

Mathematical and Numerical Methods for Non-linear Dynamics (an introduction)

Werner Herr
CERN, BE Department

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http://cern.ch/Werner.Herr/CAS2011/lectures/Chios_methods.pdf

For many more details:

<http://cern.ch/Werner.Herr/METHODS>

Werner Herr, non-linear methods, CAS 2011, Chios

Primary purpose of this lectures

- Need to introduce new tools for **non-linear** dynamics
- Give an overview of the *modern*^{*)} tools used in accelerator physics
- Necessarily incomplete
- Avoid mathematical derivations, rather give "*raison d'être*" and "*mode d'emploi*"

^{*)} *modern*: "contemporary", not "fashionable" !

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Recommended Bibliography:

E. Forest, *Beam Dynamics - A New Attitude and Framework*
Harwood Academic Publishers, 1998.

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- A. Chao and M. Tigner, *Handbook of Accelerator Physics and Engineering*
SLAC, 2001.
- A. Chao and M. Tigner, *Lecture Notes on Topics in Accelerator Physics and Engineering*
World Scientific Publishing, 1998.

Outline of this lectures

- Motivation, introduction and classical concepts
- New concepts and modern techniques
 - Maps
 - Hamiltonian theory (for our purpose)
 - Computation: maps, symplectic integration
 - Analysis: Lie transforms, normal forms
 - Analysis: Differential algebra

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Treatment of LINEAR dynamics

- Standard introduction using Hill's equation
(for simplicity: show for one dimension first):

$$\frac{d^2x(s)}{ds^2} + K(s)x(s) = 0$$

■ $K(s)$ periodic, smooth function

■ Is that true ?

■ No, is only approximate

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Treatment of LINEAR dynamics

- Used to "derive" Courant-Snyder ansatz:

$$x(s) = \sqrt{\beta(s)} \cdot \epsilon \cdot \cos(\mu(s) + \mu_0)$$

$$x'(s) = \sqrt{\frac{\epsilon}{\beta(s)}} \cdot (\sin(\mu(s) + \mu_0) + \alpha \cdot \cos(\mu(s) + \mu_0))$$

■ Is the solution to any system that is: confined and periodic !

■ Do we need Hill's equation ?

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Treatment of DISTORTED dynamics

- Hill's equation with distortions
- For both transverse dimensions:

$$\frac{d^2x(s)}{ds^2} + K(s)x(s) = \frac{F_x(x, y, s)}{v \cdot p}$$

$$\frac{d^2y(s)}{ds^2} + K(s)y(s) = \frac{F_y(x, y, s)}{v \cdot p}$$

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Treatment of DISTORTED dynamics

- What are F_x, F_y ?

$$F_x(x, y, s) = -v \cdot B_y(x, y, s)$$

$$F_y(x, y, s) = -v \cdot B_x(x, y, s)$$

- $B_x(x, y, s), B_y(x, y, s)$ can be any higher order electromagnetic field

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Treatment of DISTORTED dynamics

- We have to re-write:

$$\frac{d^2x(s)}{ds^2} + K(s)x(s) = \frac{F_x(x, y, s)}{v \cdot p}$$

- as (similar for the other plane):

$$\frac{d^2x(s)}{ds^2} + K(s)x(s) = \sum_{i,j,k,l \geq 0} p_{ijkl}(s)x^i x^j y^k y^l$$

- Very non-linear differential equation to solve ...

- Enter the field of non-linear dynamics

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Can we deal with that ?

- Under certain circumstances:

- All $p_{ijkl}(s)$ are perturbations, i.e. (very) small
- Only a few $p_{ijkl}(s)$ are non-zero
- You can avoid resonances
- Perturbations are smooth or possibly periodic

- Would you build a 3 billion Euro machine on these assumptions ?

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A disclaimer ...

- Can be useful to understand and demonstrate certain aspects
- See Oliver Brüning's lecture
- For practical work on realistic machine:
 - New tools required
 - It is much easier that you think (.. and other people tell you !)

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A better framework

Start with the differential equation:

$$\frac{d^2x(s)}{ds^2} + K(s)x(s) = \sum_{i,j,k,l \geq 0} p_{ijkl}(s)x^i x'^j y^k y'^l$$

- Bad news: We have no global analytical solution
- Good news: An analytical solution is not needed !

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A better framework

- Why not ?
- We do **not** want to know:
 - The particle's position and momentum at 2h 45min 22.3s ?
 - We do **want** to know:
 - Is the beams stable for a long time ?
 - Is the motion confined ?
 - Can we get a framework to get that (easily) ?

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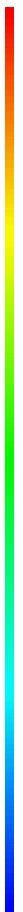


A better framework - go back 100 years ...

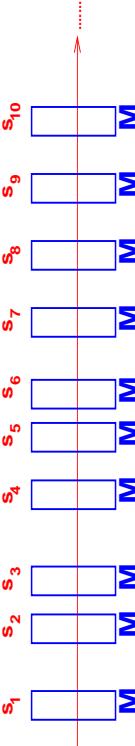


- ”Old” to ”New” classical dynamics:
 - Topology and properties of phase space (see Oliver’s lecture)
 - Chaotic motion, non-integrable systems
 - Sensitivity to initial conditions

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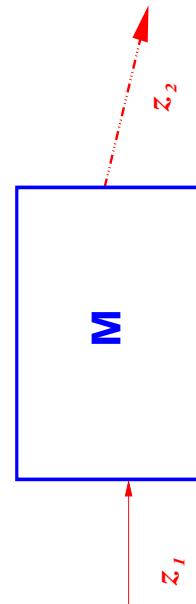
How does an **accelerator** really look like ?

- Magnetic and electric fields are (in general) not a continuous function of s
- A (finite) sequence of machine elements \mathbf{M} at longitudinal positions s_1, s_2, s_3, \dots :

- Cannot be described by a smooth differential equation

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How can an **element** really be described ?

- You need to describe what happens to the particle in M
- Assume each element \mathbf{M} (e.g. magnet) acts on the beam **locally** in a deterministic way, **functionally independent** of all other elements



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How is an element described ?

- Let \vec{z}_1, \vec{z}_2 describe a quantity (coordinates, beam sizes ...) before and after the element
- Take an machine element (e.g. magnet) and build a mathematical model \mathcal{M}
 - In general: $\vec{z}_2 = \mathcal{M}(\vec{z}_1)$
 - \mathcal{M} is a so-called map
 - Very important: no need to know what happens in the rest of the machine !!
- The complete sequence of MAPS connects the pieces together to make a ring (or beam line)

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MAPS transform coordinates through an element

- Use coordinate vector: $\vec{z} = (x, x' = \frac{\partial x}{\partial s}, y, y' = \frac{\partial y}{\partial s})^*$
- \mathcal{M} transforms the coordinates $\vec{z}_1(s_1)$ at position s_1 to new coordinates $\vec{z}_2(s_2)$ at position s_2 :

$$\vec{z}_2(s_2) = \begin{pmatrix} x \\ x' \\ y \\ y' \end{pmatrix} = \mathcal{M} \circ \begin{pmatrix} x \\ x' \\ y \\ y' \end{pmatrix}_{s_1} = \mathcal{M} \circ \vec{z}_1(s_1)$$

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*) not unique, see later



What can \mathcal{M} be ?

- Any "description" to go from \vec{z}_1 to \vec{z}_2
- This "description" can be:
 - A simple linear matrix or transformation
 - A non-linear transformation (Taylor series, Lie Transform ...)
 - High order integration algorithm
 - A computer program, subroutine etc.
- Let us look at linear theory first !

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Simple examples (one dimensional)

Focusing quadrupole of length L and strength k :

$$\begin{pmatrix} x \\ x' \end{pmatrix}_{s_2} = \begin{pmatrix} \cos(L \cdot k) & \frac{1}{k} \cdot \sin(L \cdot k) \\ k \cdot \sin(L \cdot k) & \cos(L \cdot k) \end{pmatrix} \circ \begin{pmatrix} x \\ x' \end{pmatrix}_{s_1}$$

Quadrupole with short length L (i.e.: $1 \gg L^2 \cdot k^2$)

$$\begin{pmatrix} x \\ x' \end{pmatrix}_{s_2} = \begin{pmatrix} 1 & 0 \\ k^2 \cdot L (= -\frac{1}{f}) & 1 \end{pmatrix} \circ \begin{pmatrix} x \\ x' \end{pmatrix}_{s_1}$$

They are \mathcal{M} aps, describe the movement in an element (quadrupole)

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Interlude: there was already a trap ... !

According to B. Holzer (lectures) or K. Wille (textbook):

$$k = \frac{1}{B\rho} \frac{dB_y}{dx}$$

According to "Handbook for Accelerator Physics" (A. Chao):

$$k^2 = \frac{1}{B\rho} \frac{dB_y}{dx}$$

- The lesson: check what people use !!
(remember Air Canada 143)

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Interlude: it can be worse ... !

You also find:

$$K^2 = k = \frac{1}{B\rho} \frac{dB_y}{dx}$$

Often different conventions in simulation programs !

Some programs want **fields**, not gradients !
Found this construction:

$$B_y = 0.1 \cdot k \cdot x \cdot B \cdot \rho$$

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Putting the "pieces" together

Starting from a position s_0 and applying all maps (for N elements) in sequence around a ring with circumference C to get the **One-Turn-Map** (OTM) for the position s_0 (for one dimension only):

$$\begin{pmatrix} x \\ x' \end{pmatrix}_{s_0 + C} = \mathcal{M}_1 \circ \mathcal{M}_2 \circ \dots \circ \mathcal{M}_N \circ \begin{pmatrix} x \\ x' \end{pmatrix}_{s_0}$$

$$\Rightarrow \begin{pmatrix} x \\ x' \end{pmatrix}_{s_0 + C} = \mathcal{M}_{ring}(s_0) \circ \begin{pmatrix} x \\ x' \end{pmatrix}_{s_0}$$

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Why are we interested in \mathcal{M}_{ring} ?

- In the purely linear case \mathcal{M}_{ring} corresponds to a matrix
 - \mathcal{M}_{ring} allows to derive global quantities (\mathbf{Q} , orbit, etc...)
 - Allows the analysis of imperfections (and their correction !)
 - "Straightforward" to formally extend it to complicated (e.g. non-linear) problems

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First extension: two dimensions

- Extend vectors for coordinates
 $(x, x') \rightarrow (x, x', y, y')$

- Extend transfer maps/matrices

$$\begin{pmatrix} x \\ x' \\ y \\ y' \end{pmatrix}_{s_2} = \begin{pmatrix} m_{11} & m_{12} & 0 & 0 \\ m_{21} & m_{22} & 0 & 0 \\ 0 & 0 & m_{33} & m_{34} \\ 0 & 0 & m_{43} & m_{44} \end{pmatrix} \circ \begin{pmatrix} x \\ x' \\ y \\ y' \end{pmatrix}_{s_1}$$

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Extension to two dimensions (coupling)

- The horizontal and vertical motion can be coupled:

- Additional elements in matrix

$$\begin{pmatrix} x \\ x' \\ y \\ y' \end{pmatrix}_{s_2} = \begin{pmatrix} m_{11} & m_{12} & \textcolor{red}{m_{13}} & \textcolor{red}{m_{14}} \\ m_{21} & m_{22} & \textcolor{red}{m_{23}} & \textcolor{red}{m_{24}} \\ \textcolor{red}{m_{31}} & \textcolor{red}{m_{32}} & m_{33} & m_{34} \\ \textcolor{red}{m_{41}} & \textcolor{red}{m_{42}} & m_{43} & m_{44} \end{pmatrix} \circ \begin{pmatrix} x \\ x' \\ y \\ y' \end{pmatrix}_{s_1}$$

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- Q: what about the β -functions now ??

Analysis of the One-Turn-Map

- We have obtained a map for the whole ring
- Have to get now the information we want:
 - Optics parameters (Tune, Twiss function, ..)
 - Closed orbit
 - Stability
 - etc. ...
- How to analyse a MAP (first: a matrix) ???
(see also B. Holzer lecture, but reality comes here)

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Normal forms

- Maps can be transformed into (Jordan) Normal Forms
- Original maps and normal form are equivalent, but ...
- Easily used to analyse the maps:
 - Get parameters (Q, Q' , Twiss function, ..)
 - Study invariants, etc.
 - Extract (non-linear) tune shifts
 - For resonance analysis
 - etc. ...
- Idea is to make a transformation to get a simpler form for the map

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Normal forms

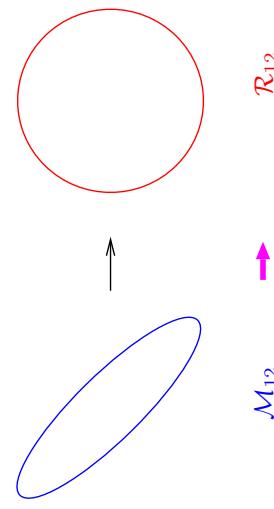
Assume the map \mathcal{M}_{12} propagates the variables from location 1 to location 2, we try to find a transformation \mathcal{A} such that:

$$\mathcal{A} \mathcal{M}_{12} \mathcal{A}^{-1} = \mathcal{R}_{12}$$

- The map \mathcal{R}_{12} is:
 - A "Jordan Normal Form", (or at least a very simplified form of the map)
 - Example: \mathcal{R}_{12} becomes a pure rotation
 - The map \mathcal{R}_{12} describes the same dynamics as \mathcal{M}_{12} , but:
 - All coordinates are transformed
 - This transformation "analyses" the non-trivial part

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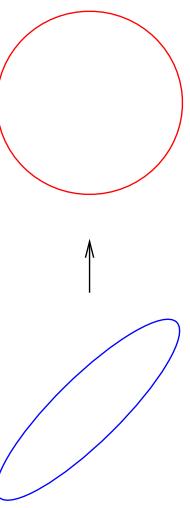
Normal forms - linear case



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- Pictorial form of the transformation
- Motion on a complicated ellipse becomes motion on a circle (i.e. a pure rotation)

Normal forms - linear case



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$$\mathcal{M} = \mathcal{A} \circ \mathcal{R}(\Delta\mu) \circ \mathcal{A}^{-1} \quad \text{or:} \quad \mathcal{R}(\Delta\mu) = \mathcal{A}^{-1} \circ \mathcal{M} \circ \mathcal{A}$$



Normal forms - linear case

Assume the one-turn-map $\mathcal{M}(s)$ at the position s is (e.g. lecture on transverse dynamics):

$$\mathcal{M}(s) = \begin{pmatrix} \cos(\Delta\mu) + \alpha(s)\sin(\Delta\mu) & \beta(s)\sin(\Delta\mu) \\ -\gamma(s)\sin(\Delta\mu) & \cos(\Delta\mu) - \alpha(s)\sin(\Delta\mu) \end{pmatrix}$$

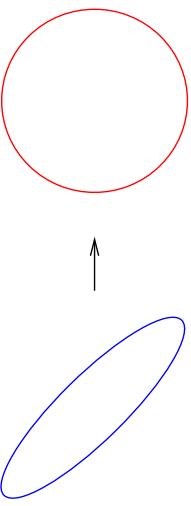
- Describes the motion on a phase space ellipse
- Re-write \mathcal{M} such that one part \mathcal{R} becomes a pure rotation (a circle), i.e.:

$$\mathcal{A}\mathcal{R}\mathcal{A}^{-1} = \mathcal{M}$$

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Normal forms - linear case



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$$\mathcal{M} = \mathcal{A} \circ \mathcal{R}(\Delta\mu) \circ \mathcal{A}^{-1} \quad \text{or:} \quad \mathcal{R}(\Delta\mu) = \mathcal{A}^{-1} \circ \mathcal{M} \circ \mathcal{A}$$

with

$$\mathcal{A} = \begin{pmatrix} \sqrt{\beta(s)} & 0 \\ -\frac{\alpha}{\sqrt{\beta}} & \frac{1}{\sqrt{\beta(s)}} \end{pmatrix} \quad \text{and} \quad \mathcal{R} = \begin{pmatrix} \cos(\Delta\mu) & \sin(\Delta\mu) \\ -\sin(\Delta\mu) & \cos(\Delta\mu) \end{pmatrix}$$



Normal forms - linear case

We had:

$$\mathcal{M} = \mathcal{A} \circ \mathcal{R}(\Delta\mu) \circ \mathcal{A}^{-1} \quad \text{or:} \quad \mathcal{R}(\Delta\mu) = \mathcal{A}^{-1} \circ \mathcal{M} \circ \mathcal{A}$$

with

$$\mathcal{A} = \begin{pmatrix} \sqrt{\beta(s)} & 0 \\ -\frac{\alpha}{\sqrt{\beta}} & \frac{1}{\sqrt{\beta(s)}} \end{pmatrix} \quad \text{and} \quad \mathcal{R} = \begin{pmatrix} \cos(\Delta\mu) & \sin(\Delta\mu) \\ -\sin(\Delta\mu) & \cos(\Delta\mu) \end{pmatrix}$$

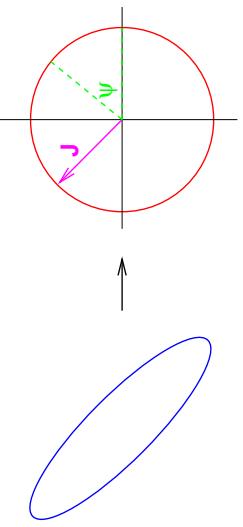
- This is just the Courant-Snyder transformation to get β, α, \dots etc., $\Delta\mu$ is the tune !

- That is: the Courant-Snyder analysis is just a **normal form transform** of the linear one turn matrix

- Works in more than one dimension

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Interlude: action - angle variables



- Once the particles "travel" on a circle:
 - Radius (say: $\sqrt{2J}$, with $J = \frac{x^2+x'^2}{2}$) is constant (invariant of motion): action J
 - Phase advances by constant amount: angle Ψ

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Another example: coupling

Assume a one-turn-matrix in 2D:

$$T = \begin{pmatrix} M & n \\ m & N \end{pmatrix}$$

M, m, N, n are 2-by-2 matrices. In case of coupling: $m \neq 0, n \neq 0$. we can try to re-write as:

$$T = \begin{pmatrix} M & n \\ m & N \end{pmatrix} = V \textcolor{red}{R} V^{-1}$$

with:

$$\textcolor{red}{R} = \begin{pmatrix} A & 0 \\ 0 & B \end{pmatrix} \quad \text{and} \quad V = \begin{pmatrix} \gamma_I & C \\ -C^t & \gamma_I \end{pmatrix}$$

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What have we obtained ?

- The matrix R is our simple rotation:
 - A and B are the one-turn-matrices for the "normal modes"
 - The matrix C contains the "coupling coefficients"
 - The matrix V transforms from the coordinates (x, x', y, y') into the "normal mode" coordinates (w, w', v, v') via the expression:
$$(x, x', y, y') = V(w, w', v, v')$$

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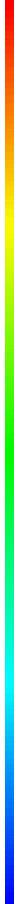
Normal forms - linear case

This is extremely useful when map is applied k times (e.g. k turns):

$$M^k(x, x') = AR^k A^{-1}(x, x') = AR^k(X, X')$$

- For multi-turns: study effect of map in normalized coordinates
 - Easier to apply k times using the simple map (e.g. a rotation of μ becomes just a rotation $k \cdot \mu$)
 - The A just transforms back to physical coordinates at the end (once !)

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The general philosophy (linear systems):

- Describe your elements by a **linear** map
- Combine all maps into a ring or beam line to get the **linear** one turn matrix
- Normal form analysis of the **linear** one turn matrix will give all the information

No need for any assumptions !

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The general philosophy (non-linear systems):

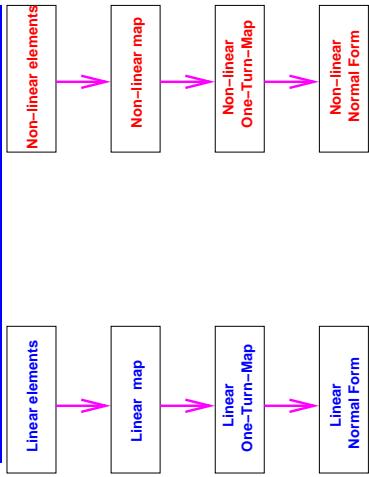
- Describe your elements by a **non-linear** map
- Combine all maps into a ring or beam line to get the **non-linear** one turn map
- Normal form analysis of the **non-linear** one turn map will give all the information

No need for any assumptions !

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The general philosophy



General formalism for all cases !

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Various types of non-linear MAPS

- Choice depends on the application
 - Taylor maps
 - Symplectic integration techniques
 - Lie transformations
 - Truncated power series algebra (TPSA), can also generate Taylor map from tracking
 - ...

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(A key concept: Symplecticity)

- Not all possible maps are allowed !
- Requires for a matrix $\mathcal{M} \rightarrow \mathcal{M}^T \cdot \mathcal{S} \cdot \mathcal{M} = \mathcal{S}$

with:

$$\mathcal{S} = \begin{pmatrix} 0 & 1 & 0 & 0 \\ -1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 1 \\ 0 & 0 & -1 & 0 \end{pmatrix}$$

- It basically means: \mathcal{M} is area preserving and

$$\lim_{n \rightarrow \infty} \mathcal{M}^n = \text{finite} \quad \Rightarrow \quad \det \mathcal{M} = 1$$

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Introducing non-linear elements

Effect of a (short) quadrupole depends **linearly** on amplitude (re-written from the matrix form):

$$\vec{z}(s_2) = \begin{pmatrix} x \\ x' \\ y \\ y' \end{pmatrix}_{s_2} = \begin{pmatrix} x \\ x' \\ y \\ y' \end{pmatrix}_{s_1} + \begin{pmatrix} 0 \\ k_1 \cdot x_{s_1} \\ 0 \\ k_1 \cdot y_{s_1} \end{pmatrix}_{s_1}$$

- $\vec{z}(s_2) = \mathbf{M} \cdot \vec{z}(s_1)$
- \mathbf{M} is a matrix

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Non-linear elements (e.g. sextupole)

Effect of a (thin) sextupole with strength k_2 is:

$$\vec{z}(s_2) = \begin{pmatrix} x \\ x' \\ y \\ y' \end{pmatrix}_{s_2} = \begin{pmatrix} x \\ x' \\ y \\ y' \end{pmatrix}_{s_1} + \begin{pmatrix} 0 \\ k_2 \cdot (x_{s_1} \cdot y_{s_1}) \\ 0 \\ \frac{1}{2}k_2 \cdot (x_{s_1}^2 - y_{s_1}^2) \end{pmatrix}_{s_1}$$

→ $\vec{z}(s_2) = \mathcal{M} \circ \vec{z}(s_1)$

- \mathcal{M} is **not** a matrix, i.e. cannot be expressed by matrix multiplication

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Non-linear elements

Cannot be written in linear matrix form !

We need something like:

$$\begin{aligned} z_1(s_2) = x(s_2) &= R_{11} \cdot \textcolor{blue}{x} + R_{12} \cdot \textcolor{blue}{x}' + R_{13} \cdot \textcolor{blue}{y} + \dots \\ &+ T_{111} \cdot \textcolor{red}{x}^2 + T_{112} \cdot \textcolor{red}{x} \textcolor{blue}{x}' + T_{122} \cdot \textcolor{red}{x}^{\prime 2} + \\ &+ T_{113} \cdot \textcolor{red}{x} \textcolor{blue}{y} + T_{114} \cdot \textcolor{red}{x} \textcolor{blue}{y}' + \dots \\ &+ U_{1111} \cdot \textcolor{red}{x}^3 + U_{1112} \cdot \textcolor{red}{x}^2 \textcolor{blue}{x}' + \dots \end{aligned}$$

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and the equivalent for all other variables ...

Higher order (Taylor -) MAPS:

We have (for: $j = 1 \dots 4$):

$$z_j(s_2) = \sum_{k=1}^4 R_{jk} z_k(s_1) + \sum_{k=1}^4 \sum_{l=1}^4 T_{jkl} z_k(s_1) z_l(s_1)$$

Let's call it : $\mathcal{A}_2 = [R, T]$ (second order map \mathcal{A}_2)

Higher orders can be defined as needed ...

$$\mathcal{A}_3 = [R, T, U] \implies + \sum_{k=1}^4 \sum_{l=1}^4 \sum_{m=1}^4 U_{jklm} z_k(s_1) z_l(s_1) z_m(s_1)$$

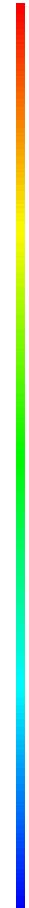


Higher order (Taylor -) MAPS:

Example: complete second order map for a (thick) sextupole with length L and strength K (in 4D):

$$\begin{aligned} x_2 &= x_1 + Lx'_1 & -K \left(\frac{L^2}{4}(x_1^2 - y_1^2) + \frac{L^3}{12}(x_1 x'_1 - y_1 y'_1) + \frac{L^4}{24}(x_1'^2 - y_1'^2) \right) \\ x'_2 &= x'_1 & -K \left(\frac{L}{2}(x_1^2 - y_1^2) + \frac{L^2}{4}(x_1 x'_1 - y_1 y'_1) + \frac{L^3}{6}(x_1'^2 - y_1'^2) \right) \\ y_2 &= y_1 + Ly'_1 & +K \left(\frac{L^2}{4}x_1 y_1 + \frac{L^3}{12}(x_1 y'_1 + y_1 x'_1) + \frac{L^4}{24}(x'_1 y'_1) \right) \\ y'_2 &= y'_1 & +K \left(\frac{L}{2}x_1 y_1 + \frac{L^2}{4}(x_1 y'_1 + y_1 x'_1) + \frac{L^3}{6}(x'_1 y'_1) \right) \end{aligned}$$

⚠ Definition of K not unique, can differ by some factor !!



Symplecticity for higher order MAPS

- Truncated Taylor expansions are not matrices !!
- It is the associated Jacobian matrix \mathcal{J} which must fulfil the symplecticity condition:

$$\mathcal{J}_{ik} = \frac{\partial z_2^i}{\partial z_1^k}$$

\mathcal{J} must fulfil: $\mathcal{J}^t \cdot \mathcal{S} \cdot \mathcal{J} = \mathcal{S}$

- In general: $\mathcal{J}_{ik} \neq \text{const}$ for truncated Taylor map can be difficult to fulfil for all z



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Thick versus thin magnets

- Real magnets have a finite length, i.e. thick magnets
- Thick magnet: field and length used to compute effect, i.e. the map
- Consequence: they are not always linear elements (even not dipoles, quadrupoles)
- For thick, non-linear magnets closed solution for maps often does not exist

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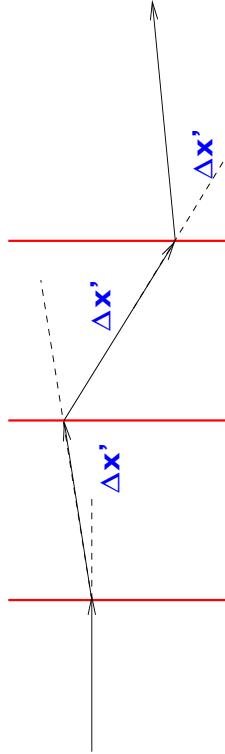
Thick versus thin magnets

- Thin "magnet": let the length go to zero, but keep field integral finite (constant)
- Thin dipoles and quadrupoles are linear elements
- Thin elements are much easier to use ...

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Moving through thin elements



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- No change of amplitudes x and y
- The momenta x' and y' receive an amplitude dependent deflection (**kick**)
 - $x' \rightarrow x' + \Delta x'$ and $y' \rightarrow y' + \Delta y'$



Using thin elements

- Can we approximate a thick element by thin element(s) ?
 - Yes, when the length is small or does not matter
 - Yes, when we can model the thick magnet correctly
 - What about accuracy, symplecticity etc. ??
 - Demonstrate with some simple examples

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Thick \Rightarrow thin quadrupole

$$\mathcal{M}_{s \rightarrow s+L} = \begin{pmatrix} \cos(L \cdot K) & \frac{1}{K} \cdot \sin(L \cdot K) \\ -K \cdot \sin(L \cdot K) & \cos(L \cdot K) \end{pmatrix}$$

- Exact map (matrix) for quadrupole
- What happens when we make it thin ?
 - Accuracy ?
 - Symplecticity ?
- (What follows is valid for all elements)

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Accuracy of thin lenses

$$\mathcal{M}_{s \rightarrow s+L} = \begin{pmatrix} \cos(L \cdot K) & \frac{1}{K} \cdot \sin(L \cdot K) \\ -K \cdot \sin(L \cdot K) & \cos(L \cdot K) \end{pmatrix}$$

► Start with exact map

► Taylor expansion in "small" length L :

$$L^0 \cdot \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix} + L^1 \cdot \begin{pmatrix} 0 & 1 \\ -K^2 & 0 \end{pmatrix} + L^2 \cdot \begin{pmatrix} -\frac{K^2}{2} & 0 \\ 0 & -\frac{K^2}{2} \end{pmatrix} + \dots$$

Slide 55



Accuracy of thin lenses (B)

► Keep up to first order term in L

$$\mathcal{M}_{s \rightarrow s+L} = L^0 \cdot \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix} + L^1 \cdot \begin{pmatrix} 0 & 1 \\ -K^2 & 0 \end{pmatrix}$$

$$\mathcal{M}_{s \rightarrow s+L} = \begin{pmatrix} 1 & L \\ -K^2 \cdot L & 1 \end{pmatrix} + \mathcal{O}(L^2)$$

► Precise to first order $\mathcal{O}(L^1)$

► $\det \mathcal{M} \neq 1$, non-symplectic

Slide 56



Accuracy of thin lenses (C)

$$\mathcal{M}_{s \rightarrow s+L} = \begin{pmatrix} 1 & L \\ -K^2 \cdot L & 1 \end{pmatrix} + \mathcal{O}(L^2)$$



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$$\mathcal{M}_{s \rightarrow s+L} = \begin{pmatrix} 1 & L \\ -K^2 \cdot L & 1 - K^2 L^2 \end{pmatrix}$$

- Precise to first order $\mathcal{O}(L^1)$
- "symplectified" by adding term $-K^2 L^2$
(it is wrong to $\mathcal{O}(L^2)$ anyway ...)



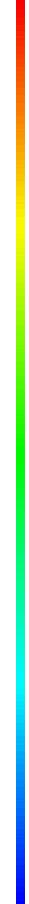
Accuracy of thin lenses

- Keep up to second order term in L

$$\mathcal{M}_{s \rightarrow s+L} = \begin{pmatrix} 1 - \frac{1}{2} K^2 L^2 & L \\ -K^2 \cdot L & 1 - \frac{1}{2} K^2 L^2 \end{pmatrix} + \mathcal{O}(L^3)$$

- Precise to second order $\mathcal{O}(L^2)$
- More accurate than (C), but not symplectic

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Accuracy of thin lenses (D)

- Symplectification like:

$$\mathcal{M}_{s \rightarrow s+L} = \begin{pmatrix} 1 - \frac{1}{2}K^2L^2 & L - \frac{1}{4}K^2L^3 \\ -K^2 \cdot L & 1 - \frac{1}{2}K^2L^2 \end{pmatrix} + \mathcal{O}(L^3)$$

- Precise to second order $\mathcal{O}(L^2)$
- Fully symplectic

Slide 59



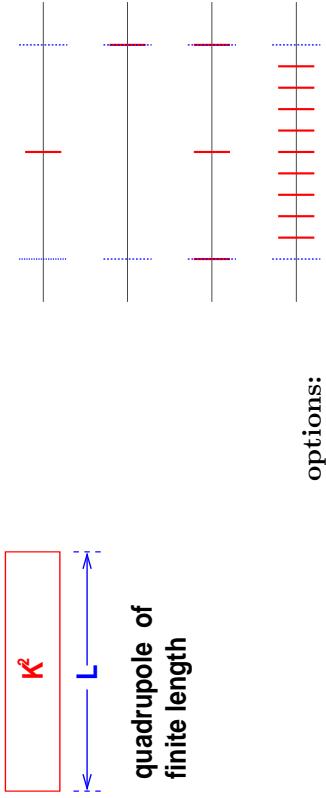
Accuracy of thin lenses

- Looks like we made some arbitrary changes and called it "symplectification"
- Is there a physical picture behind the approximations ?
- Yes, **geometry** of thin lens kicks ...
- A thick element is split into thin elements with drifts between them

Slide 60



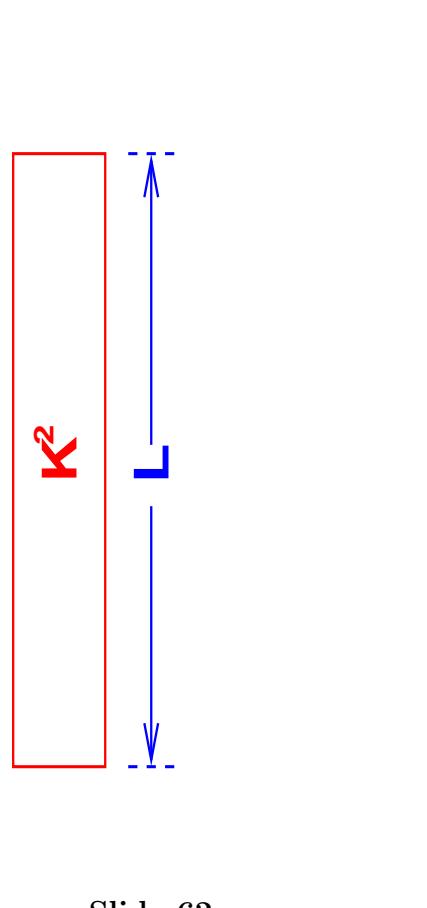
Thick \rightarrow thin quadrupole



Which is a good strategy? \rightarrow accuracy and speed

options:

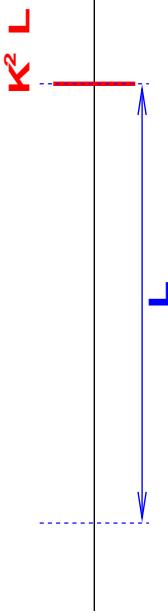
Thick quadrupole ..



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Slide 62

First order ..



➤ One thin quadrupole "kick" and one drift combined

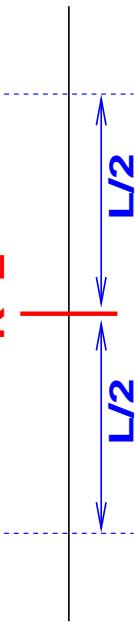
➤ Resembles "symplectification" of type (C)

$$\begin{pmatrix} 1 & 0 \\ -K^2 \cdot L & 1 \end{pmatrix} \begin{pmatrix} 1 & L \\ 0 & 1 \end{pmatrix} = \begin{pmatrix} 1 & L \\ -K^2 \cdot L & 1 - K^2 L^2 \end{pmatrix}$$



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Second order ..



➤ One thin quadrupole "kick" between two drifts

➤ Resembles more accurate, symplectic model of type (D)

$$\begin{pmatrix} 1 & \frac{1}{2}L \\ 0 & 1 \end{pmatrix} \begin{pmatrix} 1 & 0 \\ -K^2 \cdot L & 1 \end{pmatrix} \begin{pmatrix} 1 & \frac{1}{2}L \\ 0 & 1 \end{pmatrix} = \begin{pmatrix} 1 - \frac{1}{2}K^2 L^2 & L - \frac{1}{4}K^2 L^3 \\ -K^2 \cdot L & 1 - \frac{1}{2}K^2 L^2 \end{pmatrix}$$

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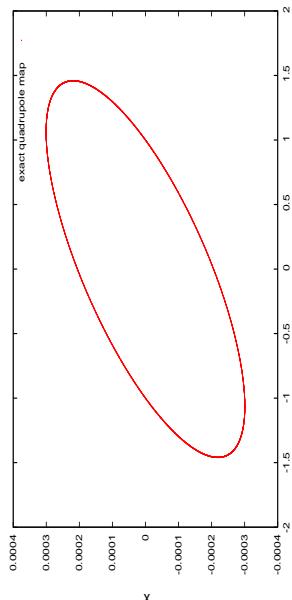
Accuracy of thin lenses

- One kick at the end (or beginning):
 - Error (inaccuracy) of first order $\mathcal{O}(L^1)$
- One kick in the centre:
 - Error (inaccuracy) of second order $\mathcal{O}(L^2)$
- It is very relevant **how** to apply thin lenses
- Aim should be to be precise and fast (and simple to implement)

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What is the point ??

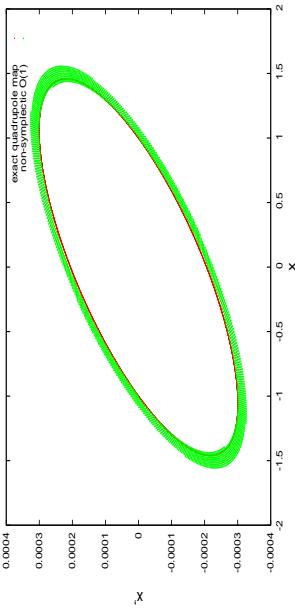


Slide 66

- Phase ellipse - quadrupole exact solution



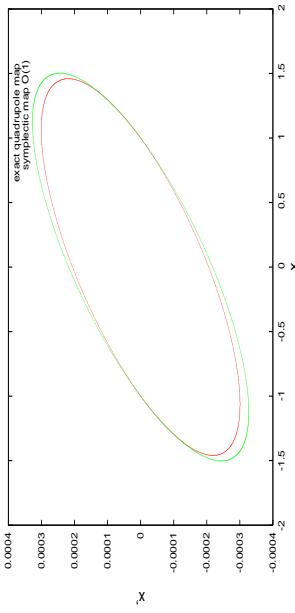
Quadrupole non-symplectic solution



► Non-symplectic: particles spiral towards outside

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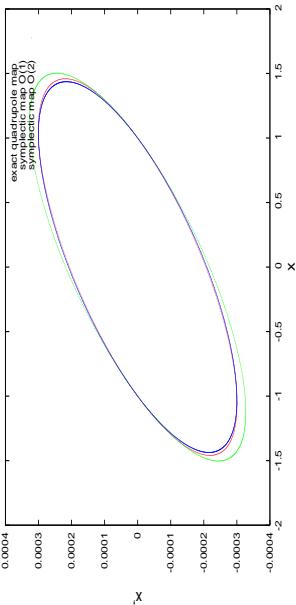
Quadrupole symplectic $O(L^1)$ solution



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► symplectic: but phase space ellipse not accurate

Quadrupole symplectic $\mathcal{O}(L^2)$ solution



➤ symplectic: phase space ellipse accurate enough

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Can we do better ?

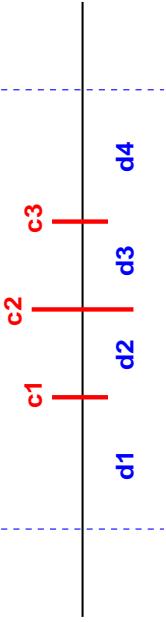
- Try more slices, e.g. 3 kicks:
- How to put them ?
- Hope you are already alerted ...
- Allow that they are at different positions **and** have different strengths
- Minimize the inaccuracy

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Question: is one of the options obviously wrong ? If yes, why ?

Can we do better ?

▷ Try a model with 3 kicks:



▶ To get best accuracy (i.e. deviation from exact solution):

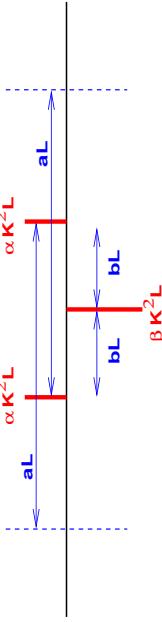
▷ Optimize kicks c_1, c_2, c_3

▷ Optimize drifts d_1, d_2, d_3, d_4

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Can we do better ?

▷ Try a model with 3 kicks:



▷ with:

$a \approx 0.6756, b \approx -0.1756, \alpha \approx 1.3512, \beta \approx -1.7024$

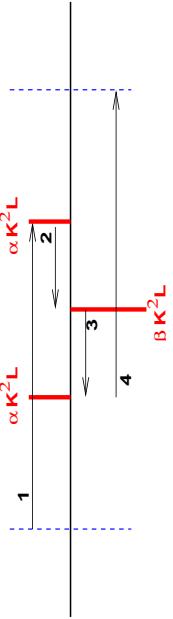
▷ we have a $\mathcal{O}(4)$ integrator ...

▷ (a $\mathcal{O}(6)$ integrator would require 9 kicks (!) ...)

Slide 72

Can we do better ?

- Try a model with 3 kicks:
 - Must track backwards ! Change interpretation !
 - Thin lenses not a new sequence of magnets (a la MAD)
 - What about space charge calculations ?



Slide 73

Symplectic integration

- What we do is **Symplectic Integration**
- From a lower order integration scheme (1 kick), construct higher order scheme
- Formally (for the formulation of $S_k(t)$ see later):
 - From a 2nd order scheme (1 kick) $S_2(t)$ we construct a 4th order scheme (3 kicks = 3 x 1 kick) like:
$$S_4(t) = S_2(x_1 t) \circ S_2(x_0 t) \circ S_2(x_1 t) \quad \text{with:}$$

$$x_0 = \frac{-2^{1/3}}{2 - 2^{1/3}} \approx -1.7024 \quad x_1 = \frac{1}{2 - 2^{1/3}} \approx 1.3512$$

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Symplectic integration

- Can be considered as an iterative scheme (see e.g.
H. Yoshida, 1990, E. Forest, 1998²⁾):
- If $S_{2k}(t)$ is a symmetric integrator of order $2k$, then:

$$S_{2k+2}(t) = S_{2k}(x_1 t) \circ S_{2k}(x_0 t) \circ S_{2k}(x_1 t) \quad \text{with:}$$

$$x_0 = \frac{-\sqrt[2k+1]{2}}{2 - \sqrt[2k+1]{2}} \quad x_1 = \frac{1}{2 - \sqrt[2k+1]{2}}$$

- Higher order integrators can be obtained in a similar way

²⁾ E. Forest, "Beam Dynamics, A New Attitude and Framework", 1998

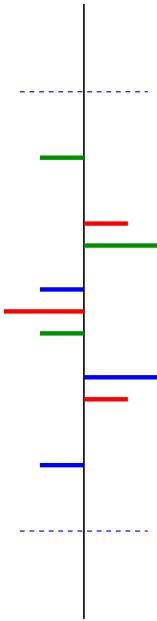
Slide 75

Symplectic integration

- Example: From a 4th order to 6th order
 $S_6(t) = S_4(x_1 t) \circ S_4(x_0 t) \circ S_4(x_1 t)$
- We get 3 times 4th order with 3 kicks each, we have the 9 kick, 6th order integrator mentioned earlier

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Integrator of order 6



- Requires 9 kicks
- We have 3 interleaved 4th order integrators
- Can be used in iterative scheme

Slide 77



Some remarks:

- We have used a linear map (quadrupole) to demonstrate the integration
- Can that be applied for other maps (solenoids, higher order, non-linear maps) ?
 - Yes !!
 - We get the same integrators !
 - Proof and systematic (and easy) extension in the form of Lie operators²⁾ (see later)
- Best accuracy for thin lenses !

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²⁾ H. Yoshida, Physics Letters A, Volume 150 (1990) 262.



Accuracy of thin lenses

What about accuracy of **non-linear** elements ?

assume a general case:

$$x'' = f(x)$$

➤ Disadvantage : usually a closed solution through the element does not exist, integration necessary

➤ Advantage : They are usually thin (thinner than dipoles, quadrupoles ...)

- Dipoles: ≈ 14.3 m

- Quadrupole: $\approx 2 - 5$ m

- Sextupoles, Octupoles: ≈ 0.30 m

➤ Can try our simplest thin lens approximation first ...

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Accuracy of thin lenses - our $\mathcal{O}(2)$ model

$$1.Step \quad \begin{pmatrix} x \\ x' \end{pmatrix}_{s_1+L/2} = \begin{pmatrix} 1 & \frac{L}{2} \\ 0 & 1 \end{pmatrix} \circ \begin{pmatrix} x \\ x' \end{pmatrix}_{s_1}$$

$$2.Step \quad \begin{pmatrix} x \\ x' \end{pmatrix}_{s_1+L/2} = \begin{pmatrix} x \\ x' + \Delta x' \end{pmatrix}_{s_1+L/2}$$

$$3.Step \quad \begin{pmatrix} x \\ x' \end{pmatrix}_{s_1+L} = \begin{pmatrix} 1 & \frac{L}{2} \\ 0 & 1 \end{pmatrix} \circ \begin{pmatrix} x \\ x' \end{pmatrix}_{s_1+L/2}$$

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Accuracy of thin lenses

Assume the general case:

$$x'' = f(x)$$

► Using this thin lens approximation (type D, $\mathcal{O}(2)$) gives:

$$x'(L) \approx x'_0 + Lf(x_0 + \frac{L}{2}x'_0)$$

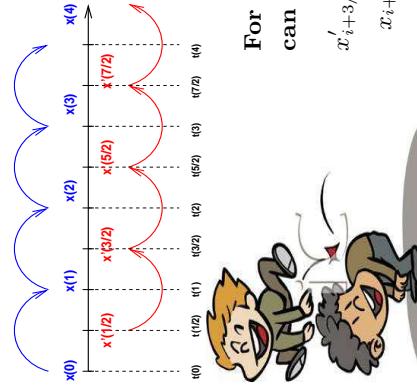
$$x(L) \approx x_0 + \frac{L}{2}(x'_0 + x'(L))$$

► This is also called "leap frog" algorithm/integration

► It is symplectic (... and time reversible) !!

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Interlude ...



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For any: $x'' = f(x, x', t)$ we
can solve it by:

$$x'_{i+3/2} \approx x'_{i+1/2} + f(x_{i+1})\Delta t$$

$$x_{i+1} \approx x_i + x'_{i+1/2}\Delta t$$

Accuracy of thin lenses

Accuracy of "leap frog" algorithm/integration"

the (exact) Taylor expansion gives:

$$x(L) = x_0 + x'_0 L + \frac{1}{2} f(x_0)L^2 + \frac{1}{6} x'_0 f'(x_0)L^3 + \dots$$

the (approximate) "leap frog" algorithm gives:

$$x(L) = x_0 + x'_0 L + \frac{1}{2} f(x_0)L^2 + \frac{1}{4} x'_0 f'(x_0)L^3 + \dots$$

➤ Errors are $\mathcal{O}(L^3)$

➤ For small L acceptable, and symplectic, extend to our symplectic integration

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Accuracy of thin lenses

For bar/coffee discussions:

why did I write:

$$x'' = f(x)$$

and not:

$$x'' = f(x, x')$$

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Accuracy of thin lenses

An application, assume a (1D) sextupole with (definition of k not unique !):

$$x'' = k \cdot x^2 = f(x)$$

using the thin lens approximation (type D) gives:

$$x(L) = x_0 + x'_0 L + \frac{1}{2} k x_0^2 L^2 + \frac{1}{2} k x_0 x'_0 L^3 + \frac{1}{8} k x_0'^2 L^4$$

$$x'(L) = x'_0 + k x_0^2 L + k x_0 x'_0 L^2 + \frac{1}{4} k x_0'^2 L^3$$

Map for thick sextupole of length L in thin lens approximation, accurate to $\mathcal{O}(L^2)$

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Some comments:

- We have interleaved kicks with drifts
- Is that always necessary ?
 - No !
 - Can be any map with an exact expression
 - In most cases the exact map is a linear map (matrix)
- We have derived element maps for tracking from the equation of motion using this technique ➔ can track now

Slide 86

The plan now ...

- Extend all concepts to non-linear dynamics
 - Lagrangian and Hamiltonian dynamics
 - How to use that \rightarrow Lie transforms
 - How to analyse that \rightarrow Non-linear normal forms
 - How to analyse that better \rightarrow Differential Algebra (DA)

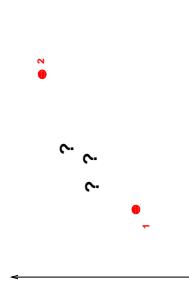
- Avoid abstract definitions and formulation, but:

- Intuitive (but correct !) treatment
- Useful formulae and examples
- Real life examples and demonstration (DA)

Slide 87

Hamilton principle

- Problem: describe the motion of a system (e.g. 1 or more particles) between times t_1 and t_2



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- Describe by coordinates q_i $(i = 1, n)$
- n are degrees of freedom of the system

Hamilton principle

- Describe motion by a function L

$$L(q_1, \dots, q_n, \dot{q}_1, \dots, \dot{q}_n, t)$$

(q_1, \dots, q_n) ... generalized coordinates

$(\dot{q}_1, \dots, \dot{q}_n)$... generalized velocities

- The function L defines the **Lagrange function**

- The integral $I = \int L(q_i, \dot{q}_i, t) dt$ defines the **action**

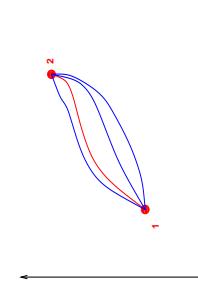
Without proof or derivation:

$$L = T - V = \text{kinetic energy} - \text{potential energy}$$

Slide 89

Hamilton principle

$$I = \int_1^2 L(q_i, \dot{q}_i, t) dt = \text{extremum}$$



Slide 90

- Hamiltonian principle: system moves such that the action I becomes an extremum

Extremum principle ?

- Not new:
 - Used in optics: Fermat principle
 - Quantum mechanics (path integrals)
 - General relativity
 - ...

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Lagrange formalism

Without proof:

$$I = \int_1^2 L(q_i, \dot{q}_i, t) dt = \text{extremum}$$

is fulfilled when:

$$\frac{d}{dt} \frac{\partial L}{\partial \dot{q}_i} - \frac{\partial L}{\partial q_i} = 0$$

(Euler - Lagrange equation)

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From Lagrangian to Hamiltonian ..

- Lagrangian $L(q_1, \dots, q_n, \dot{q}_1, \dots, \dot{q}_n, t)$ in generalized coordinates and velocities
- Provides (n) second order differential equations
- Try to get:
 - Generalized momenta instead of velocities
 - First order differential equations (always solvable)

Corresponding (so-called conjugate) momenta p_j are:

$$p_j = \frac{\partial L}{\partial \dot{q}_j}$$

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Interlude: canonical transformations

- Not treated here in detail (see literature or backup slides)
- Main ideas:
 - Find variables where equations are easier to solve
 - Action-angle variables: lead to invariants of motion (see later)

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From Lagrangian to Hamiltonian ..

■ Lagrangian:

- n second order equations
- n -dimensional coordinate space

■ Hamiltonian:

- $2n$ first order equations
- $2n$ -dimensional phase space

Slide 95



From Lagrangian to Hamiltonian ..

Once we know what the canonical momenta p_i are: the **Hamiltonian** is a transformation of the **Lagrangian**:

$$H(q_j, \mathbf{p}_j, t) = \sum_i q_i p_i - L(q_j, \dot{q}_j, t)$$

without proof.

$H = T + V$ = kinetic energy + potential energy

we obtain 2 first order equation of motion:

$$\frac{\partial H}{\partial q_j} = -\dot{p}_j = -\frac{dp_j}{dt}, \quad \frac{\partial H}{\partial p_j} = \dot{q}_j = \frac{dq_j}{dt}$$



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Hamiltonian of particle in EM fields

For the Hamiltonian of a (relativistic) particle in a electro-magnetic field we have:

$$\mathcal{H}(\vec{x}, \vec{p}, t) = c\sqrt{(\vec{p} - e\vec{A}(\vec{x}, t))^2 + m_0^2c^2} + e\Phi(\vec{x}, t)$$

where $\vec{A}(\vec{x}, t)$, $\Phi(\vec{x}, t)$ the vector and scalar scalar potential

Using canonical variables and the design path length s as independent variable (bending in x-plane):

$$\mathcal{H} = -(1 + \frac{x}{\rho}) \cdot \sqrt{(1 + \delta)^2 - p_x^2 - p_y^2} + \frac{x}{\rho} + \frac{x^2}{2\rho^2} - \frac{A_s(x, y)}{B_0\rho}$$

where $\delta = (p - p_0)/p$ is relative momentum deviation and $A_s(x, y)$ longitudinal component of the vectorpotential.

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Hamiltonian of particle in EM fields

The magnetic fields can be described with the multipole expansion:

$$B_y + iB_x = \sum_{n=1}^{\infty} (b_n + ia_n)(x + iy)^{n-1}$$

and since $\vec{B} = \nabla \times \vec{A}$:

$$A_s = \sum_{n=1}^{\infty} \frac{1}{n} [(b_n + ia_n)(x + iy)^n]$$

► For a large machine ($x \gg \rho$) we expand the root and sort the variables:

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Hamiltonian (for large machine) ..

$$\mathcal{H} = \frac{p_x^2 + p_y^2}{2(1+\delta)} - \underbrace{\frac{x\delta}{\rho}}_{\text{dispersive focusing}} + \underbrace{\frac{x^2}{2\rho^2}}_{\text{dipole}} + \underbrace{\frac{k_1(x^2 - y^2)}{2}}_{\text{quadrupole}} + \underbrace{\frac{k_2(x^3 - 3xy^2)}{3}}_{\text{sextupole}}$$

- The Hamiltonian describes exactly the motion of a particle through a magnet
- Basis to extend the linear formalism to a non-linear treatment

But how do we use it ??

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Poisson brackets

Introduce Poisson bracket for a differential operator:

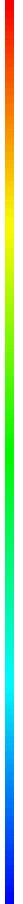
$$[f, g] = \sum_{i=1}^n \left(\frac{\partial f}{\partial q_i} \frac{\partial g}{\partial p_i} - \frac{\partial f}{\partial p_i} \frac{\partial g}{\partial q_i} \right)$$

Here the variables q_i, p_i are canonical variables, f and g are functions of q_i and p_i .
We can now write:

$$[q_i, H] = \frac{\partial H}{\partial p_i} = \frac{dq_i}{dt}$$

$$[p_i, H] = -\frac{\partial H}{\partial q_i} = \frac{dp_i}{dt}$$

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Poisson brackets

➤ Poisson bracket $[g, H]$ describes the time evolution of the system

It is a special case of:

$$\frac{du}{dt} = [u, H] + \frac{\partial u}{\partial t}$$

If H does not explicitly depend on time and:

$$[f, H] = 0$$

implies that f is an invariant of the motion !

Slide 101



Lie transformations

We can define:

$$:f:g = [f, g]$$

where $:f:$ is an operator acting on the function g :

$$:f := [f,]$$

We can define powers as:

$$(:f:)^2 g = :f:(:f:g) = [f, [f, g]] \quad \text{etc.}$$

in particular:

$$e^{:f:} = \sum_{i=0}^{\infty} \frac{1}{i!} (:f:)^i$$

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Lie transformations - example

Lie transforms acting on functions like x, p (canonical momentum, instead of x):

$$e^{-Lp^2/2}x = x - \frac{1}{2}L:p^2:x + \frac{1}{8}L^2(:p^2)^2x + \dots$$

$$= x + Lp$$

$$e^{-Lp^2/2}p = p - \frac{1}{2}L:p^2:p + \dots$$

$$= p$$

or more complicated:

$$e^{-Lp^2/2}x = x - \frac{1}{2}L:p^2:x + \frac{1}{8}L^2(:p^2)^2x + \dots$$

$$= x + Lp$$

$$e^{-Lp^2/2}p = p - \frac{1}{2}L:p^2:p + \dots$$

$$= p$$

This is the transformation of a drift space of length L !!

Lie transformations

Acting on the phase space coordinates:

$$e^{if}(x, p)_0 = (x, p)_1$$

that is:

$$e^{if}x_0 = x_1$$

$$e^{if}p_0 = p_1$$

► Lie transforms describe how to go from one point to another.

► Through a machine element (drift, magnet ...) fully described by f

► But what is f ?

Lie transformations

- The generator f is the Hamiltonian of the element !
- We use the Hamiltonian to describe the motion through an individual element
- Inside a single element the motion is "smooth" (in the full machine it is not !)

Slide 105



Some formulae for Lie transforms

With a constant, f, g, k arbitrary functions:

$$:a: = 0 \quad e^{:a:} = 1$$

$$:f:a = 0 \quad e^{:f:a} = a$$

$$e^{:f:}g(x) = g(e^{:f:}x)$$

$$e^{:f:}G(:g:)e^{-:f:} = G(:e^{:f:}g:)$$

$$e^{:f:}[g, h] = [e^{:f:}g, e^{:f:}h]$$

$$(e^{:f:})^{-1} = e^{-:f:}$$

and very important:

$$e^{:f:}e^{:g:}e^{-:f:} = e^{:e^{:f:}g:}$$



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Useful formulae for calculations

With x coordinate, p momentum, as usual:

$$\begin{aligned} :x: &= \frac{\partial}{\partial p} & :p: &= -\frac{\partial}{\partial x} \\ :x:^2 &= \frac{\partial^2}{\partial p^2} & :p:^2 &= \frac{\partial^2}{\partial x^2} \\ :x^2: &= 2x \frac{\partial}{\partial p} & :p^2: &= -2p \frac{\partial}{\partial x} \\ :xp: &= p \frac{\partial}{\partial x} - x \frac{\partial}{\partial p} & :x:p: &= :p:x: = -\frac{\partial^2}{\partial x \partial p} \end{aligned}$$


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More useful formulae for calculations

With x coordinate, p momentum, as usual:

$$\begin{aligned} :p^2:x &= \frac{\partial p^2}{\partial x} \frac{\partial x}{\partial p} - \frac{\partial p^2}{\partial p} \frac{\partial x}{\partial x} = -2p \\ :p^2:p &= \frac{\partial p^2}{\partial x} \frac{\partial p}{\partial p} - \frac{\partial p^2}{\partial p} \frac{\partial p}{\partial x} = 0 \\ (:p^2)^2x &=:p^2:(:p^2:x) =:p^2:(-2p) = 0 \\ (:p^2)^2p &=:p^2:(:p^2:p) =:p^2:(0) = 0 \end{aligned}$$


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More examples (1D):

For:

$$f = -\frac{L}{2}p^2$$

we obtained:

$$\begin{aligned} e^{if}x &= x + Lp \\ e^{if}p &= p \end{aligned}$$

➤ Drift space, seen that already

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More examples (1D):

For:

$$f = -\frac{L}{2}(k^2x^2 + p^2)$$

we would get (try it !):

$$\begin{aligned} e^{if}x &= e^{i-\frac{L}{2}(k^2x^2+p^2)}x \\ e^{if}p &= e^{i-\frac{L}{2}(k^2x^2+p^2)}p \end{aligned}$$

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Remember:

$$e^{if}g = \sum_{n=0}^{\infty} \frac{f^{(n)}}{n!} g$$

More examples (1D):

For:

$$f = -\frac{L}{2}(k^2x^2 + p^2)$$

we would get (try it !):

$$\begin{aligned} e^{i(-\frac{L}{2}(k^2x^2 + p^2))} \cdot x &= \sum_{n=0}^{\infty} \left(\frac{(-k^2L^2)^n}{(2n)!} \cdot x + L \frac{(-k^2L^2)^n}{(2n+1)!} \cdot p \right) \\ e^{i(-\frac{L}{2}(k^2x^2 + p^2))} \cdot p &= \sum_{n=0}^{\infty} \left(\frac{(-k^2L^2)^n}{(2n)!} \cdot p - k \frac{(-k^2L^2)^n}{(2n+1)!} \cdot x \right) \end{aligned}$$


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More examples (1D):

For:

$$f = -\frac{L}{2}(k^2x^2 + p^2)$$

we would get (try it !):

$$\begin{aligned} e^{if} \cdot x &= x \cos(kL) + \frac{p}{k} \sin(kL) \\ e^{if} \cdot p &= -kx \sin(kL) + p \cos(kL) \end{aligned}$$


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► Thick, focusing quadrupole !

Hamiltonians of some **thick** machine elements (2D)

dipole:

$$H = -\frac{x\delta}{\rho} + \frac{x^2}{2\rho^2} + \frac{p_x^2 + p_y^2}{2(1+\delta)}$$

quadrupole:

$$H = \frac{1}{2}k_1(x^2 - y^2) + \frac{p_x^2 + p_y^2}{2(1+\delta)}$$

sextupole:

$$H = \frac{1}{3}k_2(x^3 - 3xy^2) + \frac{p_x^2 + p_y^2}{2(1+\delta)}$$

octupole:

$$H = \frac{1}{4}k_3(x^4 - 6x^2y^2 + y^4) + \frac{p_x^2 + p_y^2}{2(1+\delta)}$$

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Many machine elements

- We can combine many machine elements by applying one transformation after the other:

$$e^{ih} = e^{if_1} \cdot e^{if_2} \cdot \dots \cdot e^{if_N}$$

- Not restricted to matrices, i.e. linear elements ...
- And arrive at a transformation for the full ring
 - a one turn map
- The one turn map is the exponential of the effective Hamiltonian:

$$\mathcal{M}_{ring} = e^{-iC\mathcal{H}_{eff}}$$

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Why all that ???

concatenation very easy:

$$e^{:h:} = e^{:f:} \cdot e^{:g:} = e^{:f+g:}$$

when f and g commute (i.e. $[f, g] = [g, f] = 0$)

■ Otherwise formalism exist

■ The miracles:

- Poisson brackets create **symplectic** maps
- Exponential form $e^{:h:}$ is **always** symplectic
- Better: the exponent is directly connected to the invariant of the transfer map !!

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Concatenation

To combine:

$$e^{:h:} = e^{:f:} \cdot e^{:g:}$$

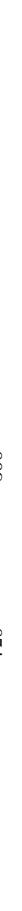
We can use the formula (Baker-Campbell-Hausdorff (BCH)):

$$\begin{aligned} h = f &+ g + \frac{1}{2}[f, g] + \frac{1}{12}[f, [f, g]] + \frac{1}{12}[g, [g, f]] \\ &+ \frac{1}{24}[f, [g, [g, f]]] - \frac{1}{720}[g, [g, [g, [g, f]]]] \\ &- \frac{1}{720}[f, [f, [f, [g, [g, f]]]]] + \frac{1}{360}[g, [f, [f, [g, [g, f]]]]] + \dots \end{aligned}$$

or :

$$\begin{aligned} h = f &+ g + \frac{1}{2}:f:g + \frac{1}{12}:f:^2g + \frac{1}{12}:g:^2f \\ &+ \frac{1}{24}:f::g:^2f - \frac{1}{720}:g:^4f \\ &- \frac{1}{720}:f:^4g + \frac{1}{360}:g::f:^3g + \dots \end{aligned}$$

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Concatenation

To combine:

$$e^{\dot{h}} = e^{\dot{f}} \cdot e^{\dot{g}}$$

if one of them (f or g) is small, can truncate the series and get a very useful formula. Assume g is small:

$$e^{\dot{f}} \cdot e^{\dot{g}} = e^{\dot{h}} = \exp \left[: f + \left(\frac{: f:}{1 - e^{-\dot{f}}} \right) g + \mathcal{O}(g^2) : \right]$$

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Non-linear kicks

General thin lens kick $f(x)$:

$$e^{\int_0^x f(x') dx'}$$

gives for the map:

$$\begin{aligned} x &= x_0 \\ p &= p_0 + f(x) \end{aligned}$$

Example: thin lens multipole of order n ($f(x) = a \cdot x^n$):

$$e^{a \cdot x^n}$$

gives for the map:

$$\begin{aligned} x &= x_0 \\ p &= p_0 + a n x^{n-1} \end{aligned}$$



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From the Hamiltonian to the map

We have seen that given the Hamiltonian f of a machine element is known, the Lie operator becomes:

$$f \rightarrow :f :$$

the corresponding map is than:

$$e^{:f:} (e^{-L_f})$$

This map is always symplectic and we have (in 1D):

$$e^{:f:} x_0 = x_1$$

$$e^{:f:} p_0 = p_1$$

or using $Z = (x, p_x, y, p_y, \dots)$:

$$e^{:f:} Z_0 = Z_1$$



From the map to the Hamiltonian

The other question \rightarrow assuming we have:

$$\mathcal{M} \equiv \begin{pmatrix} \cos(\mu) + \alpha \sin(\mu) & \beta \sin(\mu) \\ -\gamma \sin(\mu) & \cos(\mu) - \alpha \sin(\mu) \end{pmatrix}$$

i.e.:

$$\mathcal{M} Z_0 = Z_1$$

how do we find the corresponding form for f ?

$$\mathcal{M} \leftrightarrow e^{:f:} (e^{-\mu f})$$

From the map to the Hamiltonian

For the linear matrix we know that f must be a quadratic form in (x, p, \dots) .
Any quadratic form can be written as:

$$f = -\frac{1}{2}Z^*FZ \quad [= -\frac{1}{2}(a \cdot x^2 + b \cdot xp + c \cdot p^2)]$$

where F is a symmetric, positive definite (why?) matrix.
Then we can write (without proof, see e.g. Dragt):

$$: f : Z = SFZ$$

where S is the "symplecticity" matrix.

Therefore we get for the Lie transformation:

$$e^{if}Z \leftrightarrow e^{SF}Z$$

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From the map to the Hamiltonian

Since we have $n = 2$, we get (using Hamilton–Cayley theorem):

$$e^{SF} = \exp \begin{pmatrix} b & c \\ -a & -b \end{pmatrix} = a_0 + a_1 \begin{pmatrix} b & c \\ -a & -b \end{pmatrix}$$

We now have to find a_0 and a_1 !

The eigenvalues of SF are:

$$\lambda_{\pm} = \pm i\sqrt{ac - b^2}$$

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From the map to the Hamiltonian

This tells us for the coefficients the conditions:

$$e^{\lambda_+} = a_0 + a_1 \cdot \lambda_+$$

$$e^{\lambda_-} = a_0 + a_1 \cdot \lambda_-$$

and therefore:

$$a_0 = \cos(\sqrt{ac - b^2})$$

$$a_1 = \frac{\sin(\sqrt{ac - b^2})}{\sqrt{ac - b^2}}$$

and

$$e^{SF} = \cos(\sqrt{ac - b^2}) + \frac{\sin(\sqrt{ac - b^2})}{\sqrt{ac - b^2}} \begin{pmatrix} b & c \\ -a & -b \end{pmatrix}$$

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From the map to the Hamiltonian

For a general 2×2 matrix:

$$M = \begin{pmatrix} m_{11} & m_{12} \\ m_{21} & m_{22} \end{pmatrix}$$

we get by comparison:

$$\cos(\sqrt{ac - b^2}) = \frac{1}{2} \text{tr}(M)$$

and

$$\frac{a}{-m_{21}} = \frac{2b}{m_{11} - m_{22}} = \frac{c}{m_{12}} = \frac{\sqrt{ac - b^2}}{\sin(\sqrt{ac - b^2})}$$

for the quadratic form of f :

$$f = -\frac{1}{2}(a \cdot x^2 + b \cdot xp + c \cdot p^2)$$

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From the map to the Hamiltonian

For the example of a drift:

$$\mathcal{M} \equiv \begin{pmatrix} 1 & L \\ 0 & 1 \end{pmatrix}$$

we find:

$$a = 0, \quad b = 0, \quad c = L$$

and for the generator:

$$f = -\frac{1}{2}(Lp^2)$$



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From the map to the Hamiltonian

For the example of a thin quadrupole:

$$\mathcal{M} \equiv \begin{pmatrix} 1 & 0 \\ -\frac{1}{f} & 1 \end{pmatrix}$$

we find:

$$a = \frac{1}{f}, \quad b = 0, \quad c = 0$$

and for the generator:

$$f = -\frac{1}{2f}(x^2)$$



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A very important example ...

$$\mathcal{M} \equiv \begin{pmatrix} \cos \mu + \alpha \sin(\mu) & \beta \sin \mu \\ -\gamma \sin \mu & \cos \mu - \alpha \sin(\mu) \end{pmatrix}$$

corresponds to:

$$e^{:h:} = e^{:f_2:} = e^{-\mu \frac{1}{2}(\gamma x^2 + 2\alpha x p + \beta p^2)};$$

In this form f is: $-\mu \cdot (\text{Courant-Snyder invariant})$

$$e^{:h:} = e^{:f_2:} = e^{:-\mu \epsilon:}$$

► We have standard ($e^{:f_2:}$) for the linear one-turn-matrix
(a rotation)...

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Normal forms non-linear case

Normal form transformations can be generalized for non-linear maps (i.e. not matrices). If \mathcal{M} is our usual one-turn-map, we try to find a transformation:

$$\mathcal{N} = \mathcal{A} \mathcal{M} \mathcal{A}^{-1}$$

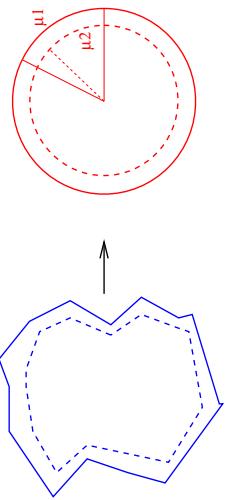
► where \mathcal{N} is a simple form (like the rotation we had before)

Of course we now do not have matrices, but find a Lie transform \mathcal{F} for:

$$\mathcal{N} = e^{-:h:} = e^{:F:} \mathcal{M} e^{-:F:}$$

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Normal forms



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- More complicated transformation required
- Transform to coordinates where map is just a rotation
- Rotation angle now amplitude dependent !



Normal forms - non-linear case

The transformation A :

$$A = e^{-iF}$$

is the canonical transformation which produces our simple form

- Works for any kind of local perturbation
- Formalism and software tools exist to find F (see e.g. Chao¹⁾ or E.Forest, M. Berz, J. Irwin, SSC-166)
- Formalism to find F is valid for any n-th order polynomial
- Once we know $h_{eff}(J_x, J_y, \delta)$ we can derive everything !

¹⁾ A. Chao, Lecture Notes on Topics in Accelerator Physics, 2001

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Normal forms - non-linear case

Once we can write the map as (now example in 3D):

$$\mathcal{N} = e^{i h_{eff}(J_x, J_y, \delta)}$$

where h_{eff} depends only on J_x, J_y , and δ , then we have the times:

$$Q_x(J_x, J_y, \delta) = \frac{1}{2\pi} \frac{\partial h_{eff}}{\partial J_x}$$

$$Q_y(J_x, J_y, \delta) = \frac{1}{2\pi} \frac{\partial h_{eff}}{\partial J_y}$$

and the change of path length:

$$\Delta s = - \frac{\partial h_{eff}}{\partial \delta} = \alpha_c \delta$$

- Dependence on J is amplitude detuning, dependence on δ are the chromaticities !



Example: sextupole (1D)

A linear map followed by a single (weak) sextupole:

$$\mathcal{M} = e^{-\frac{\mu}{2}x^2 + p^2} : e^{i f_3} : = e^{i -\mu J_x} : e^{i k x^3} :$$

using the BCH formula we get immediately for h and h_{eff} :

$$h = -\mu J + \left(\frac{i -\mu J_x}{1 - e^{i \mu J_x}} \right) \cdot k x^3$$

$$h = -\mu J - \frac{3}{8} \mu k (2\beta J)^{3/2} \cdot \left(\frac{\sin(3\Psi + \frac{3\mu}{2})}{\sin \frac{3\mu}{2}} - \frac{\sin(\Psi + \frac{\mu}{2})}{\sin \frac{\mu}{2}} \right)$$

and

$$h_{eff} = -h/C$$

Example: sextupole (3D)

When we have h_{eff} in 3D we can write (see e.g. Chao):

$$Q_x(J_x, J_y, \delta) = \frac{1}{2\pi} \frac{\partial h_{eff}}{\partial J_x} = \frac{1}{2\pi} (\mu_x - 3k|\beta_x| D \delta)$$

$$Q_y(J_x, J_y, \delta) = \frac{1}{2\pi} \frac{\partial h_{eff}}{\partial J_y} = \frac{1}{2\pi} (\mu_y + 3k|\beta_y| D \delta)$$

and the change of path length:

$$\Delta s = - \frac{\partial h_{eff}}{\partial \delta} = \alpha_c \delta - 3k D^3 \delta^2 - 3k D (\beta_x J_x - \beta_y J_y)$$

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Example: sextupole (1D)

Given the Hamiltonian h :

$$h = -\mu J - \frac{3}{8} \mu k (2\beta J)^{3/2} \cdot \left(\frac{\sin(3\Psi + \frac{3\mu}{2})}{\sin \frac{3\mu}{2}} - \frac{\sin(\Psi + \frac{\mu}{2})}{\sin \frac{\mu}{2}} \right)$$

particles move in phase space along constant h .

Back to Cartesian coordinates we get for h :

$$h = -\frac{\mu}{2} (x^2 + x'^2) \frac{3}{8} \mu \beta^{3/2} x [(3x'^2 - x^2) \cot \frac{3\mu}{2} - (x^2 + x'^2) \cot \frac{\mu}{2} - 4xx']$$

Constant h defines the trajectory in phase space !

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What about close to resonance ?

If we have $Q = \frac{\mu}{2\pi} \approx \frac{m}{3}$ (3rd order resonance). Using a "distance to resonance d^n " as:

$$Q = \frac{m+d}{3} \quad \text{where: } d \ll 1$$

The trick is to observe the motion every 3 turns:

$$\mathcal{M}^3 = (e^{i - \mu J} e^{ikx^3})^3 = e^{i3h};$$

We get a factor:

$$e^{i - 3\mu J} = e^{i - 2\pi dJ} \quad (\text{because: } e^{i - 2\pi m J} \equiv 1)$$

$$d = \frac{3\mu}{2\pi}$$

What about close to resonance ?

Without proof (but like before, see e.g. Chao), we get:

$$h = -\frac{2\pi}{3}dJ - \frac{\pi}{12}dk(2J)^{3/2} \cdot \left(\frac{\sin(3\Psi + \frac{3\mu}{2})}{\sin \frac{3\mu}{2}} - \frac{\sin(\Psi + \frac{\mu}{2})}{\sin \frac{\mu}{2}} \right)$$

For small d ($\sin \frac{3\mu}{2} \approx -\pi d$) we can simplify:

$$h \approx -\frac{2\pi}{3}dJ - \frac{1}{\sqrt{2}}k(\beta J)^{3/2} \sin(3\Psi)$$

Analysis give fixed points, i.e. (back in Cartesian again):

$$\frac{\partial h}{\partial x} = -\frac{2\pi}{3}dx - \frac{1}{4}\beta^{3/2}(3x'^2 - 3x^2) = 0$$

$$\frac{\partial h}{\partial x'} = -\frac{2\pi}{3}dx' - \frac{1}{4}\beta^{3/2}3xx' = 0$$

Normal forms - non-linear case

Assume a linear rotation (as always) followed by an octupole, the Hamiltonian is:

$$\mathcal{H} = \frac{\mu}{2}(x^2 + p^2) + \frac{x^4}{4} \quad (p = p_x)$$

With the Hamilton's equation leading to:

$$\dot{x} = \frac{\partial \mathcal{H}}{\partial p} = \mu p$$

$$\dot{p} = -\frac{\partial \mathcal{H}}{\partial x} = -\mu x - x^3$$

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Normal forms - non-linear case

The map, written in Lie representation is:

$$\mathcal{M} = e^{-\frac{\mu}{2}x^2 + p^2} e^{\frac{x^4}{4}} = Re^{\frac{x^4}{4}}$$

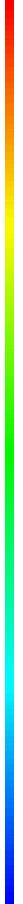
we transform by applying:

$$\begin{aligned} \mathcal{N} &= \mathcal{A} \mathcal{M} \mathcal{A}^{-1} = e^{xF} Re^{\frac{x^4}{4}} e^{-xF} = RR^{-1} e^{xF} Re^{\frac{x^4}{4}} e^{-xF} \\ &= Re^{xR^{-1}F + \frac{x^4}{4} - F + O(\epsilon^2)} = Re^{(R^{-1}-1)F + \frac{x^4}{4} + C(\epsilon^2)} \end{aligned}$$

we have now to choose F to simplify the expression:

$$= (R^{-1} - 1)F + \frac{x^4}{4}$$

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Normal forms - non-linear case

We go back to x and p coordinates and with:

$$J = (x^2 + p^2)/2$$

we can write the map:

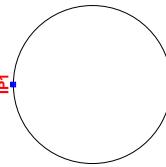
$$M = e^{-:F:} \quad e^{-\mu J + \frac{3}{8}J^2} \quad e^{:F:}$$

the term $\frac{3}{8}J^2$ is the tune shift with amplitude we know for an octupole

we have for F in x and p coordinates:

$$F = -\frac{1}{64}\{-5x^4 + 3p^4 + 6x^2p^2 + x^3p(8\cot(\mu) + 4\cot(2\mu)) + xp^3(8\cot(\mu) - 4\cot(2\mu))\}$$

A real life example: beam-beam interaction



- Localized distortion, very strong non-linearity
- Standard perturbation theory not appropriate

Effect on invariants - start with single IP

- Look for invariant \hbar
- Linear transfer e^{f_2} : and beam-beam interaction e^{F} ; i.e.:

$$e^{f_2} \cdot e^{F} = e^{\hbar}$$

with (see before):

$$f_2 = -\frac{\mu}{2} \left(\frac{x^2}{\beta} + \beta p_x^2 \right)$$

and (see before):

$$F = \int_0^x dx' f(x')$$

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Effect on invariants

using for a Gaussian beam $f(x)$ (see lecture on beam-beam effects):

$$f(x) = \frac{2}{x} \left(1 - e^{-\frac{x^2}{2\sigma^2}} \right)$$

as usual go to action angle variables Ψ, J :

$$x = \sqrt{2J\beta} \sin\Psi, \quad p = \sqrt{\frac{2J}{\beta}} \cos\Psi$$

and write $F(x)$ as Fourier series:

$$F(x) = \sum_{n=-\infty}^{\infty} c_n(J) e^{inx}$$

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We need:

REMEMBER: with this transform:

$$f_2 = -\mu J$$

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and useful properties of Lie operators (any textbook²⁾):

$$:f_2:g(J) = 0, \quad :f_2:e^{in\Psi} = in\mu e^{in\Psi}, \quad g(:f_2:)e^{in\Psi} = g(in\mu)e^{in\Psi}$$

and the formula (any textbook²⁾):

$$e^{:f_2:} e^{:F:} = e^{:h:} = \exp \left[:f_2 + \left(\frac{:f_2:}{1 - e^{-:f_2:}} \right) F + \mathcal{O}(F^2) \right]$$

²⁾ E. Forest, "Beam Dynamics, A New Attitude and Framework", 1998

gives immediately for h :

$$h = -\mu J + \sum_n c_n(J) \frac{in\mu}{1 - e^{-in\mu}} e^{in\Psi}$$

$$h = -\mu J + \sum_n c_n(J) \frac{n\mu}{2\sin(\frac{n\mu}{2})} e^{(in\Psi+i\frac{n\mu}{2})}$$

away from resonance normal form transformation gives:

$$h_n = -\mu J + c_0(J) = const.$$

$$\left[homework : \frac{dc_0(J)}{dJ} \right]$$

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Single IP - analysis of h

$$h = -\mu J + \sum_n c_n(J) \frac{n\mu}{2\sin(\frac{n\mu}{2})} e^{(in\Psi + i\frac{n\mu}{2})}$$

On resonance:

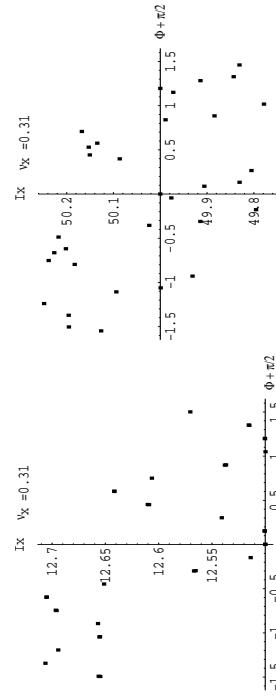
$$Q = \frac{p}{n} = \frac{\mu}{2\pi}$$

with $c_n \neq 0$:

$$\sin\left(\frac{n\pi p}{n}\right) = \sin(p\pi) \equiv 0 \quad \forall \text{ integer } p$$

and h diverges, find automatically all resonance conditions

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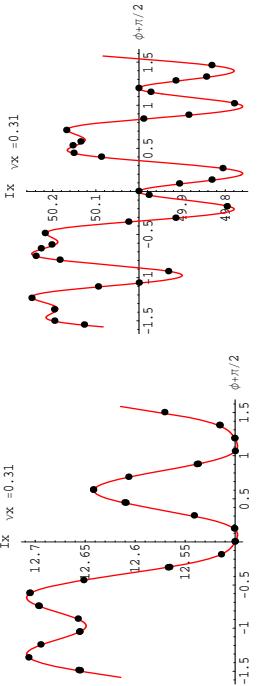


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Invariant from tracking: one IP

→ Shown for $5\sigma_x$ and $10\sigma_x$

Invariant versus tracking: one IP



► Shown for $5\sigma_x$ and $10\sigma_x$

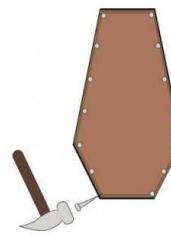
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Differential Algebra (DA)

a.k.a.

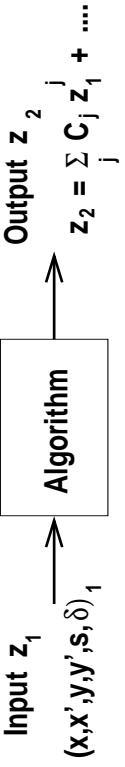
Truncated Power Series Algebra (TPSA)

- Some people like analytic methods
- Not just tracking numbers around the machine !
- Now we put the final nail into the coffin of any other approach ...



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Truncated Power Series Algebra



- The tracking of a complicated system relates the output **numerically** to the input
- Could we imagine something that relates the output **algebraically** to the input ?
- For example a Taylor series ?

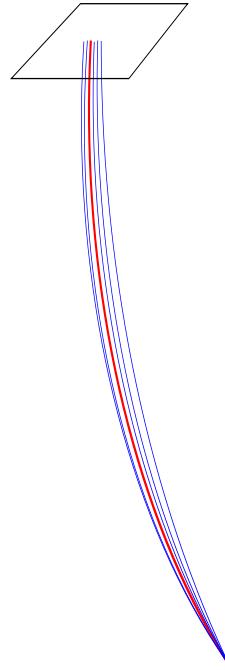
$$z_2 = \sum C_j z_1^j = \sum f^{(n)} z_1^j$$

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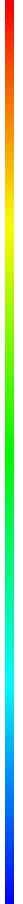
Why are Taylor series useful ?

Let us study the paraxial behaviour:



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- Red line is the ideal orbit
- Blue lines are small deviations
- If we understand how small deviations behave, we understand the system much better



Why are Taylor series useful ?

Now remember the definition of the Taylor series:

$$f(x + \Delta x) = f(x) + \sum_{n=1}^{\infty} \frac{f^{(n)}(x)}{n!} \Delta x^n$$

$$f(x + \Delta x) = f(x) + \frac{f'(x)}{1!} \Delta x^1 + \frac{f''(x)}{2!} \Delta x^2 + \frac{f'''(x)}{3!} \Delta x^3 + \dots$$

- The coefficients determine the behaviour of small deviations Δx from the ideal orbit x
- The Taylor expansion does a paraxial analysis of the system
- How to get the coefficients without extra work ?

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Numerical differentiation

The problem getting the derivatives $f^{(n)}(a)$ of $f(x)$ at a :

$$f'(a) = \frac{f(a + \epsilon) - f(a)}{\epsilon}$$

- Need to subtract almost equal numbers and divide by small number.
- For higher orders f'', f''', \dots , accuracy hopeless !
- We can use Differential Algebra (DA) (Berz, 1988)

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Differential Algebra

1. Define a pair (q_0, q_1) , with q_0, q_1 real numbers

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Differential Algebra

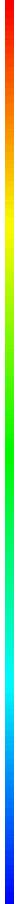
1. Define a pair (q_0, q_1) , with q_0, q_1 real numbers
2. Define operations on a pair like:

$$(q_0, q_1) + (r_0, r_1) = (q_0 + r_0, q_1 + r_1)$$

$$c \cdot (q_0, q_1) = (c \cdot q_0, c \cdot q_1)$$

$$(q_0, q_1) \cdot (r_0, r_1) = (q_0 \cdot r_0, q_0 \cdot r_1 + q_1 \cdot r_0)$$

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Differential Algebra

1. Define a pair (q_0, q_1) , with q_0, q_1 real numbers
2. Define operations on a pair like:

$$(q_0, q_1) + (r_0, r_1) = (q_0 + r_0, q_1 + r_1)$$

$$c \cdot (q_0, q_1) = (c \cdot q_0, c \cdot q_1)$$

$$(q_0, q_1) \cdot (r_0, r_1) = (q_0 \cdot r_0, q_0 \cdot r_1 + q_1 \cdot r_0)$$

3. And some ordering:

$$(q_0, q_1) < (r_0, r_1) \quad \text{if} \quad q_0 < r_0 \quad \text{or} \quad (q_0 = r_0 \quad \text{and} \quad q_1 < r_1)$$

$$(q_0, q_1) > (r_0, r_1) \quad \text{if} \quad q_0 > r_0 \quad \text{or} \quad (q_0 = r_0 \quad \text{and} \quad q_1 > r_1)$$

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Differential Algebra

1. Define a pair (q_0, q_1) , with q_0, q_1 real numbers
2. Define operations on a pair like:

$$(q_0, q_1) + (r_0, r_1) = (q_0 + r_0, q_1 + r_1)$$

$$c \cdot (q_0, q_1) = (c \cdot q_0, c \cdot q_1)$$

$$(q_0, q_1) \cdot (r_0, r_1) = (q_0 \cdot r_0, q_0 \cdot r_1 + q_1 \cdot r_0)$$

3. And some ordering:

$$(q_0, q_1) < (r_0, r_1) \quad \text{if} \quad q_0 < r_0 \quad \text{or} \quad (q_0 = r_0 \quad \text{and} \quad q_1 < r_1)$$

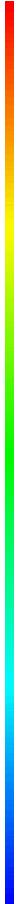
$$(q_0, q_1) > (r_0, r_1) \quad \text{if} \quad q_0 > r_0 \quad \text{or} \quad (q_0 = r_0 \quad \text{and} \quad q_1 > r_1)$$

4. This implies something strange:

$$(0, 0) < (0, 1) < (r, 0) \quad (\text{for any pos. } r)$$

$$(0, 1) \cdot (0, 1) = (0, 0) \xrightarrow{\text{?}} (0, 1) = \sqrt{(0, 0)!!}$$

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Differential Algebra

This means that $(0,1)$ is between 0 and ANY real number
→ infinitely small !!

We call this therefore "differential unit" $d = (0,1) = \delta$.

Of course $(q,0)$ is just the real number q and we define "real part" and "differential part" (a bit like complex numbers..):

$$q_0 = \mathcal{R}(q_0, q_1) \quad \text{and} \quad q_1 = \mathcal{D}(q_0, q_1)$$

With our rules we can further see that:

$$(1,0) \cdot (q_0, q_1) = (q_0, q_1)$$

$$(q_0, q_1)^{-1} = \left(\frac{1}{q_0}, -\frac{q_1}{q_0^2} \right)$$

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Differential Algebra

Of course can let a function f act on the pair (or vector) using our rules.

For example:

$$f(x) \rightarrow f(x, 0)$$

acts like the function f on the real variable x :

$$f(x) = \mathcal{R}[f(x, 0)]$$

What about the differential part \mathcal{D} ?

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Differential Algebra

For a function $f(x)$ without proof:

$$\mathcal{D}[f(x+d)] = \mathcal{D}[f((x, 0) + (0, 1))] = \mathcal{D}[f(x, 1)] = f'(x)$$

An example instead:

$$f(x) = x^2 + \frac{1}{x}$$

then using school calculus:

$$f'(x) = 2x - \frac{1}{x^2}$$

For $x = 2$ we get then:

$$f(2) = \frac{9}{2}, f'(2) = \frac{15}{4}$$

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Differential Algebra

For \mathbf{x} in:

$$f(x) = x^2 + \frac{1}{x}$$

we substitute: $\mathbf{x} \rightarrow (\mathbf{x}, 1) = (2, 1)$ and use our rules:

$$\begin{aligned} f[(2, 1)] &= (2, 1)^2 + (2, 1)^{-1} \\ &= (4, 4) + \left(\frac{1}{2}, -\frac{1}{4}\right) \\ &= \left(\frac{9}{2}, \frac{15}{4}\right) = (f(2), f'(2)) \quad !!! \end{aligned}$$

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The computation of derivatives becomes an algebraic problem, no need for small numbers, exact !

Differential Algebra - higher orders

1. The pair $(q_0, 1)$, becomes a vector of order N :
$$(q_0, 1) \xrightarrow{\quad} (q_0, 1, \mathbf{0}, \mathbf{0}, \dots, \mathbf{0}) \quad \delta = (\mathbf{0}, 1, \mathbf{0}, \mathbf{0}, \mathbf{0}, \dots)$$
2. $(q_0, q_1, q_2, \dots, q_N) + (r_0, r_1, r_2, \dots, r_N) = (s_0, s_1, s_2, \dots, s_N)$
with: $s_i = q_i + r_i$

$$3. c \cdot (q_0, q_1, q_2, \dots, q_N) = (c \cdot q_0, c \cdot q_1, c \cdot q_2, \dots, c \cdot q_N)$$

$$4. (q_0, q_1, q_2, \dots, q_N) \cdot (r_0, r_1, r_2, \dots, r_N) = (s_0, s_1, s_2, \dots, s_N)$$

with:

$$s_i = \sum_{k=0}^i \frac{i!}{k!(i-k)!} q_k r_{i-k}$$

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Differential Algebra

If we had started with:

$$x = (a, 1, 0, 0, 0, \dots)$$

we would get:

$$f(x) = (f(a), f'(a), f''(a), f'''(a), \dots, f^{(n)}(a))$$

can be extended to more variables x, y :

$$x = (a, 1, 0, 0, 0, \dots) \quad dx = (0, 1, 0, 0, 0, \dots)$$

$$y = (b, 0, 1, 0, 0, \dots) \quad dy = (0, 0, 1, 0, 0, \dots)$$

and get (with more complicated multiplication rules):

$$f((x + dx), y + dy)) = \left(f, \frac{\partial f}{\partial x}, \frac{\partial f}{\partial y}, \frac{\partial^2 f}{\partial x^2}, \frac{\partial^2 f}{\partial x \partial y}, \dots \right) (x, y)$$

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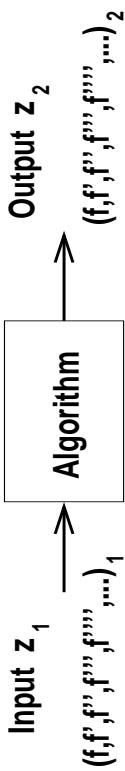
What is the use of that:



Can extract a truncated Taylor map of a beam line or ring by pushing the identity map $f(x) = (a, 1, 0, 0, \dots)$ through the algorithm as if it is a vector in phase space !
The maps are provided with the desired accuracy and to any order.

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What is the use of that:



- "Algorithm" can be a mathematical function
- "Algorithm" can be a complex computer code
- Easy using polymorphism of modern languages (see example)
- Normal form analysis on Taylor series is much easier !!
- We get a Taylor map for a computer code !!!

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What is the use of that:

- Demonstrate with simple examples (FORTRAN 95):
 - First show the concept
 - Simple FODO cell
 - Normal form analysis of the FODO cell with sextupoles
- All examples and all source code in:
[/afs/cern.ch/user/z/zwe/public/DA](http://afs/cern.ch/user/z/zwe/public/DA)

Small DA package provided by E. Forest

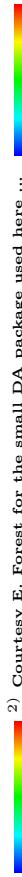


Look at this small example:

```
PROGRAM DATEST1  
use my_own_da  
real(8) x,z, dx  
my_order=3  
dx=0.0  
x=3.141592653_8/6.0_8+dx  
call track(x, z)  
call print(z,6)  
END PROGRAM DATEST1  
  
SUBROUTINE TRACK(a, b)  
use my_own_da  
real(8) a,b  
b = sin(a)  
END SUBROUTINE TRACK
```

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2) Courtesy E. Forest for the small DA package used here ..



Look at the results:

```
(0,0) 0.500000000000E+00
(1,0) 0.8660254037844E+00
(0,1) 0.000000000000E+00
(2,0) -0.250000000000E+00
(0,2) 0.000000000000E+00
(1,1) 0.000000000000E+00
(3,0) -0.1443375672974E+00
(0,3) 0.000000000000E+00
(2,1) 0.000000000000E+00
(1,2) 0.000000000000E+00
(0,0) 0.500000000000E+00
```

We have $\sin(\frac{\pi}{6}) = 0.5$ all right, but what is the rest ??



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Look at the results:

```
(0,0) 0.500000000000E+00
(1,0) 0.8660254037844E+00
(0,1) 0.000000000000E+00
(2,0) -0.250000000000E+00
(0,2) 0.000000000000E+00
(1,1) 0.000000000000E+00
(3,0) -0.1443375672974E+00
(0,3) 0.000000000000E+00
(2,1) 0.000000000000E+00
(1,2) 0.000000000000E+00
(0,0) 0.500000000000E+00
```

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$$\sin\left(\frac{\pi}{6} + \Delta x\right) = \sin\left(\frac{\pi}{6}\right) + \cos\left(\frac{\pi}{6}\right)\Delta x^1 - \frac{1}{2}\sin\left(\frac{\pi}{6}\right)\Delta x^2 - \frac{1}{6}\cos\left(\frac{\pi}{6}\right)\Delta x^3$$



What is the use of that:

- We have used a simple algorithm here (*sin*) but it can be **anything** very complex
- We can compute nonlinear maps as a Taylor expansion of **anything** the program computes
- Simply by:
 - Replacing regular (e.g. REAL) types by TPSA types (*my_taylor*) i.e. variables x, p are automatically replaced by $(x, 1, 0, \dots)$ and $(p, 0, 1, 0, \dots)$ etc.
 - Operators and functions $(+, -, *, =, \dots, exp, sin, \dots)$ automatically overloaded, i.e. behave according to new type

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What is the use of that:

Assume the *Algorithm* describes one turn, then:

- Normal tracking:
 - $X_n = (x, p_x, y, p_y, s, \delta)_n \rightarrow X_{n+1} = (x, p_x, y, p_y, s, \delta)_{n+1}$
 - Coordinates after one completed turn
- TPSA tracking:
 - $X_n = (x, p_x, y, p_y, s, \delta)_n \rightarrow X_{n+1} = \sum C_j X_n^j$
 - Taylor expansion after one completed turn
 - Automatically all X_{n+1} where it converges
 - The C_j contain useful information about behaviour
 - Taylor map directly used for normal form analysis

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Another example:

➤ Track through a FODO lattice:

QF - DRIFT - QD

Integrate 100 steps in the quadrupoles

Now we use **two** variables: $\mathbf{x}, \mathbf{p} = (\mathbf{z}(1), \mathbf{z}(2))$

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2) Courtesy E. Forest for the small DA package used here ...

Another example:

```
program fodo1
use my_own_da
use my_analysis
use my_taylor z(2)
type(normalform) NORMAL
type(my_map) M,id
real(dp) L,DL,k1,f1x(2),ma(2,2)
integer i,nstep
nstep=100
f1x=0.0_dp ! fixed point
id=1
z=f1x+id
do i=1,nstep
  z(1)=z(1)+DL/2.d0*z(2)
  z(2)=z(2)-k1*d1*z(1)
  z(1)=z(1)+DL/2.d0*z(2)
enddo
z(1)=z(1)+L*z(2)
do i=1,nstep
  z(1)=z(1)+DL/2.d0*z(2)
  z(2)=z(2)+k1*d1*z(1)
  z(1)=z(1)+DL/2.d0*z(2)
enddo
call print(z(1),6)
call print(z(2),6)
M==z
NORMAL==M
write(6,* ) normal%tune, normal%dtune,da
end program fodo1
```

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Result:

0.166310856124535, (0.000000000, 0.000000000)

2) Courtesy E. Forest for the small DA package used here ...

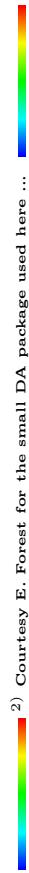
The result is:

Only linear elements in the Taylor expansion, the result is the matrix:

```
(0,0) 0.00000000000000
(1,0) -1.45367086074474
(0,1) 2.76712638076176
(2,0) 0.00000000000000
(0,2) 0.00000000000000
(1,1) 0.00000000000000
(3,0) 0.00000000000000
(0,3) 0.00000000000000
(2,1) 0.00000000000000
(1,2) 0.00000000000000
(0,0) 0.00000000000000
(1,0) -1.663241972385291
(0,1) 2.45754057212448
(2,0) 0.00000000000000
(0,2) 0.00000000000000
(1,1) 0.00000000000000
(3,0) 0.00000000000000
(0,3) 0.00000000000000
(2,1) 0.00000000000000
(1,2) 0.00000000000000
```

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2) Courtesy E. Forest for the small DA package used here ...



The output from the normal form analysis is the tune:
0.1663108 !

Modified previous example (with sextupole):

```
program fodo3
use my_own_da
use my_analysis
type(my_taylor) z(2)
type(normalform) NORMAL
type(my_map) M,id
real(dp) L,DL,k1,f(x(2)),ma(2,2)
integer i,nstep
my_order=3
nstep=100
fx=0.0,dp ! fixed point
id=1
z=f(x+id)

do i=1,nstep
  z(1)=z(1)+DL/2,d0*z(2)
  z(2)=z(2)-k1*d1*z(1)
  z(1)=z(1)+DL/2,d0*z(2)
enddo

z(2)=z(2)-z(1)**2
z(1)=z(1)+L*z(2)

do i=1,nstep
  z(1)=z(1)+DL/2,d0*z(2)
  z(2)=z(2)+k1*d1*z(1)
  z(1)=z(1)+DL/2,d0*z(2)
enddo

z(2)=z(2)-z(1)**2

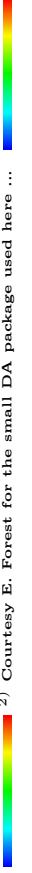
call print(z(1),6)
call print(z(2),6)
M=z
NORMAL=M
write(6,* ) normal%tune, normal%dtune_da
end program fodo3
```

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Result:

0.166310856124535, (0.000000000, -12.5243894659508)

2) Courtesy E. Forest for the small DA package used here ...



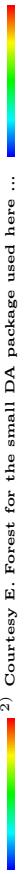
The result is:

Now non-linear elements in the Taylor expansion,

```
(0,0) 0.000000000000000
(1,0) -1.45367086074474
(0,1) 2.76712693076176
(2,0) -0.793529405369330
(0,2) -1.92470766801451
(1,1) -2.47168940711356
(3,0) 0.000000000000000
(0,3) 0.000000000000000
(2,1) 0.000000000000000
(1,2) 0.000000000000000
(0,0) 0.000000000000000
(1,0) -1.65241972385261
(0,1) 2.45754057212448
(2,0) -2.90667980020143
(0,2) -9.58167531652505
(1,1) 5.5732088161683
(3,0) -2.30706114745899
(0,3) 10.6518208440132
(2,1) -2.79446255996715
(1,2) 8.08317374091570
```

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2) Courtesy E. Forest for the small DA package used here ...



Modified previous example (with sextupole):

Remember the normal form transformation:

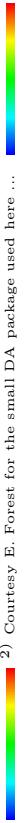
$$\mathcal{A}\mathcal{M}\mathcal{A}^{-1} = \mathcal{R}$$

The type **normalform** in the demonstration package also contains the maps \mathcal{A} and \mathcal{R} !

Question: what would be the result of something like that ?

```
r2=(x**2+p**2)**normal%at**(-1)
print sqrt(r2.sub(beta))
print -0.5*(r2.sub.twoalpha),r2.sub.gamma
```

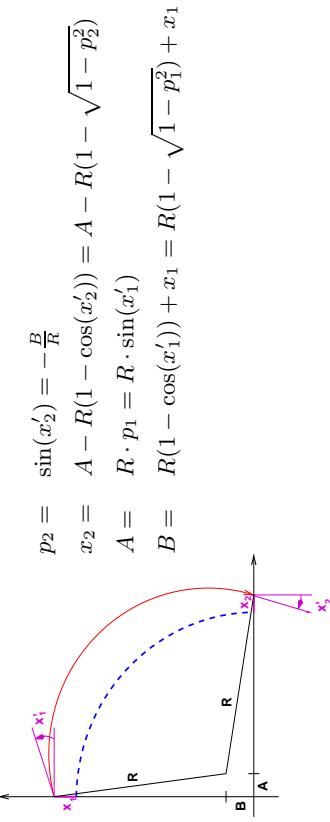
2) Courtesy E. Forest for the small DA package used here ...



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This was trivial - now a (normally) hard one

The exact map:



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A 90° bending magnet ..



How to apply Differential Algebra here ...

► Start with initial coordinates in DA style:

$$\begin{aligned} x_1 &= (0, 1, 0, \dots) \\ p_1 &= (0, 0, 1, \dots) \quad \text{and have:} \\ A &= (0, 0, R, 0, \dots) \\ B &= (0, 1, 0, 0, 0, R, 0, \dots) \end{aligned}$$

► After pushing them through the algorithm:

$$\begin{aligned} \uparrow x_2 &= (0, 0, R, -\frac{1}{R}, 0, 0, 0, \dots) = (0, \frac{\partial x_2}{\partial x_1}, \frac{\partial x_2}{\partial p_1}, \frac{\partial^2 x_2}{\partial x_1^2}, \frac{\partial^2 x_2}{\partial x_1 \partial p_1}, \dots) \\ \uparrow p_2 &= (0, -\frac{1}{R}, 0, 0, 0, -1, 0, \dots) = (0, \frac{\partial p_2}{\partial x_1}, \frac{\partial p_2}{\partial p_1}, \frac{\partial^2 p_2}{\partial x_1^2}, \frac{\partial^2 p_2}{\partial x_1 \partial p_1}, \dots) \end{aligned}$$

► Automatically evaluates all non-linearities to any desired order ..

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How to apply Differential Algebra here ...

➤ Start with initial coordinates in DA style:

$$x_1 = (0, 1, 0, \dots)$$

$$p_1 = (0, 0, 1, \dots) \quad \text{and have:}$$

$$A = (0, 0, R, 0, \dots)$$

$$B = (0, 1, 0, 0, 0, R, 0, \dots)$$

➤ After pushing them through the algorithm:

$$\rightarrow x_2 = (0, \mathbf{0}, R, -\frac{1}{R}, 0, 0, 0, \dots) = (0, \frac{\partial x_2}{\partial x_1}, \frac{\partial x_2}{\partial p_1}, \frac{\partial^2 x_2}{\partial x_1^2}, \frac{\partial^2 x_2}{\partial x_1 \partial p_1}, \dots)$$

$$\rightarrow p_2 = (0, -\frac{1}{R}, \mathbf{0}, 0, 0, -1, 0, \dots) = (0, \frac{\partial p_2}{\partial x_1}, \frac{\partial p_2}{\partial p_1}, \frac{\partial^2 p_2}{\partial x_1^2}, \frac{\partial^2 p_2}{\partial x_1 \partial p_1}, \dots)$$

➤ Automatically evaluates all non-linearities to any desired order ..

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Some we know ...

Transfer matrix of a dipole:

$$M_{dipole} = \begin{pmatrix} \cos(\frac{L}{R}) & R \sin(\frac{L}{R}) \\ -\frac{1}{R} \sin(\frac{L}{R}) & \cos(\frac{L}{R}) \end{pmatrix} = \begin{pmatrix} \frac{\partial x_2}{\partial x_1} & \frac{\partial x_2}{\partial p_1} \\ \frac{\partial p_2}{\partial x_1} & \frac{\partial p_2}{\partial p_1} \end{pmatrix}$$

For a 90° bending angle we get:

$$M_{dipole} = \begin{pmatrix} 0 & R \\ -\frac{1}{R} & 0 \end{pmatrix}$$

as computed, but we also have **all** derivatives and non-linear effects !

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What is the use of that:

- Although not strictly an analytic method in the traditional sense:
 - TPSA provide analytic expression (Taylor series) for the one turn map
 - Can be used for tracking
 - Can be analysed for dynamic behaviour of the system
 - Typical use: Normal Form Analysis discussed earlier, rather straightforward from a Taylor expansion

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Is there a summary ?

$m = z$

$\text{NORMAL} = m$

- Get the map m somehow (no matter how)
- Analyse this map (Normal form)

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And another summary

- Perturbation treatment limited to:
 - Small perturbations (not in real machines)
 - Pedagogical purpose
- For realistic machines symplectic, iterative mapping **is** appropriate, using:
 - Symplectic integration
 - Lie transformations and normal form analysis
 - Differential algebra

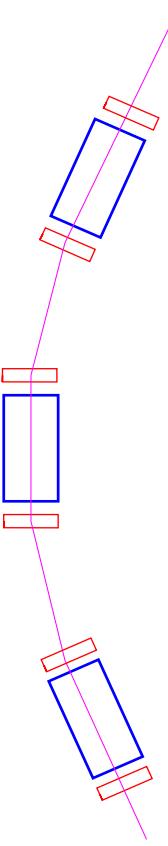
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[Back up](#)

Slide 184

- backup slides -

Where to put the elements in an accelerator ?



$$\frac{d^2x}{ds^2} + K(s) x = 0$$

- Usually use s (pathlength) along "reference path"
- "Reference path" defined geometrically by straight sections and bending magnets

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Second order MAPS concatenation

Assume now 2 maps of second order:

$$\mathcal{A}_2 = [R^A, T^A] \quad \text{and} \quad \mathcal{B}_2 = [R^B, T^B]$$

the combined second order map

$$\mathcal{C}_2 = \mathcal{A}_2 \circ \mathcal{B}_2 \quad \text{is} \quad \mathcal{C}_2 = [R^C, T^C] \quad \text{with:}$$

$$R^C = R^A \cdot R^B$$

and (after truncation of higher order terms !!):

$$T_{ijk}^C = \sum_{l=1}^4 R_{il}^B T_{ljk}^A + \sum_{l=1}^4 \sum_{m=1}^4 T_{ilm}^B R_{ij}^A R_{mk}^A$$

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Symplecticity for higher order MAPS

try truncated Taylor map in 2D, second order:

$$\begin{pmatrix} x \\ x' \end{pmatrix} = \begin{pmatrix} R_{11}x_0 + R_{12}x'_0 + T_{111}x_0^2 + T_{112}x_0x'_0 + T_{122}x_0'^2 \\ R_{21}x_0 + R_{22}x'_0 + T_{211}x_0^2 + T_{212}x_0x'_0 + T_{222}x_0'^2 \end{pmatrix}$$

The Jacobian becomes:

$$\mathcal{J} = \begin{bmatrix} R_{11} + 2T_{111}x_0 + T_{112}x'_0 & R_{12} + T_{112}x_0 + 2T_{122}x'_0 \\ R_{21} + 2T_{211}x_0 + T_{212}x'_0 & R_{22} + T_{212}x_0 + 2T_{222}x'_0 \end{bmatrix}$$

symplecticity condition requires that:

$\det \mathcal{J} = 1$ for all x_0 and all x'_0

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Symplecticity for higher order MAPS

This is only possible for the conditions:

$$\begin{cases} R_{11}R_{22} - R_{12}R_{21} = 1 \\ R_{11}T_{212} + 2R_{22}T_{111} - 2R_{12}T_{211} - R_{21}T_{112} = 0 \\ 2R_{11}T_{222} + R_{22}T_{112} - R_{12}T_{212} - 2R_{21}T_{122} = 0 \end{cases}$$

- 10 coefficients, but 3 conditions
- number of **independent** coefficients only 7 !
- Taylor map requires more coefficients than necessary
- e.g. 4D, order 4: coefficients **276** instead of 121

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Canonical transformations

- With Hamiltonian's equations, still have to solve (2n) differential equations
- Not necessarily easy, but:
 - More freedom to choose the variables q and p (because they have now "equal" status)
 - Try to find variables where they are easy to solve
- Change of variables through "canonical transformations"

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Why canonical transformations ?

- Hamiltonian have one advantage over Lagrangians:
- If the system has a symmetry, i.e. a coordinate q_i does not occur in H (i.e. $\frac{\partial H}{\partial q_i} = 0 \rightarrow \frac{dp_i}{dt} = 0$)  the corresponding momentum p_i is conserved (and the coordinate q_i can be ignored in the other equations of the set).
- Comes also from Lagrangian, but the velocities still occur in \mathcal{L} !

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Canonical transformations

Starting with $H(q, p, t)$ get new coordinates:

$$Q_i = Q_i(q, p, t)$$

$$P_i = P_i(q, p, t)$$

and new Hamiltonian $K(Q, P, t)$ with:

$$\frac{\partial K}{\partial Q_j} = -\dot{P}_j = -\frac{dP_j}{dt}, \quad \frac{\partial K}{\partial P_j} = \dot{Q}_j = \frac{dQ_j}{dt}$$

■ We can two types of canonical transformations

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Canonical transformations - type 1

➤ Ideally one would like a Hamiltonian H and coordinates with:

$$\frac{\partial H}{\partial q_j} = -\dot{p}_j = -\frac{dp_j}{dt} = 0$$

➤ Coordinate q_j not explicit in H

➤ p_j is a constant of the motion (!) and:

$$\frac{dq_j}{dt} = \frac{\partial H(p_1, p_2, .., p_n)}{\partial p_j} = F_j(p_1, p_2, .., p_n)$$

which can be directly integrated to get $q_j(t)$

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Canonical transformations - type 1, example

Harmonic oscillator:

$$H = T + V = \frac{1}{2}mv^2 + \frac{m\omega^2}{2}x^2 = \frac{p^2}{2m} + \frac{m\omega^2}{2}x^2$$

try: $x = \sqrt{\frac{2P}{m\omega}} \cdot \sin(X)$ and $p = \sqrt{2m\omega P} \cdot \cos(X)$ and we get:

$$K = \omega P \cos^2(X) + \omega P \sin^2(X) = \omega P$$

then:

$$\frac{dX}{dt} = \frac{\partial K}{\partial P} = \omega \quad \Rightarrow \quad X = \omega t + \alpha$$

back transformation to x,p:

$$x = \sqrt{\frac{2E}{m\omega^2}} \sin(\omega t + \alpha)$$

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Canonical transformations - type 2

- Find a transformation of q, p at time t to values q_0, p_0 at time $t = 0$.

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$$q = q(q_0, p_0, t)$$

$$p = p(q_0, p_0, t)$$

- The transformations ARE the solution of the problem !

For both types: how to find the transformation ?

- Without details: Hamilton-Jacobi equation ...

Extension: general monomials

Monomials in x and p of orders n and m ($x^n p^m$)

$$e^{ax^n p^m};$$

gives for the map (for $n \neq m$):

$$e^{ax^n p^m}; x = x \cdot [1 + a(n-m)x^{n-1}p^{m-1}]^{m/(m-n)}$$

$$e^{ax^n p^m}; p = p \cdot [1 + a(n-m)x^{n-1}p^{m-1}]^{n/(n-m)}$$

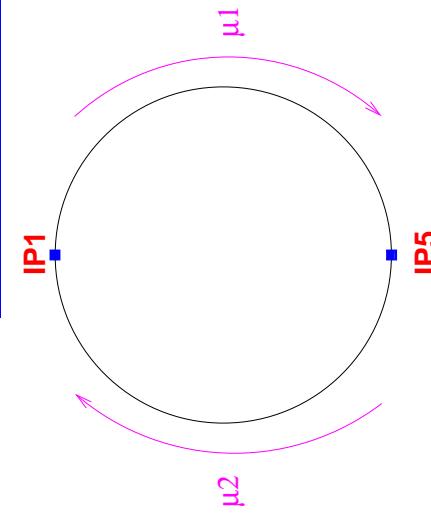
gives for the map (for $n = m$):

$$e^{ax^n p^n}; x = x \cdot e^{-anx^{n-1}p^{n-1}}$$

$$e^{ax^n p^n}; p = p \cdot e^{anx^{n-1}p^{n-1}}$$

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Collision scheme - two IPs



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Two IPs

→ two transfers f_2^1, f_2^2 and two beam-beam kicks F^1, F^2 ,
first IP at μ_1 , second IP at μ :

$$\begin{aligned} &= e^{i:f_2^1:} e^{i:F^1:} e^{i:f_2^2:} e^{i:F^2:} = e^{i:h_2:} \\ &= e^{i:f_2^1:} e^{i:F^1:} e^{-i:f_2^1:} e^{i:f_2^1:} e^{i:f_2^2:} e^{i:F^2:} = e^{i:h_2:} \\ &= e^{i:f_2^1:} e^{i:F^1:} e^{-i:f_2^1:} e^{i:f_2^1:} e^{i:F^2:} e^{-i:f_2:} e^{i:f_2:} = e^{i:h_2:} \\ &= e^{i:e^{-i:f_2^1:} F^1:} e^{i:e^{-i:f_2:} F^2:} e^{i:f_2:} = e^{i:h_2:} \end{aligned}$$

$$f_2 = -\mu A, \quad f_2^1 = -\mu_1 A, \quad \text{and} \quad f_2^2 = -\mu_2 A$$

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Two IPs

here a miracle occurs (remember $g(:f_2:)e^{in\Psi} = g(in\mu)e^{in\Psi}$):

$$e^{i:f_2^1:} e^{in\Psi} = e^{in\mu_1} e^{in\Psi} = e^{in(\mu_1 + \Psi)}$$

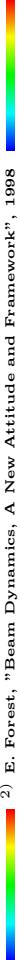
i.e. the Lie transforms of the perturbations are phase shifted²⁾. Therefore:

$$e^{i:e^{-i:f_2^1:} F^1:} e^{i:e^{-i:f_2:} F^2:} e^{i:f_2:} = e^{i:h_2:}$$

becomes simpler with substitutions of $\Psi_{\textcolor{red}{1}} = \Psi + \mu_1$ and $\Psi = \Psi + \mu$ in F^1 and F :

$$e^{i:F^1(\Psi_{\textcolor{red}{1}}):} e^{i:F(\Psi):} e^{i:f_2:} \Rightarrow e^{i:F^1(\Psi_{\textcolor{red}{1}}) + F(\Psi):} e^{i:f_2:}$$

²⁾ E. Forest, "Beam Dynamics, A New Attitude and Framework", 1998



Two IPs

gives for h_2 :

$$h_2 = -\mu A + \sum_{n=-\infty}^{\infty} \frac{n\mu c_n(A)}{2\sin(n\frac{\mu}{2})} e^{-in(\Psi+\mu/2+\mu_1)} + e^{-in(\Psi+\mu/2)}$$

$$h_2 = -\mu A + 2c_0(A) + \sum_{n=1}^{\infty} \frac{2n\mu c_n(A)}{2\sin(n\frac{\mu}{2})} \cos(n(\Psi + \frac{\mu}{2} + \frac{\mu_1}{2})) \cos(n\frac{\mu_1}{2})$$

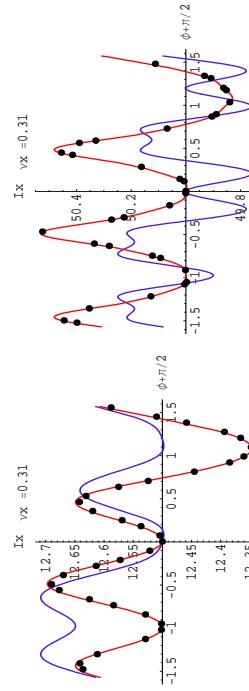
interesting part

Nota bene, because of:

$$e^{:F(\Psi):} e^{:f_2:} \xrightarrow{\quad} e^{:F^1(\Psi_1)+F(\Psi):} e^{:f_2:}$$

can be generalized to more interaction points ...

Invariant versus tracking: two IPs



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→ Shown for $5\sigma_x$ and $10\sigma_x$

Recap: Hamiltonian for a finite length element

We have from the Hamiltonian equations for the motion through an element with the Hamiltonian H for the element of length L :

$$\frac{dq}{dt} = [q, H] =: -H : q \quad (\text{from lecture 5})$$

$$\xrightarrow{\hspace{1cm}} \frac{d^k q}{dt^k} = (: -H :)^k q$$

$$\xrightarrow{\hspace{1cm}} q(t) = \sum_{k=0}^{\infty} \frac{t^k}{k!} \left(\frac{d^k q}{dt^k} \right) = \sum_{k=0}^{\infty} \frac{t^k}{k!} (- : H :)^k = e^{:-tH:}$$

with independent variable s instead of t (nota bene:

$$s_0 = 0, t_0 = 0)$$

$$\xrightarrow{\hspace{1cm}} q(s) = e^{:-LH:}$$

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Lie transformations on moments:

We have used Lie transformations mainly to propagate coordinates and momenta, i.e. like:

$$e^{\dot{f}}: x_0 = x_1$$

$$e^{\dot{f}}: p_0 = p_1$$

or using $Z = (x, p_x, y, p_y, \dots)$:

$$e^{\dot{f}}: Z_0 = Z_1$$

► Remember: can be applied to any function of x and p !!

► In particular to moments like x^2, xp, p^2, \dots

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Lie transformations on moments

Assume a matrix M of the type:

$$M = \begin{pmatrix} m_{11} & m_{12} \\ m_{21} & m_{22} \end{pmatrix}$$

described by a generator f , we have for the Lie transformation on the moment:

$$e^{if}x^2 = (e^{if}x)^2 \quad (\text{see lecture 5})$$

therefore:

$$\begin{aligned} (e^{if}x)^2 &= (m_{11}x + m_{12}p)^2 \\ (e^{if}x)^2 &= m_{11}^2x^2 + 2m_{11}m_{12}xp + m_{12}^2p^2 \end{aligned}$$

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More on moments

To summarize the moments:

$$\begin{pmatrix} x^2 \\ xp \\ p^2 \end{pmatrix}_{s_2} = \begin{pmatrix} m_{11}^2 & 2m_{11}m_{12} & m_{12}^2 \\ m_{11}m_{21} & m_{11}m_{22} + m_{12}m_{21} & m_{12}m_{22} \\ m_{21}^2 & 2m_{21}m_{22} & m_{22}^2 \end{pmatrix} \circ \begin{pmatrix} x^2 \\ xp \\ p^2 \end{pmatrix}_{s_1}$$

This is the well known transfer matrix for optical parameters

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A real life example: beam-beam interaction^{*)}

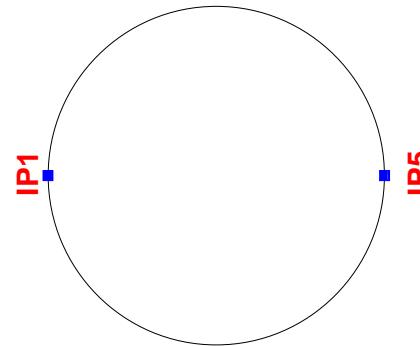
- Beam-beam interaction very non-linear
- Important to understand stability
- Non-linear effects such as amplitude detuning very important

Our questions ?

- How does the particles behave in phase space ?
 - Do we have an invariant ?
 - Can we calculate the invariant ?
- *) From: W. Herr, D. Kaltchev, LHC Project Report 1082, (2008).

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Collision scheme - two IPs



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Start with single IP

”Classic“ (B.C.) approach:

- Interaction point at beginning (end) of the ring (very local interactions, δ -functions)
- s-dependent Hamiltonian and perturbation theory:

$$\mathcal{H} = \dots + \delta(s)\epsilon V$$

■ Disadvantages:

- for several IPs endless mathematics
- conceptually and computationally easier method



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Effect on invariants - start with single IP

Look for invariants $\textcolor{red}{h}$, (see e.g. Dragt¹⁾), and evaluate for different number of interactions and phase advance.

Very well suited for local distortions (e.g. beam-beam kick)

Linear transfer e^{f_2} ; and beam-beam interaction $e^{\textcolor{red}{F}}$; i.e.:

$$e^{f_2} \cdot e^{\textcolor{red}{F}} = e^{\textcolor{red}{h}}$$

with

$$f_2 = -\frac{\mu}{2} \left(\frac{x^2}{\beta} + \beta p_x^2 \right)$$

and

$$F = \int_0^x dx' f(x')$$

¹⁾ A. Dragt, AIP Conference proceedings, Number 57 (1979)

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Effect on invariants

using for a Gaussian beam $f(x)$:

$$f(x) = \frac{2}{x} (1 - e^{\frac{-x^2}{2\sigma^2}})$$

as usual go to action angle variables Ψ, A :

$$x = \sqrt{2A}\beta \sin\Psi, \quad p = \sqrt{\frac{2A}{\beta}} \cos\Psi$$

and write $F(x)$ as Fourier series:

$$F(x) = \sum_{n=-\infty}^{\infty} c_n(A) e^{in\Psi} \quad \text{with :} \quad c_n(A) = \frac{1}{2\pi} \int_0^{2\pi} e^{-in\Psi} F(x) d\Psi$$

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We need:

REMEMBER: with this transform:

$$f_2 = -\mu A$$

and useful properties of Lie operators (any textbook²⁾):

$$:f_2:g(A) = 0, \quad :f_2:e^{in\Psi} = in\mu e^{in\Psi}, \quad g(:f_2:)e^{in\Psi} = g(in\mu)e^{in\Psi}$$

and the formula (because the beam-beam perturbation is small !):

$$e^{:f_2:F:} = e^{:h:} = \exp \left[:f_2 + \left(\frac{:f_2:}{1 - e^{-:f_2:}} \right) F + \mathcal{O}(F^2) \right]$$

²⁾ E. Forest, "Beam Dynamics, A New Attitude and Framework", 1998

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Single IP

gives immediately for h :

$$h = -\mu A + \sum_n c_n(A) \frac{in\mu}{1 - e^{-in\mu}} e^{in\Psi}$$

$$h = -\mu A + \sum_n c_n(A) \frac{n\mu}{2\sin(\frac{n\mu}{2})} e^{(in\Psi + i\frac{n\mu}{2})}$$

away from resonance, a normal form transformation takes away the pure oscillatory part and we have only:

$$h = -\mu A + c_0(A) = \text{const.}$$

$$\left[\text{homework : } \frac{dc_0(A)}{dA} \right]$$

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Single IP

If you are too lazy or too busy:

$$\Delta Q = \frac{-1}{2\pi} \frac{dc_0(A)}{dA}$$

is the detuning with amplitude, i.e. the amplitude dependent frequency change of the transformation we had before ...

We get:

$$\Delta Q = \frac{-1}{2\pi} \frac{Nr_0}{\gamma A} [1 - e^{-A\beta/2\sigma^2}] I_0(A\beta/2\sigma^2)$$

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Single IP - analysis of h

$$h = -\mu A + \sum_n c_n(A) \frac{n\mu}{2\sin(\frac{n\mu}{2})} e^{(in\Psi+i\frac{n\mu}{2})}$$

On resonance:

$$Q = \frac{p}{n} = \frac{\mu}{2\pi}$$

with $c_n \neq 0$:

$$\sin\left(\frac{n\pi p}{n}\right) = \sin(p\pi) \equiv 0 \quad \forall \text{ integer } p$$

and h diverges

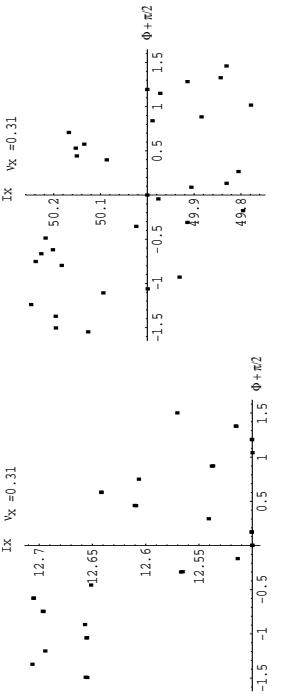
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Invariant versus tracking

- Is it useful what we obtained ?
 - ↑ Debug and compare ("benchmark")
- Compare to very simple tracking program:
 - ↑ linear transfer between interactions
 - ↑ beam-beam kick for round beam
 - ↑ compute action $I = \frac{\beta^*}{2\sigma^2} \left(\frac{x^2}{\beta^*} + p_x^2 \beta^* \right)$
 - ↑ and phase $\Psi = \arctan\left(\frac{p_x}{x}\right)$
 - ↑ compare I with h

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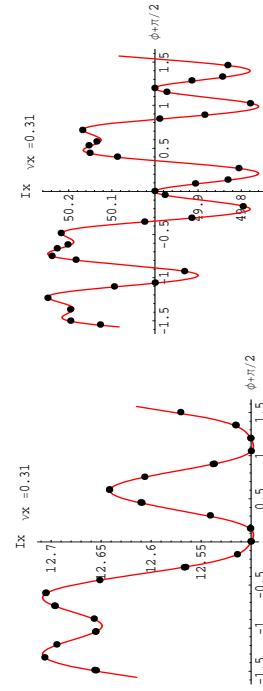
Invariant from tracking: one IP



→ Shown for $5\sigma_x$ and $10\sigma_x$

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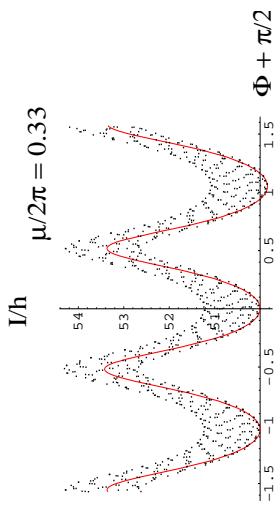
Invariant versus tracking: one IP



→ Shown for $5\sigma_x$ and $10\sigma_x$

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Invariant versus tracking:



→ Behaviour near a resonances: no more invariant possible

→ Envelope of tracking well described