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# Nonlinear cyclic reduction for the analysis of mistuned cyclic systems

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## Abstract

Predicting the vibratory response of bladed-disks in turbomachinery is of the utmost importance to design a reliable and optimized engine. Yet, such simulations are challenging mainly due to the large size of the finite-element model used to describe the system, the existence of random mistuning coming from manufacture tolerances, and the nonlinear effects arising from the different components and their coupling. As a consequence, very few reduced-order models handling nonlinear mistuned systems have yet been proposed and the existing ones try to find the best compromise between a computationally efficient simulation and a correct description of the nonlinear phenomena. In this paper, a novel approach is proposed to tackle this challenge. Its key ideas rely on the substructuring concept and the cyclic properties of the structure. A reduction basis composed of cyclic complex nonlinear normal modes is created on each sector of the system leading to a compact nonlinear superelement per sector. The system is then assembled and a synthesis simulation is performed. Due to the main concepts employed, the method is referred to as the Substructuring method based on Nonlinear Cyclic Reduction (SNCR). It is validated on a finite element model of a tuned, a randomly mistuned and an intentionally mistuned bladed-disks. It will be shown to be fast and accurate despite strong nonlinearities. Due to its flexibility and efficiency, the method is expected to have a wide range of possible applications within the community.

*Keywords:* Mistuned Structure, Friction nonlinearities, Complex nonlinear normal modes, Nonlinear reduction, Harmonic Balance Method

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## 1. Introduction

Cyclic structures are widely encountered in engineering applications: wind turbines, power stations, turboengines, etc. Determining the forced response of such systems using their cyclic symmetry properties [1] is trivial if the structure is linear: the system of equations of motion is projected on the Fourier basis and gets decoupled on its different nodal diameters [2, 3]. However, when nonlinearities are taken into account (such as friction or geometrical effects), the system of equations obtained after projection gets once again coupled and may be too large to be easily solved. Many scientists have studied the combination of the mathematical property of cyclic structures and nonlinearities [4, 5, 6]. Based on different assumptions, some numerical strategies have been proposed to compute efficiently the nonlinear response of cyclic systems [7, 8].

However, the cyclic symmetry property is generally not applicable to real-life turboengines because they contain small manufacturing imperfections. In turbomachinery, these imperfections that are tolerated in the design stage are known as random mistuning. It has been shown in linear systems that the largest vibration amplitude of mistuned structures may vary significantly with a small fluctuation of mistuning and is, for most cases, higher than the one of its tuned counterpart [9]. Furthermore, mistuning may give rise to localization phenomena [10] in which the energy of the system is mostly concentrated within a small part of the structure. In [11], it was shown that introducing intentional mistuning allows better control of the system dynamics. Strategies to account for mistuning have been developed and many of them employ the cyclic properties of the structures in their reduction basis [12, 13, 14, 15]. These reduced-order methodologies (ROM) allow fast-computation of the system dynamics and make possible detailed parametric studies.

Combining both mistuning and nonlinear effects is the new challenging problem [16] whose main difficulty is to create a ROM that provides an accurate description of the system while being sufficiently reduced to allow parametric studies. Szpempliska proposed a method to synthesize a frequency forced response from a single nonlinear normal mode (NNM) [17]. The NNM concept was first defined by Rosenberg [18] to describe the property of nonlinear systems and to gain insight into their behavior. Different definitions of the NNM have since been proposed taking into account different mechanical features such as damping [19] and friction nonlinearities [20, 21]. A review of the computation and applicability of these NNMs was proposed by Kerschen et al. [22, 23]. In addition to the work of Szpempliska [17], nonlinear forced response synthesis has also been used, for instance in [24]. Based on this method, Joannin et al. proposed the Component Nonlinear Complex Mode Synthesis method [25, 26]. This also relies on the use of NNMs but those are calculated with fixed cyclic boundaries. The modes are exploited to create a reduction basis and to simulate the nonlinear forces by substitution. Unfortunately, it may lack accuracy when the nonlinearities strongly impact

the cyclic boundaries (see Section 5.3 of [25]). Other strategies have been proposed recently [27, 28, 29], but these require to evaluate the nonlinear force for all sectors of the system with an Alternating Frequency Time procedure [30]. This operation is numerically expensive and the methodologies seem applicable neither if one intends to perform parametric studies (such as a Monte-Carlo simulation [16]) nor if the contact interface contains many degrees of freedom (DOFs).

The purpose of the work presented in this paper is to propose a new efficient method to tackle nonlinear mistuned cyclic systems. The manuscript is constructed as follows: Section 2 shows some limitations of current strategies that can be used to model mistuned cyclic structures, and illustrates the need for a new ROM. Section 3 presents the key concept of the new approach: a cyclic synthesis based on the cyclic complex NNMs of the structure. Section 4 then fully details the new method named "Substructuring method based on Nonlinear Cyclic Reduction" (SNCR). Finally, Section 5 provides extensive numerical validation on a finite-element model. In this paper, a bladed-disk model is used as a test case but the proposed methodology could be employed for any cyclic structures such as turbines, pumps, etc. Moreover, this paper handles contact interfaces with friction and possible separation but the methodology could also be extended to account for geometrical nonlinearities.

## 2. Strategies to handle nonlinear mistuned structures: application and limitations

This section presents two existing strategies that will be employed later in this paper to benchmark the new methodology developed in Section 4. A reference methodology is first proposed and then a ROM, called the CNCMS [25], is presented. The CNCMS is chosen here because it shares similar features with the new methodology presented in Section 4, such as the nonlinear modal synthesis. The purpose of this section is also to show some limitations of these possible strategies and the need to develop a more robust and less time-consuming method.

### 2.1. Common step: linear Craig-Bampton component mode synthesis

Turboengineers are nowadays designing intentionally mistuned (large level of mistuning) engines in order to overcome the difficulties associated with random mistuning and especially the energy localization on specific sectors of the cyclic system [10, 11]. An intentionally mistuned system is composed of  $M$  different kinds of sector (noted  $(S_i)_{i \in \llbracket 1, M \rrbracket}$ ). Each of these sectors  $S_i$  occurs  $m_i$  times in the structure following a

specific intentional mistuning pattern. The sectors arrangement verifies

$$\sum_{i=1}^M m_i = N, \quad (1)$$

55 where  $N$  is the total number of sectors (an example of intentional mistuning pattern is provided in Section 5.2.2). A sector can be composed of different parts such as a blade, a portion of a disk, underplatform dampers, shrouds,... Components Mode Synthesis (CMS) methods [31] are widely used to reduce the size of a model while retaining sufficient information for its accurate description. A Craig-Bampton (CB) CMS [32] is here performed on the "blade" and the "portion of disk" components forming each kind of sector  
60  $(S_i)_{i \in [1, M]}$ . The degrees of freedom (DOFs) that are kept as master DOFs  $\mathbf{u}_{\text{master},1}$  are: the right boundary DOFs (subscript b,R), the left boundary DOFs (subscript b,L) contained in  $\mathbf{u}_b = [\mathbf{u}_{b,R}^T, \mathbf{u}_{b,L}^T]^T$ , the nonlinear DOFs of both the blade and the corresponding portion of the disk (noted  $\mathbf{u}_{\text{nl}}$ ), and an observation DOF ( $\mathbf{u}_{\text{tip}}$ ) located at the tip of the blade. The internal modes of vibration are noted  $\boldsymbol{\eta}$  and their number is chosen sufficiently high to ensure convergence of the associated linearized reduced model (considered as the stuck  
65 bladed-disk). A standard CB reduction matrix is obtained and applied to both the blade and the portion of the disk. The resulting reduced mass  $\mathbf{M}$  and stiffness matrices  $\mathbf{K}$  of this superelement have a specific layout [33]. Associated with the blade, they can be written as:

$$\mathbf{M}_{\text{blade}} = \begin{bmatrix} \mathbf{M}_{\text{nl,nl}} & \mathbf{M}_{\text{nl,tip}} & \mathbf{M}_{\text{nl},\boldsymbol{\eta}} \\ \mathbf{M}_{\text{tip,nl}} & \mathbf{M}_{\text{tip,tip}} & \mathbf{M}_{\text{tip},\boldsymbol{\eta}} \\ \mathbf{M}_{\boldsymbol{\eta},\text{nl}} & \mathbf{M}_{\boldsymbol{\eta},\text{tip}} & \mathbf{I} \end{bmatrix} \quad \mathbf{K}_{\text{blade}} = \begin{bmatrix} \mathbf{K}_{\text{nl,nl}} & \mathbf{K}_{\text{nl,tip}} & \mathbf{0} \\ \mathbf{K}_{\text{tip,nl}} & \mathbf{K}_{\text{tip,tip}} & \mathbf{0} \\ \mathbf{0} & \mathbf{0} & \boldsymbol{\Omega}_{\boldsymbol{\eta}}^2 \end{bmatrix}, \quad (2)$$

where  $\boldsymbol{\Omega}_{\boldsymbol{\eta}}$  denotes the internal natural frequencies matrix derived from the CB procedure. The matrix  $\mathbf{I}$  represents the identity matrix while  $\mathbf{0}$  is the zero matrix. The layout is similar for the disk portion, simply  
70 replace the subscript "tip" with "b" for the boundary DOFs. This first CB procedure is performed on each different kind of sector  $S_i$  defining the intentional mistuning pattern. This reduction step (illustrated between States 0 and 1 in Figure 1) is common to all methodologies presented afterward.

In this paper, random mistuning is introduced by modifying the internal vibration frequencies  $\boldsymbol{\Omega}_{\boldsymbol{\eta}}$  (appearing in the reduced stiffness matrix, see Equation (2)) in this initial CB procedure. This step is also  
75 common to all methodologies and defines State 1b in Figure 1.

## 2.2. Reference methodology

For the associated linear system (stuck case), the different sectors are assembled along their boundaries as illustrated on the left column in Figure 1, and no further reduction is applied. This methodology is referred

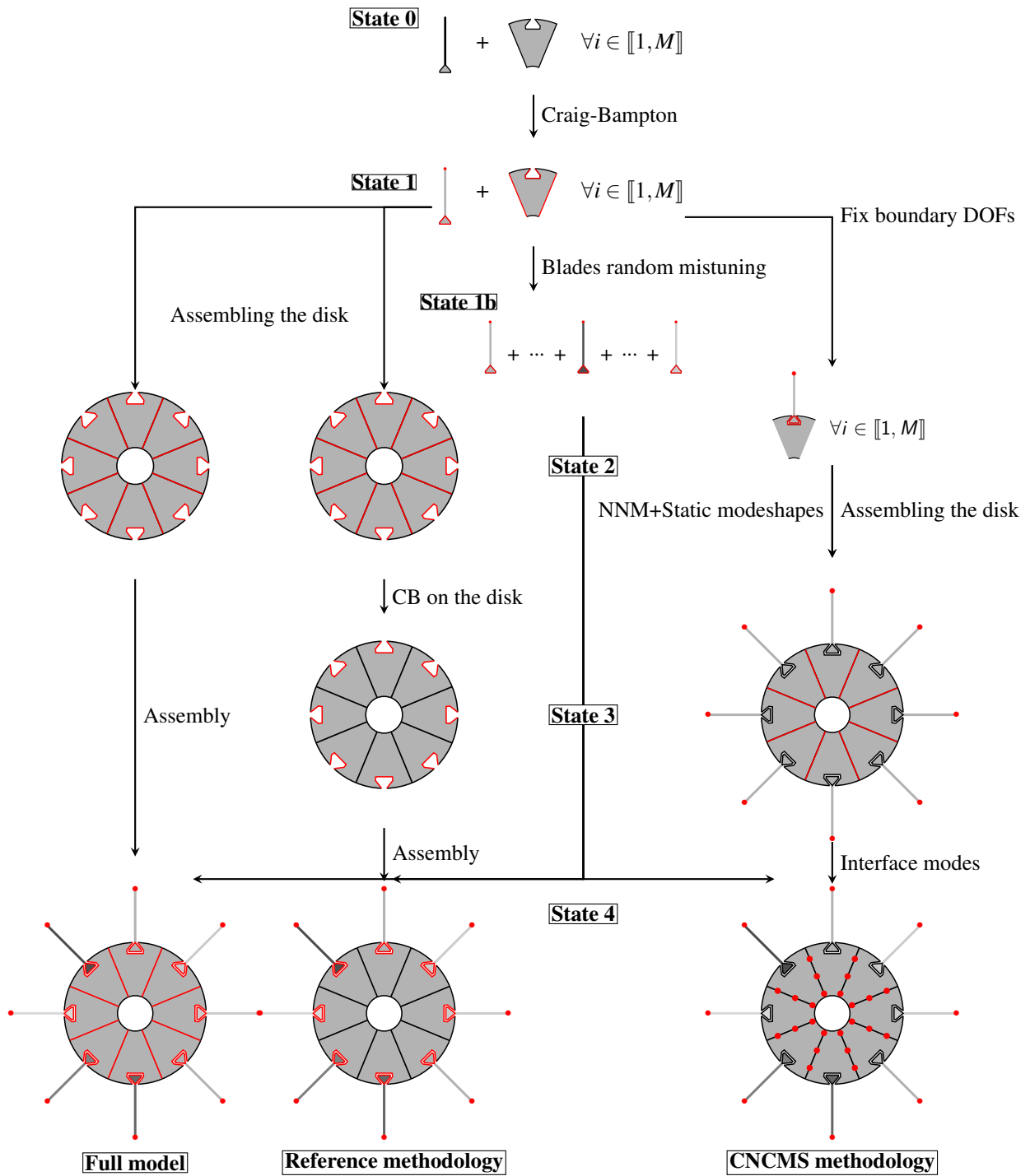


Figure 1: Description of the reference and the CNCMS methodologies.

to as the full model.

80 For a full realistic 3D model, such as the one presented in Section 2.4, using this method without further reduction is not applicable in a nonlinear analysis. A second CB reduction is therefore performed on the disk to remove the boundary DOFs. Although this reduction step creates some additional approximations, this procedure will still be referred to as the reference solution. It is illustrated in the middle column of Figure 1. Finally, the nonlinear system to solve becomes

$$\mathbf{M}_{360}\ddot{\mathbf{u}}_{360} + \mathbf{C}_{360}\dot{\mathbf{u}}_{360} + \mathbf{K}_{360}\mathbf{u}_{360} + \mathbf{f}_{\text{nl},360} = \mathbf{f}_{\text{ext},360}, \quad (3)$$

85 where  $\mathbf{M}_{360}$  is the mass matrix of the full system after the two stages of CB reduction (and similarly for the stiffness and damping matrices). A modal damping of  $5 \times 10^{-4}$  is applied to the structure. The vectors  $\mathbf{u}_{360}$ ,  $\mathbf{f}_{\text{nl},360}$  and  $\mathbf{f}_{\text{ext},360}$  represent respectively the displacement, the nonlinear forces and the external forces for the entire system. The external forces are supposed periodic in time. The force  $\mathbf{f}_{j,\text{ext}}$  applied on a particular sector  $j$  can be localized, or can follow a specific wave pattern and be defined with respect to the force  
90 applied on the first sector. This excitation pattern can then be either a traveling (forward or backward) wave pattern

$$\mathbf{f}_{j,\text{ext}} = \mathbf{f}_{1,\text{ext}} \left( t \pm \frac{2\pi h_{\text{ex}}(j-1)}{\omega N} \right), \quad j \in \llbracket 1, N \rrbracket, \quad (4)$$

or a standing wave pattern

$$\mathbf{f}_{j,\text{ext}} = \frac{1}{2} \left( \mathbf{f}_{1,\text{ext}} \left( t - \frac{2\pi h_{\text{ex}}(j-1)}{\omega N} \right) + \mathbf{f}_{1,\text{ext}} \left( t + \frac{2\pi h_{\text{ex}}(j-1)}{\omega N} \right) \right), \quad j \in \llbracket 1, N \rrbracket, \quad (5)$$

where  $h_{\text{ex}}$  is the excitation wave number,  $\omega$  is the excitation frequency that is usually the relative speed between two stages of the engines, and  $j$  is the number of sectors.

Different methodologies can be employed to solve (3), as proposed in [34]. In the case of periodic external forces, the Harmonic Balance Method (HBM) is commonly used by the community [35] as a solution procedure. In the HBM, the solution of (3) is sought after as

$$\mathbf{u}_{360} = \frac{1}{2} \left( \mathbf{c}_{360,0} + \sum_{n=1}^{N_h} \mathbf{c}_{360,n} e^{in\omega t} \right) + c.c., \quad (6)$$

where  $(\mathbf{c}_{360,n})_{n \in \llbracket 0, N_h \rrbracket}$  are the harmonic coefficients of the displacement and  $c.c.$  denotes the complex conjugate terms.  $N_h$  is the number of harmonics retained in the approximation. Substituting  $\mathbf{u}_{360}$  in (3) by (6) and

projecting the system on the exponential basis  $(e^{ni\omega t})_{n \in \llbracket 0, N_h \rrbracket}$  with the following scalar product,

$$\langle f, g \rangle = \frac{\omega}{\pi} \int_0^{\frac{2\pi}{\omega}} f(t) \bar{g}(t) dt, \quad (7)$$

where  $\bar{g}$  is the complex conjugate of  $g$ , leads to the following system

$$\mathbf{Z}_n \mathbf{c}_n + \mathbf{c}_{\text{nl},n} = \mathbf{c}_{\text{ext},n} \quad \forall n \in \llbracket 0, N_h \rrbracket. \quad (8)$$

The dynamic rigidity matrix  $\mathbf{Z}_n$  associated with the harmonic  $n$  is equal to  $(ni\omega)^2 \mathbf{M}_{360} + (ni\omega) \mathbf{C}_{360} + \mathbf{K}_{360}$ .  
 95 The projection of the nonlinear force on the exponential basis is denoted by  $\mathbf{c}_{\text{nl},n}$ . Respectively,  $\mathbf{c}_{\text{ext},n}$  refers to the excitation force. The nonlinear forces are evaluated using an AFT procedure [30]. In this paper, a Schur condensation [36] is used to describe the nonlinear phenomena (frictional contact and/or separation at interfaces) in the Dynamic Lagrangian Frequency Time procedure [37]. In the following applications, the number of unknowns thus reaches  $N \times N_{\text{r,nl}} \times (1 + 2N_h)$ , where  $N_{\text{r,nl}}$  is the relative number of nonlinear  
 100 DOFs of the FE model. Three harmonics are usually sufficient to describe accurately such systems without requiring too many computational resources [38].

### 2.3. Component Nonlinear Complex Mode Synthesis (CNCMS) methodology

The CNCMS method developed by Joannin et al. [25, 26] was mainly based on the work of Apiwat-  
 105 tanalungarn et al. [39] which assembles nonlinear superelements obtained from the invariant manifold approach, and the work of Krack et al. [24] which synthesizes a frequency forced response simulation from the computation of complex nonlinear normal modes [20]. The key concept of the method in [24] is the substitution of the nonlinear forces directly from the complex nonlinear normal modes. By this means, it avoids employing a time-consuming AFT procedure during the synthesis process.

The CNCMS combines these different methodologies. It is composed of three reduction steps. The  
 110 first step corresponds to the linear CB reduction presented in Section 2.1. Then, for each different sector  $(S_i)_{i \in \llbracket 1, M \rrbracket}$ , a nonlinear CB reduction is performed. It mainly consists in evaluating a complex nonlinear normal mode [20] with clamped boundaries and computing static modeshapes by setting successively the displacement of each boundary DOFs to unity. The unknowns of the model after this second reduction are the boundary DOFs and the control coordinates of each sector. The number of boundary DOFs is usually  
 115 large and thus a third and final reduction is employed: the interface modes [31]. This step may however be costly as the matrices associated with the boundary can be large. Finally, the unknowns of the system are the control coordinates  $(\mathbf{q}_j)_{j \in \llbracket 1, M \rrbracket}$ , which control the complex NNMs, and the interface modes control



coordinates  $\mathbf{p}$ . The equations of motion obtained are solved by the HBM as explained in Section 2.2. The first harmonic coefficients of  $\mathbf{q}_j$  and  $\mathbf{p}$  are denoted by  $\mathbf{c}_{q_j,1}$  and  $\mathbf{c}_{p,1}$ . Using a single harmonic approximation, the system to solve reduces to the algebraic system:

$$\mathbf{Z}_{r,j} \begin{bmatrix} \mathbf{c}_{q_j,1} \\ \mathbf{c}_{p,1} \end{bmatrix} + \mathbf{c}_{f_{nl,r,j}} = \mathbf{c}_{f_{ext,r,j}}, \quad (9)$$

where  $\mathbf{c}_{f_{nl,r,j}}$  is the first harmonic coefficient of the reduced nonlinear forces, and similarly for the external forces.  $\mathbf{Z}_{r,j}$  is the reduced dynamic matrix and depends on the control coordinate  $\mathbf{q}_j$ . In practice the term  $\mathbf{c}_{f_{nl,r,j}}$  is calculated by substitution via the complex nonlinear normal mode evaluated during the second stage of reduction [26, 24]. Equation (9) is then assembled along the interfaces control coordinates to account for all sectors. A semi-analytical jacobian matrix is computed, as proposed in [25], and allows fast iterations of the nonlinear solver.

#### 2.4. Description of the test case

All numerical results presented in this paper are obtained from the fundamental bladed-disk sector represented in Figure 2. The model was made with the FE-software Ansys, and its fundamental sector is described by 25920 DOFs. 180 DOFs are nonlinear (contact interfaces with friction and possible separation) and located on the dovetail attachment, and 2130 DOFs are located on the cyclic boundaries. A rotational speed of  $1000 \text{ rad s}^{-1}$  is applied and leads to static preload, centrifugal softening and geometric stiffening [33]. The gyroscopic effects are negligible for this model and are thus not accounted for. The new reduced-order methodology, presented in Section 4, could handle such effects if needed. The friction coefficient is taken equal to 0.3.

The excitation frequency range focuses on the first flexural mode. We have ensured that sufficient information were kept in the reduction steps (linear CB, complex NNM, interfaces modes) of the methodologies to accurately account for this mode. The number of internal and interface modes used are provided in Table 1.

State	Full	Reference	CNCMS
States 0-1	40 internal modes	40 internal modes	40 internal modes
States 2-3	-	600 internal modes	1 NNM with fixed boundaries
States 3-4	-	-	400 interface modes

Table 1: Components of the reduction basis for the full, the reference and the CNCMS methodologies.

140 After these reduction steps and a Schur condensation [36], the reference model contains a total of  $24 \times 90 \times 7 = 15120$  unknowns (24 sectors, 90 relative nonlinear DOFs per sector and 3 harmonics).

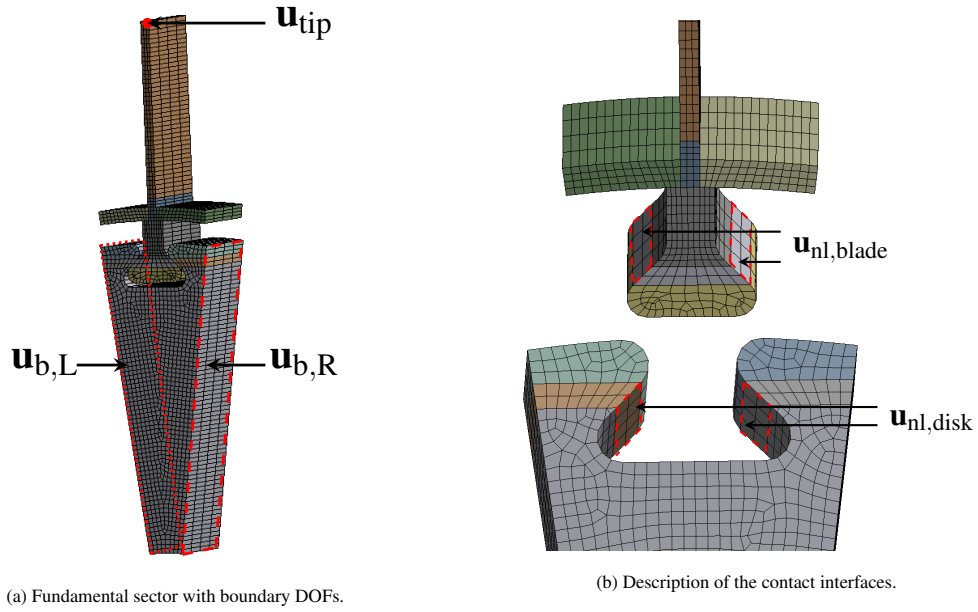


Figure 2: Description of the fundamental sector. The entire structure is made of 24 sectors.

## 2.5. Numerical simulation

An excitation of 0.25 N with a wave number equal to  $h_{ex} = 3$  is applied at the tip of the blades. The excitation follows either a traveling wave shape (see Equation (4)) or a standing wave shape (see Equation (5)).  
 145 No mistuning is considered at this point. The nonlinear solution is calculated with the reference methodology (see Section 2.2) and the CNCMS method (Section 2.3). The sectors are expected to undergo the same displacement shifted in time [7] when excited by a traveling wave, whereas a standing wave excitation

provides different levels of energy for each sector and implies different displacement amplitudes between the blades.

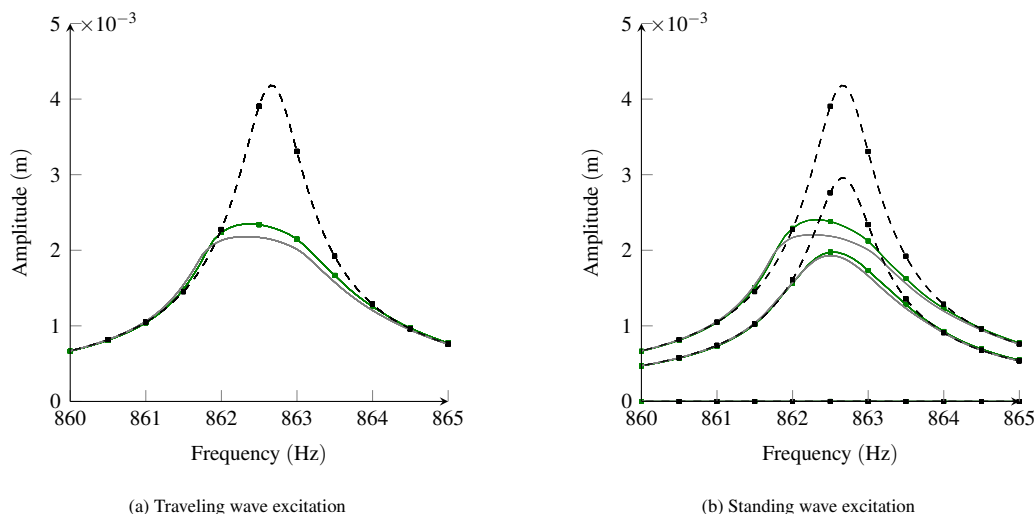


Figure 3: Frequency forced response (—■): linear response; (—■): nonlinear reference solution; (—): CNCMS solution.

150 Figure 3 illustrates the tip amplitude of all the blades. For the linear underlying system (stuck case), the full model solution, the linear reference solution, and the CNCMS method give the same results. Only the linear response obtained from the reference method is illustrated in Figure 3. As expected, the traveling wave excitation (see Figure 3a) yields an amplitude equal for all sectors. However, the response from the standing wave excitation, given in Figure 3b, shows three different levels (one of them is negligible, below  
 155  $10^{-6}$ m). For the two highest levels, the CNCMS method is not able to describe accurately the vibration amplitude and one can observe a decrease of approximately 8% at the resonant peak.

These discrepancies come from the lack of representativeness of the nonlinear CB reduction procedure in the CNCMS method. The nonlinear basis contains modes calculated with fixed boundaries and uses only linear static modeshapes for static correction. Such a basis is expected to provide good accuracy when either  
 160 the motion of the disk is small (the system mainly undergoes a blade motion mode) or when the level of nonlinearity is low and does not impact the boundary DOFs. In the case under study, the second reason is mostly responsible for the discrepancies observed.

Table 2 provides the computation time of the nonlinear simulations performed with the CNCMS and the reference methodology. In this example, the CNCMS is approximately seven times faster than the reference.  
 165 Moreover, if one is interested in performing a Monte Carlo analysis to study the impact of random mistuning

(requiring thousands of simulation), the CNCMS re-uses the pre-computed NNMs and interface modes and thus only need to relaunch the synthesis procedure.

The main objective of the paper is therefore to propose a new approach that can be more robust than the CNCMS method while still offering fast simulations.

	CNCMS			Reference
	NNM	Interface Modes	Synthesis	
Number of unknowns	630		848	15120
Computation time	10 min	11 min	40 min	7 h

Table 2: Computation times for the nonlinear simulations of the system under a traveling wave excitation. The simulations were run on a Intel(R) Core(TM) i7-7700 @ 3.6 GHz computer.

### 170 3. Nonlinear cyclic modal synthesis

The new methodology developed in this paper employs both the concept of substructuring, as explained in Section 2.1, and the concept of cyclically symmetric complex nonlinear normal modes. This latter key idea is first explained and used in a nonlinear cyclic modal synthesis to simulate the response of tuned cyclic structures.

#### 175 3.1. Cyclic complex nonlinear normal mode

One property of cyclic symmetric structures, i.e. without mistuning, is that the displacement of its  $j^{\text{th}}$  sector can be written under the Fourier form [1]

$$\mathbf{u}_j = \sum_{k=0}^{N-1} \tilde{\mathbf{u}}_k e^{-i(j-1)k\alpha}, \quad j \in \llbracket 1, N \rrbracket, \quad (10)$$

with

$$\tilde{\mathbf{u}}_{N-k} = \bar{\tilde{\mathbf{u}}}_k, \quad k \in \llbracket 1, N-1 \rrbracket \quad (11)$$

and  $\alpha = \frac{2\pi}{N}$ .  $\tilde{\mathbf{u}}_k$  are the spectral components and knowing them for  $k \in \llbracket 0, K \rrbracket$  (where  $K = \frac{N}{2}$  if  $N$  is even and  $K = \frac{N-1}{2}$  otherwise) is sufficient to recover the physical displacement  $\mathbf{u}_j$ . In practice, the components  $\tilde{\mathbf{u}}_k$  are decomposed into the internal DOFs (subscript i), and the left and right boundary DOFs (subscripts L

and R). It respects the following property [2, 40]:

$$\tilde{\mathbf{u}}_k = \begin{bmatrix} \tilde{\mathbf{u}}_{k,i} \\ \tilde{\mathbf{u}}_{k,b,L} \\ \tilde{\mathbf{u}}_{k,b,R} \end{bmatrix} = \underbrace{\begin{bmatrix} \mathbf{I} & \mathbf{0} \\ \mathbf{0} & \mathbf{I} \\ \mathbf{0} & e^{-ik\alpha\mathbf{I}} \end{bmatrix}}_{\tilde{\mathbf{B}}_k} \begin{bmatrix} \tilde{\mathbf{u}}_{k,i} \\ \tilde{\mathbf{u}}_{k,b,L} \end{bmatrix}, \quad (12)$$

where the left and right boundaries are linked by a phase change function of the nodal diameter  $k$ . By means of the matrix  $(\tilde{\mathbf{B}}_k)_{k \in \llbracket 0, K \rrbracket}$  and a projection on the Fourier basis (Equation (10)), we can rewrite the system of equations of motion for both linear [3] and nonlinear problems [7, 8]. This leads to fully decoupled equations for linear systems. However, for nonlinear systems, the different nodal diameters interact (the interacting diameters define the interaction space  $\mathcal{F}_{\text{interact}}$  [40, 8]) and the system of equations becomes,

$$\tilde{\mathbf{M}}_k \ddot{\tilde{\mathbf{u}}}_k + \tilde{\mathbf{C}}_k \dot{\tilde{\mathbf{u}}}_k + \tilde{\mathbf{K}}_k \tilde{\mathbf{u}}_k + \tilde{\mathbf{f}}_{\text{nl},k} = \tilde{\mathbf{f}}_{\text{ext},k}, \quad k \in \mathcal{F}_{\text{interact}}, \quad (13)$$

where  $\tilde{\mathbf{M}}_k$ ,  $\tilde{\mathbf{C}}_k$  and  $\tilde{\mathbf{K}}_k$  denote the mass, stiffness and damping matrices of the  $k$ -th nodal diameter. The vectors  $\tilde{\mathbf{u}}_k$ ,  $\tilde{\mathbf{f}}_{\text{nl},k}$  and  $\tilde{\mathbf{f}}_{\text{ext},k}$  represent the displacement, the nonlinear forces and the external forces on the  $k$ -th nodal diameter.  $\mathcal{F}_{\text{interact}}$  can be taken by default as the whole space  $\llbracket 0, K \rrbracket$  but it might be possible to reduce it depending on the situation [40, 8]. Solving (13) with the term  $\tilde{\mathbf{f}}_{\text{ext},k}$  allows to compute the frequency forced response of the system. The solution of the associated autonomous system (setting  $\tilde{\mathbf{f}}_{\text{ext},k} = \mathbf{0}$  in (13)) gives its cyclic nonlinear normal modes.

In this Section, we first focus on the autonomous system for a specific nodal diameter ( $k = h_{ex}$ ) in order to study one of its modes. In linear systems, the eigenfrequency of a degenerated nodal diameter ( $k \notin \{0, \frac{N}{2}\}$ ) is double and can be described by a mode pair that consists of a forward and a backward traveling modes [16]. Based on this linear description, we propose here to define the cyclic nonlinear normal modes as a sum of a forward and backward traveling modes. The computation of a forward traveling cyclic nonlinear normal modes for a particular nodal diameter  $h_{ex}$  is explained next. The computation of a backward traveling mode can be obtained similarly. Equation (13) is solved with  $\tilde{\mathbf{f}}_{\text{ext},k}$  equal to 0 (the static preload is considered but the dynamic excitation is set to 0) and by assuming that the solution can be written as a forward traveling wave [7]

$$\mathbf{u}_j(t) = \mathbf{u}_1 \left( t - \frac{\alpha(j-1)h_{ex}}{\omega} \right). \quad (14)$$

Different methodologies have been proposed to compute the nonlinear modes of a system undergoing friction nonlinearities. In [20], the dissipation introduced by friction was characterized by a parameter  $\beta$ ,

whereas it was described in [21] with an additional damping matrix. As one is here interested in synthesizing a frequency forced response based on complex nonlinear modes, it seems better-suited to parametrize the friction by a single parameter  $\beta$  instead of a full matrix. The definition of [20] is then adopted here. Using this definition, the solution on the first sector is sought after as a modified HBM form:

$$\mathbf{u}_1(t) = \frac{1}{2} \left( \mathbf{c}_0 + \sum_{n=1}^{N_h} \mathbf{c}_n e^{n(-\beta+i\omega)t} \right) + c.c. \quad (15)$$

The displacement of the remaining sectors is evaluated with Equation (14). By comparison with the traditional HBM (see Equation (6)), one can see that this expression adds an exponential term ( $e^{-\beta t}$ ) related to the dissipation effects. In the case of friction nonlinearities, the characteristic time associated with  $\beta$  is much longer than the period associated with  $\omega$ . Therefore, it may be assumed, as in [20], that  $\beta \ll \omega$ . Combining the assumption of a forward traveling wave shape (Equation (14)) with the solution (15) leads to a correspondence between the  $n$ -th harmonic and the  $k$ -th nodal diameter (governed by the number of sectors and the nodal diameter under study  $h_{ex}$ ) [7, 8]. This relationship imposes the displacement of the nodal diameter  $k$  to be governed by several harmonics (whose space will be noted  $\mathcal{N}_k$ ), whereas a given harmonic  $n$  is only present in a single nodal diameter expansion. The spectral components are hence equal to

$$\tilde{\mathbf{u}}_k = \begin{cases} \frac{\sqrt{N}}{2} \tilde{\mathbf{c}}_n e^{n\lambda_{h_{ex}}^- t} \text{ or } \frac{\sqrt{N}}{2} \mathbf{c}_n e^{n\lambda_{h_{ex}}^+ t} & \text{if } k \in \llbracket 1, M \rrbracket, n \in \mathcal{N}_k \\ \frac{\sqrt{N}}{2} \left( \mathbf{c}_n e^{n\lambda_{h_{ex}}^+ t} + \tilde{\mathbf{c}}_n e^{n\lambda_{h_{ex}}^- t} \right) & \text{if } k \in \left\{ 0, \frac{N}{2} \right\}, n \in \mathcal{N}_k \end{cases}, \quad (16)$$

where  $\lambda_{h_{ex}}^\pm = (-\beta \pm i\omega)$  and  $M = \frac{N}{2} - 1$  if  $N$  is even and  $M = \frac{N-1}{2}$  otherwise. The sign of the scalar  $\lambda_{h_{ex}}$  depends on the harmonic  $n$  which may lead to a nodal diameter aliasing.

As in the traditional HBM, the final system to solve is obtained by substituting (16) in (13) (with  $\tilde{\mathbf{f}}_{\text{ext},k} = 0$ ) and projecting the system on the exponential basis  $(e^{ni\omega t})_{n \in \llbracket 0, N_h \rrbracket}$  with the scalar product defined in (7). Since it is assumed that  $\beta \ll \omega$ , the projection of  $e^{-n\beta t}$  on the exponential basis  $(e^{ni\omega t})_{n \in \llbracket 0, N_h \rrbracket}$  is assumed to be equal to 1, and we finally have the following equations:

$$\begin{cases} \left[ \left( \lambda_{h_{ex}}^- \right)^2 \tilde{\mathbf{M}}_{h_{ex}} + \left( \lambda_{h_{ex}}^- \right) \tilde{\mathbf{C}}_{h_{ex}} + \tilde{\mathbf{K}}_{h_{ex}} \right] \frac{\sqrt{N}}{2} \tilde{\mathbf{c}}_1 + \langle \tilde{\mathbf{f}}_{\text{nl}, h_{ex}}, e^{-i\omega t} \rangle = 0 & n = 1 \\ \left[ \left( n\lambda_{h_{ex}}^\pm \right)^2 \tilde{\mathbf{M}}_k + \left( n\lambda_{h_{ex}}^\pm \right) \tilde{\mathbf{C}}_k + \tilde{\mathbf{K}}_k \right] \frac{\sqrt{N}}{2} \tilde{\mathbf{c}}_n + \langle \tilde{\mathbf{f}}_{\text{nl}, k}, e^{\pm in\omega t} \rangle = 0 & \forall n \in \llbracket 0, N_h \rrbracket \text{ and } n \neq 1 \end{cases} \quad (17a)$$

where  $\langle \tilde{\mathbf{f}}_{\text{nl}, k}, e^{\pm in\omega t} \rangle$  are the projections of the spectral nonlinear forces on the exponential basis. Details of these derivations can be found in [8] for a similar case. The nonlinear forces will be evaluated with the AFT procedure [30].

The unknowns of the system (17) are the harmonic coefficients  $(\mathbf{c}_n)_{n \in \llbracket 0, N_h \rrbracket}$ , the frequency  $\omega$  and the coefficient  $\beta$ . This system is underestimated and two additional real equations must be provided. Imposing the amplitude and phase of the first harmonic coefficient of the observation DOF (here taken as the tip of the blade) can be used to supplement the system of equation [20, 34]:

$$\Re(\bar{\mathbf{c}}_{1,\text{tip}}) = x_{\text{amp}}, \quad (18a)$$

$$\Im(\bar{\mathbf{c}}_{1,\text{tip}}) = -x_{\text{amp}} \quad (18b)$$

where  $\Re$  and  $\Im$  represent the real and imaginary parts. The term  $x_{\text{amp}} \in [0, X_f]$  is a parameter that will be gradually increased up to  $X_f$  during the evaluation of the nonlinear normal mode. The choice of the phase equation (18b) is arbitrary. Solving simultaneously (17) and (18) gives the cyclic complex NNM with traveling wave shape associated with the nodal diameter  $h_{ex}$ .

In the following applications, the pair  $(\bar{\mathbf{c}}_1, \lambda_{h_{ex}}^-)$  represents the discretized NNM (mode shape  $\bar{\mathbf{c}}_1$ , modal frequency  $\omega$  and damping coefficient  $\beta$ ) and only this information is retrieved from the solution of the system. For each level of  $x_{\text{amp}}$ ,  $\bar{\mathbf{c}}_1$  is normalized by  $\bar{\mathbf{c}}_{1,\text{tip}}$  to give the normalized nonlinear mode shape  $\Phi_{h_{ex}}^-$ :

$$\Phi_{h_{ex}}^- = \frac{\bar{\mathbf{c}}_1}{\bar{\mathbf{c}}_{1,\text{tip}}} = \frac{\bar{\mathbf{c}}_1}{x_{\text{amp}}(1-i)}, \quad x_{\text{amp}} \in [0, X_f]. \quad (19)$$

$\Phi_{h_{ex}}^-$  is numerically evaluated for discretized values of  $x_{\text{amp}}$  (a few hundred is usually sufficient to retrieve the whole NNM behavior). To describe the motion of the system for any level of amplitude, a linear interpolation of  $\Phi_{h_{ex}}^-$  is performed. The eigenvalues of the NNM,  $\lambda_{h_{ex}}^-$ , are also interpolated. In the end,  $\Phi_{h_{ex}}^-$  and  $\lambda_{h_{ex}}^-$ , are no longer discretized but are function of  $q_{\text{tip}}$  (where  $q_{\text{tip}} = |\bar{\mathbf{c}}_{1,\text{tip}}| = \sqrt{2}x_{\text{amp}}$  is the amplitude of the tip of the blade). For better readability the dependency of  $\Phi$  and  $\lambda$  on the tip displacement will be omitted. A similar development can be performed to obtain the backward traveling NNM defined with its mode shape  $\Phi_{h_{ex}}^+$  and eigenvalues  $\lambda_{h_{ex}}^+$ .

### 3.2. Cyclic synthesis

The concept of cyclic complex nonlinear normal mode explained in Section 3.1 will be largely used in Section 4 to create a reduced order model of any mistuned structure. It is first used here to simulate the forced response of a cyclic tuned structure.

Let assume that the system is excited by a forward traveling wave of frequency  $\omega$  and wave number  $h_{ex}$  ( $h_{ex} \leq K$ ). The frequency range of the excitation is supposed to be centered around a specific mode  $p$ . Based on (13), the assumption (15) and the HBM, one must solve (full details can be found in [8]):

$$\left\{ \begin{array}{l} \left[ (-i\omega)^2 \tilde{\mathbf{M}}_{h_{ex}} + (-i\omega) \tilde{\mathbf{C}}_{h_{ex}} + \tilde{\mathbf{K}}_{h_{ex}} \right] \frac{\sqrt{N}}{2} \bar{\mathbf{c}}_1 + \langle \tilde{\mathbf{f}}_{nl, h_{ex}}, e^{-i\omega t} \rangle = \langle \tilde{\mathbf{f}}_{ext, h_{ex}}, e^{-i\omega t} \rangle \quad n = 1 \quad (20a) \\ \left[ (-i\omega)^2 \tilde{\mathbf{M}}_k + (-i\omega) \tilde{\mathbf{C}}_k + \tilde{\mathbf{K}}_k \right] \frac{\sqrt{N}}{2} \bar{\mathbf{c}}_n + \langle \tilde{\mathbf{f}}_{nl, k}, e^{-i\omega t} \rangle = 0 \quad \forall n \in \llbracket 0, N_h \rrbracket \text{ and } n \neq 1. \quad (20b) \end{array} \right.$$

To solve Equation (20), we will use the strategy proposed by Krack et al. [24] (although the method was explained for a non-cyclic system, it can be easily extended to a cyclic one). The method consists in employing the cyclic complex nonlinear normal mode, solution of Equations (17) and (18) (for the nodal diameter  $h_{ex}$  and the mode  $p$ ). The Equation (20b) does not contain an excitation term and is supposed to be verified by the solution of (17b). It then remains only Equation (20a) to solve. The term  $\bar{\mathbf{c}}_1$  is substituted by  $\Phi_{h_{ex}}^- q_{h_{ex}}^-$ , where  $q_{h_{ex}}^-$  is the control coordinate of the NNM. In this specific case where only one NNM is considered we have  $q_{tip} = |q_{h_{ex}}^-|$ . Equation (20a) is then pre-multiplied by  $\Phi_{h_{ex}}^-$  and becomes

$$\left( \Phi_{h_{ex}}^- \right)^T \left( \left[ (-i\omega)^2 \tilde{\mathbf{M}}_{h_{ex}} + (-i\omega) \tilde{\mathbf{C}}_{h_{ex}} + \tilde{\mathbf{K}}_{h_{ex}} \right] \frac{\sqrt{N}}{2} \Phi_{h_{ex}}^- q_{h_{ex}}^- + \langle \tilde{\mathbf{f}}_{nl, h_{ex}}, e^{-i\omega t} \rangle = \langle \tilde{\mathbf{f}}_{ext, h_{ex}}, e^{-i\omega t} \rangle \right) = 0. \quad (21)$$

The term  $\langle \tilde{\mathbf{f}}_{nl, h_{ex}}, e^{-i\omega t} \rangle$  is substituted by the NNM (see Equation (17a)):

$$\langle \tilde{\mathbf{f}}_{nl, h_{ex}}, e^{-i\omega t} \rangle = - \left[ \left( \lambda_{h_{ex}}^- \right)^2 \tilde{\mathbf{M}}_{h_{ex}} + \left( \lambda_{h_{ex}}^- \right) \tilde{\mathbf{C}}_{h_{ex}} + \tilde{\mathbf{K}}_{h_{ex}} \right] \frac{\sqrt{N}}{2} \Phi_{h_{ex}}^- q_{h_{ex}}^-. \quad (22)$$

Finally the system to solve becomes

$$\left( \Phi_{h_{ex}}^- \right)^T \left( \left[ \left( (-i\omega)^2 - \left( \lambda_{h_{ex}}^- \right)^2 \right) \tilde{\mathbf{M}}_{h_{ex}} + \left( (-i\omega) - \left( \lambda_{h_{ex}}^- \right) \right) \tilde{\mathbf{C}}_{h_{ex}} \right] \frac{\sqrt{N}}{2} \Phi_{h_{ex}}^- q_{h_{ex}}^- - \langle \tilde{\mathbf{f}}_{ext, h_{ex}}, e^{-i\omega t} \rangle \right) = 0. \quad (23)$$

This nonlinear system of equations has only two unknowns: the real and imaginary part of  $q_{h_{ex}}^-$ , and no evaluation of the nonlinear forces is required. Therefore solving (23) is relatively fast with an iterative numerical algorithm. Although the methodology was recalled for a single mode, the basis  $\Phi$  may contain multiple modes in order to account for a wider range of excitation frequency. This method is expected to give very accurate results as long as there is no internal resonance within the modes of a given nodal diameter [24, 25]. If one is interested in a backward traveling wave excitation, the same procedure can be applied and the final equation will be similar to (23) with the quantities  $\Phi_{h_{ex}}^+$  and  $\lambda_{h_{ex}}^+$ .

Let consider a practical case to illustrate the effectiveness of this nonlinear cyclic synthesis. The tip of the blades of the structure represented in Figure 2 is excited with a forward traveling force with an amplitude of 0.1, 0.15, 0.2 or 0.25N and a wave number  $h_{ex} = 3$ . The reference solution is computed (see Section 2.2)



as well as the synthesis procedure proposed in this section. As seen and explained in Section 2.5, for a traveling wave excitation, the sectors will all undergo the same maximal amplitude shifted in time. As a consequence, the results of Figure 4 only represent the displacement of one of its sector.

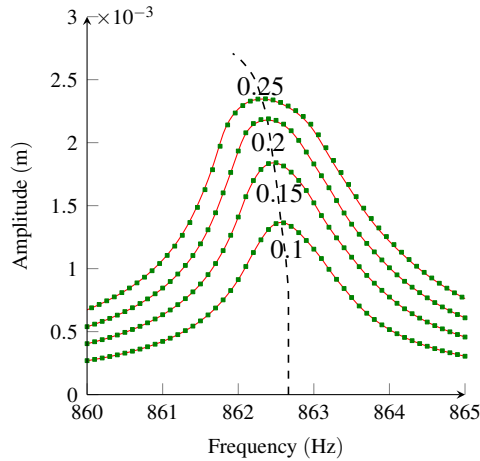


Figure 4: Frequency forced response with a forward traveling excitation of different amplitudes. (---): backbone curve of the cyclic complex nonlinear normal mode; (■): nonlinear reference solution; (—): cyclic synthesis method

Figure 4 shows that the results of the cyclic synthesis are very accurate and that the backbone curve of the NNM [23] approximately locates the peak of each frequency forced response. Furthermore, the cyclic synthesis procedure requires only a few seconds of simulation. The computation time is given in Table 3. The corresponding computation of the cyclic NNM is five times longer than the fixed NNM (see Table 2) because the boundary DOFs are kept as linear DOFs during the Schur condensation [36] in the cyclic NNM calculation.

	Cyclic synthesis		Reference
	NNM	Synthesis	
Number of unknowns	630	2	15120
Computation time	50 min	30 s	7 h

Table 3: Computation times for the cyclic synthesis simulation with a 0.25N amplitude excitation. The simulations were run on a Intel(R) Core(TM) i7-7700 @ 3.6 GHz computer.

The results of the synthesis procedure are very promising. However, as they are based on the assumption

of a traveling wave shape, they are only valid for cyclic symmetric structure undergoing traveling wave excitations. The cyclic synthesis is unable to retrieve the different levels of amplitude of the sectors under a standing wave excitation. This is illustrated in Figure 5: the cyclic synthesis only gives a single level of energy (due to the traveling wave assumption) whereas the reference solution has got three distinct energy states.

The purpose of the next section is to present a new methodology which can handle any kind of excitations (traveling, standing, localized) but can also take efficiently into account mistuned structures.

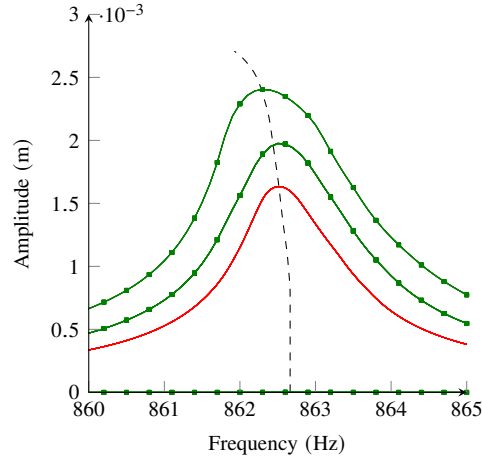


Figure 5: Frequency forced response for all sectors with a standing wave excitation (amplitude 0.25N and  $h_{ex} = 3$ ). (---): backbone curve of the cyclic complex nonlinear normal mode; (■): nonlinear reference solution; (—): cyclic synthesis method

#### 4. Substructuring method based on Nonlinear Cyclic Reduction (SNCR)

##### 4.1. Full details on the creation of the nonlinear basis

The SNCR employs as first step the linear substructuring detailed in Section 2.1. The other reduction steps are explained below and summarized in Figure 6. For each kind of sector  $(S_i)_{i \in \llbracket 1, M \rrbracket}$ , one first assumes the entire structure to be perfectly cyclic with fundamental sector  $S_i$ , and evaluates the associated cyclic matrices (State 2 in Figure 6). Both the forward and the backward traveling cyclic complex nonlinear normal modes are then computed, as explained in Section 3.1 (see State 3 in Figure 6). Since the spectral displacement can be written as the sum of the forward and backward modes over all modes [16], we assume that the spectral component of a particular nodal diameter can be written on a nonlinear modal basis as the sum of

forward and backward traveling cycling complex nonlinear modes modulated by different amplitudes:

$$\tilde{\mathbf{u}}_{k,i} = \sum_{p=1}^{N_{\text{mode}}} \left( \Phi_{k,p,i}^- (q_{\text{tip}}) q_{k,p}^- e^{-i\omega t} + \Phi_{k,p,i}^+ (q_{\text{tip}}) q_{k,p}^+ e^{i\omega t} \right), \quad i \in \llbracket 1, M \rrbracket, \quad (24)$$

where  $N_{\text{mode}}$  is the truncated number of modes kept for the analysis and  $\tilde{\mathbf{u}}_{k,i}$  represents the spectral displacement of the  $k$ -th nodal diameter for the type of sector  $i$ . The quantities  $q_{k,p}^-$  and  $q_{k,p}^+$  are the unknown generalized coordinates of the problem associated with the nodal diameter  $k$  and mode number  $p$ . Notice that for non-degenerated nodal diameters there is a single eigenvector per mode (with  $k = 0$ , all sectors move in unison, and with  $k = \frac{N}{2}$  the sectors have an alternating motion).

Let consider a sector  $j$  with the intentional mistuning type  $i \in \llbracket 1, M \rrbracket$ . Based on Equations (10) and (24), the displacement for this sector is

$$\mathbf{u}_j = \sum_{k=0}^{N-1} \tilde{\mathbf{u}}_{k,i} e^{-i(j-1)k\alpha} \quad (25a)$$

$$= \sum_{k=0}^{N-1} \left[ \sum_{p=1}^{N_{\text{mode}}} \left( \Phi_{k,p,i}^- (q_{\text{tip},j}) q_{k,p,j}^- e^{-i\omega t} + \Phi_{k,p,i}^+ (q_{\text{tip},j}) q_{k,p,j}^+ e^{i\omega t} \right) \right]. \quad (25b)$$

The term  $e^{-i(j-1)k\alpha}$  in (25a) is directly included in the amplitudes  $q_{k,p,j}^-$  and  $q_{k,p,j}^+$  in Equation (25b). Notice that the formulation in (25b) is more flexible than (10) because each sector  $j$  is here controlled by its own generalized coordinates. The displacement  $\mathbf{u}_j$  must satisfy the following equation of motion for a sector  $j$ ,

$$\mathbf{M}_j \ddot{\mathbf{u}}_j + \mathbf{C}_j \dot{\mathbf{u}}_j + \mathbf{K}_j \mathbf{u}_j + \mathbf{f}_{\text{nl},j} + \mathbf{f}_{s,j} = \mathbf{f}_{\text{ext},j}, \quad (26)$$

where  $\mathbf{M}_j$ ,  $\mathbf{C}_j$ ,  $\mathbf{K}_j$  are the mass, damping and stiffness matrices of the sector  $j$  and account for both the intentional mistuning (sector  $j$  is of type  $i \in \llbracket 1, M \rrbracket$ ) and the random mistuning. The term  $\mathbf{f}_{s,j}$  corresponds to the forces applied by the neighboring sectors on the sector  $j$  itself. The HBM with a single harmonic is applied on Equation (26) as presented in Section 2.2 and gives:

$$\mathbf{Z}_{1,j} \mathbf{c}_{1,j} + \langle \mathbf{f}_{\text{nl},j}, \mathbf{e}^{i\omega t} \rangle + \langle \mathbf{f}_{s,j}, \mathbf{e}^{i\omega t} \rangle = \langle \mathbf{f}_{\text{ext},j}, \mathbf{e}^{i\omega t} \rangle, \quad (27)$$

where  $\mathbf{Z}_{1,j}$  is the dynamic matrix associated with the first harmonic of sector  $j$  and  $\mathbf{c}_{1,j}$  is the first complex harmonic coefficient associated with  $\mathbf{u}_j$ . Based on Equation (25b), the first harmonic of  $\mathbf{u}_j$ ,  $\mathbf{c}_{1,j}$ , is written

using the nonlinear modal basis

$$\begin{aligned}
\mathbf{c}_{1,j} &= \langle \mathbf{u}_j, \mathbf{e}^{i\omega t} \rangle \\
&= \left\langle \sum_{k=0}^{N-1} \left[ \sum_{p=1}^{N_{\text{mode}}} \left( \Phi_{k,p,i}^- (q_{\text{tip},j}) \bar{q}_{k,p,j}^- e^{-i\omega t} + \Phi_{k,p,i}^+ (q_{\text{tip},j}) q_{k,p,j}^+ e^{i\omega t} \right) \right], \mathbf{e}^{i\omega t} \right\rangle \\
&= \left\langle \sum_{k=0}^K \left[ \sum_{p=1}^{N_{\text{mode}}} \left( \Phi_{k,p,i}^- (q_{\text{tip},j}) \bar{q}_{k,p,j}^- e^{-i\omega t} + \Phi_{k,p,i}^+ (q_{\text{tip},j}) q_{k,p,j}^+ e^{i\omega t} \right) \right], \mathbf{e}^{i\omega t} \right\rangle \\
&\quad + \left\langle \sum_{k=1}^M \left[ \sum_{p=1}^{N_{\text{mode}}} \left( \bar{\Phi}_{k,p,i}^- (q_{\text{tip},j}) \bar{q}_{k,p,j}^- e^{i\omega t} + \bar{\Phi}_{k,p,i}^+ (q_{\text{tip},j}) \bar{q}_{k,p,j}^+ e^{-i\omega t} \right) \right], \mathbf{e}^{i\omega t} \right\rangle \\
&= \sum_{k=0}^K \left[ \sum_{p=1}^{N_{\text{mode}}} \left( \bar{\Phi}_{k,p,i}^- (q_{\text{tip},j}) \bar{q}_{k,p,j}^- + \Phi_{k,p,i}^+ (q_{\text{tip},j}) q_{k,p,j}^+ \right) \right],
\end{aligned} \tag{28}$$

where the property (11) has been employed to simplify the initial sum (from  $K + 1$  to  $N - 1$ ) into the sum from 1 to  $M$  (with complex conjugates terms). Due to the normalization of the NNMs (see Equation (19)), the blade tip amplitude of sector  $j$  is expressed with

$$q_{\text{tip},j} = \left| \sum_{k=0}^{N-1} \left[ \sum_{p=1}^{N_{\text{mode}}} \left( \bar{q}_{k,p,j}^- + q_{k,p,j}^+ \right) \right] \right|. \tag{29}$$

In the following, the eigenvector for a given nodal diameter  $k$ , a type of sector  $i$  and a mode  $p$  that combines both forward and backward traveling modes, is noted,

$$\Phi_{k,p,i} = \left[ \bar{\Phi}_{k,p,i}^- \quad \Phi_{k,p,i}^+ \right], \tag{30}$$

where the dependency on  $q_{\text{tip},j}$  has been dropped for better readability. The complete basis is numerically created by arranging the NNMs one after the other

$$\Phi_i = \left[ \Phi_{0,1,i} \quad \dots \quad \Phi_{h_{\text{ex}},1,i} \quad \dots \quad \Phi_{K,1,i} \quad \dots \quad \Phi_{h_{\text{ex}},p,i} \quad \dots \quad \Phi_{K,N_{\text{mode}},i} \right]. \tag{31}$$

The vector of unknowns follows the same decomposition:

$$\mathbf{q}_{k,p,j} = \left[ \bar{q}_{k,p,j}^- \quad q_{k,p,j}^+ \right]^T \tag{32}$$

and

$$\mathbf{q}_j = \left[ \mathbf{q}_{0,1,j}^T \quad \dots \quad \mathbf{q}_{h_{\text{ex}},1,j}^T \quad \dots \quad \mathbf{q}_{K,1,j}^T \quad \dots \quad \mathbf{q}_{h_{\text{ex}},p,j}^T \quad \dots \quad \mathbf{q}_{K,N_{\text{mode}},j}^T \right]^T. \tag{33}$$

Equation (28) is finally written as

$$\mathbf{c}_{1,j} = \Phi(q_{\text{tip},j}) \mathbf{q}_j. \tag{34}$$

To account for the impact of the neighbouring sectors on sector  $j$  itself, one must assemble all sectors. The displacement is therefore decomposed into internal DOFs (subscript i), and the boundaries DOFs (subscript b,L and b,R for the left and right boundaries),

$$\mathbf{c}_{1,j} = \begin{bmatrix} \mathbf{c}_{1,j,i} \\ \mathbf{c}_{1,j,b,L} \\ \mathbf{c}_{1,j,b,R} \end{bmatrix} = \begin{bmatrix} \Phi_{i,i}(q_{\text{tip},j}) \\ \Phi_{i,b,L}(q_{\text{tip},j}) \\ \Phi_{i,b,R}(q_{\text{tip},j}) \end{bmatrix} \mathbf{q}_j. \quad (35)$$

This equation is then modified to control independently the internal, left and right DOFs,

$$\mathbf{c}_{1,j} = \begin{bmatrix} \mathbf{c}_{1,j,i} \\ \mathbf{c}_{1,j,b,L} \\ \mathbf{c}_{1,j,b,R} \end{bmatrix} = \underbrace{\begin{bmatrix} \Phi_{i,i}(q_{\text{tip},j}) & 0 & 0 \\ 0 & \Phi_{i,b,L}(q_{\text{tip},j}) & 0 \\ 0 & 0 & \Phi_{i,b,R}(q_{\text{tip},j}) \end{bmatrix}}_{\Phi_{i,f}} \begin{bmatrix} \mathbf{q}_j \\ \mathbf{q}_{j,b,L} \\ \mathbf{q}_{j,b,R} \end{bmatrix}. \quad (36)$$

The previous operation adds flexibility to the system with the additional unknowns  $\mathbf{q}_{j,b,L}$ ,  $\mathbf{q}_{j,b,R}$  and enables the full structure assembly. Notice that the tip amplitude of sector  $j$  is still controlled by the quantity  $\mathbf{q}_j$  through Equation (29).  $\mathbf{c}_{1,j}$  is substituted by (36) in (27) and then the system is pre-multiplied by  $\Phi_{i,f}^T$  to give

$$\mathbf{Z}_{r,j} \begin{bmatrix} \mathbf{q}_j \\ \mathbf{q}_{j,b,L} \\ \mathbf{q}_{j,b,R} \end{bmatrix} + \Phi_{i,f}^T \langle \mathbf{f}_{\text{nl},j}, \mathbf{e}^{i\omega t} \rangle + \Phi_{i,f}^T \langle \mathbf{f}_{s,j}, \mathbf{e}^{i\omega t} \rangle = \Phi_{i,f}^T \langle \mathbf{f}_{\text{ext},j}, \mathbf{e}^{i\omega t} \rangle, \quad (37)$$

The reduced dynamic rigidity matrix is equal to  $\mathbf{Z}_{r,j} = \Phi_{i,f}^T \mathbf{Z}_{1,j} \Phi_{i,f}$ . When solving (37), the different  $\mathbf{f}_{s,j}$  terms cancel each other out when assembling each sector on the  $\mathbf{q}_{j,b,L}$ ,  $\mathbf{q}_{j,b,R}$  control coordinates. As presented in Section 3.2, the nonlinear forces are substituted with the NNMs by using  $\mathbf{u}_j$  defined in Equation (25) and with the assumption that the cyclic nonlinear forces are decoupled one from another:

$$\begin{aligned} \langle \mathbf{f}_{\text{nl},j}, \mathbf{e}^{i\omega t} \rangle &= \langle \mathbf{f}_{\text{nl}}(\mathbf{u}_j), \mathbf{e}^{i\omega t} \rangle \\ &= \langle \mathbf{f}_{\text{nl}} \left( \sum_{k=0}^{N-1} \left[ \sum_{p=1}^{N_{\text{mode}}} \left( \Phi_{k,p,i}^-(q_{\text{tip},j}) q_{k,p,j}^- + \Phi_{k,p,i}^+(q_{\text{tip},j}) q_{k,p,j}^+ \right) \right] \right), \mathbf{e}^{i\omega t} \rangle \\ &= \sum_{k=0}^{N-1} \left[ \sum_{p=1}^{N_{\text{mode}}} \langle \mathbf{f}_{\text{nl}} \left( \Phi_{k,p,i}^-(q_{\text{tip},j}) q_{k,p,j}^- \right), \mathbf{e}^{i\omega t} \rangle \right] \\ &\quad + \sum_{k=0}^{N-1} \left[ \sum_{p=1}^{N_{\text{mode}}} \langle \mathbf{f}_{\text{nl}} \left( \Phi_{k,p,i}^+(q_{\text{tip},j}) q_{k,p,j}^+ \right), \mathbf{e}^{i\omega t} \rangle \right]. \end{aligned} \quad (38)$$

For each nodal diameter and mode, the term  $\langle \mathbf{f}_{\text{nl}} \left( \Phi_{k,p,i}^-(q_{\text{tip},j}) q_{k,p,j}^- \right), \mathbf{e}^{i\omega t} \rangle$  is substituted by Equation (17a)

(and similarly for the backward traveling NNM). This gives:

$$\begin{aligned} \langle \mathbf{f}_{\text{nl},j}, \mathbf{e}^{i\omega t} \rangle &= - \sum_{k=0}^K \left[ \sum_{p=1}^{N_{\text{mode}}} \left( (\bar{\lambda}_{k,p}^-)^2 \bar{\mathbf{M}}_k + (\bar{\lambda}_{k,p}^-) \bar{\mathbf{C}}_k + \bar{\mathbf{K}}_k \right) \frac{\sqrt{N}}{2} \bar{\Phi}_{k,p,i} \bar{q}_{k,p,j}^- \right] \\ &\quad - \sum_{k=0}^K \left[ \sum_{p=1}^{N_{\text{mode}}} \left( (\lambda_{k,p}^+)^2 \tilde{\mathbf{M}}_k + (\lambda_{k,p}^+) \tilde{\mathbf{C}}_k + \tilde{\mathbf{K}}_k \right) \frac{\sqrt{N}}{2} \Phi_{k,p,i}^+ q_{k,p,j}^+ \right], \end{aligned} \quad (39)$$

where the dependency of  $\Phi$  and  $\lambda$  on  $q_{\text{tip},j}$  has been omitted.

In terms of numerical implementation, one creates for each mode  $p$  and nodal diameter  $k$  the vector

$$\mathbf{z}_{\text{nl},j}^- = - \left( (\lambda_{k,p}^-)^2 \tilde{\mathbf{M}}_k + (\lambda_{k,p}^-) \tilde{\mathbf{C}}_k + \tilde{\mathbf{K}}_{\text{hex}} \right) \frac{\sqrt{N}}{2} \Phi_{k,p,i} \quad (40a)$$

$$\mathbf{z}_{\text{nl},j}^+ = - \left( (\lambda_{k,p}^+)^2 \tilde{\mathbf{M}}_k + (\lambda_{k,p}^+) \tilde{\mathbf{C}}_k + \tilde{\mathbf{K}}_k \right) \frac{\sqrt{N}}{2} \Phi_{k,p,i} \quad (40b)$$

and the different vectors are concatenated in a matrix  $\mathbf{Z}_{\text{nl},j}$  (similarly to the Equations (30) and (31)). Finally, the system to solve becomes

$$\mathbf{Z}_{\text{r},j} \begin{bmatrix} \mathbf{q}_j \\ \mathbf{q}_{j,\text{b},\text{L}} \\ \mathbf{q}_{j,\text{b},\text{R}} \end{bmatrix} + \Phi_{i,f}^T \mathbf{Z}_{\text{nl},j} \mathbf{q}_j = \Phi_{i,f}^T \langle \mathbf{f}_{\text{ext},j}, \mathbf{e}^{i\omega t} \rangle. \quad (41)$$

As the quantities  $\mathbf{Z}_{\text{r},j}$ ,  $\Phi_{i,f}^T$  and  $\mathbf{Z}_{\text{nl},j}$  depends on  $\mathbf{q}_j$  (see Equation (29)), the system (41) must be solved with a nonlinear solver. Similarly to the CNCMS method [25], a semi-analytical jacobian is provided to ensure fast computation of the solver.

Once the system is solved (the quantities  $\mathbf{q}_j$ ,  $\mathbf{q}_{j,\text{b},\text{L}}$  and  $\mathbf{q}_{j,\text{b},\text{R}}$  are known for all sectors), one can recover the displacement for all DOFs. The way is to compute the amplitude at the tip of each sector (Equation (29)), then to evaluate the mode shapes for this value (via interpolation), and finally to employ Equation (25b) to obtain the physical displacement.

#### 4.2. How to choose the nodal diameters to include in the nonlinear basis ?

The total number of unknowns of the system (41) is equal to the total number of NNMs kept in the basis multiplied by 2 (to account for the complex nature of the variables), multiplied again by 2 (for the split of unknowns introduced in Equation (36)), and finally multiplied by  $N$  (for the number of sectors). Knowing that the computation of a NNM is relatively long (see Table 3), one must keep the number of NNMs at the lowest to ensure fast computation. If the excitation frequency focuses on a specific family of mode (in our

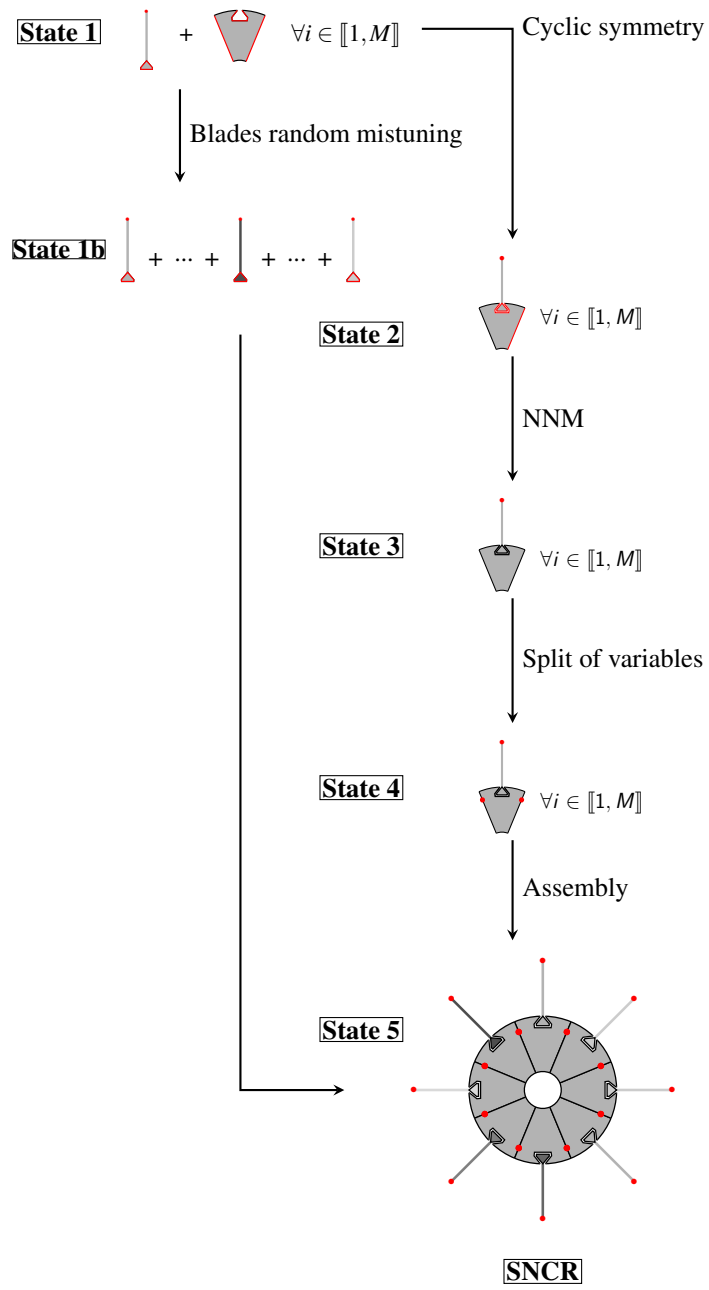


Figure 6: Schematic illustration of the SNCR.

case the first flexural mode of the different nodal diameters), the NNMs can be only computed for this family ( $N_{\text{mode}} = 1$  in (24)).

There is still to choose the nodal diameters on which one must compute the NNMs. As traveling and standing waves are the most encountered excitations in cyclic systems, the strategy adopted in this paper and  
 270 advised to potential users assumes such excitations and is explained next.

1. Retrieve the wave numbers  $h_{ex}$  of the imposed excitation forces (multiple values are possible)
2. Based on the work in [8], compute the reduced set of nodal diameters  $\mathcal{T}_{\text{interact}}$  which can get coupled by this particular excitation<sup>1</sup>.

The basis is then formed by the NNMs of these nodal diameters. Depending on the number of sectors  
 275 and the wave numbers  $h_{ex}$ , the basis may still be large. Finding a way to reduce it further is interesting as it would diminish the number of the NNMs to be calculated and reduce the final computation time of the SNCR. The following optional procedure is then proposed.

1. For each kind of sector  $i \in \llbracket 1, M \rrbracket$ , compute the reference solution for the cyclic symmetric structure of type  $i$  under a standing wave excitation with the wave numbers  $h_{ex}$ . The reference solution can either  
 280 be obtained from the full system, as presented in Section 2.2, or from the methodology presented in [8] (which was shown to be accurate for large FEM in [41]).
2. Create the basis  $\Phi$  with the  $h_{ex}$  values (solve (17) and (18) and concatenate (31))
3. Compute the frequency forced response with the SNCR using  $\Phi$  as the reduction basis (solve (41) with the system assembled)
- 285 4. Iterate between steps 2 and 3 until convergence between the SNCR and the reference solution by supplementing the basis  $\Phi$  with nodal diameters taken in  $\mathcal{T}_{\text{interact}}$ .

Once this basis is formed (following one of the procedures), one can launch simulations for both tuned and mistuned structures. One might argue that in the case of random mistuning, all nodal diameters get excited and thus the basis  $\Phi$  must contain all NNMs. However, since the SNCR parametrizes each sector  
 290 independently and that such mistuning is small, the basis is expected to be accurate enough.

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<sup>1</sup> The development presented in [8] is based on a theoretical result which determines the nodal diameters coupled by friction nonlinearities and multiple wave numbers excitations.



## 5. Numerical validation, performance and limitations

In this section, the SNCR is validated on different cases: a cyclic symmetric structure, a randomly mistuned structure and an intentionally mistuned structure. The use of various examples demonstrates the possible wide applicability of the method. In these examples, the gyroscopic effects are negligible and are thus ignored. As a consequence, the backward and forward traveling cyclic normal nonlinear modes are complex conjugates:

$$\Phi_{k,p,i}^+ = \bar{\Phi}_{k,p,i}^- \quad (42)$$

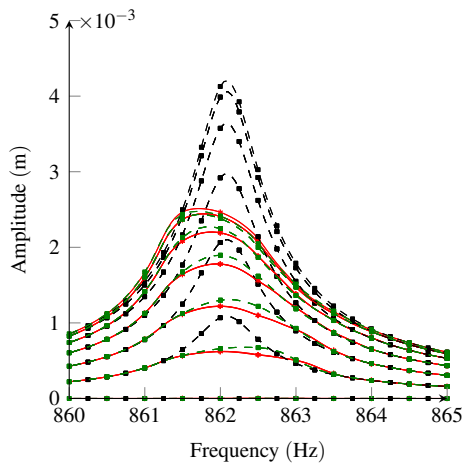
and hence only the latter are computed.

### 5.1. Numerical validation for a cyclic symmetric structure

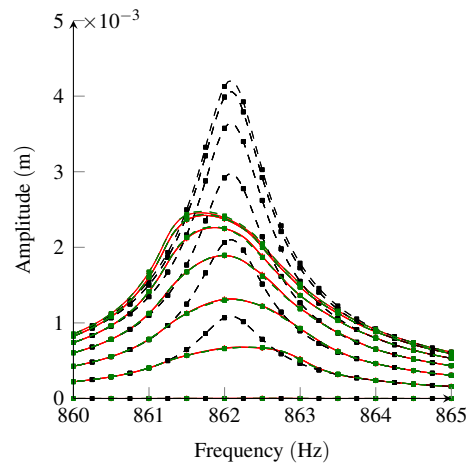
The cyclic symmetric structure is such as the one presented in Sections 2.5 and 3.2. The method has  
 295 been validated for different values of  $h_{ex}$  for both traveling and standing excitations; however, for brevity, only results associated with a standing wave excitation with two values of  $h_{ex}$  (1 and 3) are here presented. The standing wave excitation was shown to be more challenging (the sectors undergo different levels of amplitude, see Section 2.5 and more coupling occurs between the nodal diameters [8]). There is no particular difficulty in the case of a traveling wave excitation. The excitation amplitude is taken equal to 0.25 N.

300 Let first consider the case  $h_{ex} = 1$ . Based on the work in [8] (and with a structure of 24 sectors), we can predict that coupling will occur between the nodal diameters  $\{1, 3, 5, 7, 9, 11\}$  and the 0 nodal diameter excited by the static loading due to an imposed rotation. The reduction basis initially used is  $\{0, 1\}$ , composed of the zero and first nodal diameters (the numbers inside the brackets denote the nodal diameters kept for the reduction). In this section, only the first flexural mode of the different nodal diameters is studied (in  
 305 Equation (24),  $N_{mode} = 1$ ).

Figure 7 shows the amplitude of the tip of all sectors as a function of the excitation frequency, calculated for the linear and nonlinear systems, and with different methodologies. The linear response of the SNCR with basis  $\{0, 1\}$  gives the same response as the reference solution (as expected since only the first nodal diameter is excited and there is no coupling). This result is not provided in Figure 7a. The nonlinear response  
 310 of the SNCR in Figure 7a shows some discrepancies with the nonlinear reference solution. It means that the  $\{0, 1\}$  basis is not satisfactory. It is then supplemented with the third nodal diameter and is now noted  $\{0, 1, 3\}$ . Figure 7b illustrates the corresponding results. The amplitude of all sectors is now correctly obtained and shows that convergence is reached.

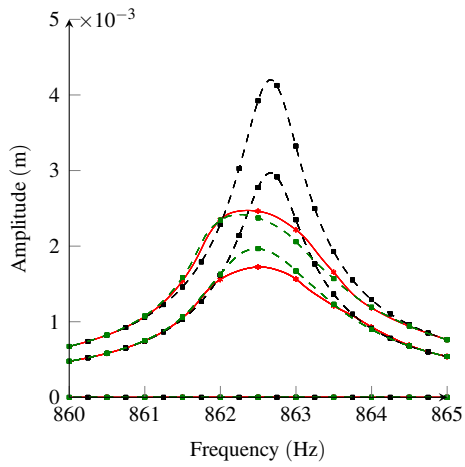


(a) Basis  $\{0, 1\}$ .

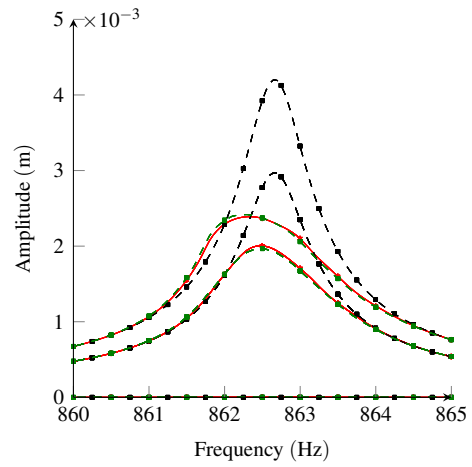


(b) Basis  $\{0, 1, 3\}$ .

Figure 7: Frequency forced responses for all sectors under a standing wave excitation and  $h_{ex} = 1$ . (—■): linear reference response; (—■): nonlinear reference solution; (—): SNCR solution.



(a) Basis  $\{0, 3\}$ .



(b) Basis  $\{0, 3, 9\}$ .

Figure 8: Frequency forced responses for all sectors under a standing wave excitation and  $h_{ex} = 3$ . (—■): linear reference response; (—■): nonlinear reference solution; (—): SNCR solution.

Similar results are found with  $h_{ex} = 3$ . Figure 8a represents the response obtained with the basis  $\{0, 3\}$ .  
 315 Adding the ninth nodal diameter [8] to form the basis  $\{0, 3, 9\}$ , the SNCR becomes able to accurately describe the response of the system, as shown in Figure 8b. Very small discrepancies are observed which may result from the substitution of the nonlinear forces (see Equation (38)) via the NNMs.

The SNCR provides very promising results as we were able to retrieve a nonlinear response under a standing wave excitation at a low cost. Moreover, the SNCR is more accurate than the CNCMS (a direct  
 320 comparison can be made between Figure 3b and Figure 8b). Details on the performance of the method will be provided in Section 5.3.

To complete the study on the tuned structure, a localized excitation is now applied on the first sector. For this specific excitation, the strategy given in Section 4.2 is no longer applicable because all nodal diameters are coupled. The basis of the SNCR was then supplemented gradually up to  $\{0, 1, 3, 5, 7\}$  (the numbers were  
 325 arbitrarily chosen to be odd) until the linear solution was considered to be accurate by comparison with the reference solution. Converged results are illustrated in Figure 9a. The nonlinear response of the system, using the same reduction basis, is presented in Figure 9b, and the differences between the reference solution and the ROM are shown to be small.

The next section focuses on the validation of the SNCR for mistuned structures.

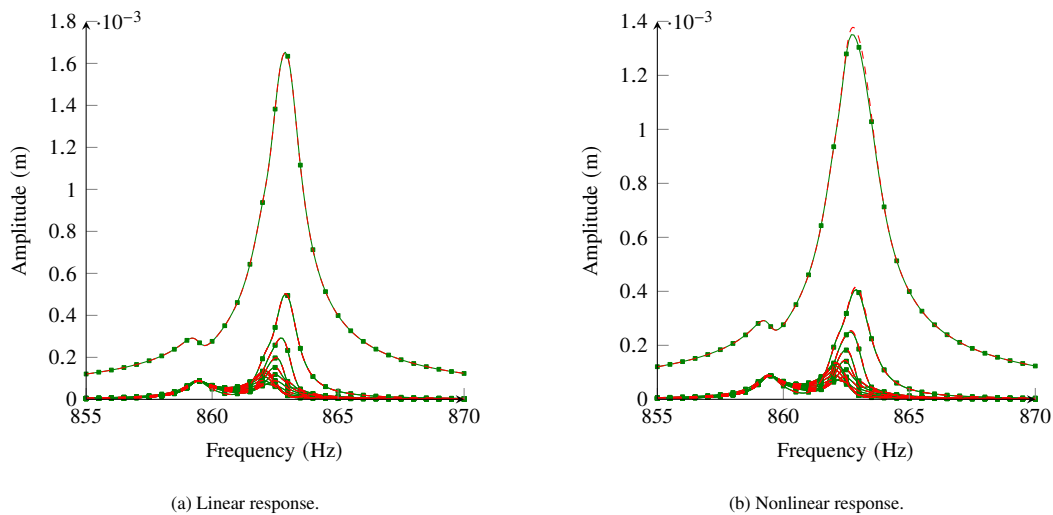


Figure 9: Frequency forced responses for all sectors under a localized excitation with an amplitude of 0.125 N. (■): reference solution; (—): SNCR solution.

## 330 5.2. Numerical validation for mistuned structures

The purpose of this section is to validate the applicability of the SNCR as a reduction method for nonlinear mistuned structures. A comparison with the CNCMS method, presented in Section 2.3, is also provided.

### 5.2.1. Random mistuning

In real engineered systems, the cyclic structure contains manufacture uncertainties [16] and exhibits a larger maximum displacement than its perfectly tuned counterpart (the ratio of both values is named the amplification factor, or AF). There exist different possibilities to take into account random mistuning in the model. For instance, in the CNCMS [26], the eigenfrequency of the fixed-NNM was shifted following a normal distribution with mean value ( $\mu$ ) equal to the natural frequency of the linear counterpart of the NNM and a variable standard deviation ( $\sigma$ ) referred to as mistuning level.

340 In this paper, the procedure to account for random mistuning modeling consists in modifying the internal natural frequencies in the initial CB procedure (see Equation (2)). For each sector, each value of  $\Omega_\eta$  is chosen randomly following a normal distribution with  $\mu$  taken as the natural frequency  $\Omega_\eta$  and a variable  $\sigma$ . This random mistuning modeling is interesting for validation and practical applications because the random matrices are common for both the reference method and the ROM. Besides, the geometry and material of the sector are not directly modified and the computation of the structural matrices only requires the original sector matrices (offering some computation gain). In most applications, as for instance in [13], only blade mistuning is considered and such practice is adopted in this paper. The choice of random mistuning modeling is method-dependent.

For both linear and nonlinear reference methodologies, the blades model is not further reduced after State 1 (see Figure 1) and the random mistuning matrices are introduced in the model assembly in State 4. The reference models retain thus their whole accuracy. In the case of the SNCR, see Figure 6, (respectively the CNCMS, see Figure 1), the blades are further reduced at step 2 with the calculation of the NNMs. Evaluating the NNMs for each new set of random mistuning would deteriorate the performance of the method. The basis  $\Phi$  is therefore assumed not to be impacted by random mistuning. Such approximation is reasonable since the mistuning is small. The modification on the stiffness matrices of the blades is nevertheless still included in the synthesis stage (Equation (41) for the SNCR and (9) for the CNCMS). As a consequence, the differences in the results between the reference method and the SNCR may be due to two approximations: the small impact of the mistuning on the NNMs that is not accounted for, and the reduction methodology itself (nonlinear basis  $\Phi$ , substitution of the nonlinear forces, etc.).

In the following numerical simulations, a traveling wave excitation with an amplitude of 0.25 N and  $h_{ex} = 3$  is applied at the tip of the blades. The  $\{0, 3, 9\}$  reduction basis is taken here as suggested in Section 4.2. Similar results are found for different values of  $h_{ex}$ .

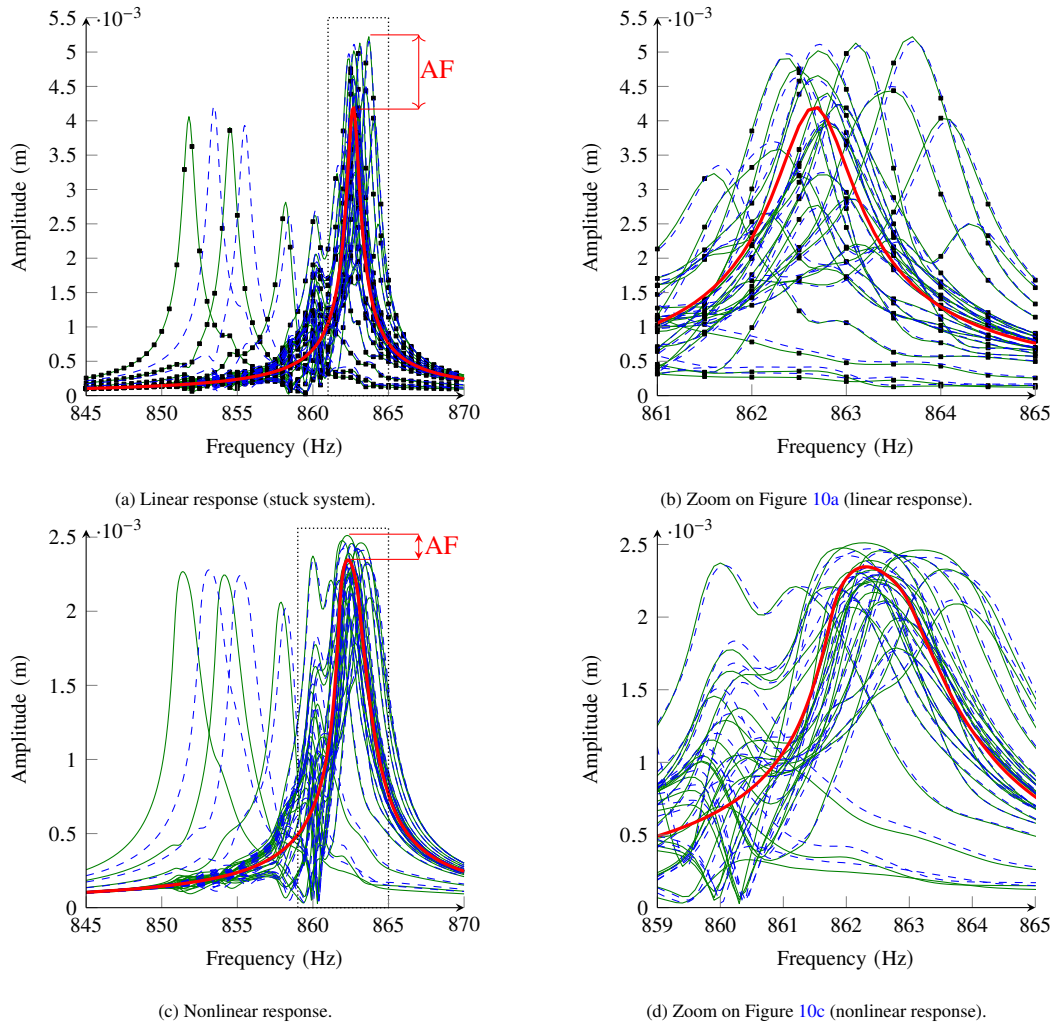


Figure 10: Frequency forced responses for all sectors with random mistuning ( $\sigma = 15\%$ ). (■): full model solution; (—): linear reference solution, (- -): SNCR solution, (—): tuned response.

Figure 10a and 10b represent the linear solution (stuck system) obtained with the full model (left column in Figure 1), the reference model (middle column in Figure 1) and the SNCR solution for a relatively large random mistuning pattern (equal to  $\sigma = 15\%$ ). The reference solution matches perfectly the full model

calculation, and it proves that the 600 internal modes kept in the nonlinear reference solution are sufficient to capture accurately the motion of the disk. For the SNCR solution, we observe that the amplitudes of the displacements are close to the reference ones although a frequency shift occurs at lower frequencies (shift of 2 Hz around 850 Hz, which corresponds to a 0.25% absolute relative error). This small discrepancy is due  
370 to the reduction procedure of the SNCR which cannot capture perfectly the flexibility of the disk. These residual flexibilities can be corrected by taking a larger base (as it was performed in Section 5.1). However, the purpose is here to show that the general behavior is correctly captured even with a very small reduced basis. It can also be noticed in Figure 10a and 10b that most of the sectors reach their maxima at higher frequencies (around 865 Hz) and the results of the SNCR are accurate in this frequency range.

375 Figure 10c and 10d represent the nonlinear response obtained with the different methodologies. Similarly to the linear case, a slight shift of frequency is observed around 850 Hz. However, the maximal displacement of the system (and thus the amplification factor) around 862 Hz is accurately captured. This ratio is of the utmost importance for the designers and we observe that the proposed ROM manages to evaluate it correctly. For this specific mistuning pattern, the SNCR gives very satisfactory results. Additional simulations are  
380 provided next to fully validate the method for different  $\sigma$  values.

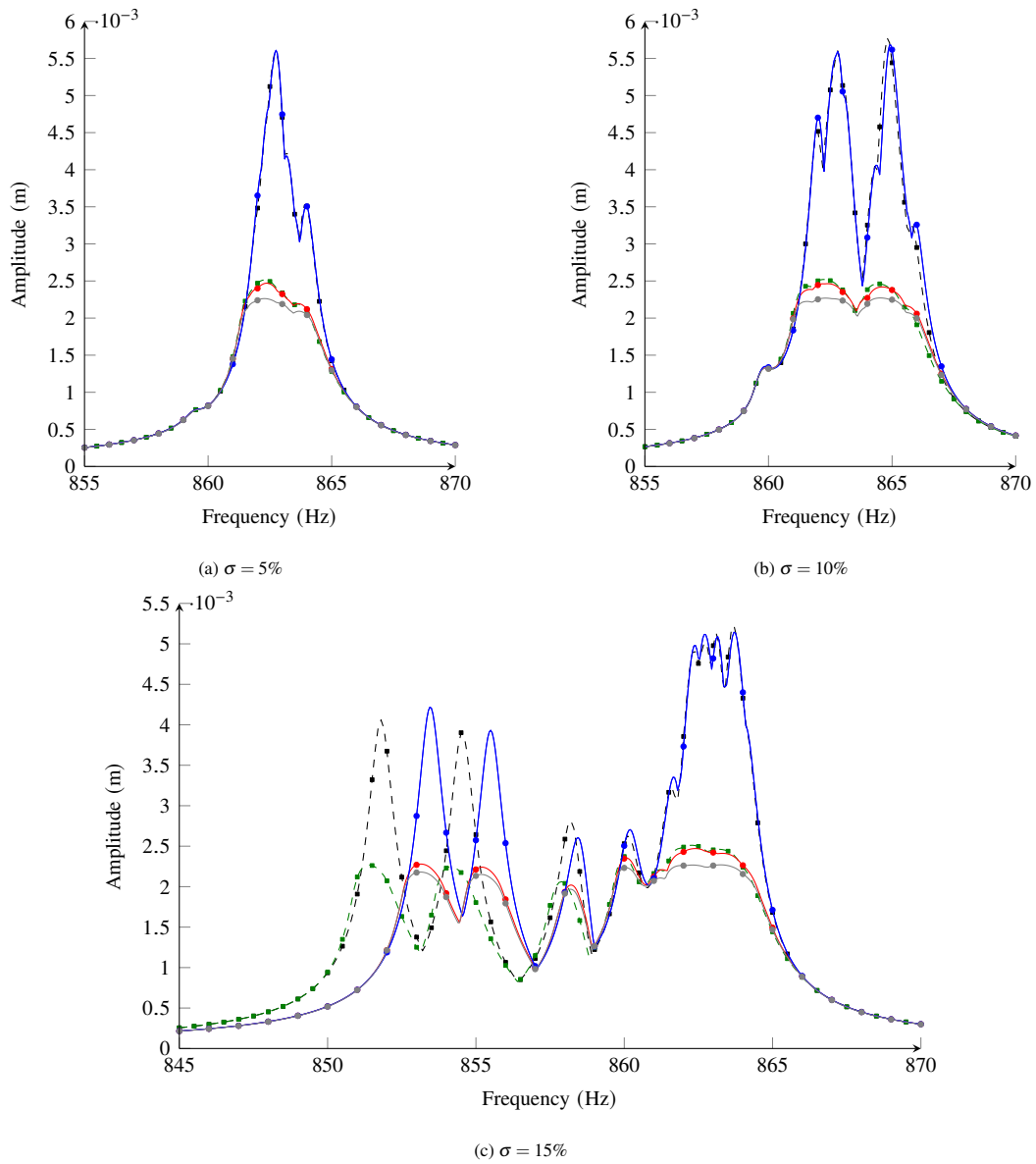


Figure 11: Frequency forced responses for different values of  $\sigma$ . (-■-): linear response for the reference solution; (-■-): nonlinear reference solution; (-●-): CNCMS solution; (-●-): linear SNCR response; (-●-): nonlinear SNCR solution.

Figure 11 represents the maximum displacement of the entire structure as a function of the excitation frequency obtained with random mistuning defined with a standard deviation equal to 5%, 10% and 15%. Once again the SNCR provides a maximum amplitude very close to the reference solution for both the

linear and nonlinear cases. For better readability, these results only represent the maximum response of the entire structure as a function of the excitation frequency (and not each blade displacement separately). The main results of Figure 10 are thus summarized in Figure 11c in which the CNCMS results are also given for comparison. Similarly to what was shown in Figure 3, the CNCMS is not able to retrieve the correct maximal displacements.

### 5.2.2. Intentional mistuning

Some studies, such as the one in [11], have shown that creating an intentional mistuning pattern may reduce the amplification factor. Intentional mistuning consists in assembling sectors which have been designed with different geometries or are made with different materials. In this study, two kinds of sectors are considered ( $M = 2$ ) arranged with the following pattern taken randomly:

$$\text{pattern} = [1, 2, 2, 1, 2, 1, 2, 1, 1, 1, 1, 2, 2, 1, 2, 1, 1, 1, 2, 2, 1, 1, 1, 2], \quad (43)$$

The number 1 denotes the standard sector used in Sections 2.5, 3.2 and 5.1, whereas 2 is a new sector derived from Sector 1 in which the length of the blade has been reduced by 0.5 mm. The reference system is obtained following the methodology given in Section 2 with  $m_1 = 14$  and  $m_2 = 10$ . For the SNCR, the cyclic matrices of the two types of sectors are evaluated, and their respective NNMs are then computed. For these two kinds of sectors ( $i \in \llbracket 1, 2 \rrbracket$ ), the basis  $\Phi_i$  in Equation (41) is formed by the NNMs corresponding to the sector of type  $i$ . In this test case, the system is excited by a standing wave excitation with  $h_{ex} = 3$  and an amplitude of 0.25 N. Hence,  $\Phi_i$  contains the NNMs of the zeroth, third and ninth nodal diameter of the sector type  $i$  (as explained in Section 4.2). Figure 12 represents the frequency forced response of the intentionally mistuned system. Once again, the results obtained with the SNCR are close to those obtained with the reference method. This validates the new methodology for intentionally mistuned structures.



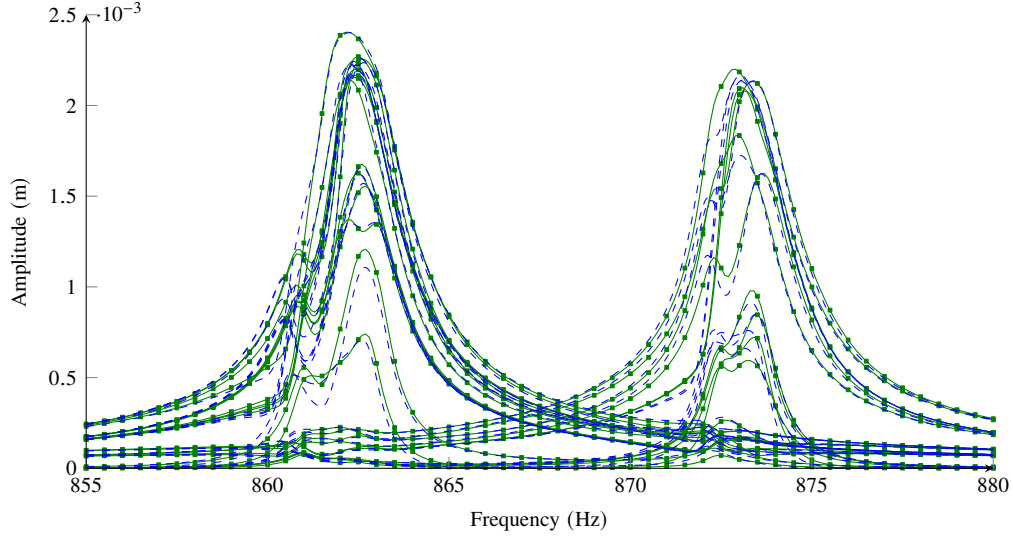


Figure 12: Nonlinear frequency forced response for all sectors of an intentionally mistuned system under a standing wave excitation  $h_{ex} = 3$ . (—■—): reference solution, (---): SNCR solution

### 400 5.3. Performances and limitations

This section provides full details on the performance of the SNCR when applied to the previous simulations. In practical applications, designers are generally interested in the highest resonant peak (its amplitude and its frequency). As a consequence we will use as quality measurement the following errors in amplitude, noted  $L_{\infty, u}$ , and frequency, noted  $L_{\infty, \omega}$ , to characterize the accuracy of the different approaches with respect to the reference.  $L_{\infty, u}$  is defined with

$$L_{\infty, u} = \frac{\left| \max_j \left[ \max_{\omega} (u_{j, \text{tip}, \text{ROM}}) \right] - \max_j \left[ \max_{\omega} (u_{j, \text{tip}, \text{Ref}}) \right] \right|}{\max_j \left[ \max_{\omega} (u_{j, \text{tip}, \text{Ref}}) \right]}, \quad (44)$$

where  $u_{j, \text{tip}, \text{ROM}}$  (respectively  $u_{j, \text{tip}, \text{Ref}}$ ) represents the displacement of sector  $j$  at the tip of the blade evaluated with the ROM (respectively the reference solution). The error  $L_{\infty, \omega}$  is evaluated with

$$L_{\infty, \omega} = \left[ \frac{\left| \max_j (\omega_{j, \text{MAX}, \text{ROM}}) - \max_j (\omega_{j, \text{MAX}, \text{Ref}}) \right|}{\max_j (\omega_{j, \text{MAX}, \text{Ref}})} \right], \quad (45)$$

where  $\omega_{j, \text{MAX}, \text{ROM}}$  (respectively  $\omega_{j, \text{MAX}, \text{Ref}}$ ) represents the frequency of the ROM (respectively the reference solution) for which the tip displacement of sector  $j$  reaches its maximum. In order to study the global

410 accuracy of the ROM we also define two mean errors, averaged over the different sectors,

$$ME_u = \frac{1}{N} \sum_{j=1}^N \left[ \max_{\omega} \left( \frac{|u_{j,\text{tip,ROM}} - u_{j,\text{tip,Ref}}|}{u_{j,\text{tip,Ref}}} \right) \right], \quad (46)$$

and,

$$ME_{\omega} = \frac{1}{N} \sum_{j=1}^N \left[ \frac{|\omega_{j,\text{MAX,ROM}} - \omega_{j,\text{Ref}}|}{\omega_{j,\text{MAX,Ref}}} \right]. \quad (47)$$

The quantity  $L_{\infty,u}$  denotes the absolute relative error between the maximum displacements while  $L_{\infty,\omega}$  corresponds to the absolute relative error between their associated frequencies.  $ME_u$  represents the average error over all the sectors and  $ME_{\omega}$  gives the average difference for their associated frequencies. Table 4 provides these information as well as the computation time and the number of unknowns for the nonlinear reference method, the SNCR and the CNCMS. The CNCMS performances are provided in Table 4 even though their associated results have not always been illustrated.

For all the test cases, the SNCR shows excellent attributes: its accuracy is very good and its computation time is at least 15 times slower than the reference solution. The highest error of 2% is reached when localization phenomena occurs (due to a localized external force). This can be explained by the fact that both mistuning and nonlinear phenomena couple many nodal diameters. As a consequence, even if the linear solution is converged, the nonlinearities may create additional couplings which will not be correctly captured by the reduction basis. However, these discrepancies are overall very small. The memory space required by the SNCR is also interesting as only the data for the  $M$  different sectors have to be saved during the solution procedure. In comparison, the CNCMS must store the model of the full disk with the boundary DOFs to perform the interface modes computation.

Overall the SNCR provides more accurate results and requires less computation time than the CNCMS method. As an example, for the simulation in Figure 8b the error level is reduced from 8.2% to 1.1% using the SNCR and the computation time is divided by 2.6. The only scenario where the CNCMS behaves better is for the localized excitation. In this particular case, the energy is mostly contained within a single sector and hence the basis of the CNCMS (fixed-boundary NNM + static modeshapes) is more accurate than the basis of the SNCR (composed of cyclic NNMs).

The main drawback of the SNCR is the computation time of the NNMs (around 50min per NNM), as shown in Table 3. This is mainly due to the inversion of the linear dynamic rigidity matrix (containing the boundary DOFs) for the Schur condensation.

Methodologies		Simulations			
		Figure 2 and 8b	Figure 9b	Figure 11c	Figure 12
<b>Reference</b>	No. of unknowns	15120	15120	15120	15120
	Computation time	7h	26h	35h	40h
<b>SNCR</b>	No. of unknowns	480	672	480	480
	Computation time	15 min	1h	2h	1h
	$L_{\infty,u}$	1.1%	2%	1.6%	0.1%
	$ME_u$	1.2%	0.8%	1.2%	2.7%
	$L_{\infty,\omega}$	< 0.01%	< 0.01%	0.011%	< 0.01%
	$ME_\omega$	< 0.01%	< 0.01%	0.02%	0.01%
<b>CNCMS</b>	No. of unknowns	848	848	848	848
	Computation time	40min	2h	4h	2h
	$L_{\infty,u}$	8.2%	0.05%	9.6%	7.3%
	$ME_u$	3.2%	0.8%	5.7%	3.9%
	$L_{\infty,\omega}$	< 0.01%	< 0.01%	0.1%	< 0.01%
	$ME_\omega$	0.08%	< 0.01%	0.02%	< 0.01%

Table 4: Performance of the different methodologies on the simulations performed in Section 5. The simulations were run on a Intel(R) Core(TM) i7-7700 @ 3.6 GHz computer.

435 To underline the performance of the SNCR, a parametric study is then performed on the intentionally  
misted structure presented in Section 5.2.2. The maximum displacements over all the sectors are repre-  
sented in Figure 13 for differents values of excitation amplitude (ranging from 0.01 to 0.25N). Such simu-  
lation lasted approximatively a day for the SNCR but would have required several weeks for the reference  
method.

440 In Figure 13, the maximal displacement over all sectors is represented as a function of the excitation

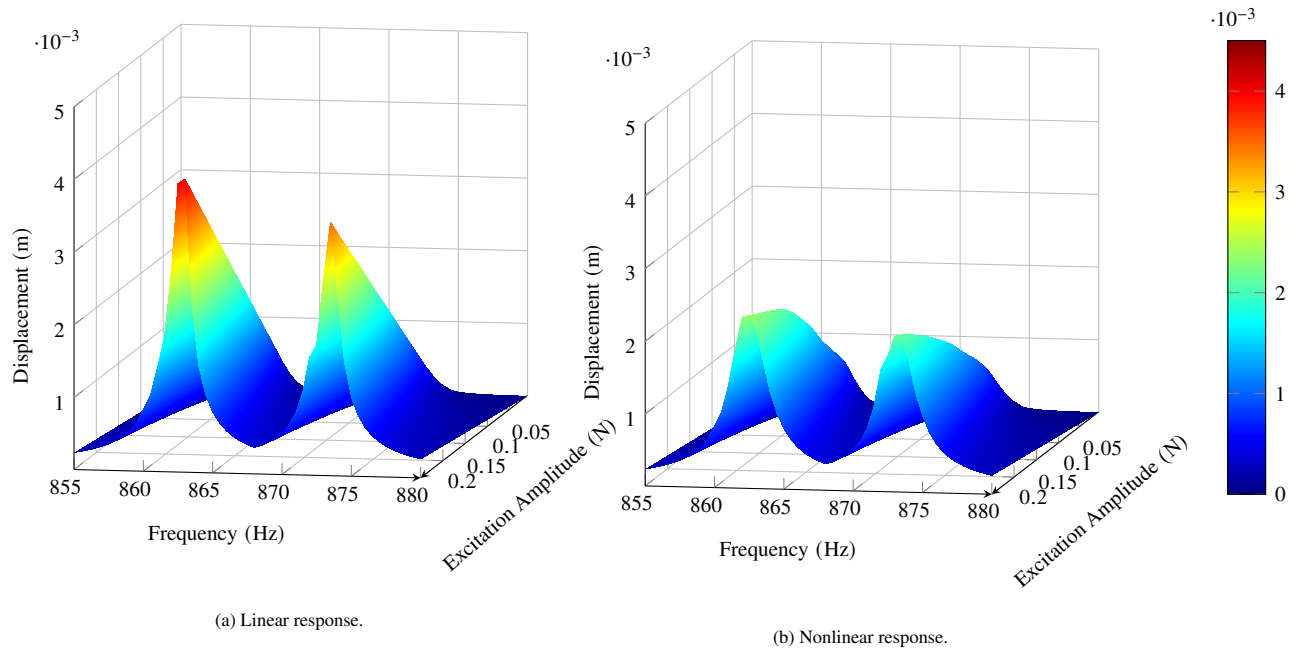


Figure 13: Parametric study on the amplitude excitation for a standing wave ( $h_{ex} = 3$ ).

frequency and the amplitude excitation for a linear and a nonlinear simulations. Using these results, different observations can be made to characterize the nonlinear system. It is for instance possible to determine the efficiency of the nonlinearity all across the excitation amplitude range. For the highest excitation amplitude simulated, the nonlinear effects assure more than a 50% displacement decrease at the resonant peaks.

Moreover, the nonlinearity only gets distinctly activated for an amplitude higher than 0.1 N. Notice also that standard features of the nonlinear system can be observed in Figure 13b: the shape of the backbone curve is similar to a quadratic shape (against a linear variation for the stuck case in Figure 13a), and the resonant peaks get flattened and widen by the nonlinearity activation level.

This result can now be compared to the nonlinear tuned case (system with only sectors of type 1) whose results are illustrated in Figure 14. To highlight the frequency split incoming from the mistuning effect, the results represent the displacement of the 8<sup>th</sup> sector in the logarithm scale. As expected, a single ray is observed for the tuned case (see Figure 14a), whereas the mistuned system contains multiple rays (three in this case, see Figure 14b). Moreover, we can see, for this sector and the mode under study that the displacement amplitude is lower for the mistuned case than its tuned counterpart.

Through this short analysis, we can highlight that the SNCR is a suited candidate to tackle a wide range

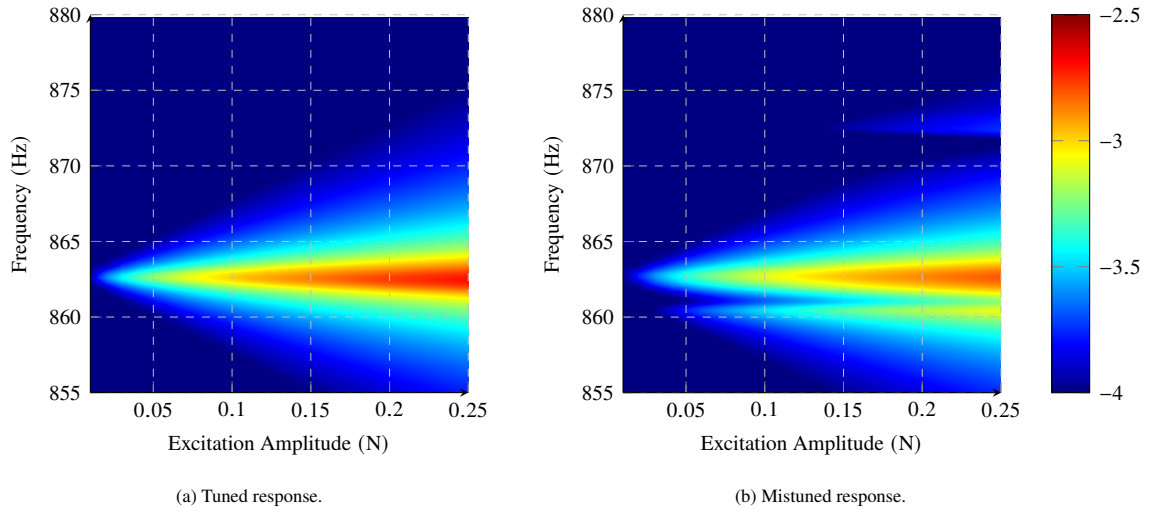


Figure 14: Comparison between the tuned and mistuned nonlinear displacement of the 8<sup>th</sup> sector in the logarithm scale.

of applications for nonlinear cyclic structures and is able to perform accurate parametric studies which were, up to now, not realistically feasible on standard computers.

## 6. Conclusion

This paper proposes a new reduced-order model to determine the steady-state vibration of nonlinear  
 460 mistuned structures. A synthesis procedure based on the cyclic complex nonlinear mode was first proposed  
 for cyclic symmetric structures. This method was shown to be very accurate but only holds for traveling  
 wave excitations. Combining a substructure technique applied to the system with the cyclic complex non-  
 linear mode, a new ROM named Substructuring method based on Nonlinear Cyclic Reduction (SNCR) was  
 proposed. For all cases simulated, the new method was shown to be very accurate (below 2% of error)  
 465 while reducing significantly the computation time (overall 15 times lower than the reference method). It  
 was also compared with a recent ROM and was proved to be a more suitable candidate for nonlinear mis-  
 tuned structures (error divided by 7 for a standard traveling wave excitation). The SNCR employs different  
 assumptions, the main ones of which are the creation of a nonlinear basis composed of only specific cyclic  
 NNMs and the substitution of the nonlinear forces by the formulation of the NNMs.

470 The flexibility of the SNCR is supposed to hold for a wide range of studies, applications and com-  
 plex phenomena: underplatform dampers, gyroscopic effects, veering area, etc. Its reduced computation

time makes it a very good candidate for nonlinear parametric studies to predict, for instance, the impact of mistuning and excitation amplitude on the amplification factor. The method is therefore expected to be of high interest for the scientific community working in the field of cyclic structures (perfectly symmetric or mistuned).

## Acknowledgements

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