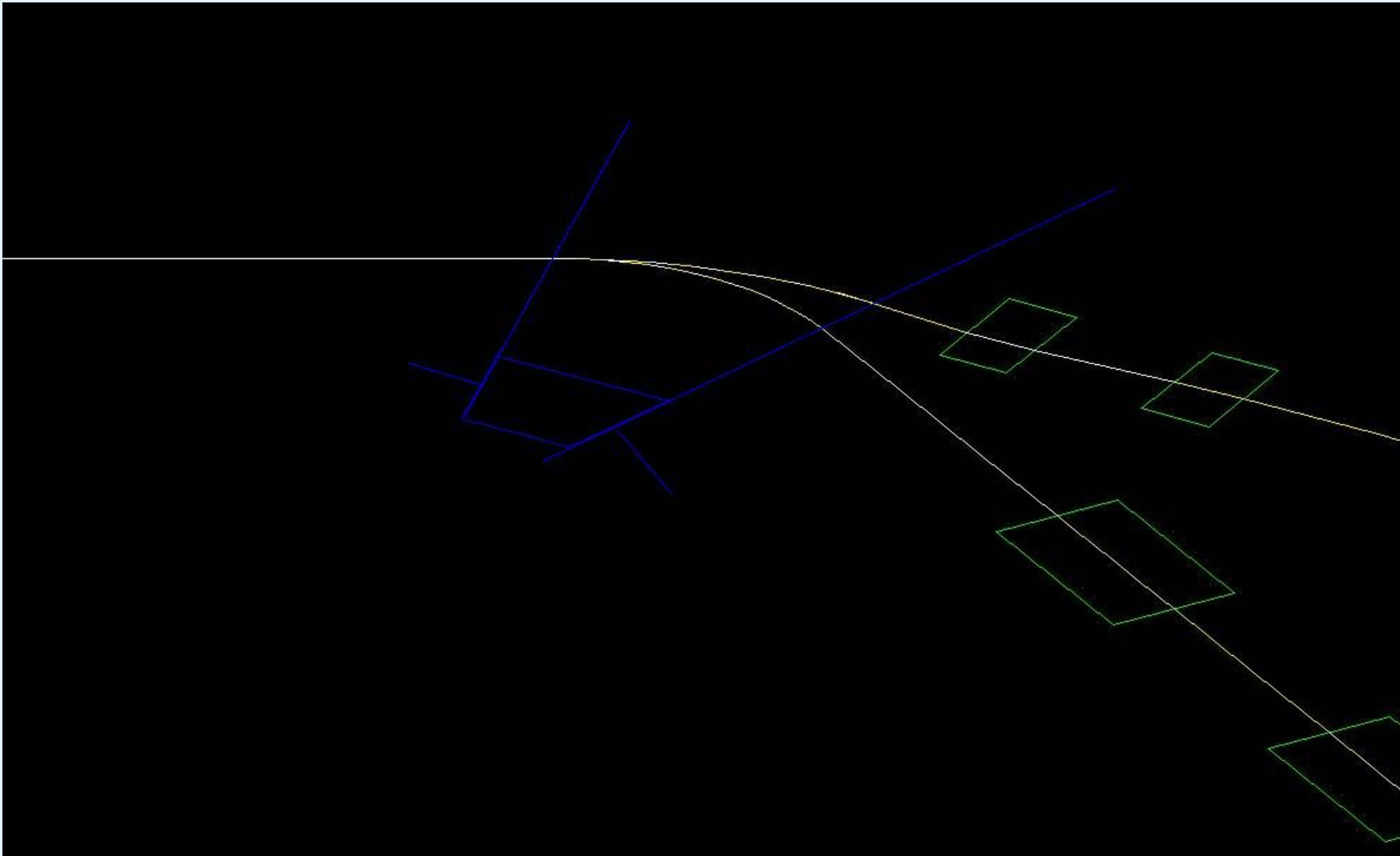
Some Strange Machida/Mohri Object



LHC



Some Recirculator



Future

- 1) Putting Laurent Deniau new TPSA in FPP
- 2) Of course anyone can incorporate FPP or PTC in their code
- 3) Perhaps one day, the full geometrical part of PTC will be integrated in a CAD environment or in a full code (BMAD is a prototype)
- 4) Always opened to perturbation theory applications: Qiang (LBL), Zhou (KEK), etc...