# Recent Advances in the Rigorous Integration of Flows of ODEs with Taylor Models

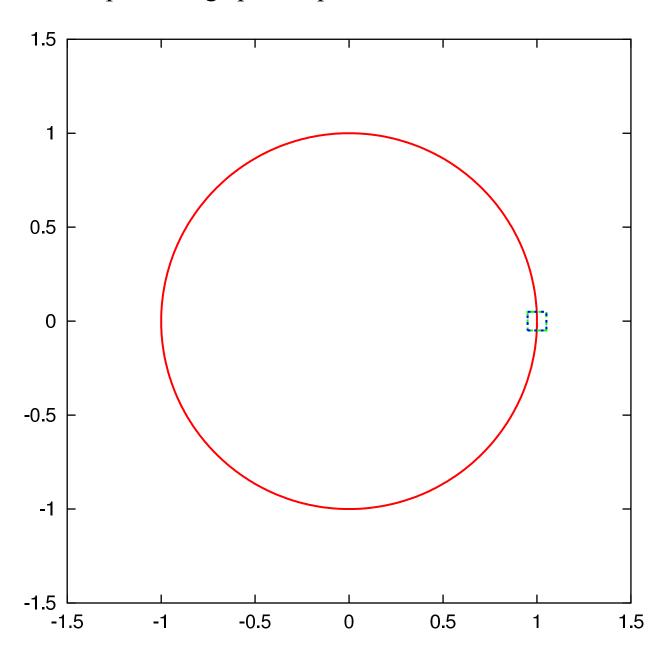
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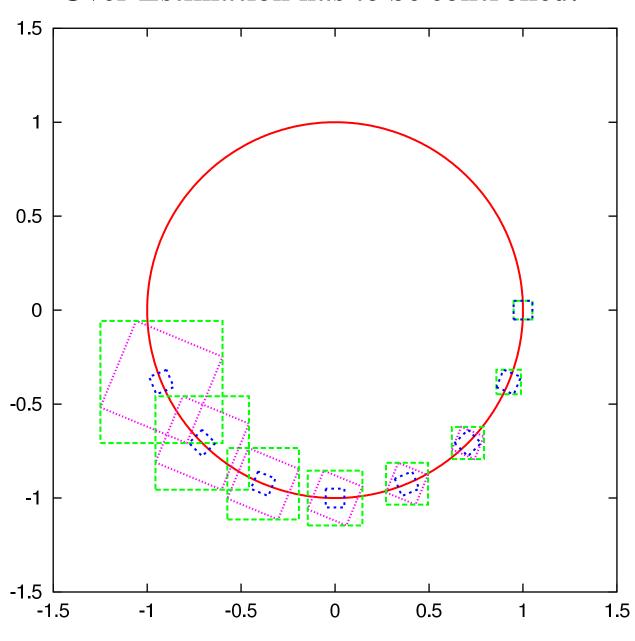
## Outline

- 1. Review of the old version of COSY-VI
- 2. The Reference Trajectory and the Flow Operator
- 3. Step Size Control
- 4. Error Parametrization of Taylor Models
- 5. Dynamic Domain Decomposition
- 6. Examples

To transport a large phase space volume with validation,



# Over Estimation has to be controlled.



# Review of the Old Version of COSY-VI

Version 2 (2004)

# Key Features and Algorithms of COSY-VI

- High order expansion not only in time t but also in transversal variables  $\vec{x}$ .
- Capability of weighted order computation, allowing to suppress the expansion order in transversal variables  $\vec{x}$ .
- Shrink wrapping algorithm including blunting to control ill-conditioned cases.
- Pre-conditioning algorithms based on the Curvilinear, QR decomposition, and blunting pre-conditioners.
- Resulting data is available in various levels including graphics output.

## The Volterra Equation

Describe dynamics of two conflicting populations

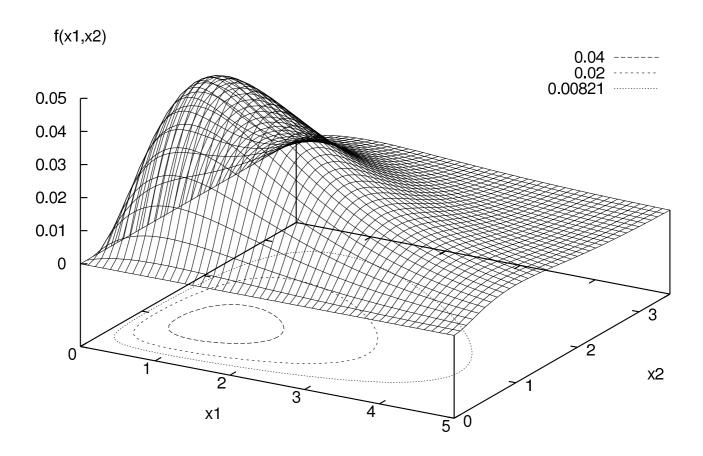
$$\frac{dx_1}{dt} = 2x_1(1-x_2), \quad \frac{dx_2}{dt} = -x_2(1-x_1)$$

Interested in initial condition

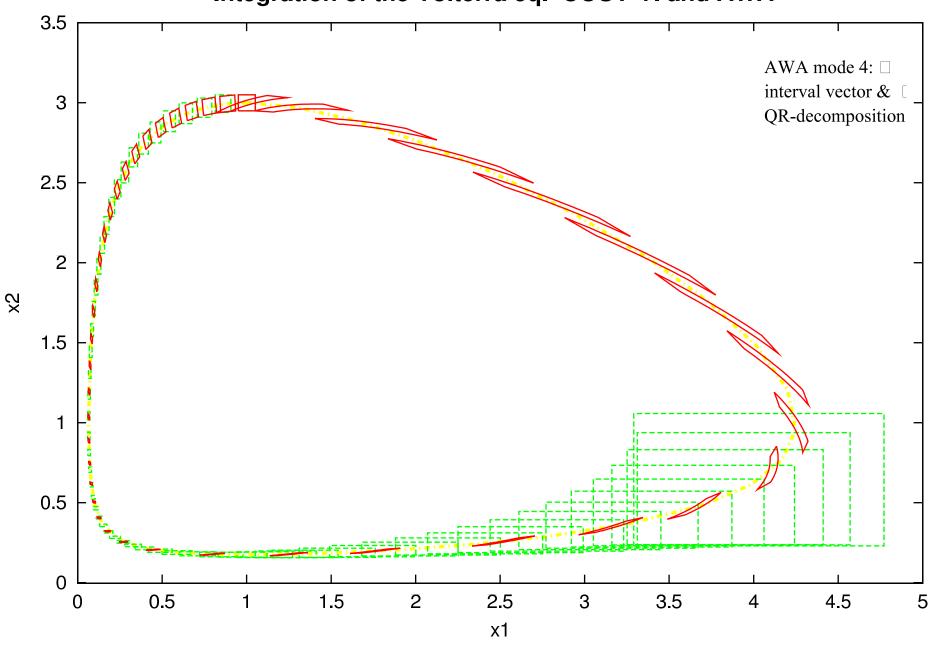
$$x_{01} \in 1 + [-0.05, 0.05], \quad x_{02} \in 3 + [-0.05, 0.05] \text{ at } t = 0.$$

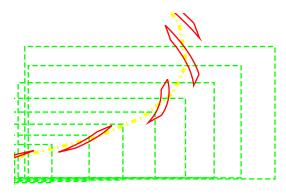
Satisfies constraint condition

$$C(x_1, x_2) = x_1 x_2^2 e^{-x_1 - 2x_2} = \text{Constant}$$



## Integration of the Volterra eq. COSY-VI and AWA





Piotr Zgliczynski, 2003

#### 2 Rössler equations

The Rössler equations are given by

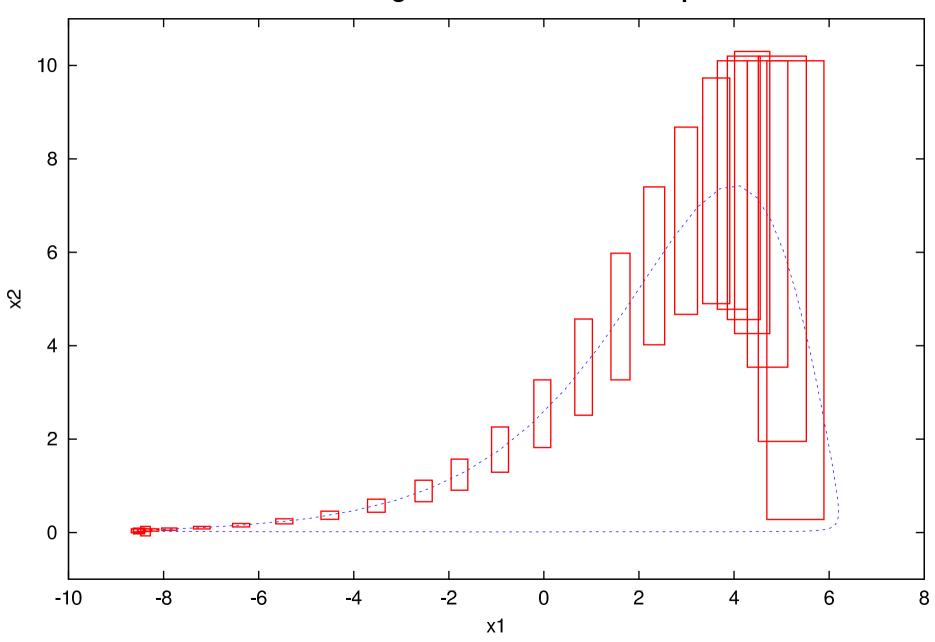
$$x' = -(y+z)$$
  
 $y' = x + 0.2y$   
 $z' = 0.2 + z(x-a),$  (4)

where a is a real parameter. We focus here at the value of a=5.7, where numerical simulations suggest an existence of a strange attractor.

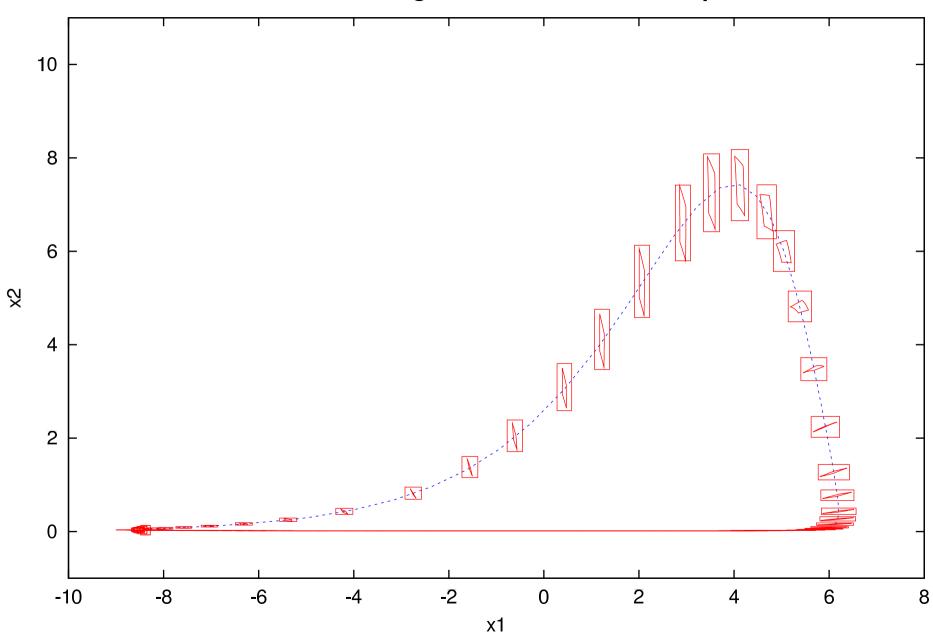
On section x=0 we consider the following initial condition  $(y,z)\in (-8.38095,0.0295902)+[-\delta,\delta]^2$ , where  $\delta$  should be considerably larger than  $10^{-3}$ . The integration time should be around T=6.

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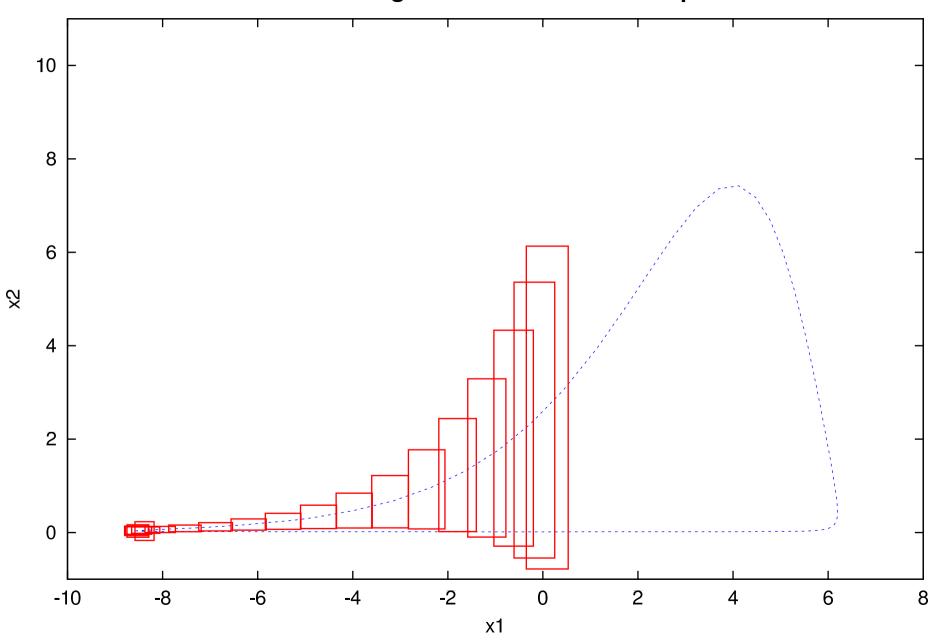
## **AWA Integration of the Roessler eqs.**



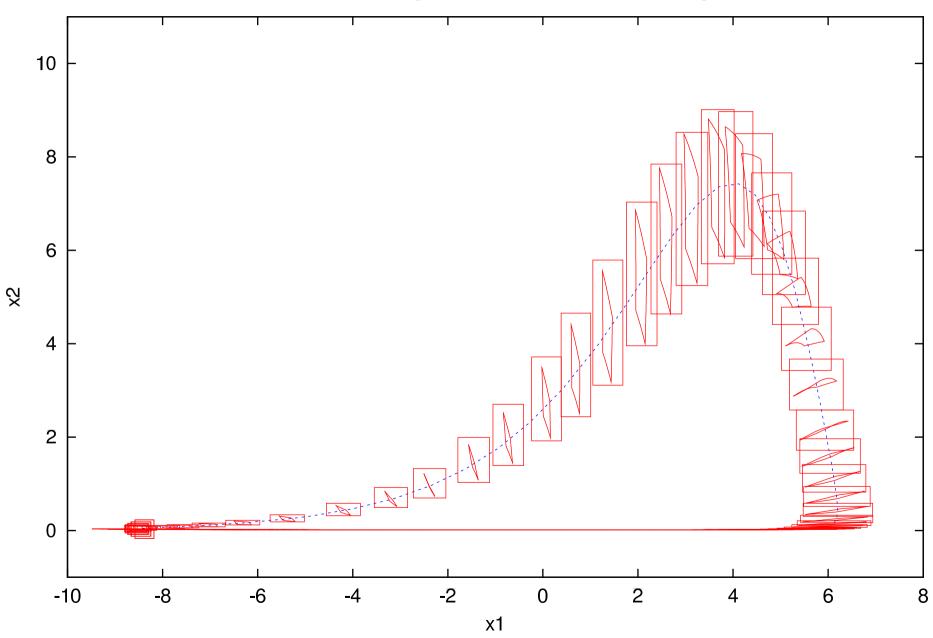
# **COSY-VI Integration of the Roessler eqs.**

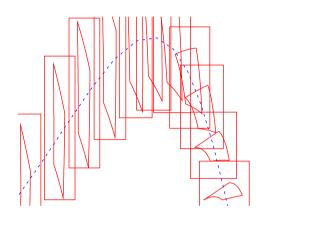


# **AWA Integration of the Roessler eqs.**



## **COSY-VI Integration of the Roessler eqs.**





## The Henon Map

Henon Map: frequently used elementary example that exhibits many of the well-known effects of nonlinear dynamics, including chaos, periodic fixed points, islands and symplectic motion. The dynamics is two-dimensional, and given by

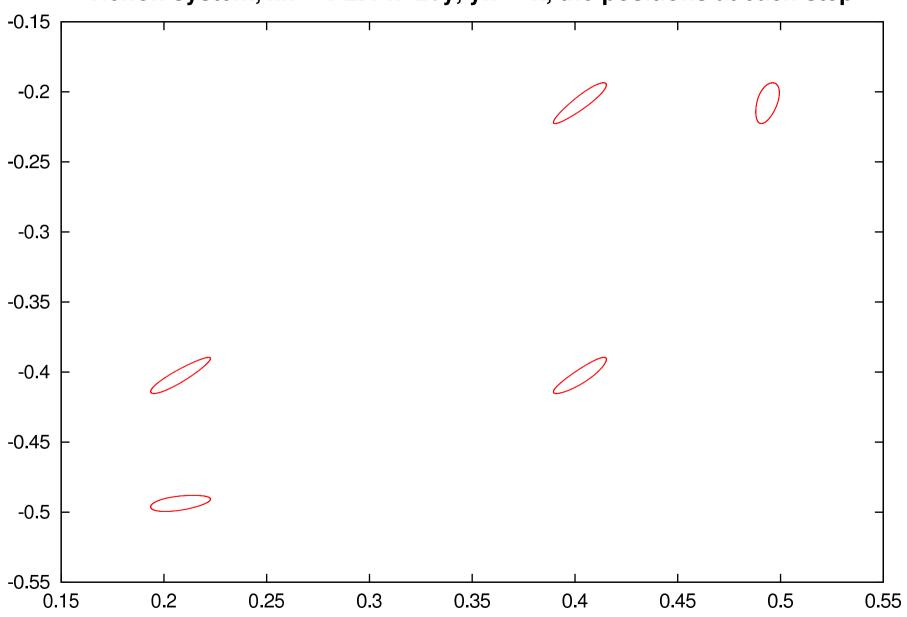
$$x_{n+1} = 1 - \alpha x_n^2 + y_n$$
$$y_{n+1} = \beta x_n.$$

It can easily be seen that the motion is area preserving for  $|\beta| = 1$ . We consider

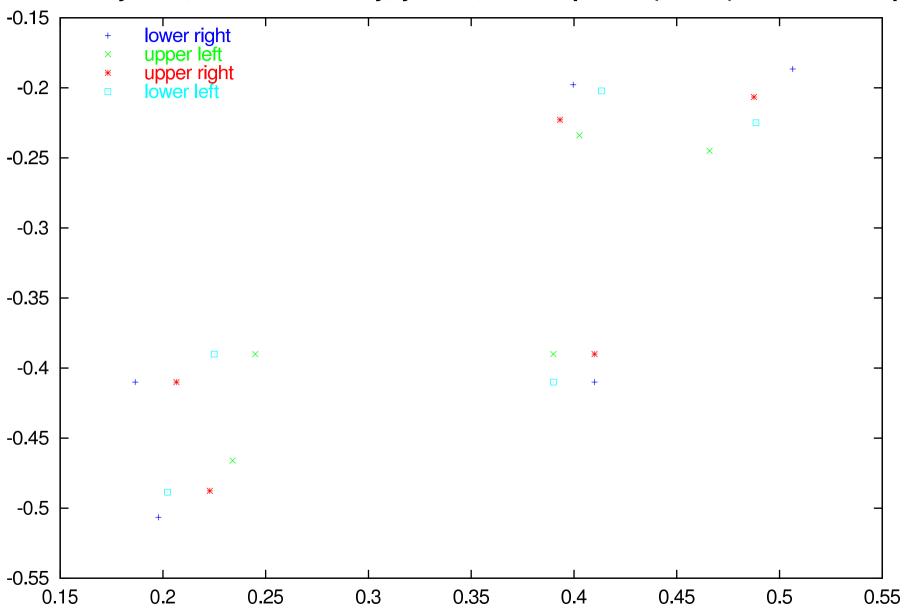
$$\alpha = 2.4$$
 and  $\beta = -1$ ,

and concentrate on initial boxes of the from  $(x_0, y_0) \in (0.4, -0.4) + [-d, d]^2$ .

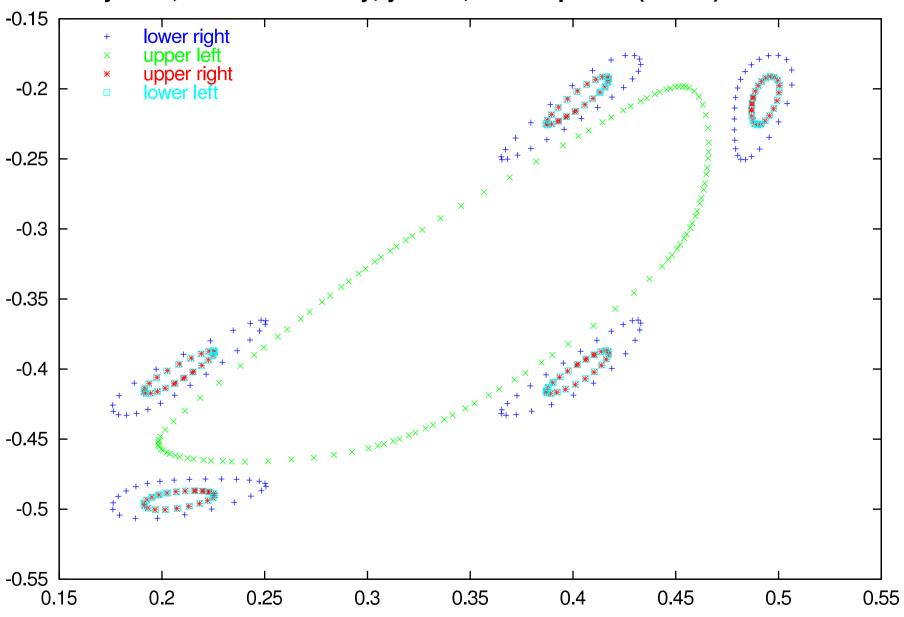
Henon system,  $xn = 1-2.4*x^2+y$ , yn = -x, the positions at each step



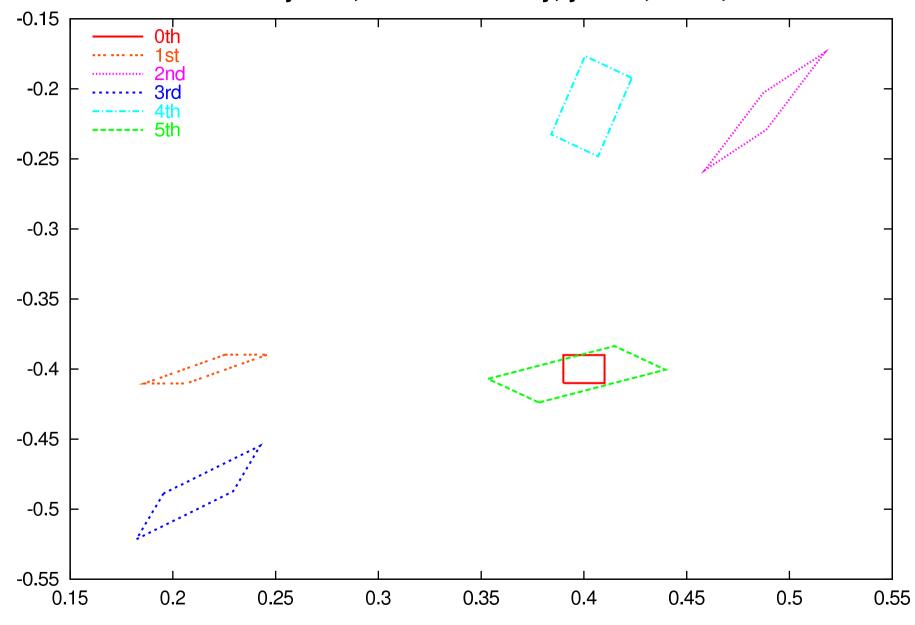
## Henon system, $xn = 1-2.4*x^2+y$ , yn = -x, corner points (+-0.01) the first 5 steps



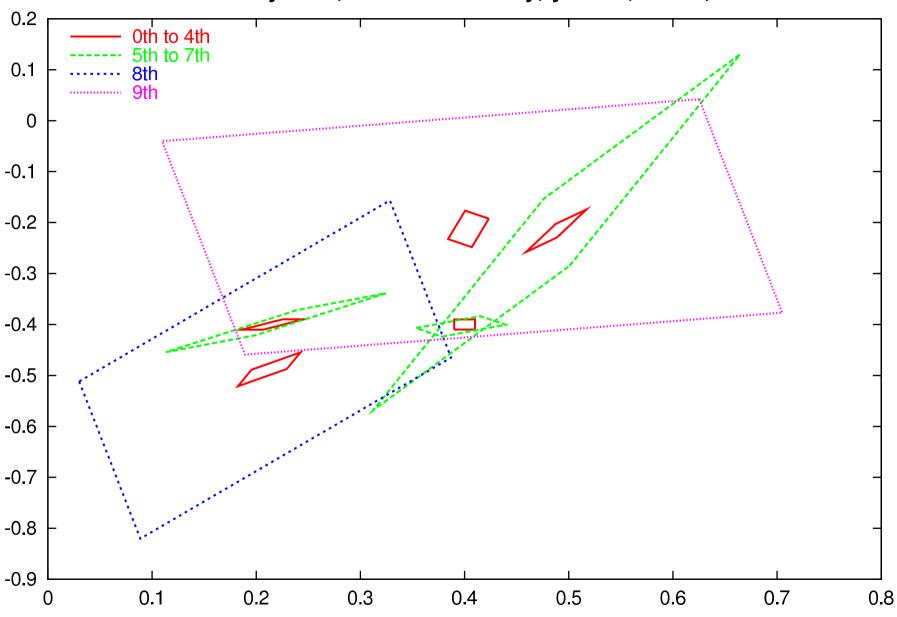
## Henon system, $xn = 1-2.4*x^2+y$ , yn = -x, corner points (+-0.01) the first 120 steps



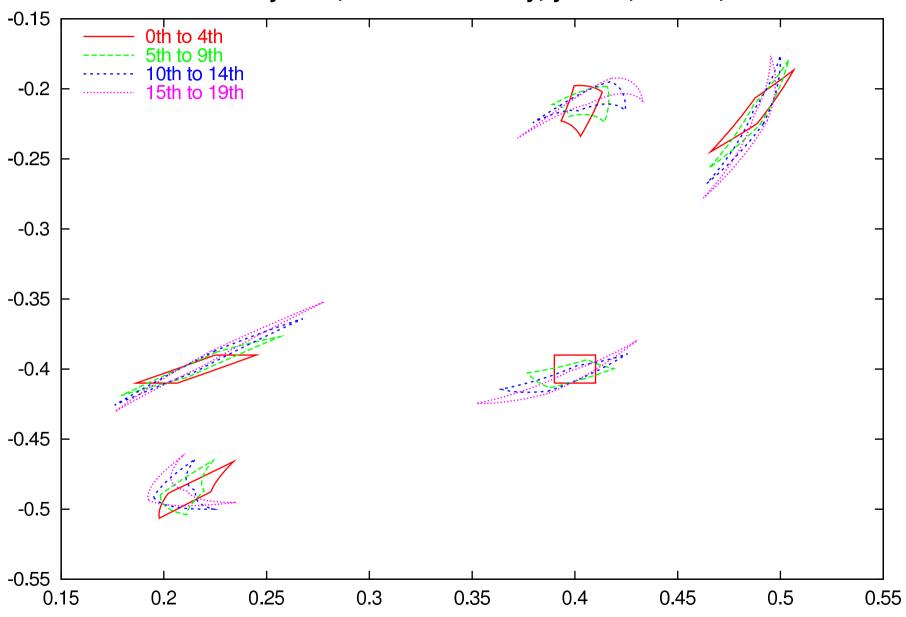
Henon system,  $xn = 1-2.4*x^2+y$ , yn = -x, NO=1, SW



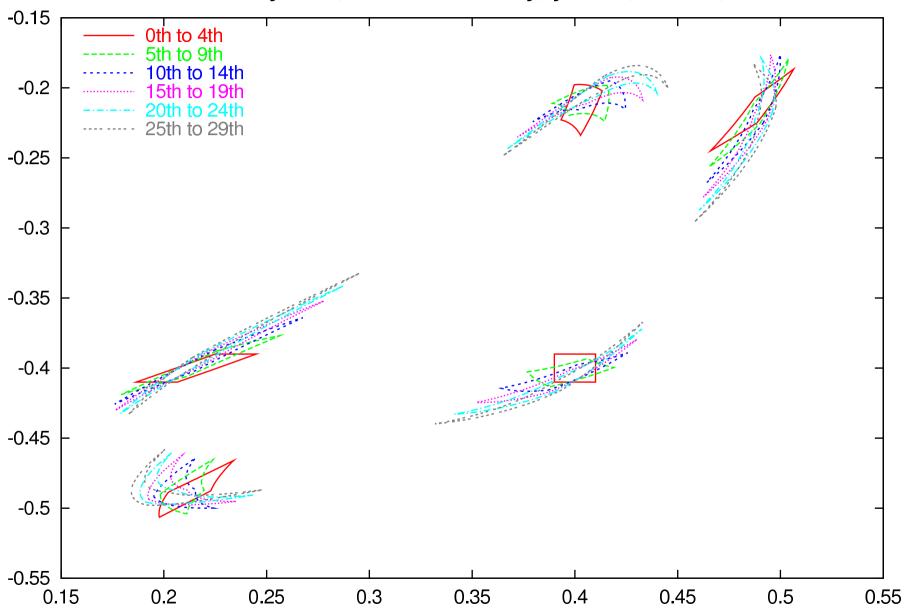
Henon system,  $xn = 1-2.4*x^2+y$ , yn = -x, NO=1, SW



Henon system,  $xn = 1-2.4*x^2+y$ , yn = -x, NO=20, SW



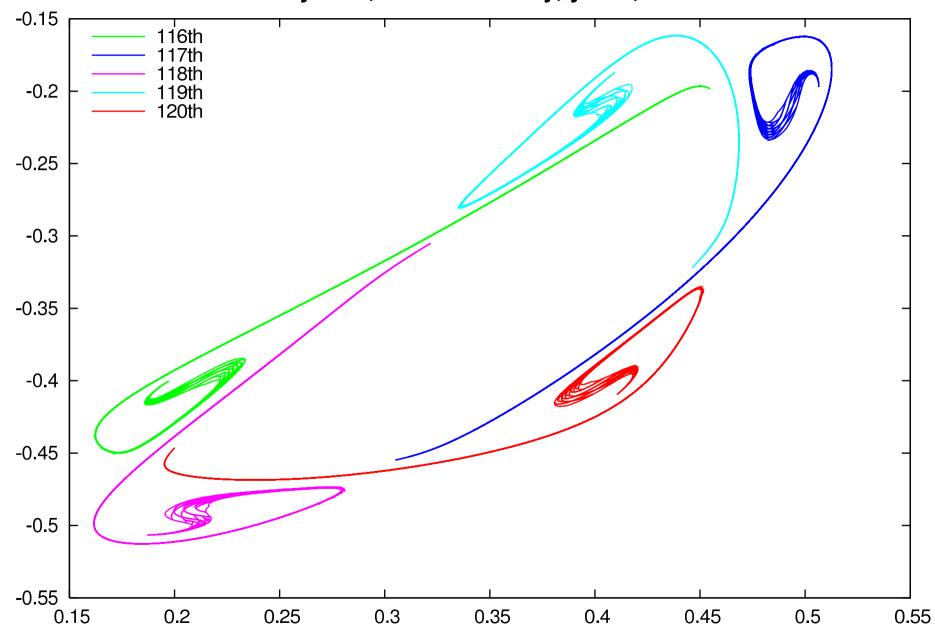
Henon system,  $xn = 1-2.4*x^2+y$ , yn = -x, NO=20, SW



## Review of the New Features

- The Reference Trajectory and the Flow Operator
- Step Size Control
- Error Parametrization of Taylor Models
- Dynamic Domain Decomposition

## Henon system, xn=1-2.4\*x^2+y, yn=-x, NO=33 w17



## The Reference Trajectory

First Step: Obtain Taylor expansion in time of solution of ODE of center point c, i.e. obtain

$$c(t) = c_0 + c_1 \cdot (t - t_0) + c_2 \cdot (t - t_0)^2 + \dots + c_n \cdot (t - t_0)^n$$

Very well known from day one how to do this with automatic differentiation. Rather convenient way: can be done by n iterations of the Picard Operator

$$c(t) = c_0 + \int_0^t f(r(t'), t)dt'$$

in one-dimensional Taylor arithmetic. Each iteration raises the order by one; so in each iteration i, only need to do Taylor arithmetic in order i. In either way, this step is **cheap** since it involves only **one-dimensional** operations.

#### The Nonlinear Flow

**Second Step:** The goal is to obtain Taylor expansion in time to order n and initial conditions to order k. Note:

- 1. This is usually the most **expensive** step. In the original Taylor model-based algorithm, it is done by n **iterations** of the Picard Operator in multi-dimensional Taylor arithmetic, where  $c_0$  is now a polynomial in initial conditions.
- 2. The case k = 1 has been known for a long time. Traditionally solved by setting up **ODEs for sensitivities** and solving these as before.
- 3. The case of higher k goes back to Beam Physics (M. Berz, Particle Accelerators 1988)
- 4. Newest Taylor model arithmetic naturally supports different expansions orders k for initial conditions and n for time.

Goal: Obtain flow with one single evaluation of right hand side.

#### The Nonlinear Relative ODE

We now develop a better way for second step.

**First:** introduce new "perturbation" variables  $\tilde{r}$  such that

$$r(t) = c(t) + A \cdot \tilde{r}(t).$$

The matrix A provides **preconditioning**. ODE for  $\tilde{r}(t)$ :

$$\tilde{r}' = A^{-1} [f(c(t) + A \cdot \tilde{r}(t)) - c'(t)]$$

**Second:** evaluate ODE for  $\tilde{r}'$  in Taylor arithmetic. Obtain a Taylor expansion of the ODE, i.e.

$$\tilde{r}' = P(\tilde{r}, t)$$

up to order n in time and k in  $\tilde{r}$ . Very important for later use: the polynomial P will have no constant part, i.e.

$$P(0,t) = 0.$$

#### Reminder: The Lie Derivative

Let

$$r' = f(r, t)$$

be a dynamical system. Let g be a variable in state space, and let us study g(r(t)), i.e. along a solution of the ODE. We have

$$\frac{d}{dt}g(t) = f \cdot \nabla g + \frac{\partial g}{\partial t}$$

Introducing the **Lie Derivative**  $L_f = f \cdot \nabla + \partial/\partial t$ , we have

$$\frac{d^n}{dt^n}g = L_f^n g \text{ and } g(t) \approx \sum_{i=0}^n \frac{(t-t_0)^i}{i!} L_f^i g /_{t=t_0}$$

## Differential Algebras on Taylor Polynomial Spaces

Consider space  $_nD_v$  of Taylor polynomials in v variables and order n with truncation multiplication. Formally: introduce **equivalence relation** on space of smooth functions

$$f =_n g$$

if all derivatives from 0 to n agree at 0. Class of f is denoted [f]. This induces addition, multiplication and scalar multiplication on classes. The resulting structure forms an algebra.

An algebra is a **Differential Algebra** if there is an operation  $\partial$ , called a derivation, that satisfies

$$\partial(s \cdot a + t \cdot b) = s \cdot \partial a + t \cdot \partial b$$
 and  $\partial(a \cdot b) = a \cdot (\partial b) + (\partial a) \cdot b$ 

for any vectors a and b and scalars s and t. Unfortunately, the **natural partial derivative** operations  $[f] \rightarrow [\partial_i f]$  does **not** introduce a differential algebra, because of loss of highest order.

## Differential Algebras on Taylor Polynomial Spaces

However, consider the modified operation

$$\partial_f$$
 with  $\partial_f g = f \cdot \nabla g$ 

If f is origin preserving, i.e. f(0) = 0, then  $\partial_f$  is a derivation on the space  ${}_nD_v$ . Why?

- Each derivative operation in the gradient  $\nabla g$  looses the highest order;
- but since f(0) = 0, the missing order in  $\nabla g$  does not matter since it does not contribute to the product  $f \cdot \nabla g$ .

## Polynomial Flow from Lie Derivative

Remember the ODE for  $\tilde{r}'$ :

$$\tilde{r}' = P(\tilde{r}, t)$$

up to order n in time and k in  $\tilde{r}$ . And remember P(0, t) = 0. Thus we can obtain the n-th order expansion of the flow as

$$\tilde{r}(t) = \sum_{i=0}^{n} \frac{(t - t_0)^i}{i!} \cdot \left( P \cdot \nabla + \frac{\partial}{\partial t} \right)^i \tilde{r}_0 \bigg/_{t=t_0}$$

- The fact that P(0,t)=0 restores the derivatives lost in  $\nabla$
- The fact that  $\partial/\partial t$  appears without origin-preserving factor limits the expansion to order n.

#### Performance of Lie Derivative Flow Methods

Apparently we have the following:

- Each term in the Lie derivative sum requires v + 1 derivations (very cheap, just re-shuffling of coefficients)
- $\bullet$  Each term requires v multiplications
- We need **one** evaluation of f in  ${}_{n}D_{v}$  (to set up ODE)

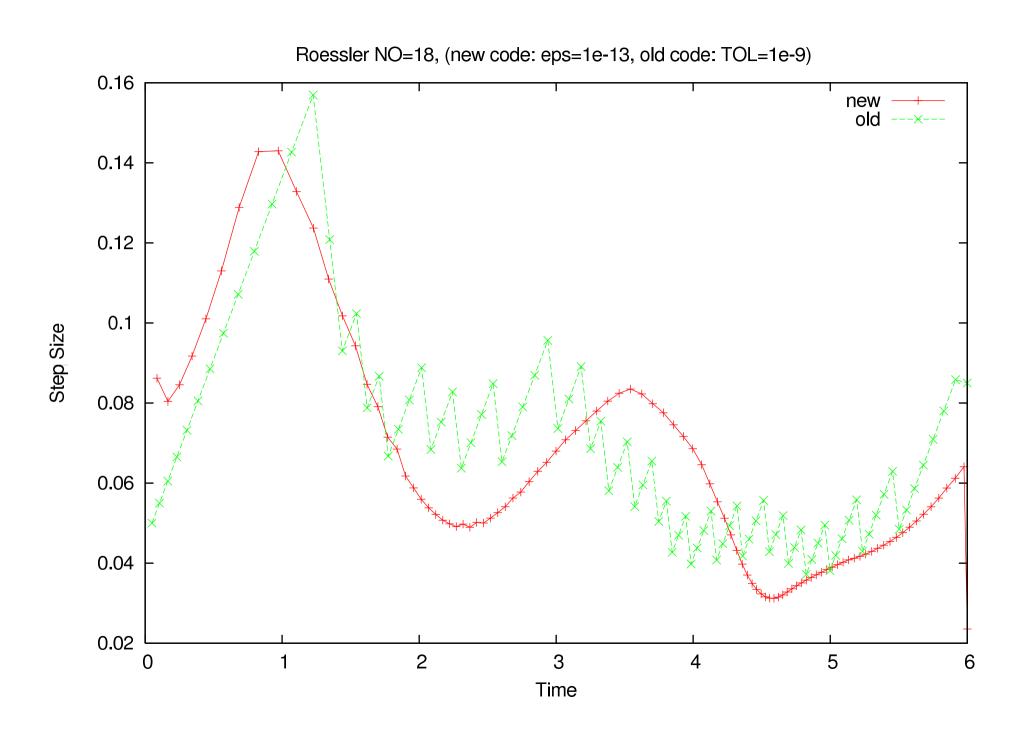
Compare this with the conventional algorithm, which requires n evaluations of the function f of the right hand side. Thus, roughly, if the evaluation of f requires more than v multiplications, the new method is more efficient.

- $\bullet$  Many practically appearing right hand sides f satisfy this.
- But on the other hand, if the function f does not satisfy this (for example for the linear case), then also P will be simple (in the linear case: P will be linear), and thus less operations appear

## Step Size Control

Step size control to maintain approximate error  $\varepsilon$  in each step. Based on a suite of tests:

- 1. Utilize the **Reference Orbit.** Extrapolate the size of coefficients for estimate of remainder error, scale so that it reaches and get  $\Delta t_1$ . Goes back to Moore in 1960s. This is one of conveniences when using Taylor integrators.
- 2. Utilize the **Flow.** Compute flow time step with  $\Delta t_1$ . Extrapolate the contributions of each order of flow for estimate of remainder error to get update  $\Delta t_2$ .
- 3. Utilize a **Correction factor** c to account for overestimation in TM arithmetic as  $c = \sqrt[n+1]{|R|/\varepsilon}$ . Largely a measure of complexity of ODE. Dynamically update the correction factor.
- 4. Perform verification attempt for  $\Delta t_3 = c \cdot \Delta t_2$



COSY-VI Roessler until Break-down, Step Size, April 13 2007 0.16 Step Size 0.14 0.12 0.1 Step Size 0.08 0.06 0.04 0.02 0 2 8 4 6 10 12 14 16 18 20 0 Time

## Error Parametrization of Taylor models

**Motivation:** Is it possible to absorb the remainder error bound intervals of Taylor models into the polynomial parts using additional parameters?

Phrase the question as the following problem:

1. Have Taylor models with 0 remainder error interval, which depend on the independent variables  $\vec{x}$  and the parameters  $\vec{\alpha}$ .

$$\vec{T}_0 = \vec{P}_0(\vec{x}, \vec{\alpha}) + \overrightarrow{[0, 0]}.$$

2. Perform Taylor model arithmetic on  $\vec{T}_0$ , namely  $\vec{F}(\vec{T}_0)$ 

$$\vec{F}(\vec{T}_0) = \vec{P}(\vec{x}, \vec{\alpha}) + \vec{I}_F$$
, where  $\vec{I}_F \neq [0, 0]$ .

3. Try to absorb  $\vec{I}_F$  into the polynomial part that depends on  $\vec{\alpha}$ 

$$\vec{P}(\vec{x}, \vec{\alpha}) + \vec{I}_F \subseteq \vec{P}'(\vec{x}, \vec{\alpha}) + [0, 0].$$
 (A)

#### Observe

$$\vec{P}(\vec{x}, \vec{\alpha}) = \underbrace{\vec{P}(\vec{x}, 0)}_{\vec{\alpha} \text{-indep.}} + \underbrace{\vec{P}(\vec{x}, \vec{\alpha}) - \vec{P}(\vec{x}, 0)}_{\vec{\alpha} \text{-dependent}} = \vec{P}(\vec{x}, 0) + \vec{P}_{\alpha}(\vec{x}, \vec{\alpha})$$

The size of  $\vec{P}(\vec{x},0)$  is much larger than the rest, because the rest is essentially errors. The process of (A) does not alter  $\vec{P}(\vec{x},0)$ , so set the  $\vec{\alpha}$ -independent part  $\vec{P}(\vec{x},0)$  aside from the whole process, which helps the numerical stability of the process.

The task is now

$$\vec{P}_{\alpha}(\vec{x}, \vec{\alpha}) + \vec{I}_F \subseteq \vec{P}'_{\alpha}(\vec{x}, \vec{\alpha}) + [0, 0].$$

We limit  $\vec{P}_{\alpha}(\vec{x}, \vec{\alpha})$  to be only **linearly** dependent on  $\vec{\alpha}$ .

$$\vec{P}_{\alpha}(\vec{x}, \vec{\alpha}) + \vec{I}_F = \left(\widehat{M} + \widehat{M}(\vec{x})\right) \cdot \vec{\alpha} + \vec{I}_F.$$

Express  $\vec{I}_F$  by the matrix form using additional parameters  $\vec{\beta}$ 

$$\vec{I}_F \subseteq \left(\widehat{I}_F + \widehat{\bar{I}}_F(\vec{x})\right) \cdot \vec{\beta}.$$

where  $\widehat{\overline{I}}_F(\vec{x}) = 0$  and  $(\widehat{I}_F)_{ii} = |I_{Fi}|$ .

$$\vec{P}_{\alpha}(\vec{x}, \vec{\alpha}) + \vec{I}_F \subseteq \left(\widehat{M} + \widehat{M}(\vec{x})\right) \cdot \vec{\alpha} + \left(\widehat{I}_F + \widehat{\bar{I}}_F(\vec{x})\right) \cdot \vec{\beta}.$$

View this as a collection of  $2 \cdot v$  column vectors associated to  $2 \cdot v$  parameters  $\vec{\alpha}$  and  $\vec{\beta}$ . Recall a matrix, or a collection of v column vectors, represent a parallelepiped. The problem is now to find a set sum of two parallelepipeds.

## Psum Algorithm for choosing column vectors

**Task**: Choose v vectors out of n vectors  $\vec{s_i}$ ,  $i = 1, ..., n, n \ge v$ .

- 1. Choose the longest vector  $\vec{s}_k$ , and assign it as  $\vec{t}_1$ . Normalize it as  $\vec{e}_1 = \vec{t}_1 / |\vec{t}_1|$ .
- 2. Out of the remaining vectors  $\vec{s}_i$ , choose the j-th vector  $\vec{t}_j = \vec{s}_k$  such that

$$\frac{|\vec{s}_k|^2 - \sum_{m=1}^{j-1} |\vec{s}_k \cdot \vec{e}_m|^2}{|\vec{s}_k|^{2p}}$$

is largest. Compute  $\vec{e_j}$ , the orthonormalized vector of  $\vec{t_j}$  to  $\vec{e_1}, ..., \vec{e_{j-1}}$ . (Gram-Schmidt)

3. Repeat the process 2 until j = v.

Experimentally, p = 0.5 is found to be efficient and robust for obtaining a set sum of two parallelepipeds

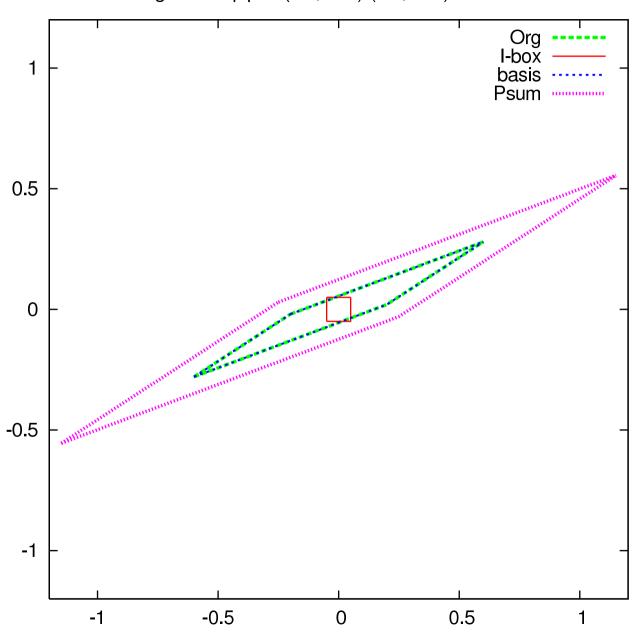
## Psum Algorithm for two parallelepipeds

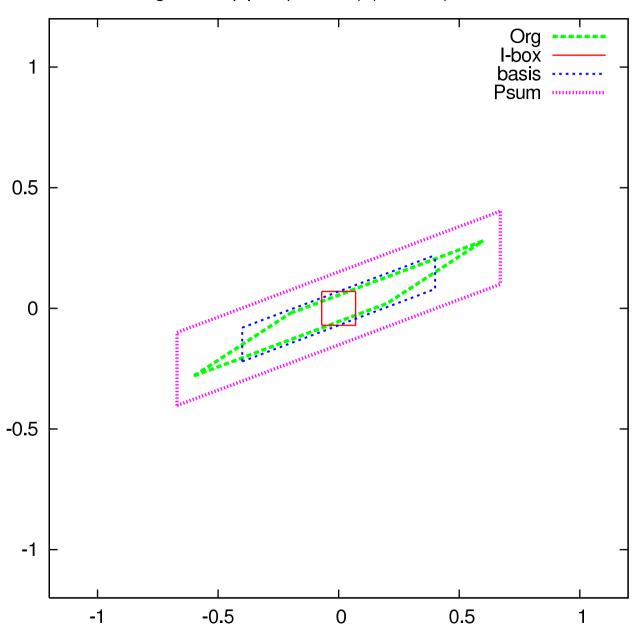
**Task**: Obtain a set sum of two parallelepipeds  $\widehat{M}_1$  and  $\widehat{M}_2$ .

- 1. Prepare the basis  $\widehat{M}_b$  using the Psum algorithm for choosing v column vectors out of  $2 \cdot v$  column vectors from  $\widehat{M}_1$  and  $\widehat{M}_2$ .
- 2. Compute conditioned parallelepipeds  $\widehat{M}_b^{-1} \cdot \widehat{M}_1$  and  $\widehat{M}_b^{-1} \cdot \widehat{M}_2$ .
- 3. Confine the conditioned parallelepipeds by bounding them.

$$\vec{B}_1 = \text{bound}\left(\widehat{M}_b^{-1} \cdot \widehat{M}_1\right) \text{ and } \vec{B}_2 = \text{bound}\left(\widehat{M}_b^{-1} \cdot \widehat{M}_2\right).$$

- 4. Compute the interval sum  $\vec{B} = \vec{B_1} + \vec{B_2}$ .  $\vec{B}$  confines the conditioned set sum of the conditioned parallelepipeds.
- 5. From  $\vec{B}$ , set up a parallelepiped as a box  $\hat{B} = \begin{pmatrix} |B_1| & 0 \\ & \ddots & \\ 0 & |B_v| \end{pmatrix}$ .
- 6. Compute  $\widehat{M}_b \cdot \widehat{B}$ , which is a set sum of  $\widehat{M}_1$  and  $\widehat{M}_2$  under  $\widehat{M}_b$ .





#### Error Absorption

We now chose a favoured collection of v column vectors  $\widehat{L} + \widehat{\overline{L}}(\vec{x})$  using the Psum algorithm. Collect the left over v column vectors to  $\widehat{E} + \widehat{\overline{E}}(\vec{x})$ . Associate them to  $2 \cdot v$  parameters  $\vec{\alpha}'$  and  $\vec{\beta}'$ .

$$\vec{P}_{\alpha}(\vec{x}, \vec{\alpha}) + \vec{I}_F \subseteq \left(\widehat{L} + \widehat{\bar{L}}(\vec{x})\right) \cdot \vec{\alpha}' + \left(\widehat{E} + \widehat{\bar{E}}(\vec{x})\right) \cdot \vec{\beta}'.$$

Since  $\vec{\alpha}$  and  $\vec{\beta}$  do not appear anymore, we can rename  $\vec{\alpha}'$  and  $\vec{\beta}'$  as  $\vec{\alpha}$  and  $\vec{\beta}$  for the simplicity.

$$\vec{P}_{\alpha}(\vec{x}, \vec{\alpha}) + \vec{I}_{F} \subseteq \left(\widehat{L} + \widehat{\bar{L}}(\vec{x})\right) \cdot \vec{\alpha} + \left(\widehat{E} + \widehat{\bar{E}}(\vec{x})\right) \cdot \vec{\beta} 
= \widehat{L} \circ \left[\widehat{L}^{-1} \circ \left(\widehat{L} + \widehat{\bar{L}}(\vec{x})\right) \cdot \vec{\alpha} + \widehat{L}^{-1} \circ \left(\widehat{E} + \widehat{\bar{E}}(\vec{x})\right) \cdot \vec{\beta}\right] 
\subseteq \widehat{L} \circ \left[\left(\widehat{I} + \widehat{L}^{-1} \circ \widehat{\bar{L}}(\vec{x})\right) \cdot \vec{\alpha} + \widehat{B} \cdot \vec{\beta}\right]$$

where  $\widehat{B}$  is a diagonal matrix with the *i*-th element is  $|B_i|$  and  $\overrightarrow{B} = \text{bound}(\widehat{L}^{-1} \circ (\widehat{E} + \widehat{E}(\overrightarrow{x})) \cdot \overrightarrow{\beta})$ .

If the diagonal terms of  $(\widehat{I} + \widehat{L}^{-1} \circ \widehat{\overline{L}}(\vec{x}))$  are positive,

$$\vec{P}_{\alpha}(\vec{x}, \vec{\alpha}) + \vec{I}_{F} \subseteq \widehat{L} \circ \left[ \left( \widehat{I} + \widehat{L}^{-1} \circ \widehat{\bar{L}}(\vec{x}) \right) \cdot \vec{\alpha} + \widehat{B} \cdot \vec{\alpha} \right] 
= \widehat{L} \circ \left( \widehat{I} + \widehat{L}^{-1} \circ \widehat{\bar{L}}(\vec{x}) \right) \cdot \vec{\alpha} + \widehat{L} \circ \widehat{B} \cdot \vec{\alpha} 
= \left( \widehat{L} + \widehat{\bar{L}}(\vec{x}) + \widehat{L} \circ \widehat{B} \right) \cdot \vec{\alpha}.$$

**Note**: A modification to use  $\widehat{A}$  instead of  $\widehat{L}$ , when  $\widehat{A} \approx \widehat{L}$ , is done easily. This involves bounding of  $\widehat{A}^{-1} \circ \left(\widehat{L} - \widehat{A}\right) \cdot \overrightarrow{\alpha}$  and the diagonal terms to be checked positive are those of  $\left(\widehat{I} + \widehat{A}^{-1} \circ \widehat{L}(\overrightarrow{x})\right)$ .

henon (area preserving). Performance Comparison. TM order 13, IC width 4e-3 0.01 0.0001 1e-006 Error 1e-008 1e-010 1e-012 Error parametrization w Psum old PreConditioning w QR old PreConditioning w blunting old naive 1e-014 1e-016 50 100 300 150 200 250 Step

#### Cost of Additional Parameters

For a v dimensional system, we need v parameters  $\vec{\alpha}$  to absorb Taylor model remainder error bound intervals. The dependence on  $\vec{\alpha}$  is limited to **linear**. So, we use weighted DA. Choose an appropriate weight order w for  $\vec{\alpha}$ .

• The dependence on  $\vec{\alpha}$  has to be kept linear. Namely  $2 \cdot w > n$ , where n is the computational order of Taylor models. Choose

$$w = \operatorname{Int}\left(\frac{n}{2}\right) + 1.$$

Maximum size necessary for DA and TM for v=2.

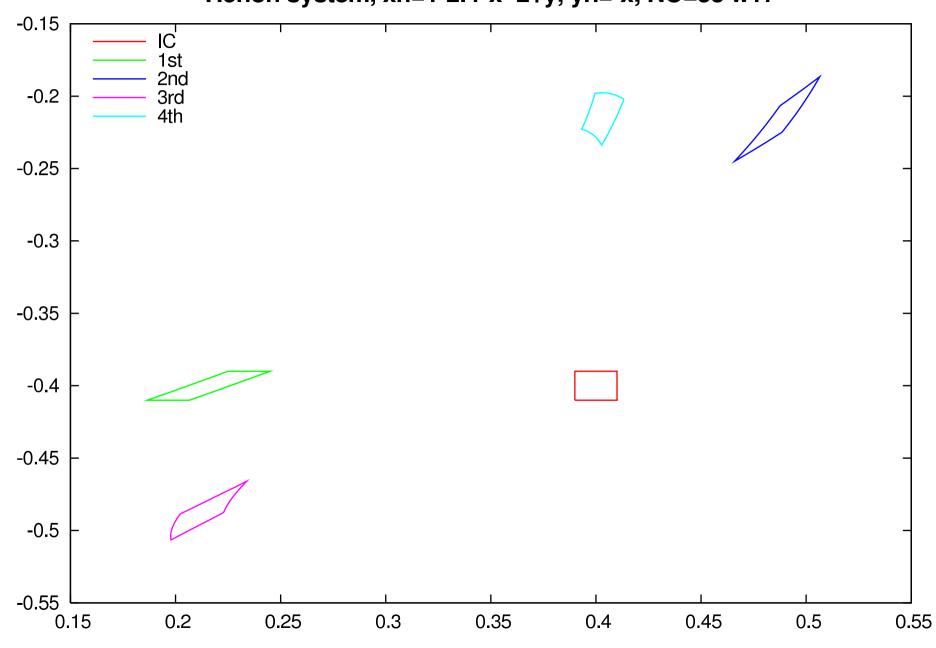
## Dynamic Domain Decomposition

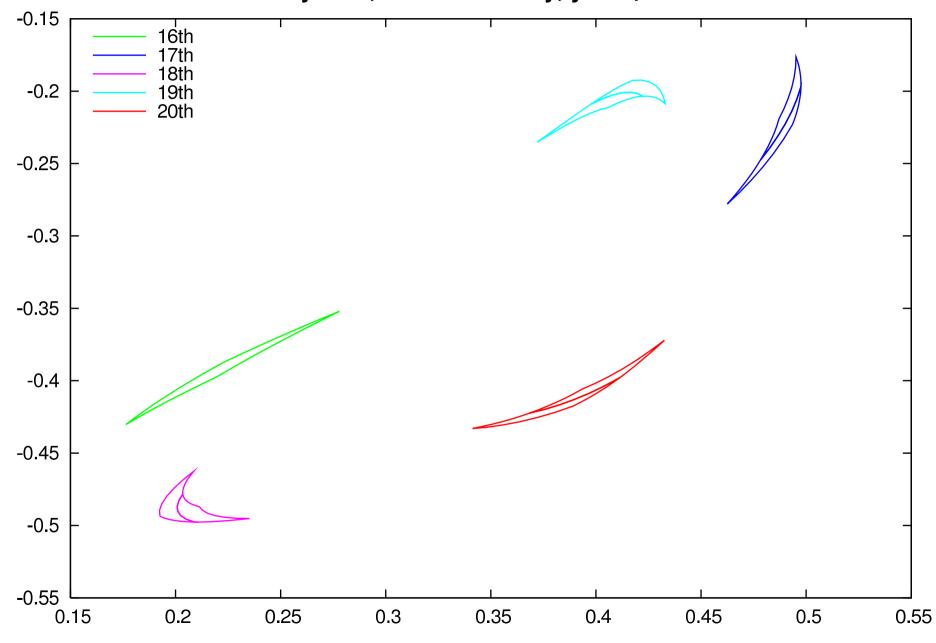
For extended domains, this is **natural equivalent** to step size control. Similarity to what's done in global optimization.

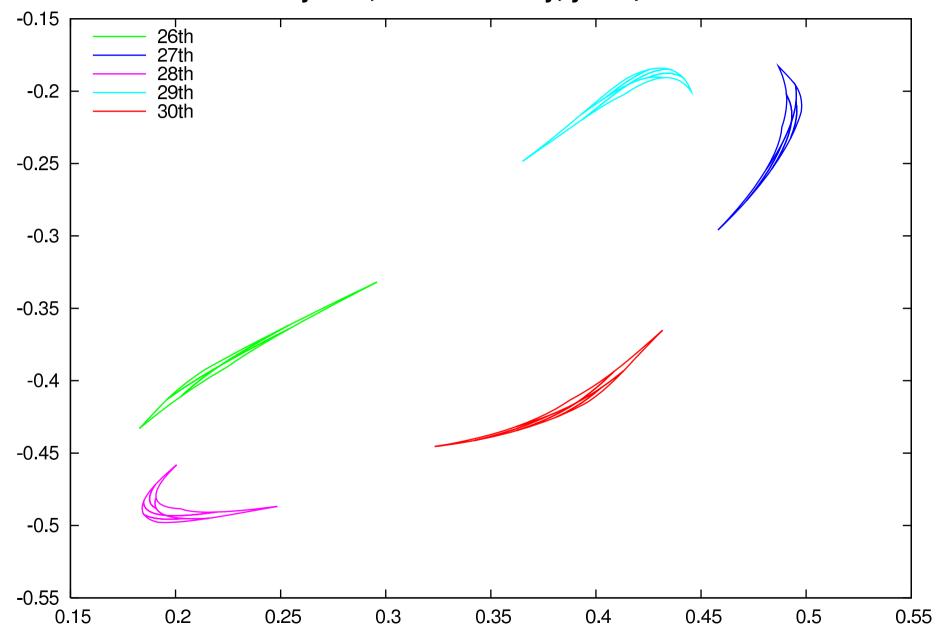
- 1. Evaluate ODE for  $\Delta t = 0$  for current flow.
- 2. If resulting remainder bound R greater than  $\varepsilon$ , split the domain along variable leading to longest axis.
- 3. Absorb R in the TM polynomial part using the error parametrization method. If it fails, split the domain along variable leading to largest x dependence of the error.
- 4. Put one half of the box on stack for future work.

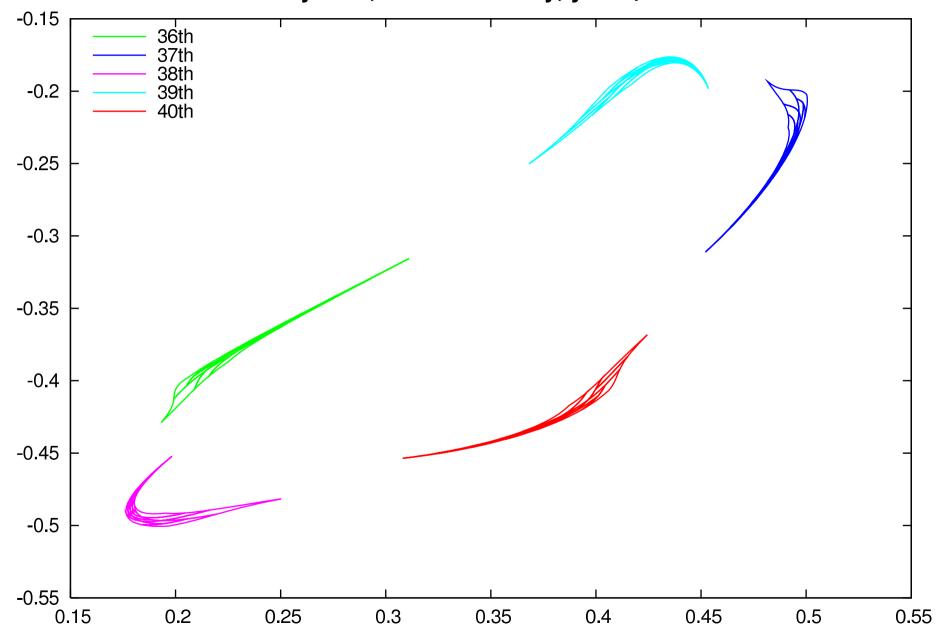
Things to consider:

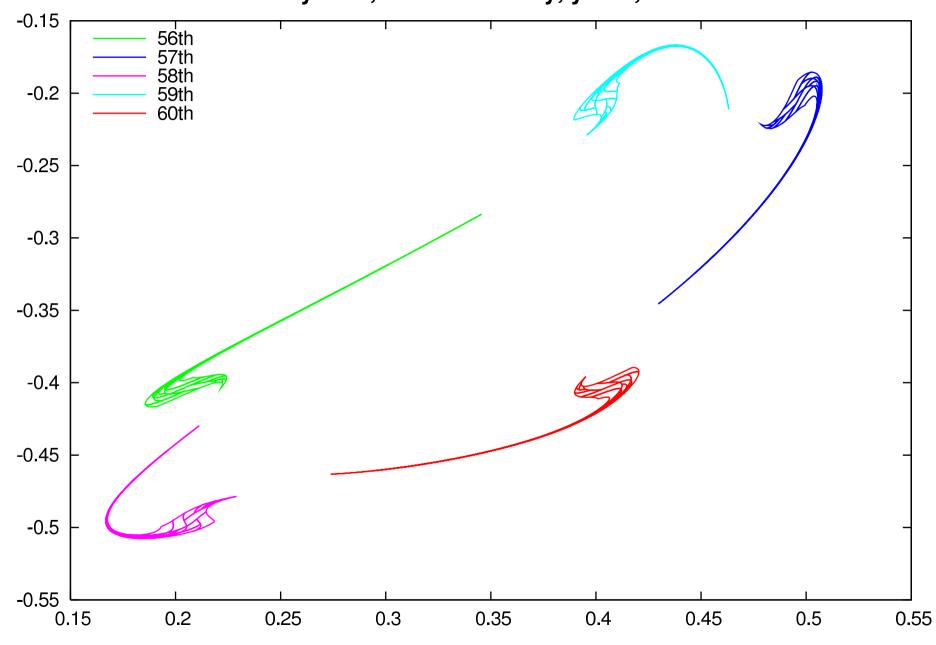
- Utilize "First-in-last-out" stack; minimizes stack length. Special adjustments for stack management in a parallel environment, including load balancing.
- Outlook: also dynamic order control for dependence on initial conditions

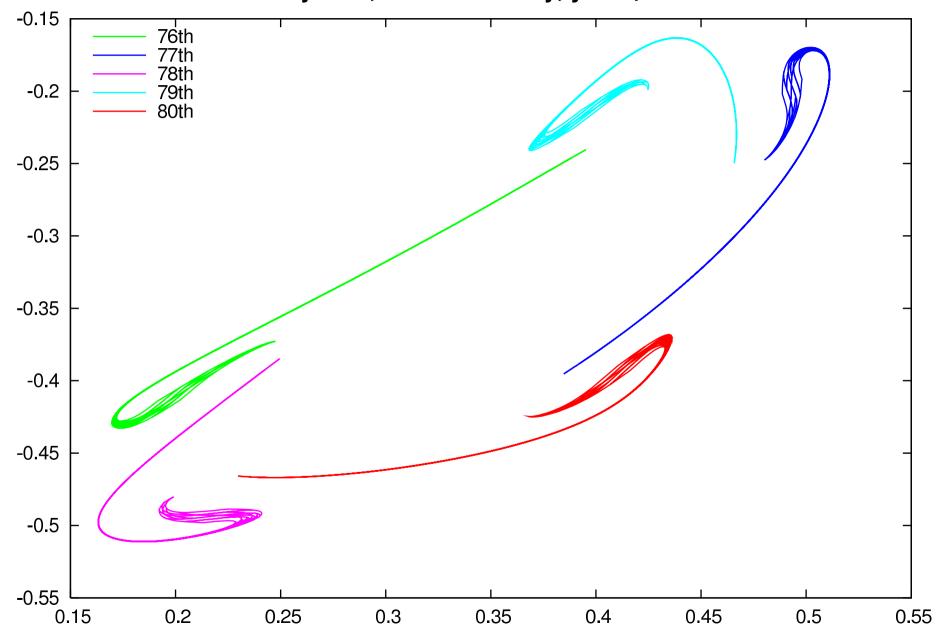


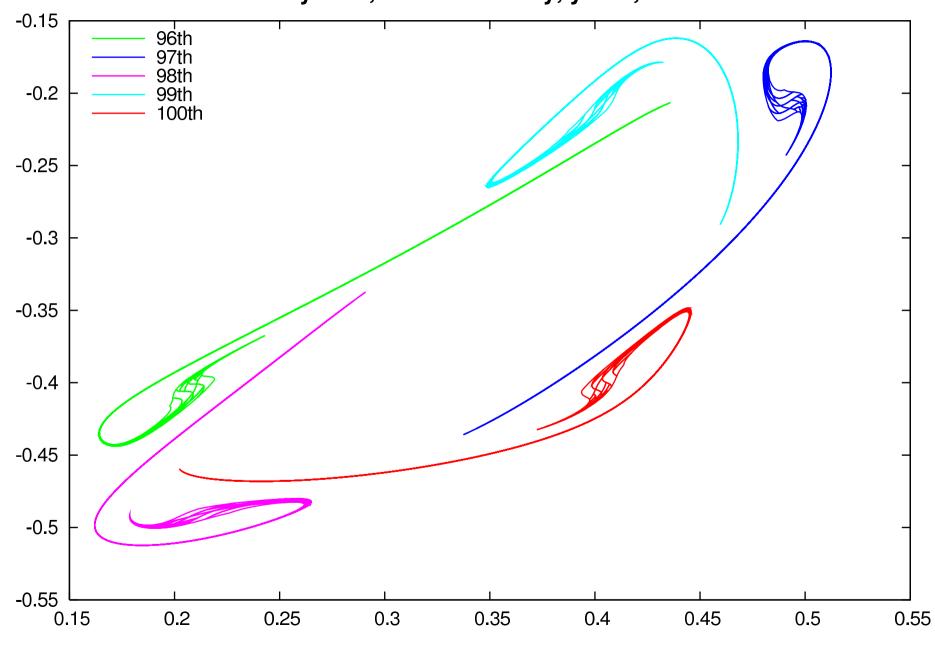


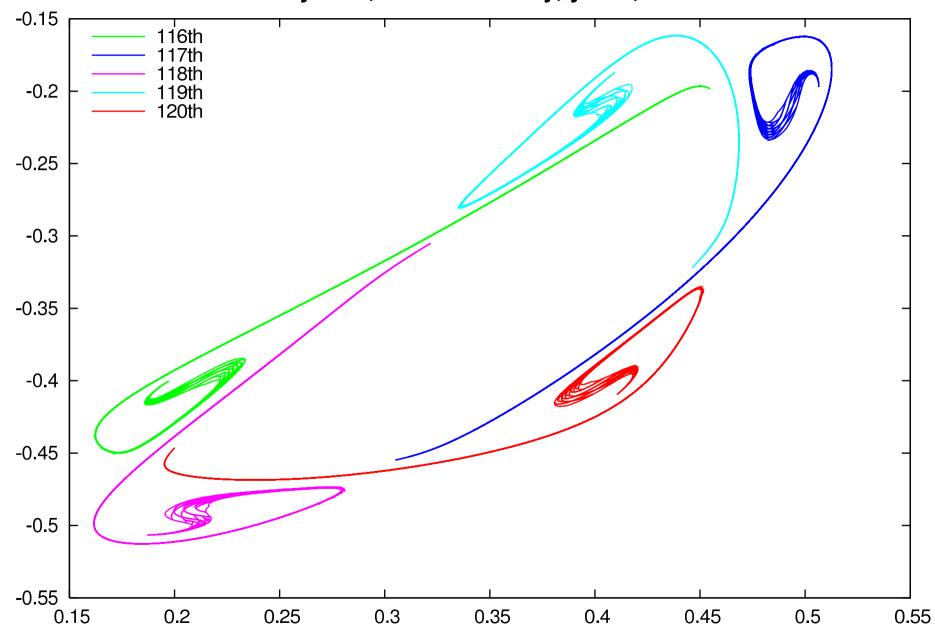


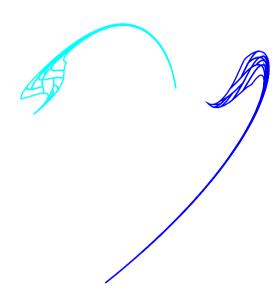




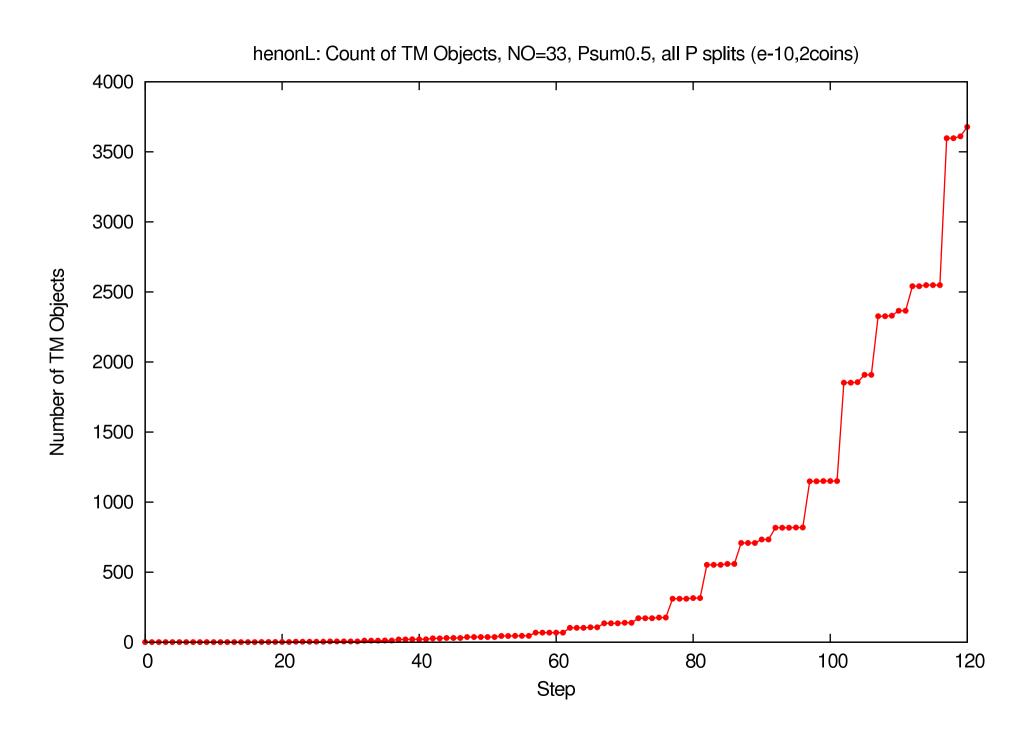


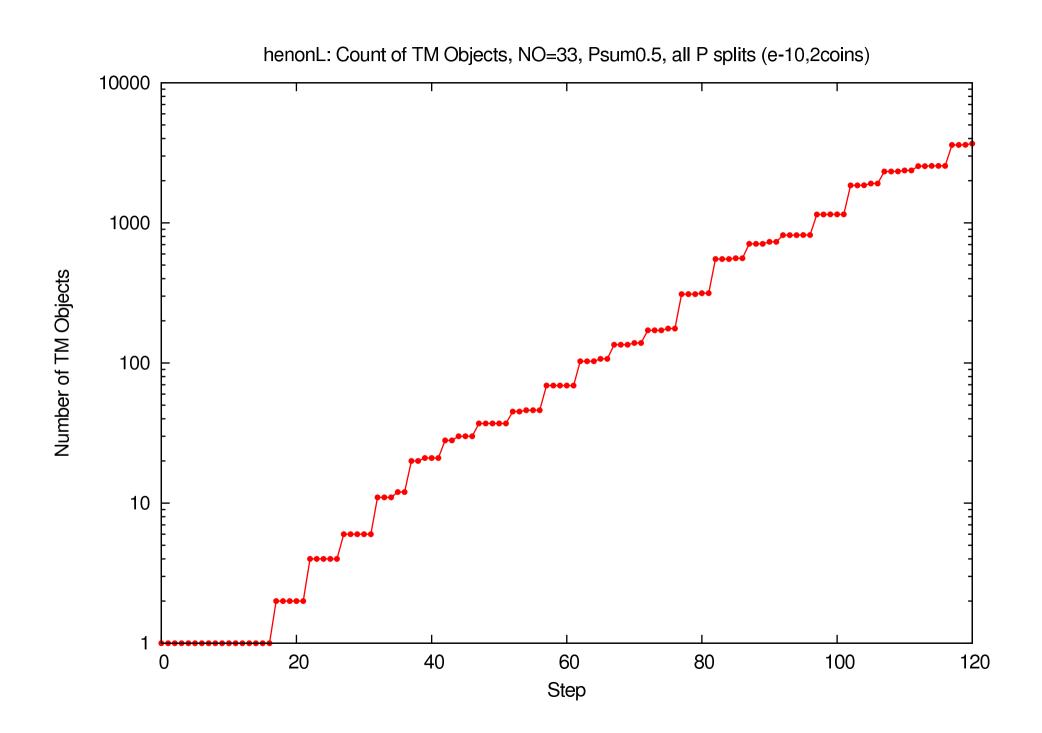




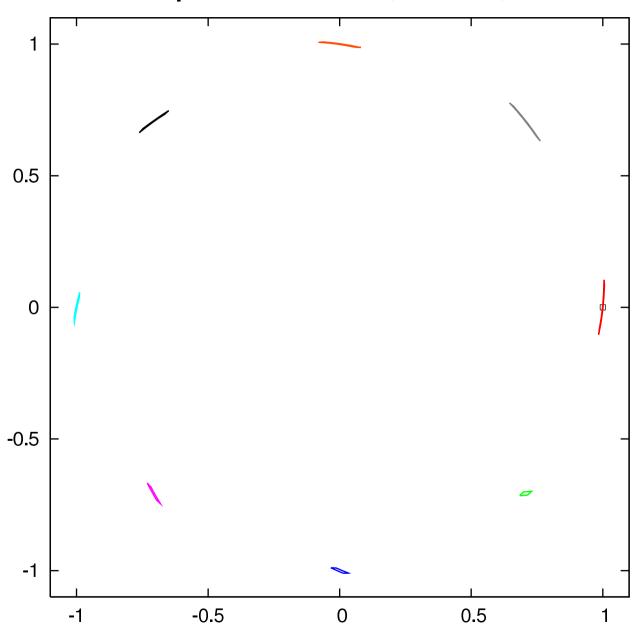




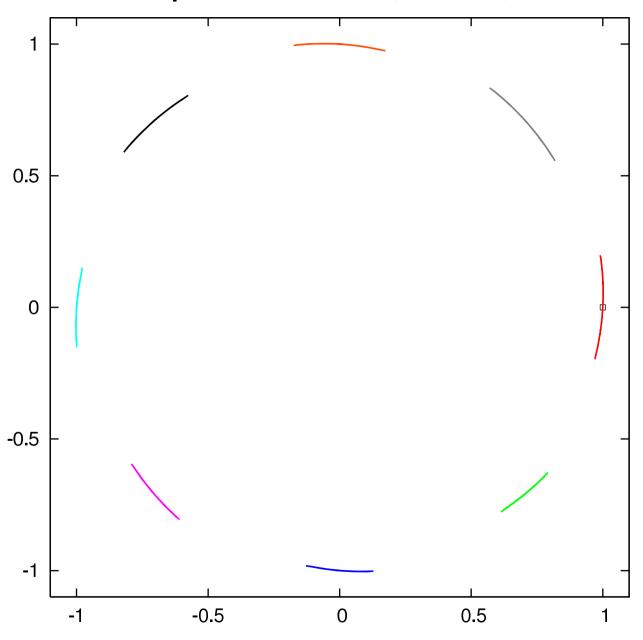




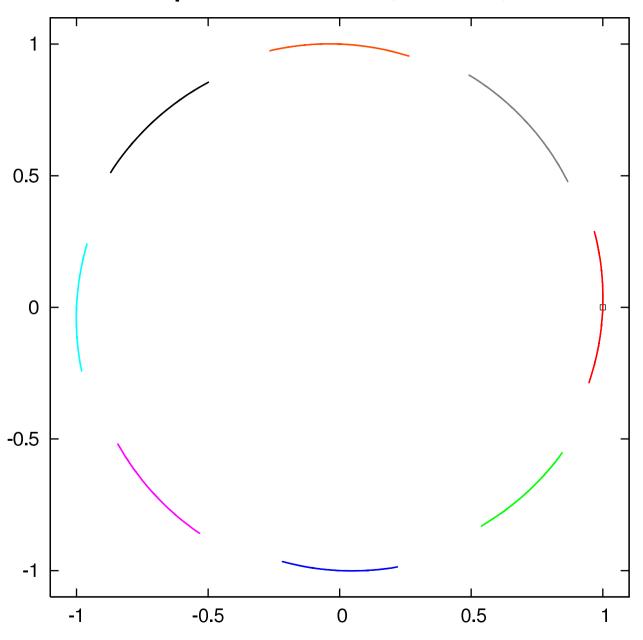
## discrete kepler. 1st revolution, ICw 0.02, NO=13 w7



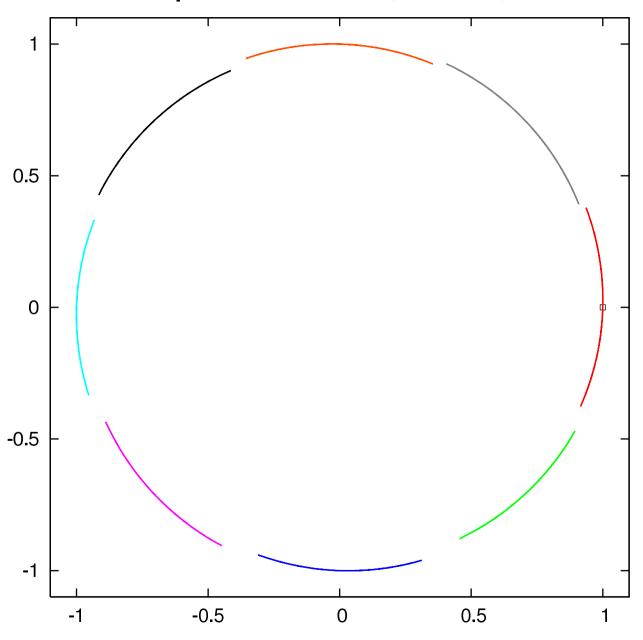
## discrete kepler. 2nd revolution, ICw 0.02, NO=13 w7



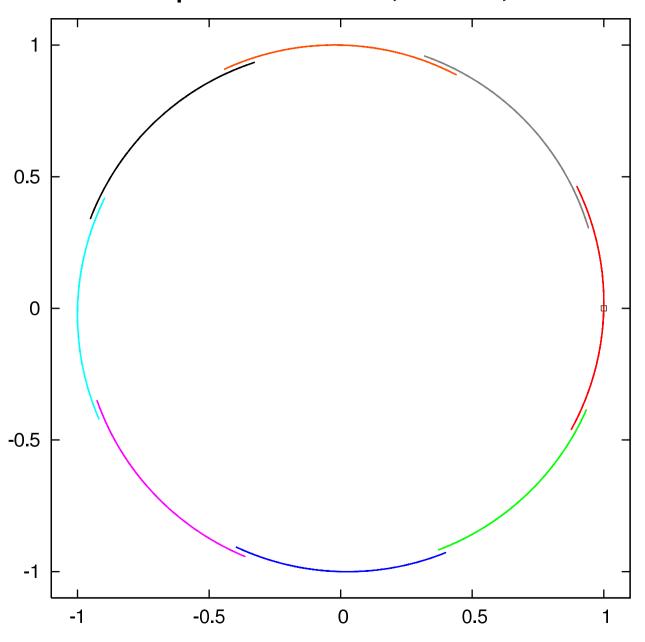
## discrete kepler. 3rd revolution, ICw 0.02, NO=13 w7



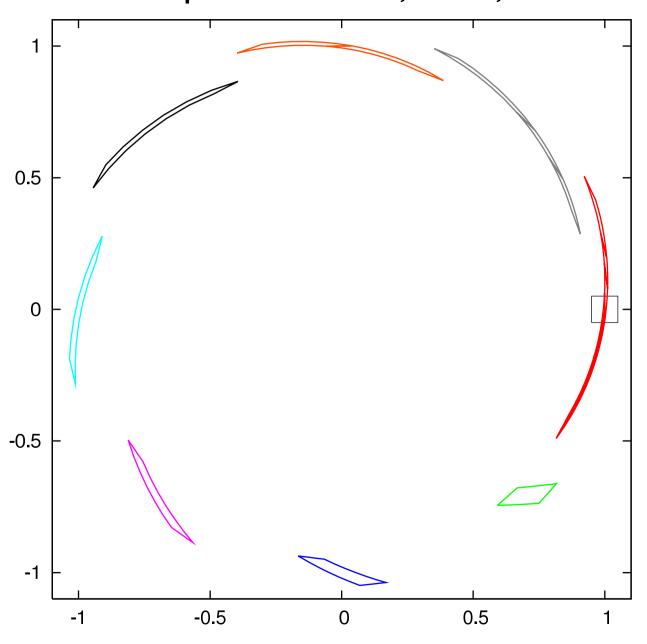
## discrete kepler. 4th revolution, ICw 0.02, NO=13 w7



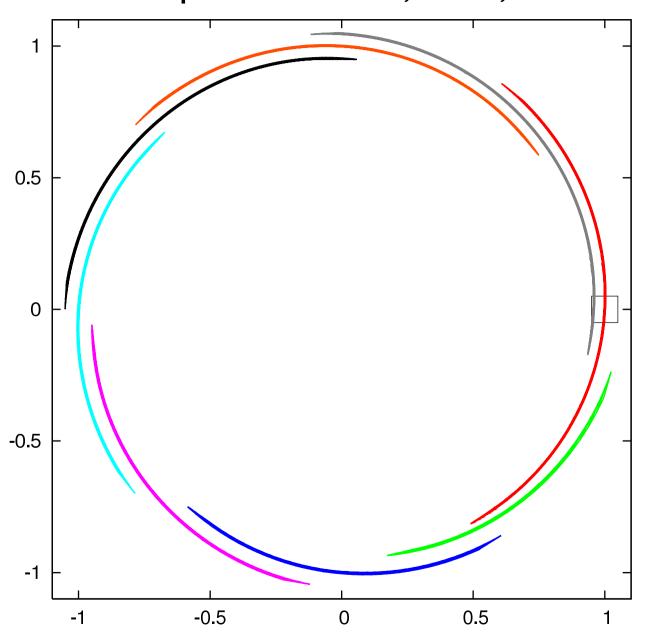
## discrete kepler. 5th revolution, ICw 0.02, NO=13 w7



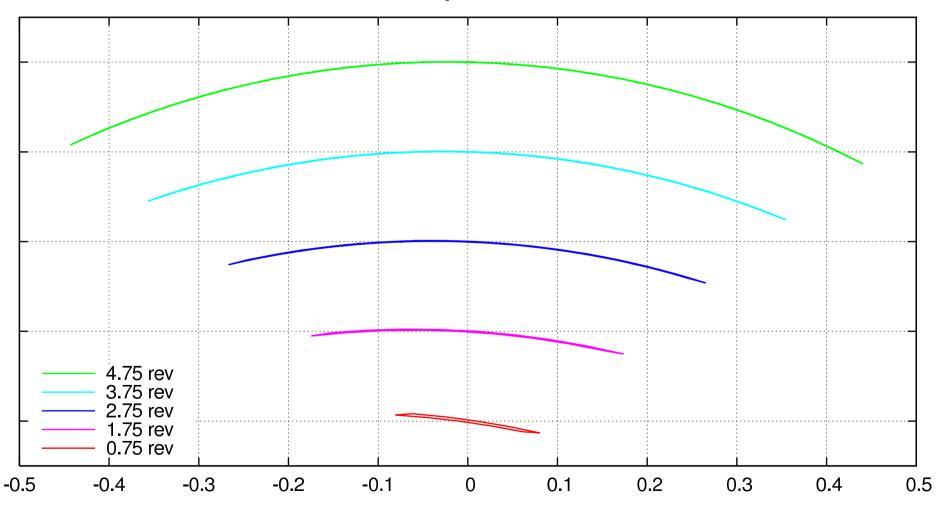
discrete kepler. 1st revolution, ICw 0.1, NO=13 w7



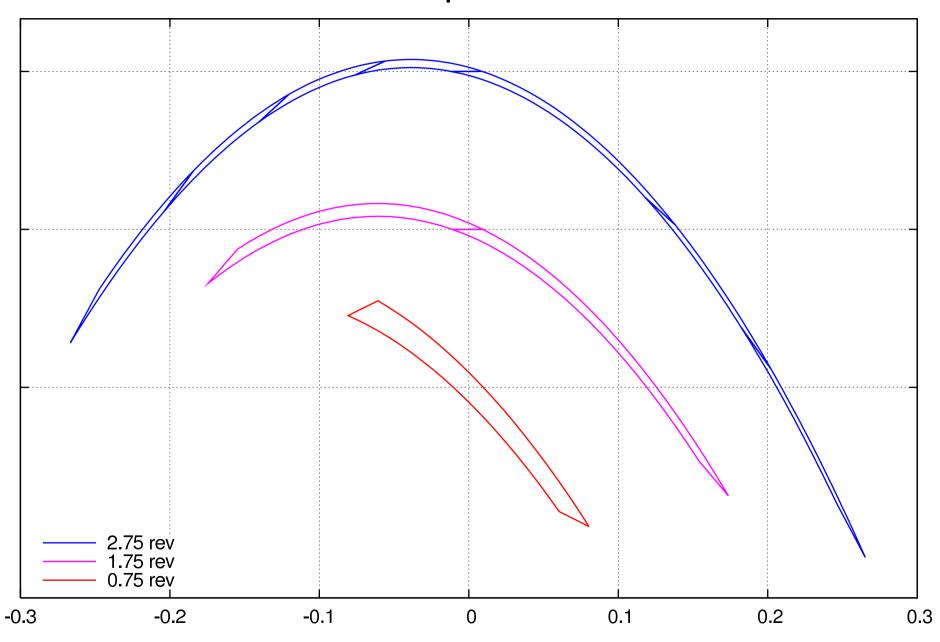
## discrete kepler. 2nd revolution, ICw 0.1, NO=13 w7



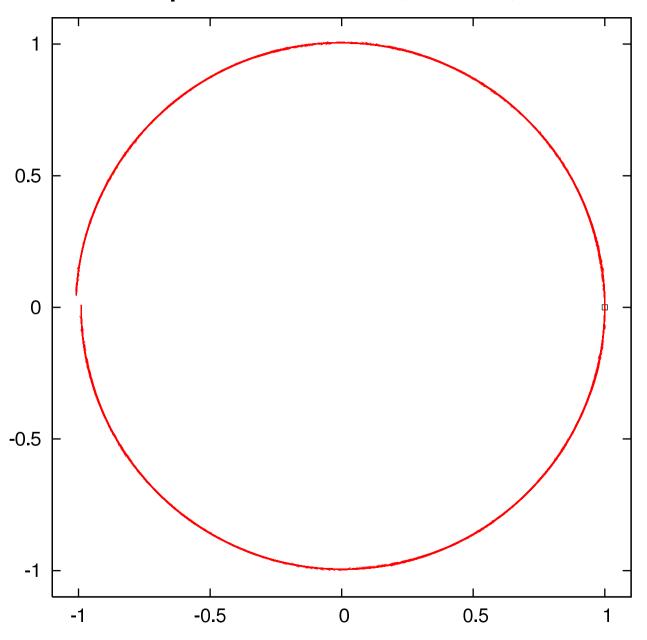
## discrete kepler. NO=13 w7



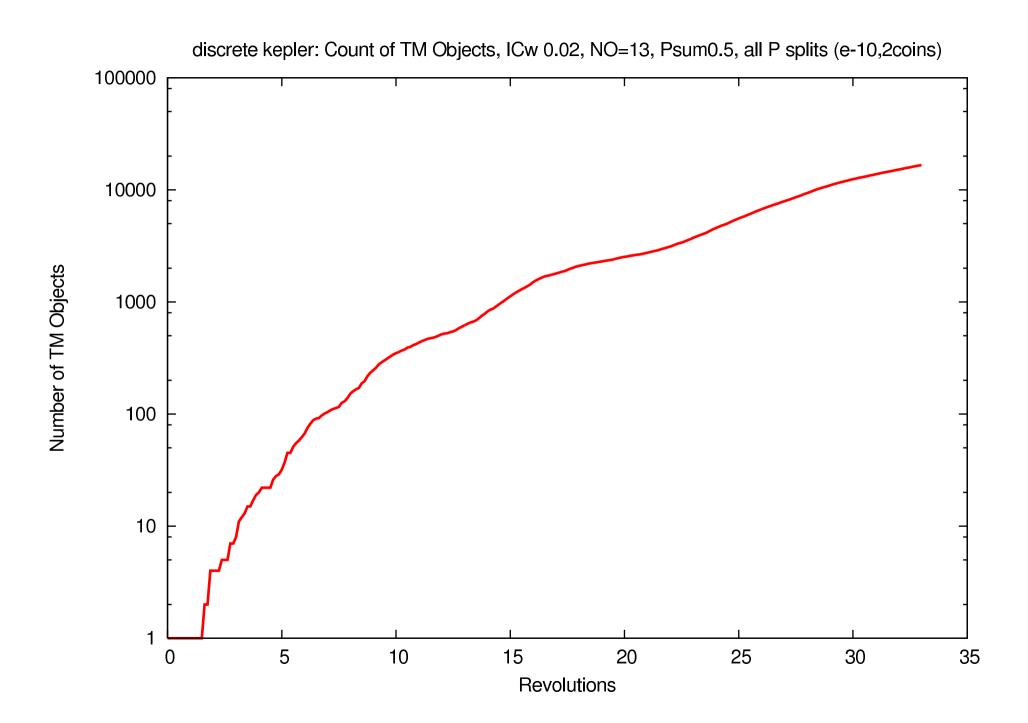
# discrete kepler. NO=13 w7



# discrete kepler. 33rd revolution, ICw 0.02, NO=13 w7



discrete kepler: Count of TM Objects, ICw 0.02, NO=13, Psum0.5, all P splits (e-10,2coins) Number of TM Objects Revolutions



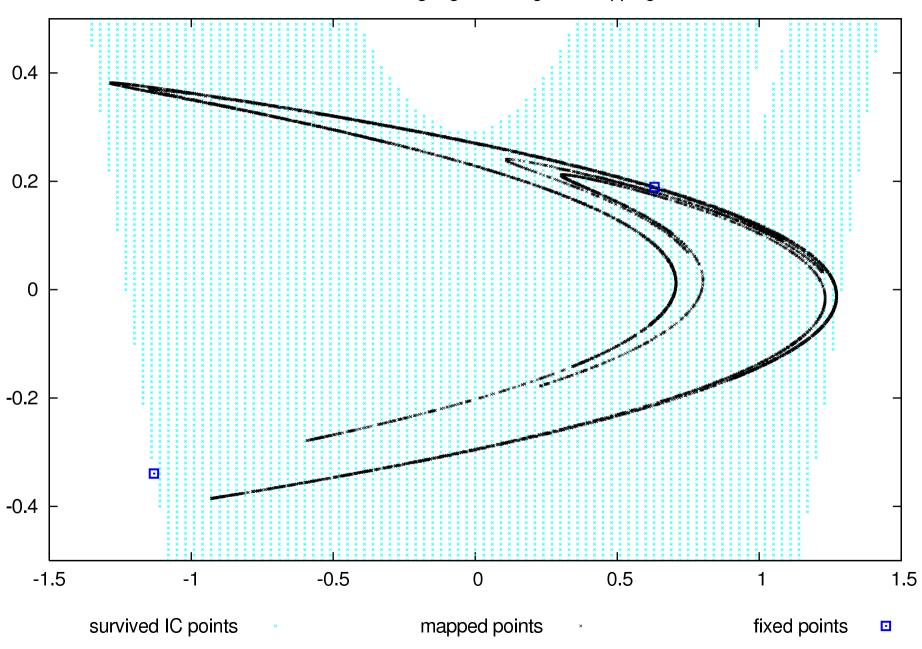
## The Henon Map

$$H(x,y) = (1 - ax^2 + y, bx).$$

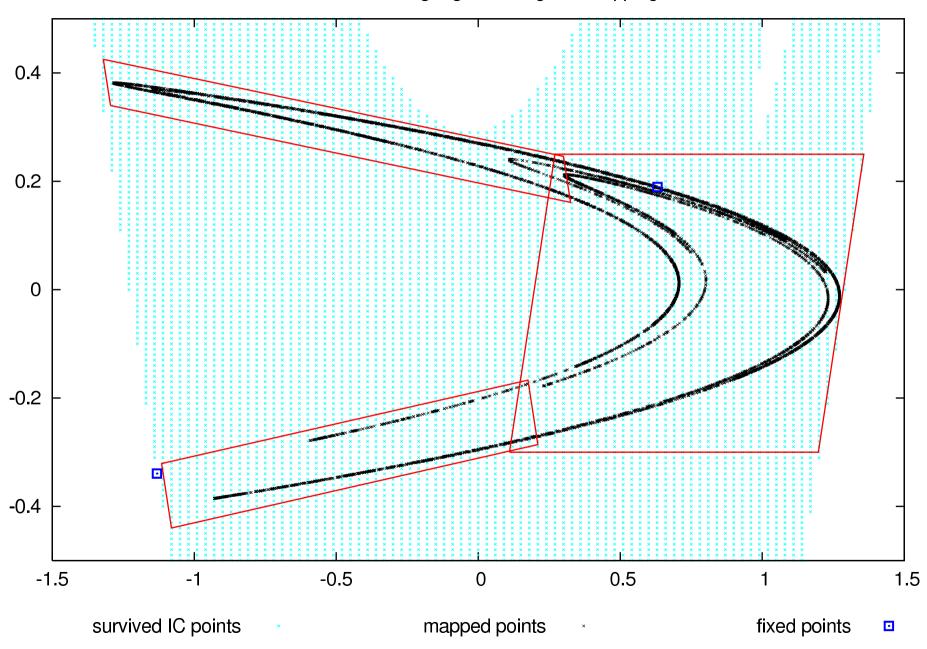
We set the parameters a=1.4 and b=0.3, which are originally considered by Henon. The map H has two fixed points.

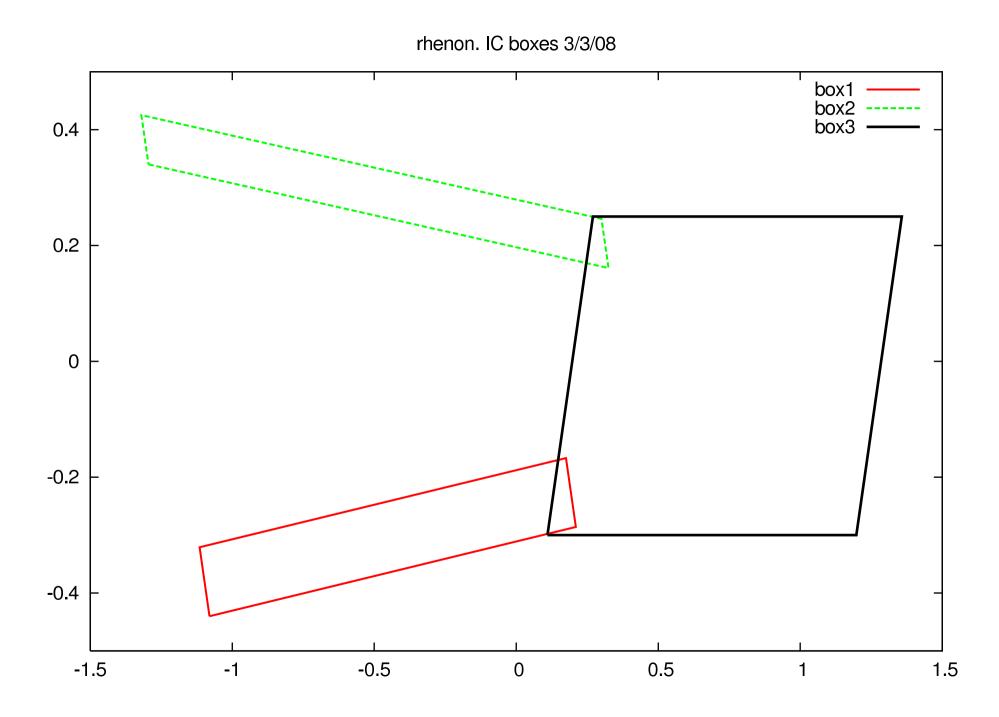
$$\vec{p}_1 = (0.63135, 0.18940)$$
 and  $\vec{p}_2 = (-1.13135, -0.33941)$ .

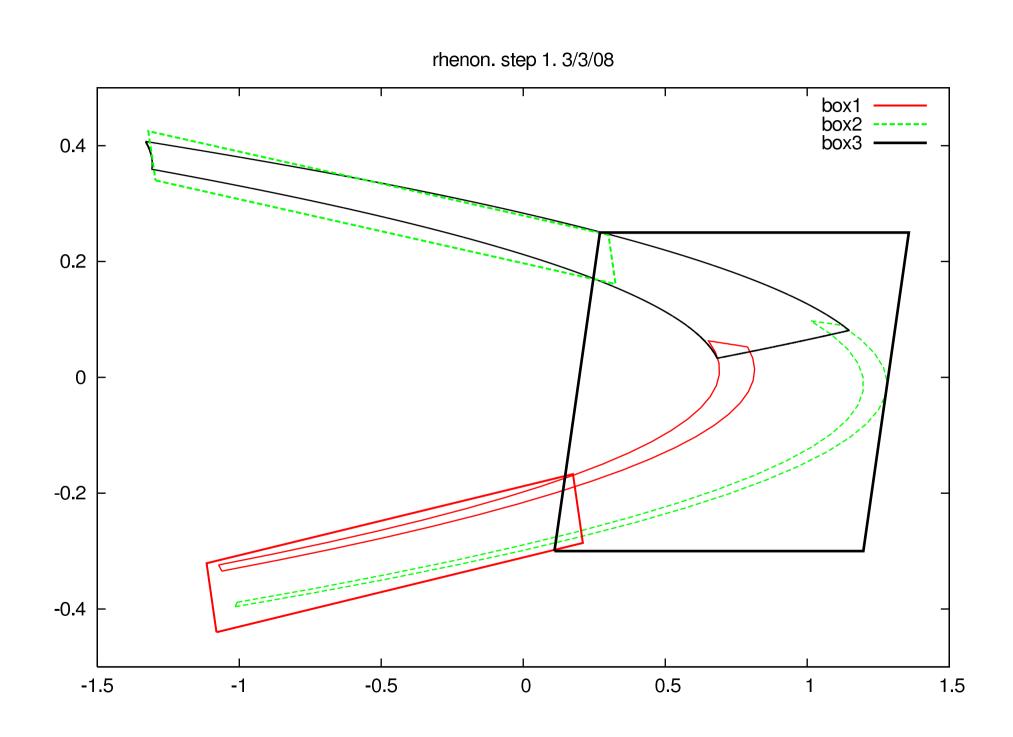
rhenon. surviving region through 12 mappings

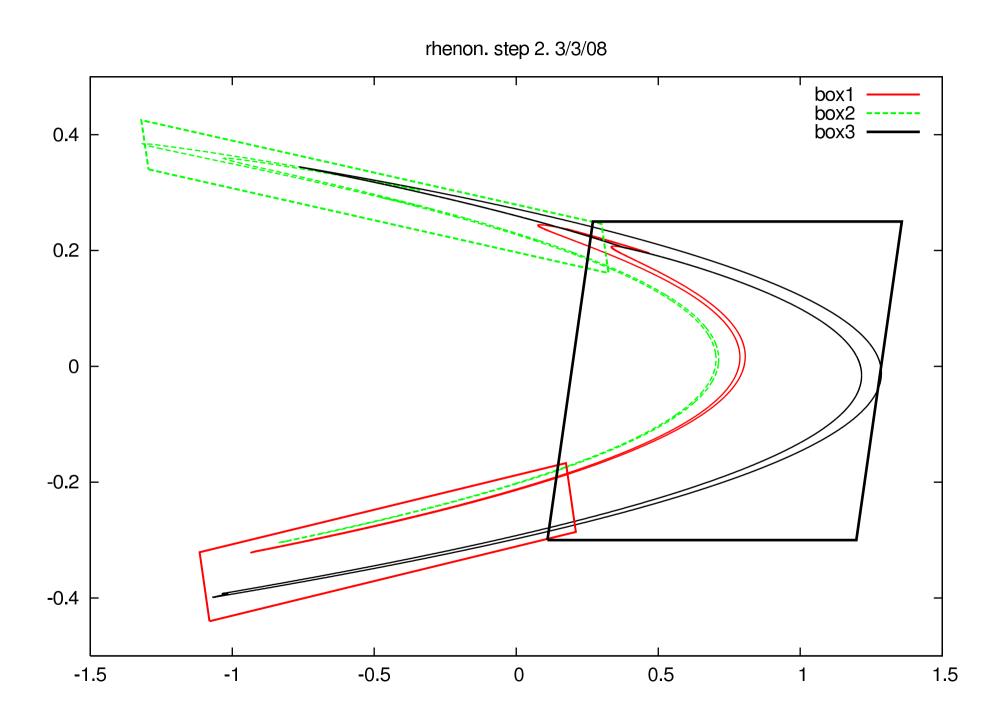


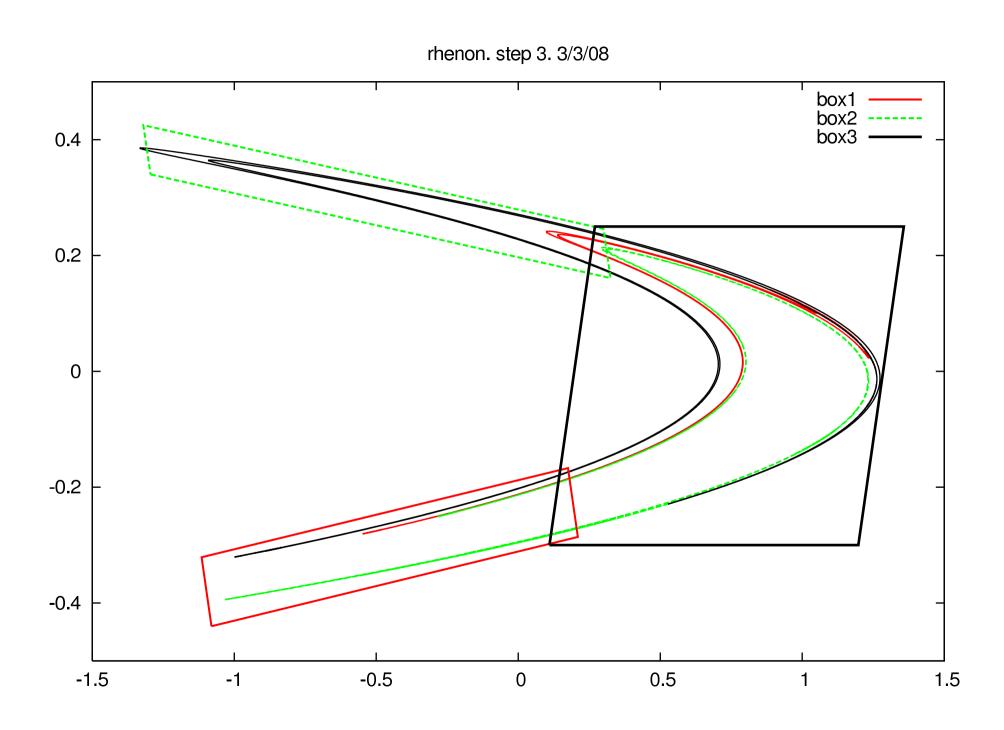
rhenon. surviving region through 12 mappings

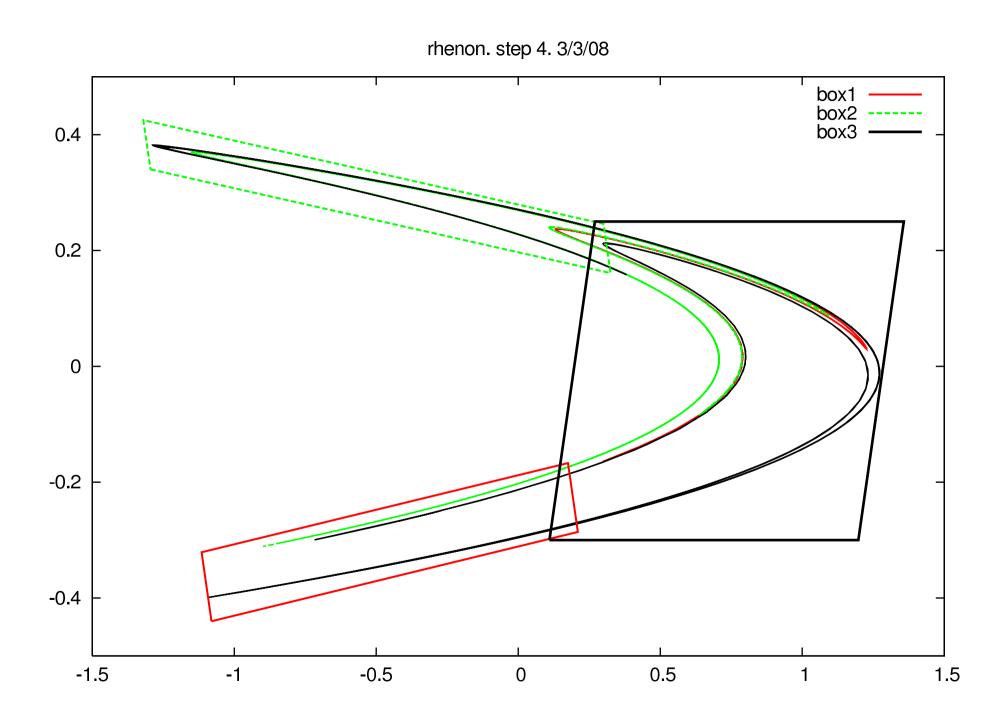








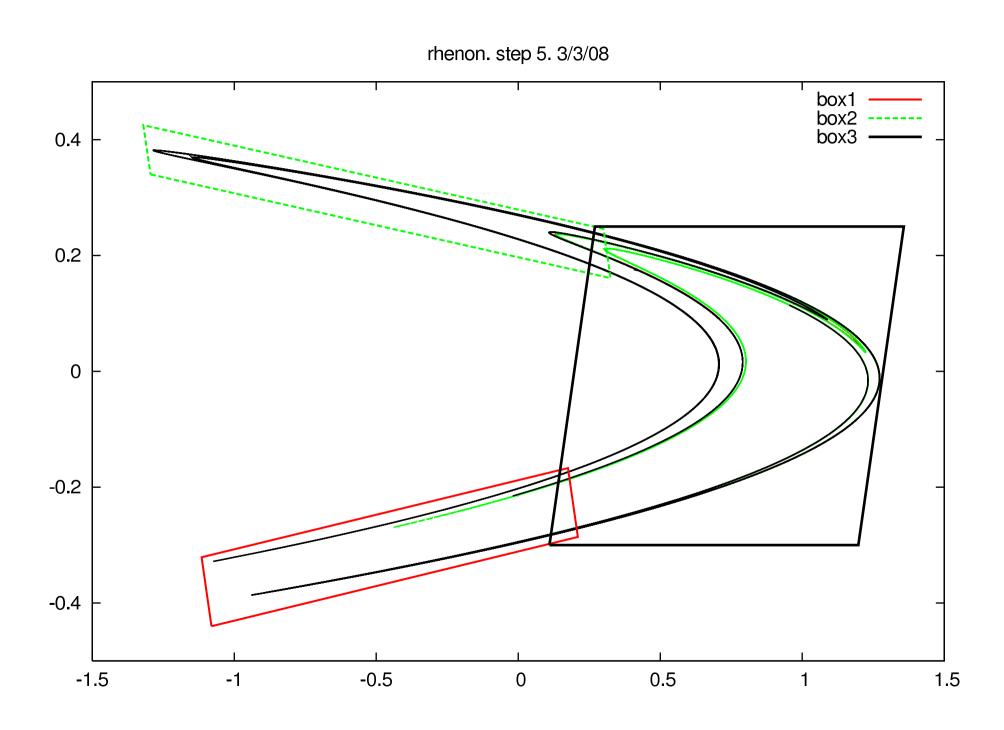




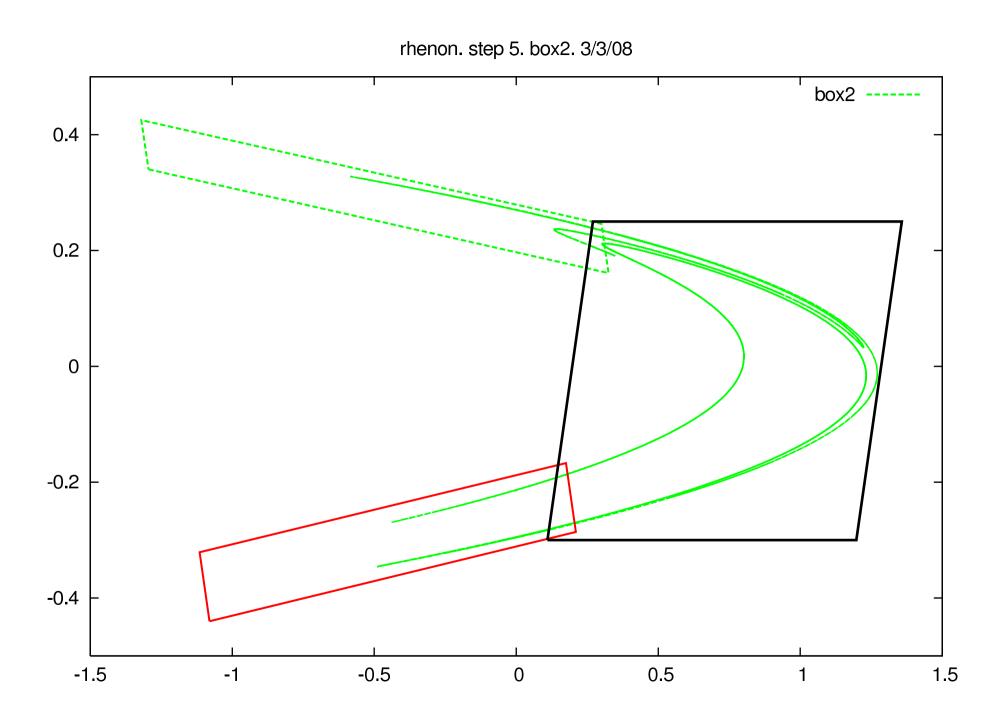
rhenon. step 4. box1. 3/3/08 box1 0.4 0.2 0 -0.2 -0.4 0.5 -1.5 -1 -0.5 1.5 0

rhenon. step 4. box2. 3/3/08 box2 -0.4 0.2 0 -0.2 -0.4 0.5 -1.5 -1 -0.5 1.5 0

rhenon. step 4. box3. 3/3/08 box3 0.4 0.2 0 -0.2 -0.4 0.5 -1.5 -1 -0.5 1.5 0



rhenon. step 5. box1. 3/3/08 box1 0.4 0.2 0 -0.2 -0.4 0.5 -1.5 -1 -0.5 1.5 0



rhenon. step 5. box3. 3/3/08 box3 0.4 0.2 0 -0.2 -0.4 0.5 -1.5 -1 -0.5 1.5 0

## rhenon: Number of Objects

To carry out multiple mappings of the Henon map, Taylor model objects underwent the domain decomposition.

Number of Taylor model objects used for multiple mappings:

|      | n  | w  | for 5 steps | for 7 steps |
|------|----|----|-------------|-------------|
| box1 | 33 | 17 | 3           | 1386        |
| box2 | 21 | 11 | 148         | 1691        |
| box3 | 33 | 17 | 8           | 2839        |

Coming very soon...

Dynamic Domain Decomposition for the ODE integrator