

MANLAB 4 theory and tutorial

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MANLAB

- ▶ An interactive path-following and bifurcation analysis software based on the Asymptotic numerical method;
- ▶ developed by the Bruno COCHELIN and Christophe VERGEZ team of Laboratoire de Mécanique et d'Acoustique (LMA) de Marseille since 2004:
<http://manlab.lma.cnrs-mrs.fr/>
- ▶ addition of stability computation by Arnaud LAZARUS (Institut d'Alembert, Paris) and Olivier THOMAS (Arts et Métiers, Lille)

MANLAB versions

- ▶ Manlab 1 (2009, R. Arquier PhD) \rightsquigarrow a continuation kernel for quadratic algebraic systems of equations;
- ▶ Manlab 2 (2010, S. Karkar PhD) \rightsquigarrow addition of Harmonic balance method for the continuation of periodic orbits;
- ▶ Manlab 2 (2010, S. Karkar PhD) \rightsquigarrow addition of Harmonic balance method for the continuation of periodic orbits + stability computation by the Hill method + Fortran acceleration;
- ▶ Manlab 3 (2011, 2014, S. Karkar) \rightsquigarrow Tensor acceleration
- ▶ **Manlab 4 (2018, L. Guillot PhD) \rightsquigarrow complete rewriting of the code with high acceleration + quasi-periodic solutions ...**

Outline

▷ MANLAB

▷ The kernel: continuation of algebraic systems

Theory

Examples

▷ Continuation of periodic solutions

Theory

Results visualizations

General framework

We consider the N -dimensional algebraic system

$$\mathbf{R}(\mathbf{U}, \lambda) = \mathbf{0},$$

with

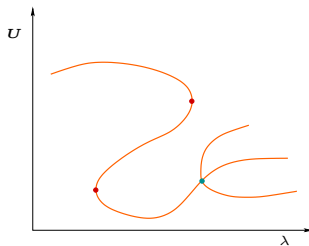
$$\mathbf{R} : \left\{ \begin{array}{ll} \mathbb{R}^N \times \mathbb{R} & \longrightarrow \mathbb{R}^N \\ (\mathbf{U}, \lambda) & \longmapsto \mathbf{R}(\mathbf{U}, \lambda) \end{array} \right.$$

The implicit functions theorem says that if \mathbf{R} is continuously differentiable with respect to \mathbf{U} and λ and if $\partial \mathbf{R} / \partial \mathbf{U}$ is invertible, there exists a continuously differentiable function \mathbf{g} such that:

$$\mathbf{g} : \left\{ \begin{array}{ll} \mathbb{R} & \longrightarrow \mathbb{R}^N \\ \lambda & \longmapsto \mathbf{g}(\lambda) = \mathbf{U} \end{array} \right.$$

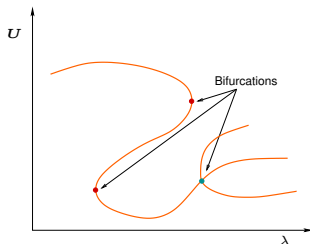
General framework

We can then wind a network of curves of U as a function of λ :



General framework

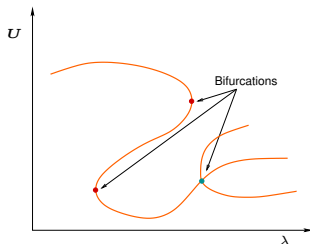
We can then wind a network of curves of U as a function of λ :



But in some **bifurcation points**, the curves can cross or have a vertical tangent with respect to λ .

General framework

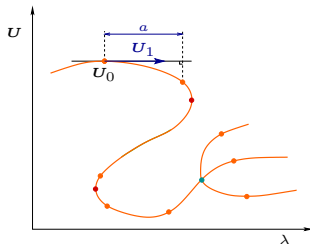
We can then wind a network of curves of U as a function of λ :



But in some **bifurcation points**, the curves can cross or have a vertical tangent with respect to λ .

A continuation (or path following) method is a numerical method that computes U for several values of λ in a given set and that manages the bifurcation points.

Path parametrization

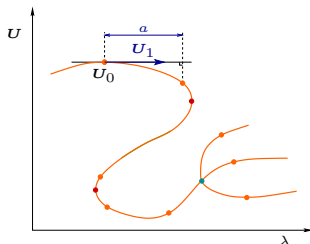


A way of dealing with the bifurcation points is to add a path parametrization. For instance, define a scalar a such that:

$$\begin{cases} U = U(a) \\ \lambda = \lambda(a) \\ f(U, \lambda, a) = 0 \end{cases} \Rightarrow \begin{cases} R(U(a), \lambda(a)) = 0 \\ f(U, \lambda, a) = 0 \end{cases}$$

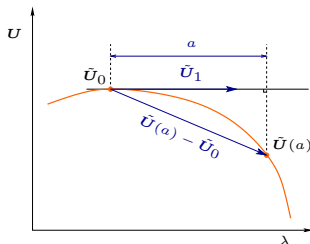
A new set of algebraic equations is defined, of size \mathbb{R}^{N+1} which is not singular as a function of a .

Path parametrization



Depending on the definition $f(U, \lambda, a) = 0$ of the path parameter a , several parametrization are available (arclength, pseudo-arclength, secant. . .) (see for instance [Seydel 2010]).

Path parametrization

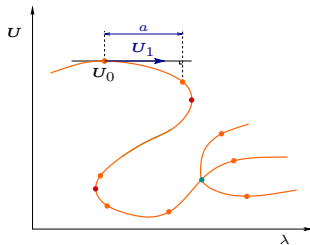


In Manlab, a pseudo-arclength parametrization is used. Let us first define $\tilde{U} = [U^t \lambda]^t$ such that $R(U, \lambda) = R(\tilde{U})$. Then, a is defined by:

$$f(U, \lambda, a) = 0 \quad \Rightarrow \quad a = \left[\tilde{U}(a) - \tilde{U}_0 \right]^t \tilde{U}_1$$

with $\tilde{U}_0 = \tilde{U}(a = 0)$ and $\tilde{U}_1 = \partial \tilde{U} / \partial a|_{a=0}$.

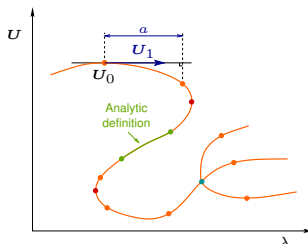
Numerical solving



▷ Newton-Raphson method

- ▶ the most used;
- ▶ enables the computation of a discrete set of solutions such that $\mathbf{R}(\mathbf{U}(a), \lambda(a)) = 0$ by sweeping the a parameter;
- ▶ one has to control the a stepping. . .

Numerical solving



► The asymptotic numerical method (ANM)

- Méthode Asymptotique Numérique (MAN) in french;
- Initiated by M. Potier-Ferry in Metz, France in the 1990';
- Power series expansions of the unknowns as a function of a : the branches are continuous function of a ;
- Automatic stepping, very few control parameters;

Asymptotic Numerical Method

- The unknowns U and λ are expanded as power series of $a \in \mathbb{R}$ of order $n \in \mathbb{N}$:

$$U(a) = U_0 + aU_1 + a^2U_2 + \dots + a^nU_n,$$

$$\lambda(a) = \lambda_0 + a\lambda_1 + a^2\lambda_2 + \dots + a^n\lambda_n.$$

- For **efficiency** of the numerical procedure, we write $R(\tilde{U})$ quadratically:

$$R(\tilde{U}) = C + L(\tilde{U}) + Q(\tilde{U}, \tilde{U})$$

where $C \in \mathbb{R}^N$, $L \in \mathbb{R}^N$ and $Q \in \mathbb{R}^N$ are constant, linear and quadratic functions of \tilde{U} .

- Introducing the above equations into $R(\tilde{U}) = 0$ **for all** $a \in \mathbb{R}$ leads to a cascade of successive problems:

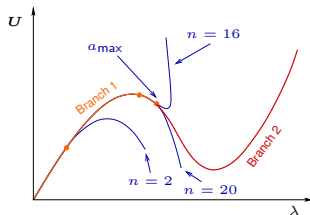
$$\text{order 0:} \quad R(\tilde{U}_0) = 0 \quad \leftarrow \text{NL initial system}$$

$$\text{order 1:} \quad J_0\tilde{U}_1 = 0 \quad \leftarrow \text{Linear system}$$

$$\text{order } p = 2, 3, \dots, n: \quad J_0\tilde{U}_p = - \sum_{i=1}^{p-1} Q(\tilde{U}_i, \tilde{U}_{p-i}) \quad \leftarrow \text{Linear system}$$

n linear systems with the same “stiffness” matrix $J_0 = \partial R / \partial \tilde{U}|_{\tilde{U}_0}$.

Asymptotic Numerical Method



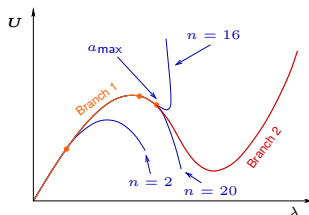
- The power series above defined have a radius of convergence such that the range of utility of a given series is defined by the value of a_{\max} such that:

$$\forall a \in [0 a_{\max}], \quad ||R(\tilde{U}(a))|| \leq \varepsilon$$

where ε is a user tolerance parameter.

- a_{\max} is automatically computed.
- The complete branch of solutions is obtained by successive power series. The $k + 1$ series is initiated such that its first point equals the last point of the previous one: $\tilde{U}_0^{k+1} = \tilde{U}(a_{\max})^k$.

Asymptotic Numerical Method



- The tolerance ε does not grow from branches to branches.
- In Manlab, the first point of the first branch \tilde{U}_0 is obtained by a Newton-Raphson method. Its accuracy conditions the one of the whole branche. Local Newton-Raphson corrections are possible with the interface.
- A bifurcation detector, special to the ANM, is also implemented in Matlab 4.

Outline

▷ MANLAB

▷ The kernel: continuation of algebraic systems

Theory

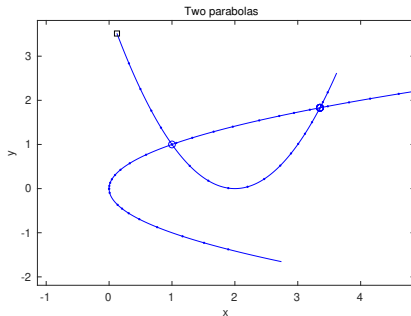
Examples

▷ Continuation of periodic solutions

Theory

Results visualizations

Two parabolas



- We consider two parabolas:

$$\begin{cases} x = y^2 \\ y = (x - 2)^2 + a; \end{cases}$$

where (x, y) are two variables and a a scalar parameter.

- Those two parabolas are solution of the implicit equation:

$$R(x, y) = (x - y^2)[y - (x - 2)^2 - a] = 0$$

In Manlab

$$R(x, y) = (x - y^2)[y - (x - 2)^2 - a] = 0$$

- Add some auxiliary variables to obtain a quadratic system:

$$\begin{cases} u_1 = x - y^2 \\ u_2 = y - (x - 2)^2 - a \end{cases}$$

- Implement the following system, quadratic in the variables $U = [x \ y \ u_1 \ u_2]^t$:

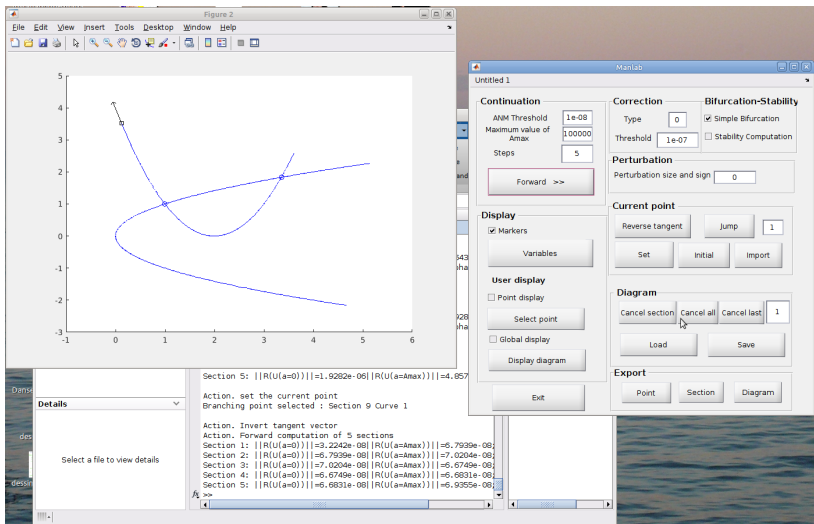
Primary system: $R_1 = u_1 u_2 = 0$

Auxiliary system: $\begin{cases} R_{1\text{aux}} = u_1 - x + y^2 = 0 \\ R_{2\text{aux}} = u_2 - y + (x - 2)^2 + a = 0 \end{cases}$

- Four .m files:

- the main file: `your_name.m`: contains all the parameters of the simulation; Its name can be chosen by the user; this file has to be executed to launch the interactive simulation;
- the equation file: `equation.m`: where the algebraic system is coded;
- two optional files for automatic display: `point_display.m` and `global_display.m`

Manlab screenshot



Outline

- ▷ **MANLAB**
- ▷ **The kernel: continuation of algebraic systems**
 - Theory
 - Examples
- ▷ **Continuation of periodic solutions**
 - Theory
 - Results visualizations

Framework

- We consider a N -dimensional first order dynamical system:

$$\dot{\mathbf{x}} = \mathbf{f}(\mathbf{x}, \lambda, t),$$

with

$$\mathbf{f} : \begin{cases} \mathbb{R}^N \times \mathbb{R} \times \mathbb{R} & \longrightarrow \mathbb{R}^N \\ (\mathbf{x}, \lambda, t) & \longmapsto \mathbf{f}(\mathbf{x}, \lambda, t) \end{cases}$$

where t is the time and λ a control parameter.

- We look for T -periodic solutions (at frequency $\omega = 2\pi/T$) of this system:

$$\forall T \in \mathbb{R} \quad \mathbf{x}(t+T) = \mathbf{x}(t).$$

- For this, we expand \mathbf{x} in Fourier series:

$$\mathbf{x}(t) = \mathbf{x}^{(0)} + \sum_{h=1}^H \left(\mathbf{x}^{(hc)} \cos \omega t + \mathbf{x}^{(hs)} \sin \omega t \right),$$

and we use the the harmonic balance method (HBM), which leads to replace the initial dynamical system by an algebraic dynamical system $\mathbf{R}(\mathbf{X}, \lambda, \omega) = \mathbf{0}$ where \mathbf{X} collects all the Fourier series coefficients $\mathbf{x}^{(0)}$, $\mathbf{x}^{(hc)}$, $\mathbf{x}^{(hs)}$ of \mathbf{x} .

In MANLAB, four cases

► Autonomous system (ex: Van der Pol)

$$\ddot{u} - \lambda(1 - u^2)\dot{u} + u = 0$$

– N primary unknowns + $\lambda + \omega \Rightarrow N + 2$ unknowns

– N primary equations + 1 phase condition $\Rightarrow N + 1$ equations

```
sys=SystHBQ(nz,nz_aux,H,@equations,@point_display,@global_display,  
            parameters,'autonomous','standard');  
sys.zi_phase = 1; % Indice of component for the phase condition
```

► Autonomous conservative system (Hamiltonian, ex: free Duffing)

$$\ddot{u} + \lambda \dot{u} + u + u^3 = 0$$

– N primary unknowns + $\lambda + \omega \Rightarrow N + 2$ unknowns

– N primary equations + phase condition $\Rightarrow N + 1$ equations

```
sys=SystHBQ(nz,nz_aux,H,@equations,@point_display,@global_display,  
            parameters,'autonomous','standard');  
sys.zi_phase = 1; % Indice of component for the phase condition
```

In MANLAB, four cases

- Forced system at Ω and Ω is the bifurcation parameter (ex: Duffing)

$$\ddot{u} + 2\mu\dot{u} + u + u^3 = F \cos \Omega t$$

- N primary unknowns + $\lambda = \Omega \Rightarrow N + 1$ unknowns
- N primary equations $\Rightarrow N$ equations

```
parameters.angfreq = 'omega';  
sys=SystHBQ(nz,nz_aux,H,@equations,@point_display,@global_display,  
parameters,'forced','standard')
```

- Forced system at fixed Ω with a bifurcation parameter (ex: Duffing)

$$\ddot{u} + 2\mu\dot{u} + u + u^3 = \lambda \cos \Omega t$$

- N primary unknowns + $\lambda \Rightarrow N + 1$ unknowns
- N primary equations $\Rightarrow N$ equations

```
parameters.angfreq = 2; % fixed value of the angular frequency  
sys=SystHBQ(nz,nz_aux,H,@equations,@point_display,@global_display,  
parameters,'forced','standard')
```


The phase condition in Manlab

- For forced systems, the time reference is given by the forcing signal, so that the phase of the periodic solution $\mathbf{x}(t)$ is referenced with respect to the phase of the forcing signal. No phase condition is needed.
- For autonomous systems, there is no reference for the phase of the periodic solution of the system and if $\mathbf{x}(t)$ is solution, $\mathbf{x}(t + \tau)$ for any $\tau < T$ is also solution. One has then to impose the phase of the seeked periodic solution to make it unique.
- In harmonic balance methods, a way is to set to zero a given harmonic of a given component of $\mathbf{x}(t)$ (for instance, the sine component of the first harmonics of the i -th. component: $x_i^{(1s)} = 0$. This is the phase condition.
- In Manlab 4.1.5, it is the time derivative at $t = 0$ of the i -th. component of $\mathbf{x}(t)$ which is set to zero:

$$\frac{d x_i}{d t}(t = 0) = 0 \quad \Rightarrow \quad \sum_{h=1}^H h x_i^{(hs)} = 0$$

- The component number i is set by `sys.zi_phase = i;` in Manlab.

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Results visualizations

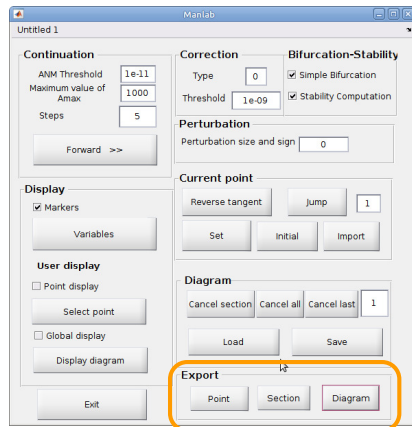
Storage of data in Manlab

- In Manlab, the basic object is a power series of \tilde{U} :

$$\tilde{U}(a) = \tilde{U}_0 + a\tilde{U}_1 + a^2\tilde{U}_2 + \dots + a^n\tilde{U}_n,$$

valid for $a \in [0, a_{\max}]$. For graphical purpose, it is computed at several points between $a = 0$ and $a = a_{\max}$. All those data (\tilde{U}_0 , \tilde{U}_1 , a_{\max} etc.) are stored in a section object. An ensemble of sections is a diagram object.

- For a given simulation, you have access to a particular section or the whole diagram by clicking on the buttons of the Export section of the Manlab interface.



Basic export

- If you click on the **Point** button, and you select a given point of the diagram, you simply obtain a vector U that contains all the components of \tilde{U} .
- If you click on the **Section** button, and you select a given section of the diagram, you obtain an object `Section` that contains all the data of the section. In particular, `Section.Upp` contains the components of \tilde{U} for all the points of the section.
- If you click on the **Diagram** button, you obtain a `cell` array of section objects. In particular, `Diagram{3}.Upp` contains the components of \tilde{U} for all the points of the 3rd. section of the Diagram.

$$U = \begin{bmatrix} U \\ \lambda \\ U_{\text{aux}} \end{bmatrix} \quad \text{Section.Upp} = \underbrace{\begin{bmatrix} \dots & U & \dots \\ \dots & \lambda & \dots \\ \dots & U_{\text{aux}} & \dots \end{bmatrix}}_{\text{number of computation points in the section}}$$

A convenient object for Diagrams

- Diagram objects can be converted to a simpler object:

`[Diag] = calcdiagUpp(sys,Diagram)`

$$\text{Diag.DiagUpp} = \underbrace{\begin{bmatrix} \dots & U & \dots \\ \dots & \lambda & \dots \\ \dots & U_{\text{aux}} & \dots \end{bmatrix}}_{\substack{\text{number of computation} \\ \text{points in the diagram}}}$$

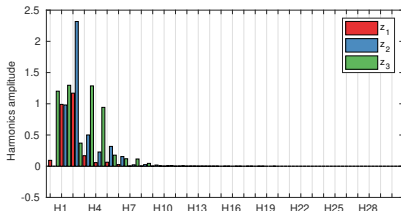
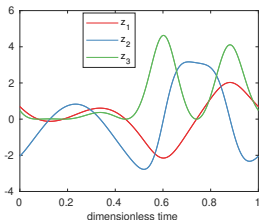
where `Diag.DiagUpp` contains \tilde{U} for all the points of the diagram in a single matrix. `Diag.Stabinfo` and `Diag.change` contains data relative to the stability of the branches.

Structure of U_{pp} for HBM simulations

For HBM simulations, U and U_{aux} contains all the Fourier series coefficients of $X(t)$ as well as ω , ω^2 and λ^2 in the following way:

$$\mathbf{X}_i = \begin{bmatrix} x_i^{(0)} \\ x_i^{(1c)} \\ \vdots \\ x_i^{(Hc)} \\ x_i^{(1s)} \\ \vdots \\ x_i^{(Hs)} \end{bmatrix} \quad \text{Diag.DiagUpp} = \underbrace{\begin{bmatrix} \dots & \mathbf{X}_1 & \dots \\ \dots & \vdots & \dots \\ \dots & \mathbf{X}_N & \dots \\ \dots & \omega & \dots \\ \dots & \lambda & \dots \\ \dots & \mathbf{X}_{1aux} & \dots \\ \dots & \vdots & \dots \\ \dots & \mathbf{X}_{Naux} & \dots \\ \dots & \omega^2 & \dots \\ \dots & \lambda^2 & \dots \end{bmatrix}}_{\text{number of computation points in the diagram}}$$

Display / computation functions for single point U



Let U be the result of a single point export.

- ▶ `Utime = calcperiodHBM(sys,U,Icalc,time)`
`plotperiodHBM(sys,U,Idisp)`
 computes / plots the periodic time evolution of several variables.
- ▶ `plotbarHBM(sys,U,Idisp)`
`plotbarsincoshBM(sys,U,Idisp)`
 plots the harmonics content of several variables as a bargraph

Display / computation functions for Sections / Diagram

Let `Diag` be a Section or a Diagram / `Diag` object.

- ▶ `plotdiagnormHBMbif(sys,Diag,Idisp,bifpara_str)`
`plotdiagHBMbif(sys,Diag,Idisp,Hdisp,bifpara_str)`
`plotdiagHBM(sys,Diag,Idisp,Hdisp,bifpara_str)`
 plots the amplitude or the L_2 norm of the variables in `Idisp` as a function of ω or λ (`bifpara_str='omega'` or `bifpara_str='lambda'`). For the first two functions, the type of bifurcations are specified by letters ('B': branch point; 'PD': period doubling; 'NS': Neimark-Sacker);
- ▶ `plotdiagYHBM(sys,Diag,Y,bifpara_str)`
`plotdiagXYHBM(sys,Diag,X,Y)`
 The same as before with y -axis (or both axis) plotted with the stability. X and Y can be obtained by
`calcdiagHBM(sys,Diag,Icalc,Hcalc)`
`calcdiagUpp(sys,Diagram)`
- ▶ Other functions are under addition, for Manlab 4.1.6...

During simulation plots

Don't hesitate to include the previous defined functions in:

- ▶ `point_display.m` → applies to a single point U
- ▶ `global_display.m` → applies to a section or the diagram.

Doing plots and post-treatments

- ▶ do your Manlab simulations. Check them during the simulation by programming `point_display.m` and `global_display.m`;
- ▶ when you know your bifurcation diagram, compute and export using the “Diagram” button of the interface. It creates a global variable called `Diagram` that contains all the results of the simulation.
- ▶ Save **all** the variables (`Diagram`, `sys` and the others).
`> save simulation.mat`
- ▶ Apply all the previous defined display functions.

