

John Talman

Unified Accelerator
Libraries
(UAL)

ETEAPOT Status
With EDM Experiment
Emphasis

E(lectric)**TEAPOT** Partially
Cloned From(magnetic)TEAPOT

(Thin Element Accelerator
Program For Optics And
Tracking)

Little electric/magnetic distinction for
Quads and Sexts (e.g. fringe field).

“Exactly” solve Bends (treating them as
thick elements) in ETEAPOT(as contrasted
with modeling them as kicks in UAL/TEAPOT).
Correct for deviant radial field dependence
using kicks.

Status:

Physical Ring	—	Stable
Lattice	—	Stable
Survey	—	Finished
Orientation	—	Stabilizing
Equations	—	Stabilizing
Code	—	Stabilizing

Probably biased to the
optimistic side!

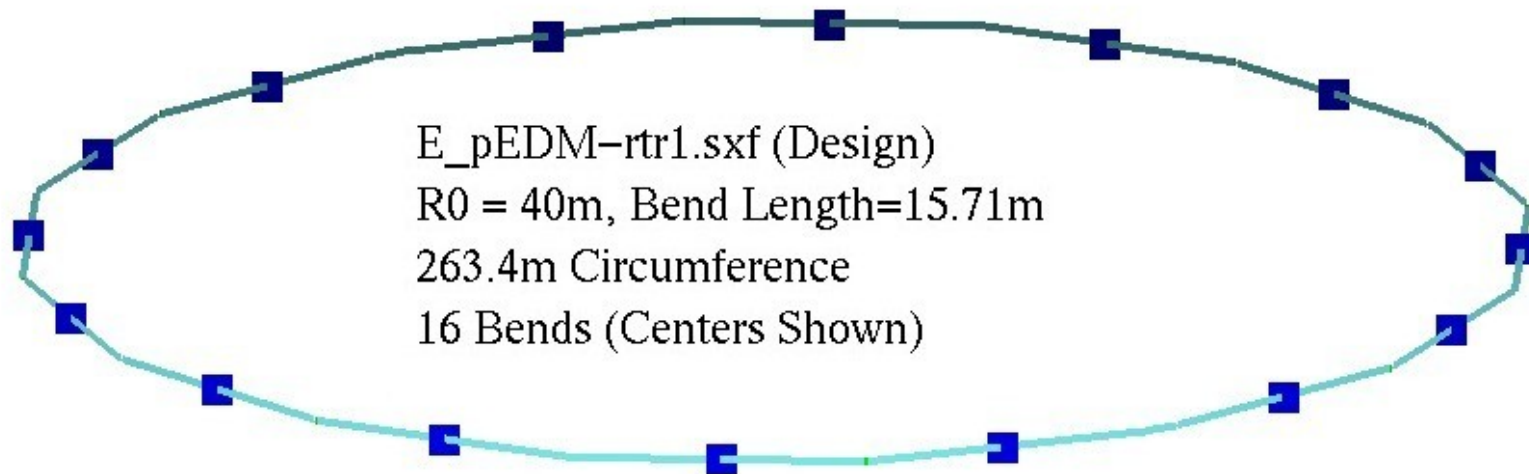
Status: Physical Ring



Status: Physical Ring

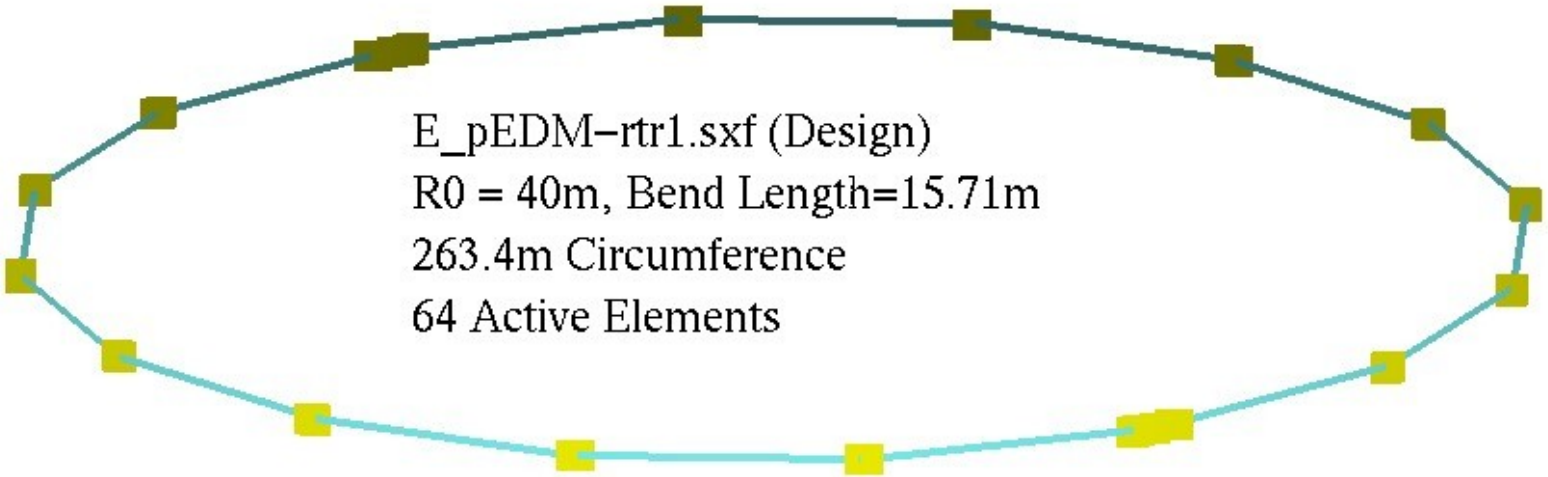
Stable

Status: Lattice



Drift, Kick, Drift --->>> Exact, Split Boundary Kick, Exact

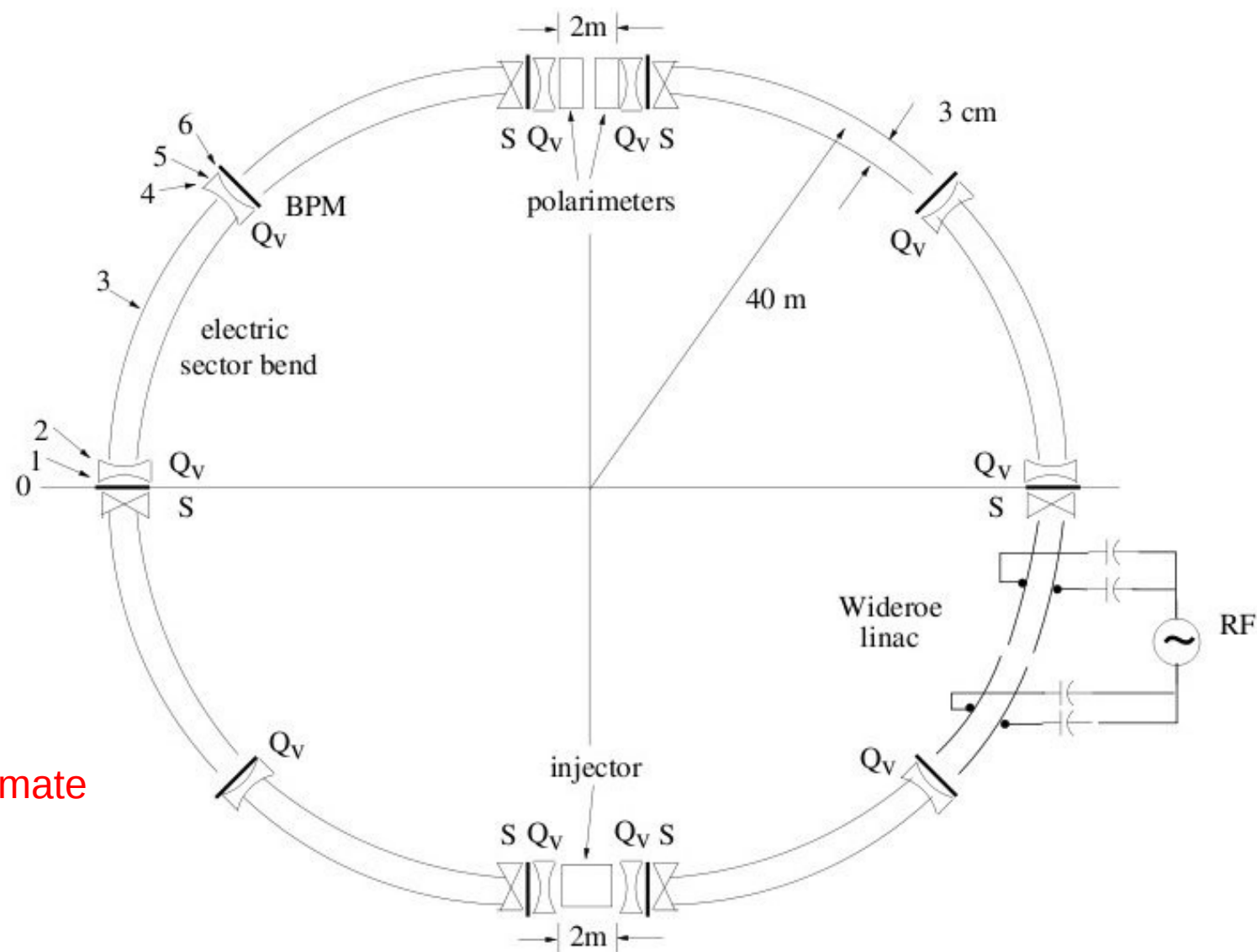
Status: Lattice



The diagram illustrates a particle accelerator lattice design. It features two concentric elliptical paths. The outer path is represented by a dark teal line with 16 square markers, and the inner path is represented by a light blue line with 16 square markers. The markers are distributed evenly along the perimeter of each ellipse. In the center of the diagram, there is a text block providing design specifications.

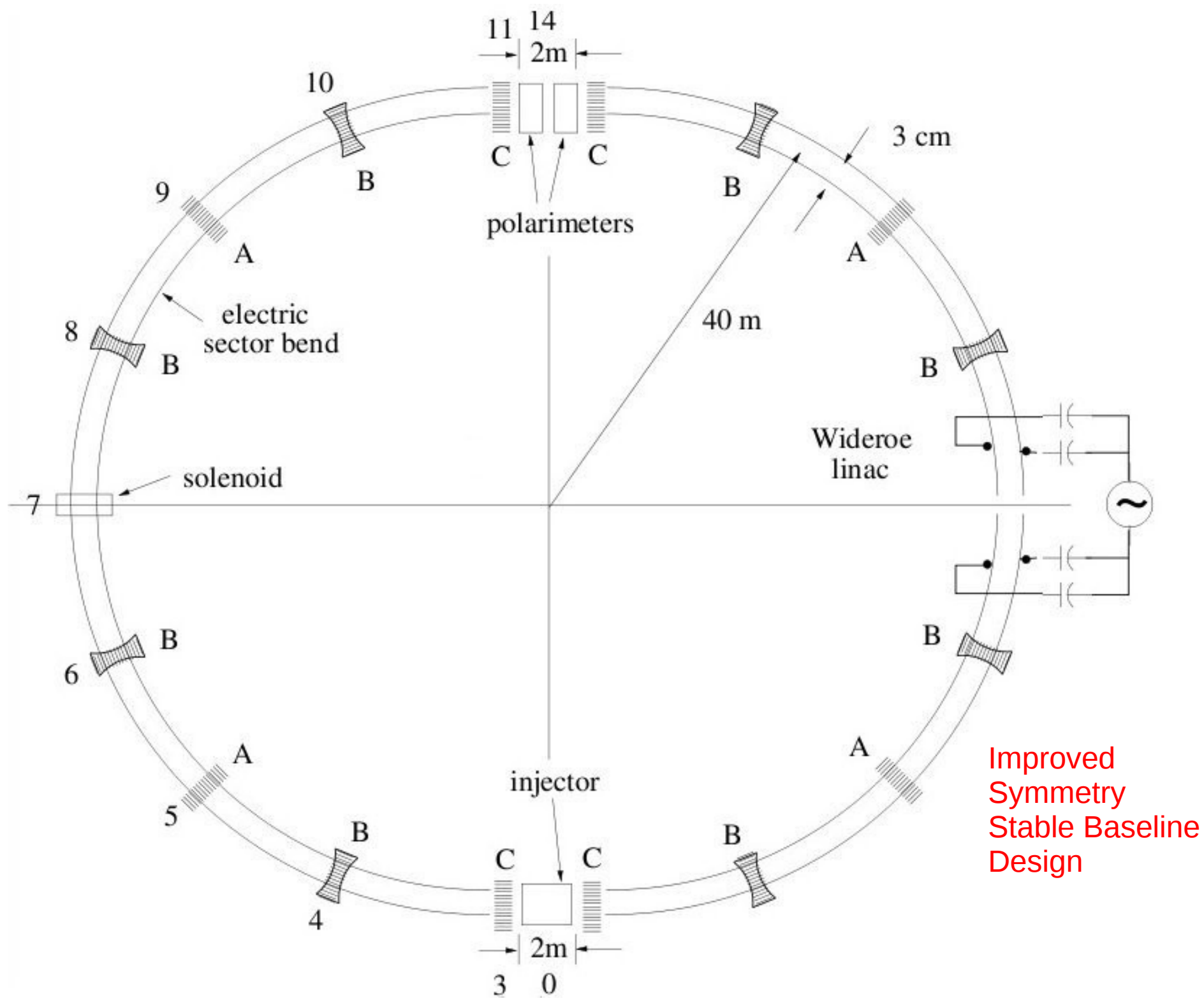
E_pEDM-rtr1.sxf (Design)
R0 = 40m, Bend Length=15.71m
263.4m Circumference
64 Active Elements

Penultimate
Design



Scale in straight sections is expanded

Figure 1: Full ring of the wCFSD, (weak) Combined-Focusing, Separate-Defocusing, lattice. The horizontal focusing is provided by the (combined function) saddle-shaped ($m = -1.2$) electrodes. (Except for the extremely short straightaways of the racetrack, the ring is made from nothing but eight of these cells. Not all elements are shown.



Status: Lattice

Stable

Status: Survey

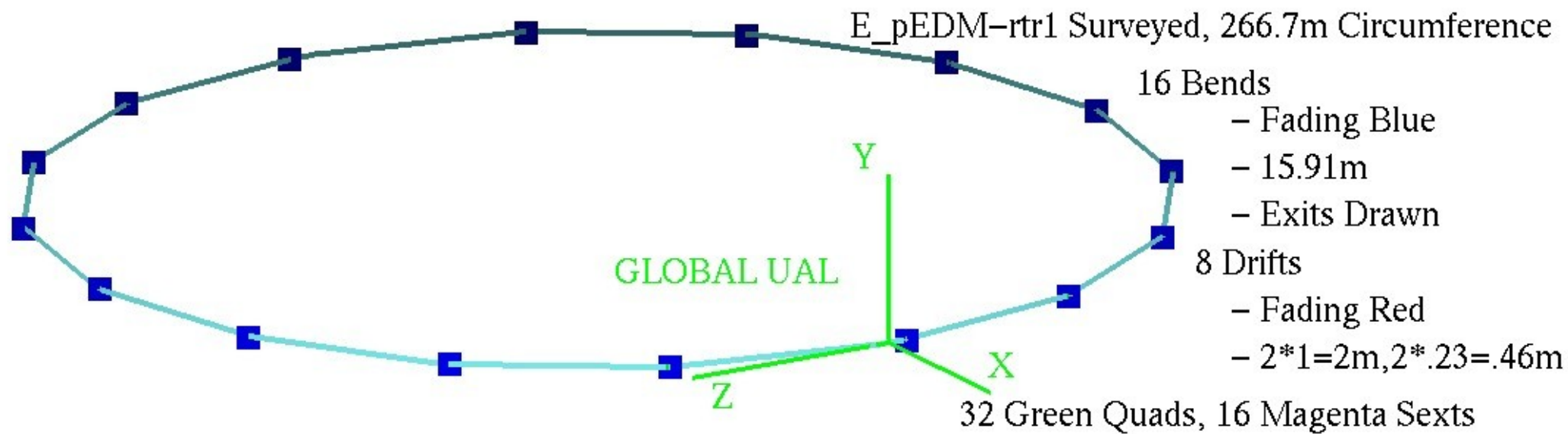
Very Good Corroboration

Fully Arc Length Sequenced

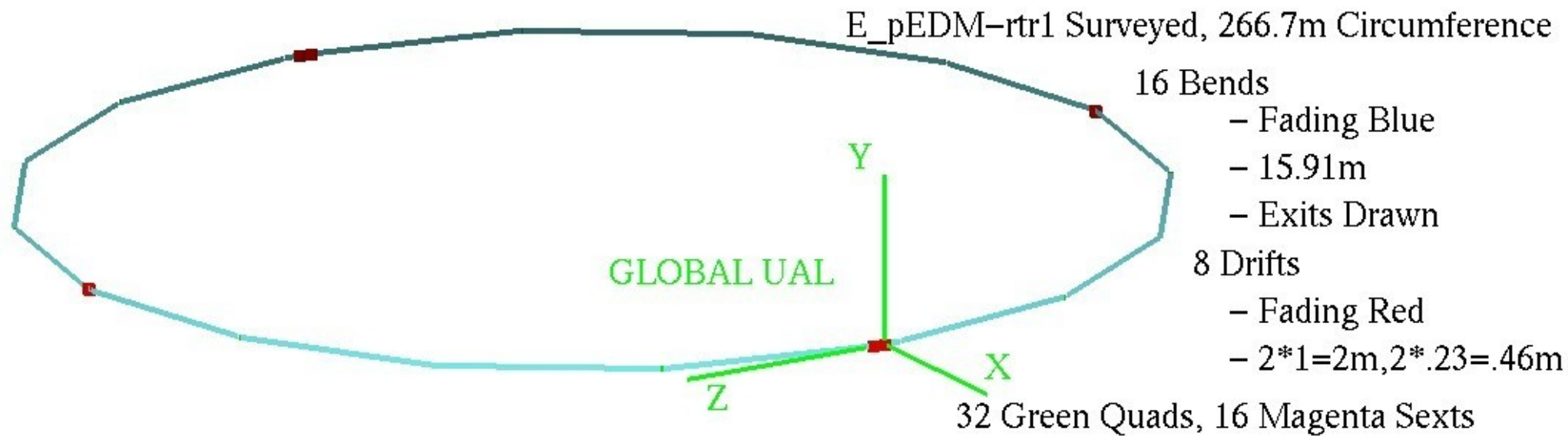
Absolute Coordinates

(Not Strictly Necessary)

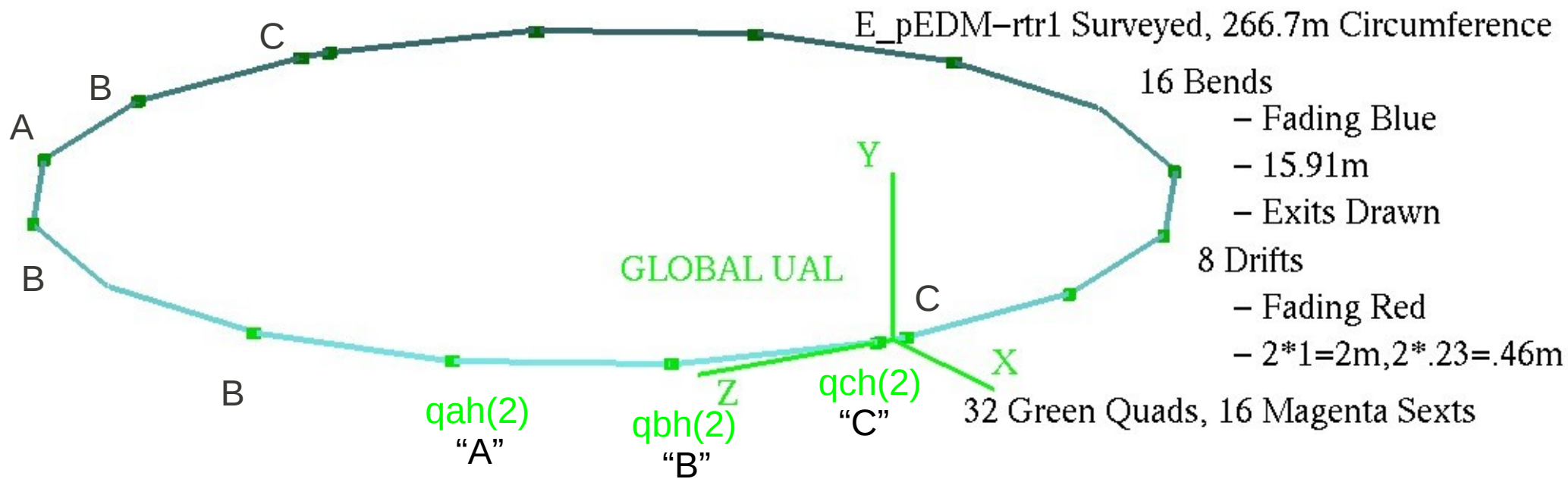
Status: Survey



Status: Survey



Status: Survey



Fundamental Symmetries:

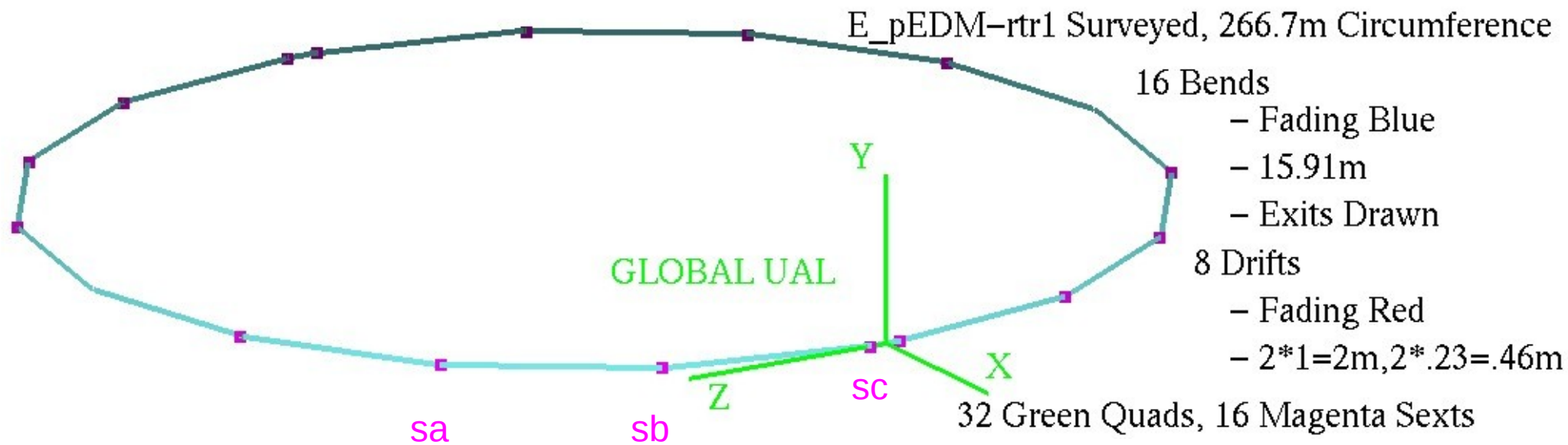
Superperiodicity – 2

Mirror:

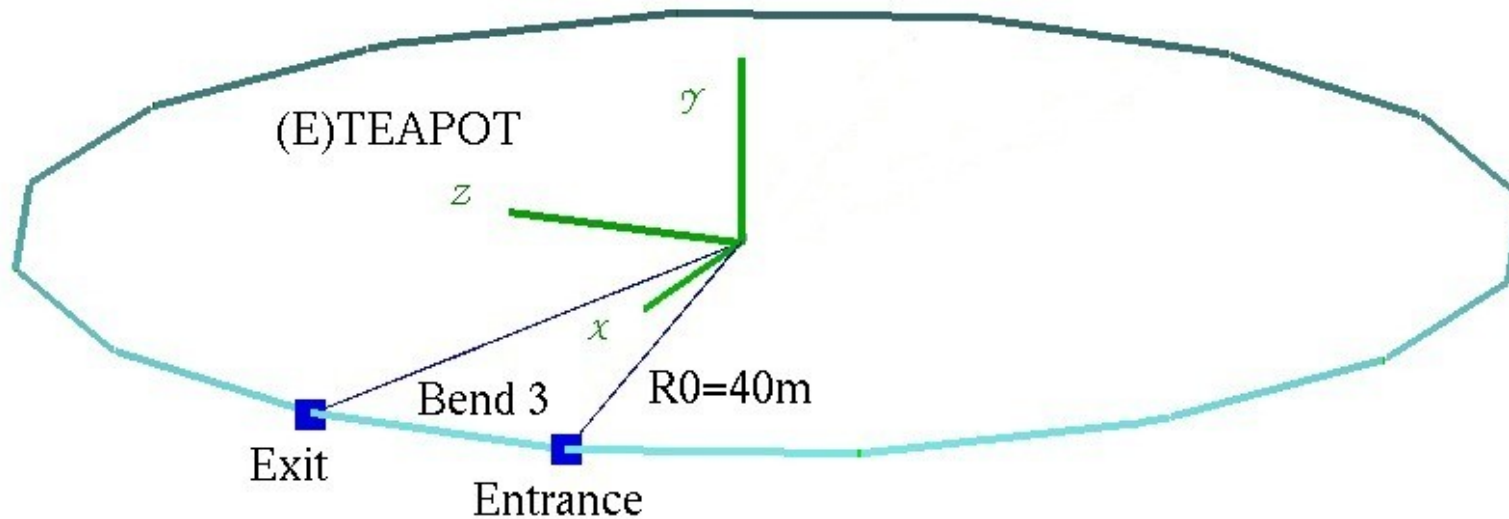
East – West (except for RF \leftrightarrow Solenoid)

North – South (except for Injection \leftrightarrow Polarimeter)

Status: Survey



Status: Survey



Not Strictly Necessary

Status: Survey

Finished

Status: Orientation

Only Need Local
Coordinates Relative
To The Appropriate
Reference/Design Orbit!

New Initial Value
Problem ($F=dp/dt$)
At Each Interval

Status: Orientation

Reference Orbit:

Drift	—	Straight Line
Bend	—	Circular Arc
Quad	—	Straight Line
Sext	—	Straight Line

Status: Orientation

Reference Orbit:

$p[0]$, $p[2]$, $p[4]$ are
coordinates referred to
the reference orbit of the
particular element.

$p[1]$, $p[3]$, $p[5]$ are the
canonical conjugates.

Status: Orientation

Reference Orbit:

$p[0] = x$

$p[2] = y$

$p[4] = \text{time delta}$

$p[1] = px/p0$

$p[3] = py/p0$

$p[5] = \text{energy delta}$

(Deviations From Reference Orbit)

Reference Orbit - Bend

$$\mathbf{E}_m = -E_0 \frac{R_0^{1+m}}{r^{1+m}} \hat{\mathbf{r}}$$

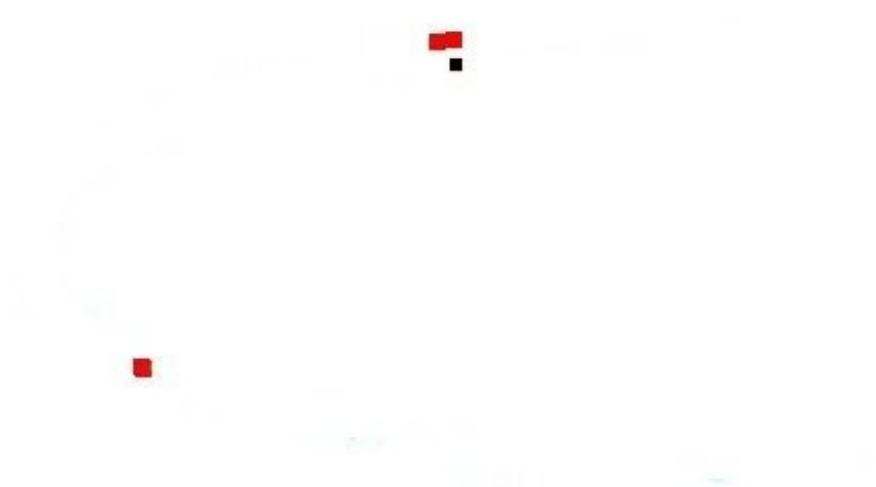


- particle/probe has R_0 available
- $r = r(x,y)$
- r close to $R_0 + p[0]$
- $V(r) = \ln(r)$
- or $1/r$
- or $(1/r)^m$
- ...
- treat “exactly” (thick element, precision of machine, analytical solutions), correct with kick

$$V(r) = -\frac{E_0 R_0}{m} \left(\frac{R_0^m}{r^m} - 1 \right)$$

Reference Orbit - Drift, Quad, Sext

- supposed to indicate straight sections!
- same as magnetic
- probe $x = p[0]$
- $px = p[1] * p0$
- $y = p[2]$
- ...
- same as TEAPOT



Status: Orientation

Element – Algorithm – Probe

UAL algorithms are developed with respect to reference system of the particular element. In short, we (beam physics) need to know only relative positions represented by PAC::Position (probe).

GOOD SOFTWARE ARCHITECTURE!

Status: Orientation

Stabilizing

Status: Equations

The cylindrical coordinate ρ is approximately equal to the spherical coordinate r (nearly co-planar orbits, very close to the horizontal plane)

Conservation of energy,
differentiation with respect to θ , ...

Status: Equations

$$\xi = 1 - \frac{R_0}{r} = \frac{x/R_0}{1 + x/R_0}, \quad \text{and} \quad \frac{d\xi}{d\theta} = \frac{R_0}{r^2} \frac{dr}{d\theta}.$$

$$(\mathcal{E} + eE_0R_0 \ln(1 - \xi))^2 = \frac{L_y^2 c^2}{R_0^2} \left(\frac{d\xi}{d\theta} \right)^2 + \frac{L_y^2 c^2}{R_0^2} (1 - \xi)^2 + p_y^2 c^2 + m^2 c^4$$

$$\begin{aligned} \frac{d^2 \xi}{d\theta^2} &= 1 - \xi - \mathcal{E} \frac{eE_0 R_0^3}{L_y^2 c^2} \frac{1}{1 - \xi} - \frac{e^2 E_0^2 R_0^4}{L_y^2 c^2} \frac{\ln(1 - \xi)}{1 - \xi} \\ &= 1 - \xi - \frac{\mathcal{E}/e}{L_y c/(eR_0)} \frac{E_0 R_0}{L_y c/(eR_0)} \frac{1}{1 - \xi} - \left(\frac{E_0 R_0}{L_y c/(eR_0)} \right)^2 \frac{\ln(1 - \xi)}{1 - \xi} \\ &= 1 - \xi - \frac{\mathcal{E}}{\mathcal{E}_0} \frac{L_0^2}{L_y^2} \frac{1}{1 - \xi} - \frac{L_0^2}{L_y^2} \beta_0^2 \frac{\ln(1 - \xi)}{1 - \xi}. \end{aligned}$$

Why ξ ? It permits analytical solution in special cases, and $\xi = x/R_0$ in linear approximation. Thus, ...

Status: Equations

$$(1 - \xi)^{-1} = 1 + \xi + \xi^2 + \dots, \quad \frac{\ln(1 - \xi)}{1 - \xi} = -\xi - \frac{3}{2}\xi^2 + \dots$$

The orbit equation becomes

$$\begin{aligned} \frac{d^2\xi}{d\theta^2} &= 1 - \frac{L_0^2}{L_y^2} \frac{\mathcal{E}}{\mathcal{E}_0} - \left(1 + \frac{L_0^2}{L_y^2} \left(\frac{\mathcal{E}}{\mathcal{E}_0} - \beta_0^2\right)\right) \xi - \frac{L_0^2}{L_y^2} \left(\frac{\mathcal{E}}{\mathcal{E}_0} - \frac{3}{2}\beta_0^2\right) \xi^2 + \dots \\ &\approx -Q^2(\xi - \xi_0), \quad \text{Nice/Convenient Form!} \end{aligned}$$

$$Q^2 = 1 + \frac{L_0^2}{L_y^2} \left(\frac{\mathcal{E}}{\mathcal{E}_0} - \beta_0^2\right), \quad \text{and} \quad \xi_0 = \frac{1 - \frac{L_0^2}{L_y^2} \frac{\mathcal{E}}{\mathcal{E}_0}}{Q^2}.$$

Status: Equations

$$C_{\xi}(\theta) = \cos(Q\theta)(1 - \xi_0) + \xi_0$$

$$C'_{\xi}(\theta) = -\sin(Q\theta) Q(1 - \xi_0)$$

$$S_{\xi}(\theta) = \frac{\sin(Q\theta)}{Q} - \cos(Q\theta) \xi_0 + \xi_0$$

$$S'_{\xi}(\theta) = \cos(Q\theta) + \sin(Q\theta) Q\xi_0$$

cosine-like orbit/trajectory

sine-like orbit/trajectory

Status: Equations

$$\xi_{\text{in}} = \frac{x_{\text{in}}}{R_0 + x_{\text{in}}}, \quad \xi'_{\text{in}} = \frac{R_0 x'_{\text{in}}}{(R_0 + x_{\text{in}})^2}.$$

$$\frac{x/R_0}{1 + x/R_0} = C_\xi(\theta) \xi_{\text{in}} + S_\xi(\theta) \xi'_{\text{in}}.$$

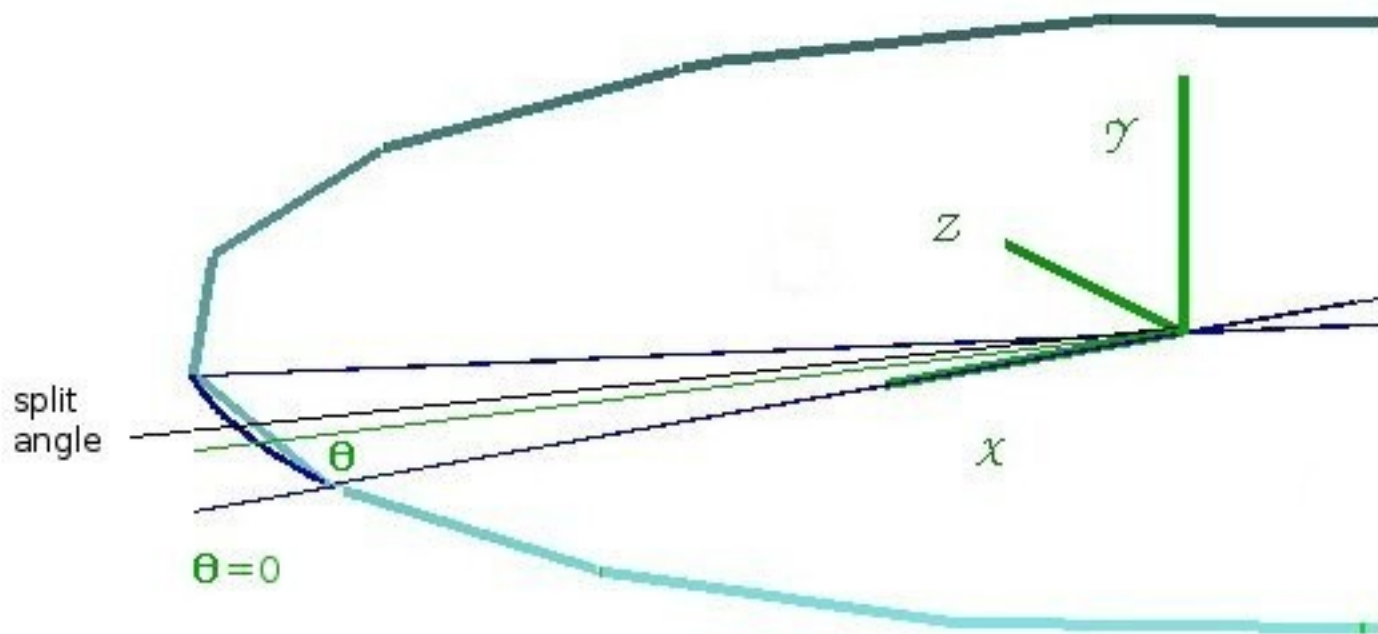
$$\frac{x}{R_0} = \frac{C_\xi(\theta) \xi_{\text{in}} + S_\xi(\theta) \xi'_{\text{in}}}{1 - C_\xi(\theta) \xi_{\text{in}} - S_\xi(\theta) \xi'_{\text{in}}}$$

$$\frac{x'}{R_0} = \frac{C'_\xi(\theta) \xi_{\text{in}} + S'_\xi(\theta) \xi'_{\text{in}}}{\left(1 - C_\xi(\theta) \xi_{\text{in}} - S_\xi(\theta) \xi'_{\text{in}}\right)^2}.$$

“Exact”
Output in terms of
Input

Stable!

Status: Equations



Status: Equations

Stabilizing

Status: Code

Bunch loop

```
PAC::Position& p =  
bunch[ip].getPosition();
```

(currently 1 probe in bunch)

Splitting

```
simplest split (into 2 equal  
halves)
```

Status: Code

```
const PAC::BeamAttributes cba
```

```
double e0=cba.getEnergy();  
double m0=cba.getMass();  
double q0=cba.getCharge();  
double t0=cba.getElapsedTime();  
double f0=cba.getRevfreq();  
double M0=cba.getMacrosize();  
double G0=cba.getG();  
double L0=cba.getL();           // hybrid? p0 beam, R0 element?  
double E0=cba.getE();           // more like an element property  
double R0=cba.getR();           // more like an element property  
                                // needs to be reworked
```

```
    "p0"=sqrt(e0*e0-m0*m0); // design momentum  
        =L0/R0;  
    "g0"=e0/m0;             // design gamma
```

g also stands for gap

Status: Code

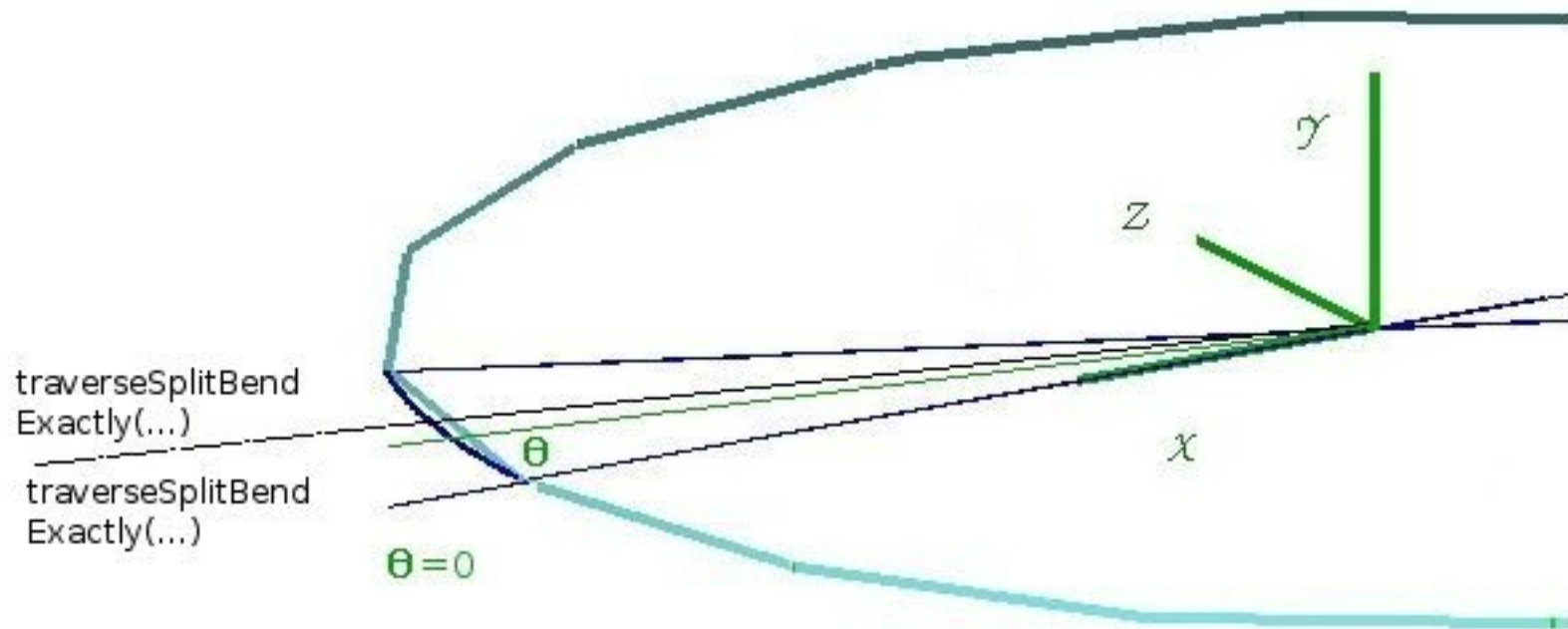
```
// Drift, Kick, Drift --->>> Exact, Split Boundary Kick, Exact

// passBendSlice          (data.m_slices[0], p, tmp, v0byc);
  traverseSplitBendExactly(data.m_slices[0], p, tmp, v0byc,
                          cba,data.m_angle/2);

// applyThinBendKick(data, edata, 1, p, v0byc);
  handleSplitBendBoundary(p,cba);

// passBendSlice(data.m_slices[1], p, tmp, v0byc);
  traverseSplitBendExactly(data.m_slices[1], p, tmp, v0byc,
                          cba,data.m_angle/2);
```

Status: Code



Plus Corrections at Entrance, Split, and Exit

Status: Code

```
enterBendCorrection(p,cba);

// passBendSlice          (data.m_slices[0], p, tmp, v0byc);
  traverseSplitBendExactly(data.m_slices[0], p, tmp, v0byc,
                          cba,data.m_angle/2);

// applyThinBendKick(data, edata, 1, p, v0byc);
  handleSplitBendBoundary(p,cba);

// passBendSlice(data.m_slices[1], p, tmp, v0byc);
  traverseSplitBendExactly(data.m_slices[1], p, tmp, v0byc,
                          cba,data.m_angle/2);

leaveBendCorrection(p,cba);
```

Status: Code

```
enterBendCorrection(Coordinates& p,  
                    const PAC::BeamAttributes cba){  
  
    #include "getDesignBeam.h"  
  
    //          eMD=e0+p0*p[5]          // Mechanical energy - "Drift" (non Bend)  
    //          eTD=eMD                // Total energy - Drift  
    //          eTB=eTD                // Total energy conserved  
    //          =eMB+V(r)              // Total energy - Bend  
    //          eMD=eMB+V(r)          // substitute  
    //  
    //      (eMD-e0)/p0=(eMB-e0)/p0+V(r)/p0      // rearrange  
    //      p[5]Drift = p[5]Bend +V(r)/p0         // p[5]Bend is output  
  
    double q      = UAL::elemCharge;          // units  
    double GeVperJ = 1/q/1e9;                 // units  
    double r      = R0+p[0];  
    double p0     = sqrt(e0*e0-m0*m0);        // derived beam momentum  
  
    p[5] = p[5]-GeVperJ*getPotentialEnergy(q,E0,R0,r)/p0;  
}
```

Status: Code

```
double getPotentialEnergy(double q0,double E0,double R0,double r){  
    return q0*E0*R0*log(r/R0);  
}
```

```
double Cxi(double Q,double theta,double xi0){  
    double value=cos(Q*theta)*(1-xi0)+xi0;  
    return value;  
}
```

```
double CxiP(double Q,double theta,double xi0){  
    double value=-sin(Q*theta)*Q*(1-xi0);  
    return value;  
}
```

```
double Sxi(double Q,double theta,double xi0){  
    double value=sin(Q*theta)/Q-cos(Q*theta)*xi0+xi0;  
    return value;  
}
```

```
double SxiP(double Q,double theta,double xi0){  
    double value=cos(Q*theta)+sin(Q*theta)*Q*xi0;  
    return value;  
}
```

Convenient
Equations!

Status: Code

- good start on stubs with doubles
- need better time of flight ($p[4]$)
- need better $p[2]$, $p[3]$ (y and p_y)
- TODO:
 - testing
 - benchmarking
 - spin
 - (vector of) truncated power series (polymorphism)

Status: Code

Stabilizing

Status: Summary

- in control of geometry,
element sequencing,
code sequencing,
probe parameters
- good UAL software architecture allows
the algorithm to be developed
independently
- probably about $\frac{1}{2}$ way to code
content

Thanks To

Nikolay Malitsky

Richard Talman

Yannis Semertzidis

Bill Morse

Selcuk Haciomeroglu

Alfredo Luccio

...

Status:

Physical Ring	—	Stable
Lattice	—	Stable
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Code	—	Stabilizing

Probably biased to the
optimistic side!

Free Associating

Flight Time

- Flight time is given by an integral which is not in the spirit of the TEAPOT approach.
- Flight time should be viewed as a split operation; The higher the split # (finer the splitting), the more accurate the flight time.

$$mc^2\gamma^I(\theta) = \mathcal{E} - eE_0(C_\xi(\theta)x_0 + S_\xi(\theta)x'_0).$$

From this formula one can obtain β^I using

$$\beta^{I^2}(\theta) = 1 - \frac{1}{\gamma^{I^2}(\theta)}.$$

For motion in the horizontal plane, the angular velocity can be obtained from the y -component of Eq. (9);

$$\frac{d\theta}{dt} = \frac{L}{mr^2\gamma^I}.$$

The right hand side of this equation can be expressed in terms of θ , invariant and initial conditions using Eq. (48). This relation is useful primarily to obtain the flight time through bend elements. After rearrangement, the flight time $t_o - t_i$ from input to output of a bend element is given by

$$t_o - t_i = \frac{m}{L} \int_{\theta_i}^{\theta_o} r^2(\theta')\gamma^I(\theta') d\theta' \approx \frac{mR_0^2}{L} \int_{\theta_i}^{\theta_o} (1 + 2x(\theta')/R_0)\gamma^I(\theta') d\theta'.$$

Acceptance/ Admittance/ Emittance

- E0 central/design/ideal/reference Electric Field
1.7E7 V/m
- g design gap
3cm?

Acceptance/ Admittance/ Emittance

Are "admittance" and "acceptance" more or less synonymous?

Yes. Furthermore "emittance" (which applies to a beam) and "admittance" (which applies to a lattice) are commensurate (same units) quantities. The lattice admittance has to exceed the beam emittance in order for the whole beam to be captured. "Dynamic aperture" is similar to "admittance" though it has different units since it is just the maximum stable particle transverse displacement from the design orbit.


```

$UAL/codes/ETEAPOT/src/ETEAPOT/Integrator/DipoleTracker.cc
93     for(int ip = 0; ip < bunch.size(); ip++) {
    if(bunch[ip].isLost()) continue;
    PAC::Position& p = bunch[ip].getPosition();      <<<----
    tmp = p;
    s_algorithm.passEntry(m_edata, p);
    s_algorithm.makeVelocity(p, tmp, v0byc);
    s_algorithm.makeRV(p, tmp, e0, p0, m0);

    s_algorithm.passBend(m_data, m_edata, p, tmp, v0byc, cba);
//          [          original interface          ] [central orbit ]
    s_algorithm.passExit(m_edata, p);
    // testAperture(p);
}

```

```

$UAL/examples/ETEAPOT/tracker.cc
50     UAL::Shell shell;

    // *****
    std::cout << "\nDefine the space of Taylor maps." << std::endl;
    // *****

    shell.setMapAttributes(UAL::Args() << UAL::Arg("order", 5));  <<<----

```

Simulation Initialization

Beam -

```
shell.setBeamAttributes(UAL::Args() << UAL::Arg("energy",          e0));
shell.setBeamAttributes(UAL::Args() << UAL::Arg("mass",            m0));
shell.setBeamAttributes(UAL::Args() << UAL::Arg("charge",          q0));
//shell.setBeamAttributes(UAL::Args() << UAL::Arg("elapsedTime",    t0));
//shell.setBeamAttributes(UAL::Args() << UAL::Arg("frequency",      f0));
shell.setBeamAttributes(UAL::Args() << UAL::Arg("macrosize",        M0));
//shell.setBeamAttributes(UAL::Args() << UAL::Arg("gyromagnetic",    G0));
shell.setBeamAttributes(UAL::Args() << UAL::Arg("designAngularMomentum",L0));
shell.setBeamAttributes(UAL::Args() << UAL::Arg("designElectricField", E0));
shell.setBeamAttributes(UAL::Args() << UAL::Arg("designRadius",      R0));
```

Simulation Initialization

Probe -

```
for(int ip=0;ip<bunch.size();ip++){  
    bunch[ip].getPosition().set  
        (probe__dx0,probe_dpx0,probe__dy0,probe_dpy0,probe_cdt0,probeEscr0);  
    bunch[ip].setSpin(spin);  
}
```

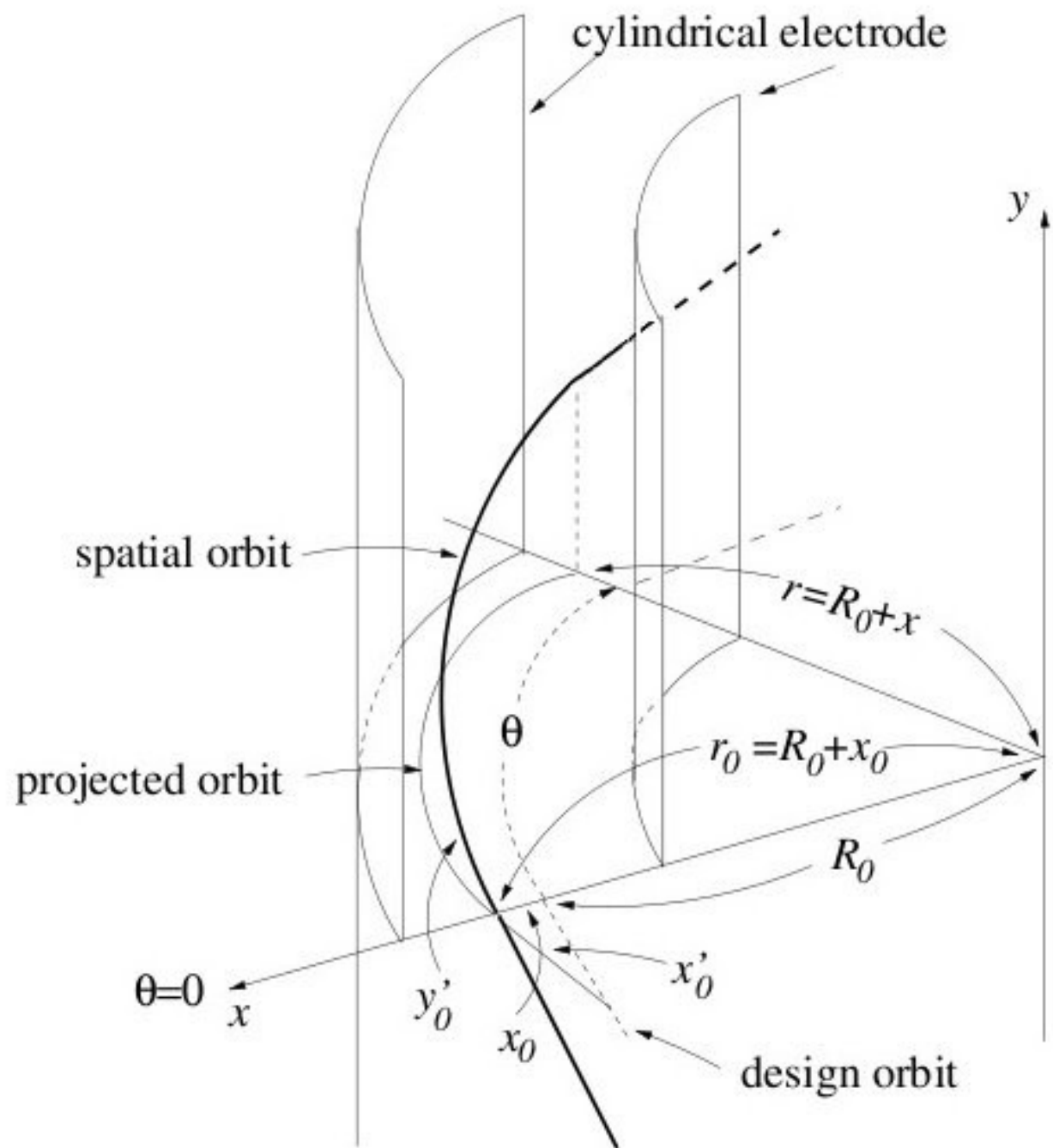


Figure 1: The bold curve shows a proton orbit passing through a curved-planar cylindrical electrostatic bending element. The electrode spacing is g and the design orbit is centered between the electrodes.

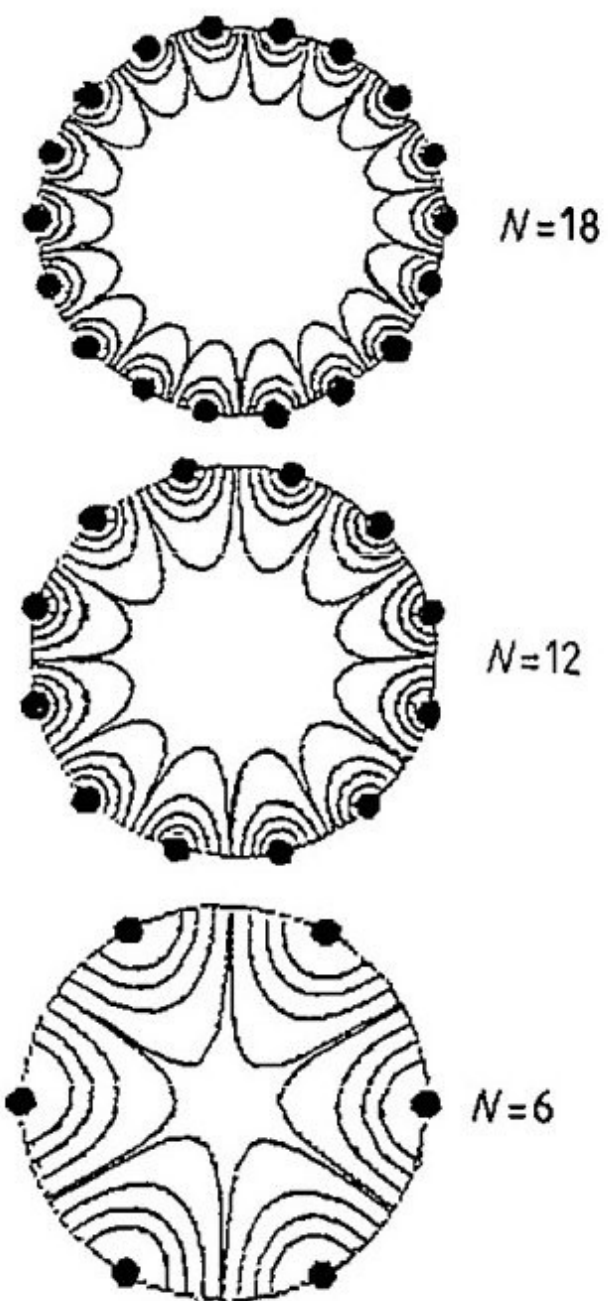


Figure 3. Configuration of multipole magnetic field in a 'squirrel cage' scheme.

