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Reduction of cyclic symmetric mechanical systems subject to blade-tip/casing contact interactions

Thibaut Vadcard, Samuel Quaegebeur, Fabrice Thouverez

Abstract

This paper proposes a novel reduction technique for industrial bladed disks featuring blade-tip/casing contacts. By analyzing a regularized contact law, a criterion allowing to drastically reduce the size of the problem is derived. It consists in restricting solutions to only a subset of nodal diameters identified in advance. An additional parity rule allowing to reduce the size of problem by half in certain configurations when solving through the harmonic balance method (HBM). These rules are tested on an academic cyclic model featuring unilateral contacts. The reduction is exact both for a regularized law but also for a nonsmooth contact law. The results match perfectly with the reference calculations led on the full system. The previously published traveling wave hypothesis (deducing the full response from a single sector) also provides exact results when considering traveling wave excitations. All the previously mentioned observations remain true for the simulations led on the industrial fan stage ECL5/Catana under rubbing interactions. After including friction forces, the reduction technique is still exact for the industrial model. The traveling wave hypothesis also remains valid. Several loading scenarios demonstrate the applicability of the strategy in numerous configurations. Localized responses are accurately captured by the reduction method in the case of excitation through perturbed traveling waves. This contribution allows to take a major step forward in the numerical modeling of bladed disks featuring blade-tip/casing interactions by taking advantage of cyclic symmetry. The proposed strategy yields an average reduction of 37.6% for the HBM and 25.2% for other methods. For traveling wave excitations, the reduction is rapidly over 90% and goes up as the number of sectors increases.

Keywords

harmonic balance method, unilateral contact, blade-tip/casing interactions, nonlinear dynamics, computational mechanics, cyclic symmetry, reduction method

1 - École Centrale de Lyon, Laboratoire de Tribologie et Dynamique des Systèmes, UMR CNRS 5513, 36 avenue Guy de Collongue, Écully, 69134, France

Réduction de systèmes mécaniques sujets à des contacts aube/carter en symétrie cyclique

Thibaut Vadcard, Samuel Quaegebeur, Fabrice Thouverez

Résumé

Cet article propose une nouvelle technique de réduction pour les roues aubagées industrielles subissant des contacts aube/carter. En analysant une loi de contact régularisée, un critère permettant de réduire considérablement la taille du problème est obtenu. Il consiste à restreindre les solutions à un sous-ensemble de diamètres nodaux identifiés à l'avance. Une règle de parité supplémentaire permet de réduire de moitié la taille du problème dans certaines configurations lors de la résolution par la méthode de l'équilibre harmonique (HBM). Ces règles sont testées sur un modèle cyclique académique comportant des contacts unilatéraux. La réduction est exacte à la fois pour une loi régularisée et pour une loi de contact non régulière. Les résultats correspondent parfaitement aux calculs de référence effectués sur le système complet. L'hypothèse d'onde tournante publiée précédemment (déduisant la réponse complète d'un seul secteur) fournit également des résultats exacts lorsque l'on considère des excitations tournantes. Toutes les observations mentionnées précédemment restent vraies pour les simulations menées sur l'étage de soufflante industrielle ECL5/Catana en cas de contact aube/carter. Après l'inclusion des forces de frottement, la technique de réduction reste toujours exacte pour le modèle industriel. L'hypothèse de l'onde tournante reste également valable. Plusieurs scénarios de chargement démontrent l'applicabilité de la stratégie dans de nombreuses configurations. Des réponses localisées sont détectées avec précision par la méthode de réduction dans le cas d'excitations par des ondes tournantes perturbées. Cette contribution constitue une avancée majeure dans la modélisation numérique des roues aubagées en situation de contact rotor/stator en tirant parti de la symétrie cyclique. La stratégie proposée permet une réduction moyenne de 37,6% pour le HBM et de 25,2% pour les autres méthodes. Pour les excitations par ondes tournantes, la réduction est rapidement supérieure à 90% et augmente avec le nombre de secteurs.

Mots-clés

méthode d'équilibrage harmonique, contact unilatéral, contact aube/carter, dynamique non linéaire, mécanique numérique, symétrie cyclique, méthode de réduction

1 - École Centrale de Lyon, Laboratoire de Tribologie et Dynamique des Systèmes, UMR CNRS 5513, 36 avenue Guy de Collongue, Écully, 69134, France

1 Introduction

In order to comply with international regulations that are aiming towards the achievement of net carbon neutrality [1], aircraft engine manufacturers are committed to design continuously more efficient turbomachines. An aeronautical engine is a highly nonlinear and multiphysic system that should be tackled by using advanced modeling techniques. From a structural viewpoint, the nonlinearities play a pivotal role in the dynamics of bladed disks. Notably, the blade-tip/casing contact interface is among the nonlinearities that are the most studied. This interface is mainly addressed for safety concerns about the engine. Indeed, the gaps between the rotating and fixed parts of the engine, here blades and the casing, are decreased in new engines as a way of increasing the aerodynamic efficiency of the turbomachine. This has the impact of yielding structural contacts in nominal operating conditions $[2, 3]$. These contacts at high speeds are examined with great care by designers since their consequences can threaten the integrity of engine itself, the plane and thus the safety of the passengers.

The realization of full scale experiments of blade-tip/casing contacts in order to evaluate the dynamics response of the engine are oftentimes prohibitively expensive for the manufacturers. In addition, this type of experiments can not be carried out until the final stage of the design cycle. Such tests can not serve to actually design the system, and are generally just about safety verifications. Hence, there is a need to setup efficient simulation tools that are able to account for the nonlinear behavior of the engine at early stage of design. This is why many researchers are working towards industry-ready, robust, predictive and accurate numerical strategies in order to characterize these sophisticated and highly nonlinear systems.

Among the strategies used to reduce computational times without hurting the accuracy of the results, the cyclic symmetry [4, 5] is widely used in a linear framework. This formulation takes advantage of the angular periodicity of cyclic systems, such as bladed disks. In a perfectly tuned configuration, the dynamics of the bladed disk can be expressed as a sum of decoupled nodal diameters (spatial spectral components). Hence, by analyzing the spectral components present in the loading applied to the system, one can only keep the subset of nodal diameters that are undergoing a loading and thus drastically reduce the size of the system to be solved at no cost in terms of accuracy. However, in a nonlinear framework, the different nodal diameters are not uncoupled anymore and specific strategies should be used. Petrov [6] formulated an hypothesis that states that in the case of a traveling wave excitation on friction-damping applications, the displacement field and the nonlinear forces have the same shape as the excitation. This so-called traveling wave hypothesis greatly improves the performances of the simulations since the results on a single sector can be used to deduce the response of the full bladed disk [7]. Following this, Petrov also tackled standing wave excitations that can be considered to be traveling wave on a supersector composed of several actual sectors of the bladed disk. Siewert *et al.*[8] also tackled the problem of friction between shrouds in a cyclic symmetric framework with the traveling wave hypothesis. Quaegebeur [9] went further and derived a criterion allowing to determine the couplings that occur between the different nodal diameters by analyzing the Taylor expansion of the nonlinear forces. This last strategy allows to compute the response of the full system for any excitation without making any approximation on the solution. However, these strategies were designed and tested to be applied only on systems featuring friction interfaces.

Indeed, when considering systems featuring other nonlinearities than friction, such as unilateral contacts or polynomial nonlinearities, no specific criterion exists to determine how the nodal diameters are coupled with each other. Moreover, the traveling wave hypothesis remains untested on cyclic systems featuring unilateral contacts and blade-tip/casing contacts. Until now, simulations of blade-tip/casing contacts are carried out either on a single blade [10, 11, 12, 13, 14, 15, 16] or on a full bladed disk [17, 18, 19]. While simulations on single blades are able to obtain results in a reasonable time, they are unable to reflect the dynamics of the full bladed disk and thus yield an approximation on the vibrations of the system. On the contrary, studies on the full bladed disk do not make any assumption on the solutions but come at an extremely prohibitive computational cost. This contribution aims to tackle this modeling issue and propose a method that allows to model the full bladed disk without having to solve the problem for all unknowns.

The harmonic balance method (HBM) has recently known a growth in popularity when it comes to the characterization of blade-tip/casing interactions $[11, 20, 21]$. Its ability to gather information about nonlinear solutions is particularly powerful. Indeed, it is able to state on the stability of periodic orbits $[21, 22]$, the connectivity of the solutions $[23, 24]$ and it allows to compute nonlinear modal parameters of a system $[25, 26]$. Moreover, it was shown to achieve a similar level of accuracy than the reference explicit time-marching strategies [11] on industrial models at a lower computational cost. For these reasons, it is generally considered that the HBM outperforms the time integration methods for the computation of periodic orbits of a nonlinear system. This article is then centered around the HBM but some of its theoretical elements can still be applied to time-marching simulations.

This paper aims to tackle the challenging subject of reduction of bladed disks in a nonlinear framework. Firstly, analytical derivations are made in order to determine a method of computing a reduction basis on which the solutions of the problem should be projected. This reduction basis is actually composed of a subset of nodal diameters that are expected to be coupled through the unilateral nonlinearity. This analysis relies on the Taylor expansion of a regularized penalty law [20]. Another reduction of the size of the problem is proposed when considering a resolution with the HBM based on the spatio-temporal couplings of nodal diameters with temporal harmonics. Then, the validity of these reductions are assessed on an academic model both for the regularized law that served to derive the reduction rules, but also on a nonsmooth law [27, 28] relying on augmented lagrangians. The verification on the latter non-differentiable contact law allows to extend the reduction rule to any contact law that models unilateral laws regardless of the contact formulation that it relies on. Several contact scenarios are studied in order to give an exhaustive view of the reduction capabilities of the strategy. Then, the simulations are taken a step further to characterize a full industrial bladed disk through the cyclic symmetry based reduction strategy, on both contact laws again. In all configurations, the traveling wave hypothesis is also shown to hold true on these applications which yields a significant reduction in the case of traveling wave excitations.

The physical framework relying on the cyclic symmetry is described in Sec. 2: it includes the formulation of the problem and the derivation of the reduction criterion. Then, the HBM-related resolution strategies are described in Sec. 3, it also includes the formulation of the supplementary reduction criterion based on the HBM. Then, numerous simulations are led to validate the reduction strategy on a simplified model in cyclic symmetry in Sec. 4. Finally, all the presented developments are applied to a representative fan blade model in Sec. 5. It is shown that the procedure is industry-ready and allows to accurately model the dynamics of the full bladed disk at a low computational cost.

2 Cyclic symmetry and reduction strategy

2.1 Equations of motion in a cyclic symmetric framework

In this section, the equations of motion for a cyclic symmetric mechanical system are derived in terms of spectral generalized coordinates. For each sector $j \in [1, N]$ where N is the number of sectors of the system, the equations of motion read

$$
\mathbf{M}_0 \ddot{\mathbf{x}}_j + \mathbf{C}_0 \dot{\mathbf{x}}_j + \mathbf{K}_0 \mathbf{x}_j = \mathbf{f}_{\mathrm{s},j}^{\mathrm{p}} + \mathbf{f}_{\mathrm{s},j}^{\mathrm{n}} + \mathbf{f}_{\mathrm{nl},j}(\mathbf{x}_j(t), \dot{\mathbf{x}}_j(t)) + \mathbf{f}_{0,j} + \mathbf{f}_{\mathrm{ex},j}(\omega, t)
$$
(1)

where M_0 , C_0 and K_0 are respectively the mass, damping and stiffness matrix of a single sector of the structure. The full structure is considered to be tuned, meaning that all sectors are identical. x_i is the displacement field of the n degrees of freedom (dof) of the jth sector. In order to represent the full system, the influence of neighboring sectors is accounted for by the forces applied to the cyclic boundary of the sector $\mathbf{f}_{s,j}^{\mathrm{p}}$ and $\mathbf{f}_{s,j}^{\mathrm{n}}$, related to the forces applied by the previous and the next sector respectively. Note that sector number 1 is adjacent to sectors N and 2 , and that sector number N is adjacent to sectors $N-1$ and 1 so that the cyclic nature is preserved. The system also undergoes nonlinear forces, represented by $\mathbf{f}_{n,l,j}$, external pre-loading forces $\mathbf{f}_{0,j}$ and excitation forces $\mathbf{f}_{ex,j}$.

The physical displacement x_i can then be expressed as a function of the real spectral components of the cyclic system thanks to the theory of linear representation of cyclic groups of order N [5]. It follows that

$$
\mathbf{x}_{j} = \frac{1}{\sqrt{N}}\hat{\mathbf{x}}_{0} + \sqrt{\frac{2}{N}}\sum_{h=1}^{\frac{N}{2}-1} \left[\hat{\mathbf{x}}_{h,c}\cos\left(hj\alpha\right) + \hat{\mathbf{x}}_{h,s}\sin\left(hj\alpha\right)\right] + \frac{(-1)^{j}}{\sqrt{N}}\hat{\mathbf{x}}_{N/2} \quad \text{if } N \text{ is even},\tag{2}
$$

$$
\mathbf{x}_{j} = \frac{1}{\sqrt{N}} \hat{\mathbf{x}}_{0} + \sqrt{\frac{2}{N}} \sum_{h=1}^{\frac{N-1}{2}} \left[\hat{\mathbf{x}}_{h,c} \cos(hj\alpha) + \hat{\mathbf{x}}_{h,s} \sin(hj\alpha) \right]
$$
 if *N* is odd (3)

where $\alpha = \frac{2\pi}{N}$ is the intersector angle and $\hat{\mathbf{x}}_{h,c}$ and $\hat{\mathbf{x}}_{h,s}$ are the real spectral displacements of order h. Following the definition of the quality Equation patrix \mathbf{S} in A, physical displacemen definition of the cyclic Fourier matrix **S** in A, physical displacements of all sectors $\mathbf{x} = [\mathbf{x}_1, \dots, \mathbf{x}_j, \dots, \mathbf{x}_N]$ can be substituted by the spectral components such as

$$
\mathbf{x} = \mathbf{S}^{\top} \hat{\mathbf{x}} \tag{4}
$$

with $\hat{\mathbf{x}} = \begin{bmatrix} \hat{\mathbf{x}}_0^{\top}, \ldots, \hat{\mathbf{x}}_h^{\top}, \ldots, \hat{\mathbf{x}}_m^{\top} \end{bmatrix}^{\top}$ with $m = \frac{N}{2}$ if N is even or $m = \frac{N-1}{2}$ if N is odd. All spectral components of order $0 < h < \frac{N}{2}$ contain both cosine and sine contributions such as $\hat{\mathbf{x}}_k = \left[\hat{\mathbf{x}}_{h,c}^{\top}, \hat{\mathbf{x}}_{h,s}^{\top}\right]^{\top}$. The notation \bullet^{\top} is the transposition operation.

In this work the cyclic boundaries of the sectors remain linear, thus the different cyclic boundaries are assembled [4, 5] leading to a system of equations where all matrices are block-diagonal accounting for the fact that all cyclic indexes are uncoupled. In the nonlinear framework studied, a specific coupling is appearing between the cyclic indexes due to the nonlinear forces applied on the system. This coupling is the main topic of this contribution and is thoroughly addressed throughout the article. In the end, the problem associated to the full system is then expressed such as

$$
\widehat{\mathbf{M}}\ddot{\hat{\mathbf{x}}}(t) + \widehat{\mathbf{C}}\dot{\hat{\mathbf{x}}}(t) + \widehat{\mathbf{K}}\hat{\mathbf{x}}(t) = \widehat{\mathbf{f}}_{\text{ex}}(\omega, t) + \widehat{\mathbf{f}}_{0} + \widehat{\mathbf{f}}_{\text{nl}}(\mathbf{x}(t), \dot{\mathbf{x}}(t))
$$
\n(5)

with

$$
\widehat{\mathbf{M}} = \text{blkdiag}\left(\left\{\widehat{\mathbf{M}}^h\right\}_{h=0}^m\right), \widehat{\mathbf{C}} = \text{blkdiag}\left(\left\{\widehat{\mathbf{C}}^h\right\}_{h=0}^m\right), \widehat{\mathbf{K}} = \text{blkdiag}\left(\left\{\widehat{\mathbf{K}}^h\right\}_{h=0}^m\right), \tag{6}
$$
\n
$$
\widehat{\mathbf{f}}_{\mathbf{C}}^0
$$

$$
\hat{\mathbf{f}}_{\text{ex}}(\omega, t) = \begin{bmatrix} \mathbf{f}_{\text{ex}}^0 \\ \vdots \\ \hat{\mathbf{f}}_{\text{ex}}^h \\ \vdots \\ \hat{\mathbf{f}}_{\text{ex}}^m \end{bmatrix}, \qquad \hat{\mathbf{f}}_0 = \begin{bmatrix} \mathbf{f}_0^0 \\ \vdots \\ \hat{\mathbf{f}}_0^h \\ \vdots \\ \hat{\mathbf{f}}_0^m \end{bmatrix} \text{ and } \qquad \hat{\mathbf{f}}_{\text{nl}}(\mathbf{x}(t), \dot{\mathbf{x}}(t)) = \begin{bmatrix} \mathbf{f}_{\text{nl}}^0 \\ \vdots \\ \hat{\mathbf{f}}_{\text{nl}}^h \\ \vdots \\ \hat{\mathbf{f}}_{\text{nl}}^m \end{bmatrix}.
$$
\n(7)

The notation $\hat{\bullet}^h$ refers to spectral components of order h, oftentimes referred to as nodal diameters (ND) of order h. It is worth noting that the nonlinear forces expressed in terms of spectral components \hat{f}_{nl} still depends on physical displacements $\mathbf{x}(t)$ and velocities $\dot{\mathbf{x}}(t)$. Indeed, the evaluation of nonlinear forces, presented further in the article in Sec. 2.2 should be made on physical coordinates. For numerical conditioning purposes, a space and time normalization of the unknown $\hat{\mathbf{x}}(t)$ can be introduced by means of scalar coefficients [24]. The latter do not affect the solutions of the problem but allow better convergence of nonlinear solvers. However, these normalizations are not included in the equations of this article for the sake of simplifying the expressions.

This general formulation in cyclic symmetry is actually equivalent of solving the full problem with all sectors and thus yields no computational reduction with respect to the full physical problem. Nevertheless, this strategy enables to perform full computations without defining the whole problem as a single matrix. This particular property is helpful when generating reduced-order models of industrial size models. These models generally cause issues in terms of memory that can be avoided through the use of cyclic symmetry.

Since the nodal diameters are linearly decoupled (*i.e.* \hat{M} , \hat{C} and \hat{K} are block-diagonal), the only link between the different nodal diameters in Eq. (5) is made through the nonlinear forces. Hence, it is possible to analyze the analytical expressions of the nonlinear forces in order to determine the couplings that occur between nodal diameters. By doing so, a subset of nodal diameters $\mathcal{D} \subset [0, m]$ is determined for a specific loading scenario. The latter can be used for a nonlinear calculation instead of the whole range $\mathcal{D}^f = [0, m]$. Indeed, for certain loading scenarios, some nodal diameters are not involved at all in the dynamics and can thus be withdrawn from the unknown vector without affecting the accuracy of the solutions. This generally yields a significant reduction in computation time and memory usage. Equation 5 becomes

$$
\widehat{\mathbf{M}}^{\mathcal{D}} \ddot{\hat{\mathbf{x}}}^{\mathcal{D}}(t) + \widehat{\mathbf{C}}^{\mathcal{D}} \dot{\hat{\mathbf{x}}}^{\mathcal{D}}(t) + \widehat{\mathbf{K}}^{\mathcal{D}} \widehat{\mathbf{x}}^{\mathcal{D}}(t) = \widehat{\mathbf{f}}_{\text{ex}}^{\mathcal{D}}(\omega, t) + \widehat{\mathbf{f}}_{0}^{\mathcal{D}} + \widehat{\mathbf{f}}_{\text{nl}}^{\mathcal{D}}(\mathbf{x}(t), \dot{\mathbf{x}}(t))
$$
\n(8)

with

$$
\widehat{\mathbf{M}}^{\mathcal{D}} = \text{blkdiag}\left(\left\{\widehat{\mathbf{M}}^h\right\}_{h\in\mathcal{D}}\right), \widehat{\mathbf{C}}^{\mathcal{D}} = \text{blkdiag}\left(\left\{\widehat{\mathbf{C}}^h\right\}_{h\in\mathcal{D}}\right), \widehat{\mathbf{K}}^{\mathcal{D}} = \text{blkdiag}\left(\left\{\widehat{\mathbf{K}}^h\right\}_{h\in\mathcal{D}}\right),\tag{9}
$$

$$
\widehat{\mathbf{f}}_{\text{ex}}^{\mathcal{D}} = \left[\left\{ \widehat{\mathbf{f}}_{\text{ex}}^{h} \right\}_{h \in \mathcal{D}} \right], \qquad \widehat{\mathbf{f}}_{0}^{\mathcal{D}} = \left[\left\{ \widehat{\mathbf{f}}_{0}^{h} \right\}_{h \in \mathcal{D}} \right] \qquad \text{and } \widehat{\mathbf{f}}_{\text{nl}}^{\mathcal{D}} = \left[\left\{ \widehat{\mathbf{f}}_{\text{nl}}^{h} \right\}_{h \in \mathcal{D}} \right]. \tag{10}
$$

In the general case, the choice of an appropriate set of nodal diameter is a difficult task. However, by analyzing the nonlinearities at stake, it is possible to derive rules that provide ways of determining the right nodal diameters to keep in D. For instance, Quaegebeur *et al.* derived a specific rule allowing to determine the couplings between the spectral components of a cyclic structure with friction nonlinearities [9]. This rule analyzes the spectral components of the excitations and the number of sectors of the system in order to determine the subset D of nodal diameters allowing for eventual reduction of the system to be solved. This subset is used by only accounting for the nodal diameter that can be coupled with the excitation as well as with each other by the nonlinearity. In this work, the same type of rule is derived for nonlinearities that model unilateral contacts, it is presented in Sec. 2.3.

2.2 Evaluation of nonlinear forces in the physical space

It is reminded that the nonlinear forces oftentimes depend on physical displacements and should be evaluated in the physical space. In order to go from the spectral (cyclic) variables to physical (sectors) variables, one should involve the spatial Fourier matrix S and its inverse, defined in A. As a way of computing the spectral components of the nonlinear forces, the cyclic variables linked to the displacements $\hat{\mathbf{x}}^{\mathcal{D}}(t)$ are firstly transformed to physical
unrighted on all sectors $\mathbf{x}(t)$ than the nonlinear forces are computed in the physical g variables on all sectors $\mathbf{x}(t)$, then the nonlinear forces are computed in the physical space on each sector $\mathbf{f}_{nl}(t)$ and these physical variables are then transformed back to the cyclic space $\hat{\mathbf{f}}_{nl}^{\mathcal{D}}(t)$.

Firstly, let's assume that D is chosen to be the full set $D = D^f$. Then, for sampling purposes, it is necessary to evaluate the nonlinear forces on all sectors $S^f = \{1, 2, ..., N\}$ in order to be able to reconstruct the whole spectral decomposition.

However, given Eq. (8), the unknown vectors $\hat{\mathbf{x}}^{\mathcal{D}}(t)$ and $\hat{\mathbf{f}}_{\text{nl}}^{\mathcal{D}}(\mathbf{x}(t), \dot{\mathbf{x}}(t))$ are composed of all spectral components to reacting the subset $\mathcal{D} \subset \mathcal{D}^{\text{f}}$. Then, if this time $\mathcal{D$ that are retained in the subset $D \subset \mathcal{D}^f$. Then, if this time $D \neq \mathcal{D}^f$ (does not contain every spectral component), the equivalence between physical and cyclic spaces yields that the less nodal diameters are considered, the less sectors are necessary for the computation of the spectral components retained in \mathcal{D} . Indeed, by suppressing some spectral components of the full basis, the information over all sectors becomes redundant.

As a way of determining the set of sectors $S \subset S^f$, over which the nonlinear forces should be sampled (composed of a number of j^{\max} sectors), it is necessary to analyze the content of D . For each non-degenerated spectral component in D (*i.e.* $h = 0$ and $h = \frac{N}{2}$ if N is even and only $h = 0$ if N is odd), a single sector should be evaluated. For each degenerated spectral components $0 < h < \frac{N}{2}$ in D , two sectors should be evaluated. For instance, a structure composed of $N = 24$ sectors for which D is arbitrarily set to $\mathcal{D} = \{0, 2, 4, 6, 8, 10, 12\}$, the number of sectors needed for the evaluation of all spectral components of the nonlinear forces is $j^{\max} = 2 \times 1 + 5 \times 2 = 12$ (2 non-degenerated and 5 degenerated spectral components). Then, one should evaluate the nonlinear forces on all sector numbers belonging to $S = \{1, 2, \ldots, j^{\max}\}.$

Following these guidelines, it is possible to establish a relation between spectral components restricted to D and the physical information over the sectors of S. In order to clarify the notations, complementary sets $\overline{\mathcal{D}} = \mathcal{D}^f \setminus \mathcal{D}$ and $\overline{S} = S^f \setminus S$ are defined. The relation is obtained by partitioning $\hat{\mathbf{x}} = \mathbf{S} \mathbf{x}$ over the different sets such as

$$
\begin{bmatrix} \hat{\mathbf{x}}^{\mathcal{D}} \\ \hat{\mathbf{x}}^{\mathcal{D}} \end{bmatrix} = \begin{bmatrix} \mathbf{S}_{\mathcal{S}}^{\mathcal{D}} & \mathbf{S}_{\mathcal{S}}^{\mathcal{D}} \\ \mathbf{S}_{\mathcal{S}}^{\mathcal{D}} & \mathbf{S}_{\mathcal{S}}^{\mathcal{D}} \end{bmatrix} \begin{bmatrix} \mathbf{x}^{\mathcal{S}} \\ \mathbf{x}^{\mathcal{S}} \end{bmatrix} . \tag{11}
$$

By definition, restricting the spectral components to a specific set \mathcal{D} means that $\hat{\mathbf{x}}^{\mathcal{D}} = \mathbf{0}$. Then, the inverse transformation is expressed as $\mathbf{x} = \mathbf{S}^\top \hat{\mathbf{x}}$. Since the Fourier matrix **S** is orthogonal, it yields

$$
\begin{bmatrix} \mathbf{x}^{\mathcal{S}} \\ \mathbf{x}^{\overline{\mathcal{S}}} \end{bmatrix} = \begin{bmatrix} \left(\mathbf{S}_{\mathcal{S}}^{\mathcal{D}} \right)^{\top} & \left(\mathbf{S}_{\mathcal{S}}^{\overline{\mathcal{D}}} \right)^{\top} \\ \left(\mathbf{S}_{\mathcal{S}}^{\mathcal{D}} \right)^{\top} & \left(\mathbf{S}_{\mathcal{S}}^{\overline{\mathcal{D}}} \right)^{\top} \end{bmatrix} \begin{bmatrix} \widehat{\mathbf{x}}^{\mathcal{D}} \\ \mathbf{0} \end{bmatrix}
$$
(12)

and it comes that

$$
\mathbf{x}^{\mathcal{S}} = \left(\mathbf{S}_{\mathcal{S}}^{\mathcal{D}}\right)^{\top} \widehat{\mathbf{x}}^{\mathcal{D}}.\tag{13}
$$

Since the set S was built to be equivalent to D in the physical space, $(S_S^{\mathcal{D}})^{\top}$ is square and invertible, yielding the direct transformation from $\mathcal S$ to $\mathcal D$ is

$$
\widehat{\mathbf{x}}^{\mathcal{D}} = \left[\left(\mathbf{S}_{\mathcal{S}}^{\mathcal{D}} \right)^{\top} \right]^{-1} \mathbf{x}^{\mathcal{S}}.
$$
\n(14)

This shows that restricting the number of spectral components in D yields a double advantage: firstly it decreases the overall number of unknowns of the problem and secondly it reduces the computational cost of evaluating nonlinear forces since it requires their evaluation on less sectors than for the full system.

2.3 Selection of nodal diameters for unilateral contacts

In this section, a complex formulation is employed for the sake of simplifying the developments, however both complex and real cyclic symmetry feature the same properties. In this work, the resolution is carried out with real cyclic symmetry. The complex formulation is only used in this section and it relies on the complex counterpart of the spectral decomposition in Eq. (3).

In the complex formalism, the spectral decomposition of the displacement field of each sector x_i reads

$$
\mathbf{x}_{j} = \frac{1}{\sqrt{N}} \sum_{h=0}^{N-1} \left(\widehat{\mathbf{x}}^{h} e^{i\alpha h(j-1)} \right) \text{ for } j \in [1, N], \tag{15}
$$

with $\hat{\mathbf{x}}^h = \bar{\hat{\mathbf{x}}}^{N-h}$ for $h \in [0, m]$. Similarly as in the real formalism, a block-diagonal system can be built for all $h \in \mathcal{D}$ such sections $h \in \mathcal{D}^f$ such as

$$
\widehat{\mathbf{M}}^{h}\widetilde{\mathbf{x}}^{h}(t) + \widehat{\mathbf{C}}^{h}\dot{\mathbf{x}}^{h}(t) + \widehat{\mathbf{K}}^{h}\widetilde{\mathbf{x}}^{h}(t) = \widehat{\mathbf{f}}_{\text{ex}}^{h}(\omega, t) + \widehat{\mathbf{f}}_{0}^{h} + \widehat{\mathbf{f}}_{\text{nl}}^{h}(\mathbf{x}(t), \dot{\mathbf{x}}(t)).
$$
\n(16)

As stated, when nonlinearities are taken into account the nonlinear forces create interactions between the different nodal diameters and Eq. (16) must be solved for all nodal diameter at once.

In this work, unilateral contact nonlinearities are considered. For theoretical purposes, a regularized penalty law [20] is considered throughout these developments, but it is shown in this article that these results also apply to non-regularized laws. Furthermore, a single nonlinear dof per sector is assumed, its displacement is noted $x_{nl,j}$. The extension to other dof is straightforward, based on the developments that follow. The nonlinear forces for sector j are evaluated with the following

$$
f_{\mathrm{nl},j}(t) = -\kappa \frac{\left(g_j(t) - x_{\mathrm{nl},j}(t)\right)}{2} + \sqrt{\left(\kappa \frac{\left(g_j(t) - x_{\mathrm{nl},j}(t)\right)}{2}\right)^2 + \gamma^2},\tag{17}
$$

where κ is the contact stiffness and γ the regularization parameter.

The gap function can be written in terms of cyclic components such as

$$
g_j = \frac{1}{\sqrt{N}} \sum_{h=0}^{N-1} \left(\widehat{g}^h e^{i\alpha h(j-1)} \right) \text{ for } j \in [1, N]. \tag{18}
$$

Th gap is a known quantity, as such all the components $(\hat{g}^h)_{h \in \mathcal{D}^{\text{f}}}$ are also known. As an example, for a perfectly constant and over one period, only the component \hat{g}_h is different from 0. To simplify t constant gap over one period, only the component \hat{g}_0 is different from 0. To simplify the following derivations, the new unknown of the problem is set to

$$
\widehat{u}^h = \widehat{g}^h - \widehat{x}_{\rm nl}^h. \tag{19}
$$

The equation of motion, accounting for the nonlinearities reads

$$
\widehat{\mathbf{M}}^h \ddot{\hat{\mathbf{u}}}^h + \widehat{\mathbf{C}}^h \dot{\hat{\mathbf{u}}}^h + \widehat{\mathbf{K}}^h \mathbf{u}^h = \widehat{\mathbf{f}}^h_{\text{ext}} + \widehat{\mathbf{f}}^h_0 + \widehat{\mathbf{f}}^h_{\text{nl}} - \widehat{\mathbf{K}}^h \mathbf{g}^h \text{ for } h \in \mathcal{D}^f,
$$
\n(20)

where the force applied to the only nonlinear dof of the system inside $\hat{\mathbf{f}}_{\text{nl}}^h$ is expressed as

$$
\widehat{f}_{\rm nl}^h = \frac{1}{\sqrt{N}} \sum_{j=1}^N \left(f_{\rm nl,j} e^{-i\alpha h(j-1)} \right) \text{ for } h \in \mathcal{D}^{\rm f}.
$$
\n(21)

Substituting the term $f_{n,l,j}$ in Eq. (21) by Eq. (17) (where the physical quantities have been replaced by the cyclic ones) for a given $q \in \mathcal{D}^f$ reads

$$
\widehat{f}_{\rm nl}^{q} = \frac{1}{\sqrt{N}} \sum_{j=1}^{N} \left(\left[-\frac{\kappa}{2} \sum_{h=0}^{N-1} \widehat{u}^{h} e^{\mathrm{i}h(j-1)\alpha} + \frac{\kappa}{\gamma} \sqrt{1 + \frac{1}{\gamma^{2}} \left(\sum_{h=0}^{N-1} \widehat{u}^{h} e^{\mathrm{i}h(j-1)\alpha} \right)^{2}} \right] e^{-\mathrm{i}q(j-1)\alpha} \right). \tag{22}
$$

The second term is expanded as a Taylor series at 0

$$
\sqrt{1+y^2} = 1 + \sum_{l=0}^{\infty} \beta_l y^{2l}.
$$
\n(23)

This leads to

$$
\widehat{f}_{\rm nl}^{q} = \frac{1}{\sqrt{N}} \sum_{j=1}^{N} \left(\left[-\frac{\kappa}{2} \sum_{h=0}^{N-1} \widehat{u}^{h} e^{i(h-q)(j-1)\alpha} + \frac{\kappa}{\gamma} \left(1 + \sum_{l=0}^{\infty} \left(\frac{\beta_{l}}{\gamma^{2}} \left(\sum_{h=0}^{N-1} \widehat{u}^{h} e^{ih(j-1)\alpha} \right)^{2l} \right) \right) \right] e^{-iq(j-1)\alpha} \right). \tag{24}
$$

Notice that the first term in \hat{u}_h is different from 0 when $h = q$. This means that if a nodal diameter responds, it will provide a component to the nonlinear forces on the same nodal diameter. Moreover the constant term $\frac{\kappa}{\gamma}$ provides a nonlinear force component on its 0th nodal diameter. Finally, the following expression explicits the last term (noted $\widehat{f}_{\text{nl},3}^q$) of Eq. (24)

$$
\widehat{f}_{\mathrm{nl},3}^{q} = \frac{\kappa}{\sqrt{N}\gamma^{3}} \sum_{j=1}^{N} \left(\sum_{l=0}^{\infty} \left(\beta_{l} \left(\sum_{h=0}^{N-1} \widehat{u}_{h} e^{\mathrm{i}h(j-1)\alpha} \right)^{2l} \right) \right) e^{-\mathrm{i}q(j-1)\alpha}
$$
\n(25)

$$
= \frac{\kappa}{\sqrt{N}\gamma^3} \sum_{l=0}^{\infty} \sum_{j=1}^{N} \left(\left(\beta_l \left(\sum_{h_1=0}^{N-1} \hat{u}_{h_1} e^{ih_1(j-1)\alpha} \right) \cdots \left(\sum_{h_{2l}=0}^{N-1} \hat{u}_{h_{2l}} e^{ih_{2l}(j-1)\alpha} \right) e^{i\alpha(j-1)-q} \right) \right).
$$
 (26)

Developing the sums in Eq. (26) gives

$$
\widehat{f}_{\mathrm{nl},3}^{q} = \frac{\kappa}{\sqrt{N}\gamma^{3}} \sum_{l=0}^{\infty} \left(\beta_{l} \sum_{k \in \mathcal{T}_{l}} \left(\alpha_{k} \prod_{p_{k}=1}^{Q_{k}} \left(\widehat{u}_{h_{p_{k}}} \right)^{\beta_{p_{k}}} \left(\sum_{j=1}^{N} e^{i\alpha(j-1)\left(\sum_{p_{k}=1}^{Q_{k}} h_{p_{k}} \beta_{p_{k}} - q\right)} \right) \right) \right). \tag{27}
$$

The set \mathcal{T}_l describes all the possible combinations of the sums' multiplications. Each combination k involves one or multiple nodal diameters noted h_{p_k} and their multiplicity β_{p_k} . The terms α_k denotes the number of occurrence of such a term. Since the degree of the polynomial in Eq. (27) is $\sum_{p_k=1}^{Q_k} \beta_{p_k}$. This degree must be equal to the sum of the degrees of all monomials associated with the 2l products of Eq. (26) (from h_1 to h_{2l}). Thus, the following relation must be verified

$$
\sum_{p_k=1}^{Q_k} \beta_{p_k} = 2l. \tag{28}
$$

It corresponds to the sum of the degrees of the different monomials in the combination k considered. The nodal diameter couplings that are created through all terms of Eq. (27) are now analyzed in Sec. 2.4.

2.4 Determination of the nodal diameter coupling

This section proposes an algorithm to determine the nodal diameter coupling occurring when contact nonlinearities are taken into account, it relies on the equations set up is Sec. 2.3. Firstly, \mathcal{D}_q denotes the set of nodal diameters that include the nodal diameters corresponding to \hat{g} and the external forces, that is known data. Based on Eq. (20), all these nodal diameters are expected to respond and are thus kept in the basis. Moreover, based on Eq. (24), due to the first term, the 0th nodal diameter also has to be included in the basis, now noted $\mathcal{D}_{q,0}$. This basis is being completed through an iterative process with Eq. (25) : a new nodal diameter noted p must be included in the basis if

$$
\sum_{i=1}^{2l} h_i - p \equiv 0 \,[N] \text{ with } h_i \in \mathcal{D}_{q,0} \text{ and } l \in \mathbb{N}.
$$

As l goes up to infinity, such an iterative process is cumbersome. However, a few mathematical considerations allow to perform such procedure in an efficient way. First, it is assumed that the nodal diameter basis only contains the diameters 0 and h_{ex} (the nodal diameter associated with the external loading). The 0th nodal diameter is not involved in the iterative process as it does not create further coupling. Then, Eq. (29) becomes

$$
2lh_{\text{ex}} \equiv p[N] \quad l \in \mathbb{N}.\tag{30}
$$

There exists an integer K such that the remainder of the euclidian division of $2h_{\text{ex}}$ by l (noted p), for $l \geq K$, is the same as the one of the euclidian division of $2h_{\rm ex}$ by l with $l \in [1, K]$. Then, it comes that

$$
2(l+K)h_{\text{ex}} \equiv p[N].\tag{31}
$$

As $2lh_{\text{ex}} \equiv p[N]$, one obtains the following

$$
2Kh_{\text{ex}} \equiv 0\,[N] \tag{32}
$$

Therefore, K is the first integer for which $2Kh_{ex}$ is a multiple of N. The determination of K is then made thanks to the implementation of

$$
K = \frac{\text{lcm}(2h_{\text{ex}}, N)}{2h_{\text{ex}}}
$$
\n(33)

with $\text{lcm}(\bullet,\bullet)$ returns the lowest common multiple of the two arguments. Note that Eq. (33) can not be used when $h_{\text{ex}} = 0$, but $K = 1$ and only the nodal diameter 0 is involved in the response.

Based on this result, the updated set $\mathcal{D}_{q,0}$ must contains 0, h_{ex} , and all even multiples of h_{ex} up to 2K. However, these new nodal diameters can get coupled once again with h_{ex} and thus the set $\mathcal{D}_{q,0}$ must also contain all the uneven multiple of h_{ex} up to 2K. However thanks to Eq. (32), the remainder of the euclidian division of $(2K - l) h_{ex}$ by N is equal to the remainder of the euclidian division of lh_{ex} by N for $l \in [1, K]$. As a consequence, in order to determine the nodal diameter coupling when the initial set $\mathcal{D}_{q,0}$ contains the nodal diameter $h_{\rm ex}$, one must first determine K through Eq. (32) and then include all the multiples of h_{ex} up to K only. Then, the final set D is built following these three steps

1. the list $\mathbf{v}_{K,h_{ex}}$ of all multiples of h_{ex} up to K is firstly built such as

$$
\mathbf{v}_{K,h_{\text{ex}}}=h_{\text{ex}}\times\underbrace{[0,1,2,\ldots,q,\ldots,K]}_{\mathbf{v}_{K}}=[0,h_{\text{ex}},2h_{\text{ex}},\ldots,qh_{\text{ex}},\ldots,Kh_{\text{ex}}],
$$
\n(34)

- 2. then, the remainder of each element of $\mathbf{v}_{K,h_{\text{ex}}}$ by the euclidian division by N, noted $r_q \forall q \in \mathbf{v}_K$, is stored in the vector $\mathbf{v}_r = [r_0, r_1, r_2, \dots, r_q, \dots, r_K],$
- 3. finally, the nodal diameters associated with each r_q are computed and stored in $\mathcal{D} = [d_0, d_1, \ldots, d_q, \ldots, d_K]$ with $d_q = r_q$ if $r_q < m$ and $d_q = N - r_q$ if $r_q > m \quad \forall q \in \mathbf{v}_K$

Examples of application of this strategy are found throughout this article: each time a computation in cyclic symmetry is conducted with this criterion, the value of K and the final basis used is given.

In the case where the initial set $\mathcal{D}_{q,0}$ contains several nodal diameters, two cases arise. First, if the included nodal diameters share a common divider strictly greater than 1, then the previous procedure can be employed to

obtained a reduction by using $h_{\text{ex}} = \text{gcd}(h_{\text{ex}}^1, h_{\text{ex}}^2)$, where $\text{gcd}(\bullet, \bullet)$ returns the greater common divider of the two arguments. The other case corresponds to the situation where the nodal diameters are prime with one another $(\gcd(h_{\text{ex}}^1, h_{\text{ex}}^2) = 1)$. In this case, unfortunately all nodal diameters respond and no reduction can be achieved.

An additional case where no reduction can be achieved by this criterion is the configuration where N is a prime number. Then, no engine order (except the 0th) are subject to any reduction and full calculations should be made. This property is a consequence of the fact that $K = N$ when N is prime and it yields that $\mathcal{D} = \mathcal{D}^{\text{f}}$.

3 Numerical framework : harmonic balance method with nodal diameters selection

The resolution of Eq. (8) can be carried out indifferently with time domain or frequency domain methods. The determination of the set D does not make any assumption on the shape of the solution, thus the strategy presented in Sec. 2.4 is valid for all numerical strategies. In this work, the retained solving method for the application of the diameter selection strategy method is the HBM. Firstly the general formulation of the HBM is applied to Eq. (5), then a particular form of the HBM is used to take advantage of the selection of certain nodal diameters. A new criterion is also proposed in addition to the one in Sec. 2.4. It takes advantage of the coupling between nodal diameters and temporal harmonic to halve the size of the system. Finally, specific developments concerning the response to traveling wave excitations are carried out.

3.1 General harmonic balance method

The HBM is used to solve Eq. (5) for periodic solutions and allows fast solving and flexibility. It relies on the assumption that, under an excitation of angular frequency ω , the time-dependant unknown $\hat{\mathbf{x}}(t)$ is written as a truncated Fourier series of order N_h and fundamental angular frequency ω

$$
\widehat{\mathbf{x}}(t) \simeq \frac{1}{2}\widehat{\mathbf{a}}_0 + \sum_{k=1}^{N_{\rm h}} \left(\widehat{\mathbf{a}}_k \cos(k\omega t) + \widehat{\mathbf{b}}_k \sin(k\omega t) \right)
$$
(35)

This decomposition allows to perform a Fourier-Galerkin projection on Eq. (5) in order to obtain an algebraic system of equations related to cyclic coordinates in the frequency domain

$$
\mathbf{Z}(\omega)\hat{\tilde{\mathbf{x}}} - \hat{\tilde{\mathbf{f}}}_{\text{nl}}\left(\hat{\tilde{\mathbf{x}}}\right) - \hat{\tilde{\mathbf{f}}}_{\text{ex}} - \hat{\tilde{\mathbf{f}}}_{0} = \mathbf{0}.\tag{36}
$$

The new unknown of the problem is then the so-called multiharmonic vector $\hat{\tilde{\mathbf{x}}}$ containing the real Fourier decomposition of the spectral components of the displacement field $\mathbf{x}(t)$, written as in Eq. (37a). The vectors associated with external and nonlinear forces are also organized as multiharmonic vector which expressions are given in Eq. (37b), Eq. (37c) and Eq. (37d)

$$
\begin{cases} \n\hat{\mathbf{x}} = [\frac{1}{2}\hat{\mathbf{a}}_0^\top \ \hat{\mathbf{a}}_1^\top \ \hat{\mathbf{b}}_1^\top \dots \ \hat{\mathbf{a}}_{N_{\rm h}}^\top \ \hat{\mathbf{b}}_{N_{\rm h}}^\top]^\top,\n\end{cases} \tag{37a}
$$

$$
\begin{cases}\n\mathbf{x} = \left[\frac{1}{2}\mathbf{a}_0 \mathbf{a}_1 \mathbf{b}_1 \dots \mathbf{a}_{N_{\rm h}} \mathbf{b}_{N_{\rm h}}\right], & (37a) \\
\hat{\mathbf{f}}_{\rm nl} = \left[\frac{1}{2}(\hat{\mathbf{a}}_0^{\rm nl})^\top (\hat{\mathbf{a}}_1^{\rm nl})^\top (\hat{\mathbf{b}}_1^{\rm nl})^\top \dots (\hat{\mathbf{a}}_{N_{\rm h}}^{\rm nl})^\top (\hat{\mathbf{b}}_{N_{\rm h}}^{\rm nl})^\top \right],\n\end{cases}
$$
\n(37b)

$$
\begin{cases}\n\widehat{\widetilde{\mathbf{f}}}_{\text{ex}} = [\mathbf{0} (\widehat{\mathbf{a}}_1^{\text{ex}})^{\top} (\widehat{\mathbf{b}}_1^{\text{ex}})^{\top} \mathbf{0} \dots \mathbf{0} \mathbf{0}]^{\top}, \\
\widehat{\widetilde{\mathbf{f}}}_{0} = [(\widehat{\mathbf{a}}_0^0)^{\top} \mathbf{0} \dots \mathbf{0} \mathbf{0}]^{\top}.\n\end{cases}
$$
\n(37c)

$$
\widetilde{\mathbf{f}}_0 = [(\widehat{\mathbf{a}}_0^0)^\top \ \mathbf{0} \ \dots \ \mathbf{0} \ \mathbf{0}]^\top. \tag{37d}
$$

Z is a block-diagonal matrix accounting for the linear behavior of the system and is composed of $(N_h + 1)$ blocks, defined such as

$$
\mathbf{Z}_0 = \hat{\mathbf{K}} \text{ and } \mathbf{Z}_k(\omega) = \begin{bmatrix} \hat{\mathbf{K}} - (k\omega)^2 \widehat{\mathbf{M}} & k\omega \widehat{\mathbf{C}} \\ -k\omega \widehat{\mathbf{C}} & \hat{\mathbf{K}} - (k\omega)^2 \widehat{\mathbf{M}} \end{bmatrix} \quad \forall k \in [1, N_h].
$$
 (38)

3.2 Modifications induced by the selection of specific nodal diameters

However, the HBM allows a specific methodology that consists in selecting a subset of all spectral components, or nodal diameters, to reduce the size of the system that is actually solved. While in the time domain equation the restriction of the unknown to a set of nodal diameters $\mathcal D$ affects the whole equation in the same way, it is not necessarily the case for HBM. Indeed, since the blocks \mathbf{Z}_k of the dynamic stiffness matrix are independent, it is possible to use different subsets \mathcal{D}^k specific to each temporal harmonic k as a way of taking advantage of the spatio-temporal couplings that occur in cyclic structures. Then, these sets \mathcal{D}^k are used to build specific constitutive matrices $\widehat{\mathbf{M}}_k$, $\widehat{\mathbf{C}}_k$ and $\widehat{\mathbf{K}}_k$ and corresponding loading and unknown vectors of each kth harmonic. Then, Eq. (38) is rewritten such as

$$
\mathbf{Z}_0 = \widehat{\mathbf{K}}_0 \text{ and } \mathbf{Z}_k(\omega) = \begin{bmatrix} \widehat{\mathbf{K}}_k - (k\omega)^2 \widehat{\mathbf{M}}_k & k\omega \widehat{\mathbf{C}}_k \\ -k\omega \widehat{\mathbf{C}}_k & \widehat{\mathbf{K}}_k - (k\omega)^2 \widehat{\mathbf{M}}_k \end{bmatrix} \forall k \in [1, N_h].
$$
 (39)

with

$$
\widehat{\mathbf{M}}_k = \text{blkdiag}\left(\left\{\widehat{\mathbf{M}}^h\right\}_{h \in \mathcal{D}^k}\right), \widehat{\mathbf{C}}_k = \text{blkdiag}\left(\left\{\widehat{\mathbf{C}}^h\right\}_{h \in \mathcal{D}^k}\right), \widehat{\mathbf{K}}_k = \text{blkdiag}\left(\left\{\widehat{\mathbf{K}}^h\right\}_{h \in \mathcal{D}^k}\right) \tag{40}
$$

The different loading vectors of Eq. (7) should also be rewritten as

$$
\widehat{\mathbf{f}}_{\mathrm{ex},k} = \left[\left\{ \widehat{\mathbf{f}}_{\mathrm{ex}}^h \right\}_{h \in \mathcal{D}^k} \right] \qquad \text{and} \qquad \widehat{\mathbf{f}}_{0,k} = \left[\left\{ \widehat{\mathbf{f}}_0^h \right\}_{h \in \mathcal{D}^k} \right] \tag{41}
$$

depending on the temporal harmonic number k they're related to. Following these notations, the unknown vector is also reduced and defined as $\widehat{\mathbf{x}}_k = \left[\left\{\widehat{\mathbf{x}}^h\right\}_{h\in\mathcal{D}^k}\right]$ and $\widehat{\mathbf{f}}_{\text{nl},k} = \left[\left\{\widehat{\mathbf{f}}_{\text{nl}}^h\right\}_{h\in\mathcal{D}^k}\right]$ i .

If all subsets \mathcal{D}^k for $k \in [0, N_h]$ are $\mathcal{D}^k = \mathcal{D}^f$, then this strategy is equivalent to solving the full system with all core sectors.

In the literature, Petrov showed that for cyclic symmetric systems undergoing friction nonlinearities and traveling wave excitations, solution displacement fields can be searched as traveling waves of the same shape as the excitation [6]. This particular hypothesis allows to select only a single index in each of the \mathcal{D}^k , significantly reducing the computational costs without accuracy loss. Knowing the direction of propagation of the traveling wave also allows to divide the size of the system by two. The strategy used for nodal diameter selection and system condensation are explained in Sec. 3.3 in the case of traveling wave excitations.

At this point, it is possible to perform a condensation of the linear dof by taking advantage of the algebraic nature of the HBM equation and the fact that contact nonlinearities are localized [28]. Then, the system is restricted to nonlinear dof and the computations are significantly sped up. The reduced system of equations formulated on sets of nodal diameters \mathcal{D}^k finally reads

$$
\mathbf{Z}_{k,\mathbf{r}}\widehat{\widetilde{\mathbf{x}}}_{k,\mathbf{r}} - \widehat{\widetilde{\mathbf{f}}}_{\mathbf{n},k,\mathbf{r}} - \widehat{\widetilde{\mathbf{f}}}_{\mathbf{ex},k,\mathbf{r}} - \widehat{\widetilde{\mathbf{f}}}_{0,k,\mathbf{r}} = \mathbf{0} \quad \forall k \in [\![0, N_{\mathbf{h}}]\!]
$$
(42)

that is aggregated into a single equation

$$
\mathbf{Z}_{\mathrm{r}}\widehat{\widetilde{\mathbf{x}}}_{\mathrm{r}} - \widehat{\widetilde{\mathbf{f}}}_{\mathrm{nl},\mathrm{r}} - \widehat{\widetilde{\mathbf{f}}}_{\mathrm{ex},\mathrm{r}} - \widehat{\widetilde{\mathbf{f}}}_{0,\mathrm{r}} = \mathbf{0}.
$$
 (43)

This equation is solved by means of a Newton-Raphson nonlinear iterative solver. The jacobian matrix of Eq. (43) should thus be computed at each iteration.

3.2.1 Parity criterion for unilateral nonlinearities

For unilateral nonlinearities such as the one studied in Sec. 2.3 and Sec. 2.4, the procedure of selection of diameters harmonic-wise can help further reducing of the size of the problem. The additional criterion and its implementation are provided in this section, however the underlying mathematical developments are left in B.

The additional reduction consists in taking advantage of how the different harmonics of the nonlinear forces couple themselves. Indeed, when certain conditions described hereafter are met, it is observed that all diameters among the subset D determined in Sec. 2.4 do not respond on all harmonics. A distinction comes from the parity of the harmonic number: a subset \mathcal{D}^e responds only on even harmonics while a subset \mathcal{D}^o responds only on odd harmonics. The sets \mathcal{D}^e and \mathcal{D}^o respect the relations

$$
\mathcal{D}^{\mathbf{e}} \cup \mathcal{D}^{\mathbf{o}} = \mathcal{D} \text{ and } \mathcal{D}^{\mathbf{e}} \cap \mathcal{D}^{\mathbf{o}} = \emptyset. \tag{44}
$$

This means that all diameters in $\mathcal D$ are needed, but only a subset $\mathcal D^e$ or $\mathcal D^o$ can be selected according to the parity of the harmonic.

In order to know if this parity criterion can be used, one should examine the value of Kh_{ex} where the definition of K is given in Eq. (33) and h_{ex} is the nodal diameter of the excitation. Two cases arise:

if $Kh_{\text{ex}} \equiv 0$ [N]: no reduction is expected and all harmonics should use the whole D set such as

$$
\mathcal{D}^k = \mathcal{D} \qquad \forall k \in [0, N_h], \tag{45}
$$

if $Kh_{\text{ex}} \equiv \frac{N}{2}$ [N]: the reduction according to the parity of the harmonic number can be used and it comes that

$$
\int \mathcal{D}^k = \mathcal{D}^e \text{ if } k \text{ is even, } \quad \forall k \in \mathbb{I} \quad N \text{-} \mathbb{I}
$$
\n
$$
(46)
$$

$$
\begin{cases}\n\mathcal{D}^k = \mathcal{D}^{\text{o}} \text{ if } k \text{ is odd,} \\
\mathcal{D}^k = \mathcal{D}^{\text{o}} \text{ if } k \text{ is odd,}\n\end{cases} \tag{17}
$$

The sets \mathcal{D}^e and \mathcal{D}^o are built such as \mathcal{D}^e contains all d_q associated to even values of q in \mathbf{v}_K while \mathcal{D}^o is composed of all d_q associated to odd values of q in \mathbf{v}_K .

Overall, this parity criterion allows to further divide the size of the system by a factor 2 when the condition $Kh_{\text{ex}} \neq 0$ [N] is respected. It is noted that the demonstration of B shows that this condition can only occur when N is even.

Due to the necessity of splitting the diametral basis among different harmonics, this method only applies to an HBM solving process. Throughout the rest of the article, this criterion is used for the simulations when it is relevant and applicable.

3.3 Traveling wave excitations

In this section, the particular case of traveling wave excitation is dealt with. For this type of loading, further reductions are possible and the associated developments are provided. This idea was initially proposed by Petrov [6].

In the following, a traveling wave excitation of order h_{ex} of the form

$$
\mathbf{f}_{\text{ex},j}^{\text{TW}}(t) = f_{\text{ex}} \cos\left(t + \frac{\alpha(j-1)h_{\text{ex}}}{\omega}\right) \quad \forall j \in [\![1, N]\!]
$$
\n(48)

is applied to a cyclic system. The excitation is defined in the fixed frame such as it can model stator wakes. Then, assuming the bladed disk is rotating in the same direction as blades are numbered, the fixed frame excitation propagates backwards in terms of blade numbering: 1, N , $N - 1$, ...

In the case of such an excitation, the subsets \mathcal{D}^k should be chosen by reading the zig-zag diagram, such as

$$
\mathcal{D}^k = \{ \mathcal{Z}(kh_{\text{ex}}) \} \tag{49}
$$

where $\mathcal{Z}(h)$ is the function that stands for the zig-zag diagram. In the particular case of traveling waves, the spectral indexes in \mathcal{D}^k are signed, allowing to know the specific propagation direction of the traveling waves. This is necessary for further condensation of the problem. Zig-zag diagram are illustrated in Fig. 1 for both even and odd cases.

If there exists a spectral index $-m \leq q < 0$ such that $(q-h) \equiv 0$ [N] then $\mathcal{Z}(h) = q$, otherwise $\mathcal{Z}(h) = q$ where q respects $(q + h) \equiv 0 |N|$. Both conditions are exclusive and only a single spectral component $-m \le q \le m$ can be obtained by this method.

The selection of nodal diameters through Eq. (49) already yields a significant reduction in the system's size and computational burden as each harmonic only contains one specific nodal diameter. In addition, it is possible to go further in the reduction of the size of the system by imposing the direction of propagation of the different traveling

Figure 1. Examples of zig-zag diagrams, backward traveling wave $(-)$ $(q < 0)$, forward traveling wave $(-)$ $(q > 0)$.

waves. Traveling waves only make sense on so-called degenerated nodal diameters (size doubled with respect to the fundamental sector size), *i.e.* $0 < q < \frac{N}{2}$. So, this condensation is only useful on harmonics where $0 < |\mathcal{Z}(kh_{ex})| < \frac{N}{2}$. In this case, for a physical dof numbered d in the fundamental sector on harmonic $k > 0$ (harmonic 0 can only be composed of nodal diameter 0 in the definition of Eq. (49)) where $q = |Z(kh_{ex})|$, its displacements on this specific harmonic are given by

$$
\begin{cases} \hat{\mathbf{x}}_{k,c}^d = a_{k,q}^d \cos(k\omega t) + c_{k,q}^d \sin(k\omega t) \end{cases} \tag{50}
$$

$$
\hat{\mathbf{x}}_{k,s}^d = b_{k,q}^d \cos(k\omega t) + d_{k,q}^d \sin(k\omega t) \tag{51}
$$

as a function of the unknown vector $\hat{\tilde{\mathbf{x}}}^d = \left[a_{k,q}^d, b_{k,q}^d, c_{k,q}^d, d_{k,q}^d\right]^\top$. The general expression of a forward traveling wave is of the shape

$$
\int \widehat{\mathbf{x}}_{k,c}^d = a_{k,q}^d \cos(k\omega t) + c_{k,q}^d \sin(k\omega t)
$$
\n(52)

$$
\hat{\mathbf{x}}_{k,s}^d = -c_{k,q}^d \cos(k\omega t) + a_{k,q}^d \sin(k\omega t)
$$
\n(53)

where $b_{k,q}^d$ and $d_{k,q}^d$ are actually directly linked to $c_{k,q}^d$ and $a_{k,q}^d$. By imposing these links between the four coefficients, the direction of propagation of the wave is also imposed, but it yields a reduction in terms of computational cost. The expression of a backward traveling wave is given by

$$
\begin{cases} \widehat{\mathbf{x}}_{k,c}^d = a_{k,q}^d \cos(k\omega t) + c_{k,q}^d \sin(k\omega t) \end{cases} \tag{54}
$$

$$
\hat{\mathbf{x}}_{k,s}^d = c_{k,q}^d \cos(k\omega t) - a_{k,q}^d \sin(k\omega t). \tag{55}
$$

This new expression causes two signs to be swapped in Eq. (55), justifying the need to retrieve the signed nodal diameter when reading the zig-zag diagram. The so-called traveling wave hypothesis, relies on this link by substituting $\widehat{\widetilde{\mathbf{x}}}_k^d$ \bar{k} in the equations following the relation

$$
\widehat{\widetilde{\mathbf{x}}}_{k}^{d} = \mathbf{H}_{k}^{\mathrm{TW},1} \widehat{\widetilde{\mathbf{x}}}_{k}^{d,\mathrm{TW}} \tag{56}
$$

where

$$
\mathbf{H}_{k}^{\mathrm{TW},1} = \begin{bmatrix} 1 & 0 \\ 0 & \mathrm{sign}\left(\mathcal{Z}(kh_{\mathrm{ex}})\right) \\ 0 & 1 \\ -\mathrm{sign}\left(\mathcal{Z}(kh_{\mathrm{ex}})\right) & 0 \end{bmatrix} \tag{57}
$$

is the traveling hypothesis transform matrix for a single dof and $\hat{\tilde{\mathbf{x}}}_k^{d,\text{TW}} = \begin{bmatrix} a_{k,q}^d, c_{k,q}^d \end{bmatrix}^\top$. The transform matrix is extended to the size of the harmonic by performing a Kronecker product such as $H_k^{TW} = H_k^{TW,1} \otimes I_n$ and

$$
\widehat{\widetilde{\mathbf{x}}}_k = \mathbf{H}_k^{\mathrm{TW}} \widehat{\widetilde{\mathbf{x}}}_k^{\mathrm{TW}}.
$$
\n(58)

The harmonic balance equations corresponding to the kth harmonic

$$
\mathbf{Z}_{k}\widehat{\widetilde{\mathbf{x}}}_{k} - \widehat{\widetilde{\mathbf{f}}}_{\mathrm{nl},k} - \widehat{\widetilde{\mathbf{f}}}_{\mathrm{ex},k} - \widehat{\widetilde{\mathbf{f}}}_{0,k} = \mathbf{0},\tag{59}
$$

can then be rewritten using the change of basis Eq. (58) such as

$$
\left(\mathbf{H}_{k}^{\mathrm{TW}}\right)^{\top} \mathbf{Z}_{k} \mathbf{H}_{k}^{\mathrm{TW}} \widehat{\widetilde{\mathbf{x}}}_{k}^{\mathrm{TW}} - \left(\mathbf{H}_{k}^{\mathrm{TW}}\right)^{\top} \mathbf{H}_{k}^{\mathrm{TW}} \widehat{\widetilde{\mathbf{f}}}_{\mathrm{nl},k}^{\mathrm{TW}} - \left(\mathbf{H}_{k}^{\mathrm{TW}}\right)^{\top} \widehat{\widetilde{\mathbf{f}}}_{\mathrm{ex},k} - \left(\mathbf{H}_{k}^{\mathrm{TW}}\right)^{\top} \widehat{\widetilde{\mathbf{f}}}_{0,k} = \mathbf{0}.\tag{60}
$$

Following the definition of \mathbf{H}_k^{TW} , the product with its transposed is

$$
\left(\mathbf{H}_k^{\text{TW}}\right)^{\top} \mathbf{H}_k^{\text{TW}} = 2\mathbf{I}_{2n}.\tag{61}
$$

The Eq. (60) is then multiplied by $\frac{1}{2}$ as a way of including this factor in the matrices and vectors that define the problem rather than in the unknowns $\hat{\tilde{\mathbf{x}}}_k^{\text{TW}}$ and $\hat{\tilde{\mathbf{f}}}$ TW $_{\mathrm{nl},k}$. The final step of the reduction of the *k*th harmonic then yields

$$
\mathbf{Z}_{k}^{\text{TW}} \widehat{\widetilde{\mathbf{x}}}_{k}^{\text{TW}} - \widehat{\widetilde{\mathbf{f}}}_{\text{nl},k}^{\text{TW}} - \widehat{\widetilde{\mathbf{f}}}_{\text{ex},k}^{\text{TW}} - \widehat{\widetilde{\mathbf{f}}}_{0,k}^{\text{TW}} = \mathbf{0}.
$$
 (62)

with

$$
\mathbf{Z}_{k}^{\text{TW}} = \frac{1}{2} \left(\mathbf{H}_{k}^{\text{TW}} \right)^{\top} \mathbf{Z}_{k} \mathbf{H}_{k}^{\text{TW}}, \qquad \widehat{\widetilde{\mathbf{f}}}_{\text{ex},k}^{\text{TW}} = \frac{1}{2} \left(\mathbf{H}_{k}^{\text{TW}} \right)^{\top} \widehat{\widetilde{\mathbf{f}}}_{\text{ex},k} \quad \text{and} \quad \widehat{\widetilde{\mathbf{f}}}_{0,k}^{\text{TW}} = \frac{1}{2} \left(\mathbf{H}_{k}^{\text{TW}} \right)^{\top} \widehat{\widetilde{\mathbf{f}}}_{0,k} \tag{63}
$$

It is noted that, the determination of j^{\max} (see Sec. 2.2) for this case is immediate and a single sector is necessary for the evaluation of nonlinear forces: $j^{\text{max}} = 1$. Moreover, similarly as for the classical HBM, the condensation of linear dof [28] is also possible in Eq. (62) and allows for savings in computing time.

3.4 Evaluation of nonlinear forces

This section details the strategy retained for the evaluation of the nonlinear forces throughout the HBM resolution. Two different contact laws are used in the article and both rely on the alternating frequency/time procedure (AFT) [29]. A modification of the classical AFT scheme introduced to deal with the alternation between the cyclic and physical space is also proposed. These developments are presented in Sec. 3.4.1. Then the regularized penalty law (referred to as Regularized-HBM (R-HBM)) used for the analytical developments in Sec. 2.3 is presented in Sec. 3.4.2. A stiffer contact law relying on augmented lagrangian [27, 28], denominated Dynamic Lagrangian Frequency/Time (DLFT-HBM), is also presented in Sec. 3.4.3. Both contact laws constitute the state-of-the-art in terms of computing unilateral contact dynamics through HBM $[20, 28]$, especially for blade-tip/casing contacts $[11, 20, 30]$

3.4.1 Alternating Frequency/Time-Cyclic/Physical

In order to compute the nonlinear forces in a HBM framework, it is oftentimes necessary to resort to the AFT scheme [29]. In this work, a supplementary step is required since the HBM unknown is composed of spectral components. Indeed, the cyclic components should be transformed to the physical space for the computation of nonlinear forces and back to the cyclic space for the evaluation of the residual at each iteration. This new alternation type of is denominated Alternating Frequency/Time-Cyclic/Physical (AFT-CP).

The transformation between the frequency and time domain is done with by using the real inverse discrete Fourier transform (IDFT) matrix \mathcal{F}^{-1} so that $\{ \bullet (t_i) \}_{i=1...N_t} = \mathcal{F}^{-1} \widetilde{\bullet}$. Whereas, the transformation from the time domain to the frequency domain is made by means of the real discrete Fourier transform (D relation between both domains is $\widetilde{\bullet} = \mathcal{F}\{\bullet(t_i)\}_{i=1...N_t}$. A time period is sampled into N_t (user-defined parameter) instants. It should be chosen as a compromise between computation time and accuracy of the procedure: increasing N_t favors convergence and prevents aliasing problems but increases computational costs $[31]$.

In this work, the link between the cyclic and physical spaces is made in the frequency domain. It relies on \tilde{G} and \widetilde{G} that are built out of sub-matrices of the spatial Fourier transform matrix **S**. The construction of these matrices requires to determine in advance the sets \mathcal{S}^k (defined by the integers j_k^{\max}) associated with each \mathcal{D}^k following the rule presented in Sec. 2.2. In practice, since the evaluation of nonlinear forces in the time domain on a set of sectors requires all harmonics for all sectors, the set of sectors over which nonlinear forces are computed is actually

 $\mathcal{S}=\left(\begin{array}{c} N_{\rm h}\ \mid \end{array}\right)$ $k=0$ s^k . Then, the matrices ${\bf G}$ and ${\bf G}$ are defined as

$$
\widetilde{\mathbf{G}} = \text{blkdiag}\left(\widetilde{\mathbf{G}}^k\right) \otimes \mathbf{I}_{n_{\text{nl}}} \text{ with } \text{ and } \widetilde{\mathbf{G}} = \text{blkdiag}\left(\overline{\widetilde{\mathbf{G}}}^k\right) \otimes \mathbf{I}_{n_{\text{nl}}} \text{ with } k \in [0, N_{\text{h}}] \tag{64}
$$

where n_{nl} is the number of nonlinear dof of a single sector. Hereafter, the notations are consistent with what is defined in Sec. 2.2. The inverse transformation (from cyclic to physical variables) is straightforward and

$$
\overline{\widetilde{\mathbf{G}}}^{k} = \begin{bmatrix} \left(\mathbf{S}_{\mathcal{S}}^{\mathcal{D}^{k}}\right)^{\top} & \mathbf{0} \\ \mathbf{0} & \left(\mathbf{S}_{\mathcal{S}}^{\mathcal{D}^{k}}\right)^{\top} \end{bmatrix} \text{ if } k > 0 \text{ or } \widetilde{\mathbf{G}}^{k} = \left(\mathbf{S}_{\mathcal{S}}^{\mathcal{D}^{k}}\right)^{\top} \text{ if } k = 0. \tag{65}
$$

However, for the forward transformation, since S contains a larger number of sectors than each individual harmonic set \mathcal{S}^k , some of the information contained over sectors of S is redundant and should be truncated to switch from physical to cyclic variables. This truncation from S to S^k is made by the matrix \mathbf{R}_{S^k} defined as

$$
\mathbf{R}_{\mathcal{S}^k} = \begin{bmatrix} \mathbf{I}_{j_k^{\max}} & \mathbf{0} \end{bmatrix} \text{ if } \mathcal{S}^k \neq \mathcal{S} \quad \text{ and } \quad \mathbf{R}_{\mathcal{S}^k} = \mathbf{I}_{j_k^{\max}} \text{ if } \mathcal{S}^k = \mathcal{S}. \tag{66}
$$

Following this definition, it comes that the direct transformation \tilde{G} is built out of blocks \tilde{G}^k such as

$$
\widetilde{\mathbf{G}}^{k} = \begin{bmatrix} \left[\left(\mathbf{S}_{\mathcal{S}^{k}}^{\mathcal{D}^{k}} \right)^{\top} \right]^{-1} \mathbf{R}_{\mathcal{S}^{k}} & \mathbf{0} \\ \mathbf{0} & \left[\left(\mathbf{S}_{\mathcal{S}^{k}}^{\mathcal{D}^{k}} \right)^{\top} \right]^{-1} \mathbf{R}_{\mathcal{S}^{k}} \end{bmatrix} \text{ if } k > 0 \text{ or } \widetilde{\mathbf{G}}^{k} = \left[\left(\mathbf{S}_{\mathcal{S}^{k}}^{\mathcal{D}^{k}} \right)^{\top} \right]^{-1} \mathbf{R}_{\mathcal{S}^{k}} \text{ if } k = 0. \tag{67}
$$

Note that the block $\left[\left(\mathbf{S}_{\mathcal{S}^k}^{\mathcal{D}^k}\right)^{\top}\right]$ is invertible since sets \mathcal{S}^k and \mathcal{D}^k were chosen so that this sub-matrix of **S** is square. The distinction with respect to k comes from the presence of cosine and sine terms in real HBM for non-zero harmonic numbers k.

Now that all matrices are defined for the alternation between the different spaces, the AFT-CP methodology is presented hereafter. Say that an estimate of the solution is known in the shape of a multiharmonic vector corresponding to spectral components $\hat{\tilde{\mathbf{x}}}$, the overall alternating strategy is composed of five steps:

- 1. the vector $\hat{\tilde{\mathbf{x}}}$ is transformed to the physical space in the frequency domain by means of $\tilde{\mathbf{x}} = \overline{\tilde{\mathbf{G}}}\hat{\tilde{\mathbf{x}}}$. At this stage, $\tilde{\mathbf{x}}$ contains all real Fourier coefficients of the nonlinear dof of the system on sectors $j \in [1, j^{\max}]$ where $j^{\max} = \max$ $k \in \llbracket 0, N_{\rm h} \rrbracket$ $j_k^{\max},$
- 2. the multiharmonic vector of displacements is transformed to the time domain by using the IDFT matrix such that $\mathbf{x}(t) = \mathcal{F}^{-1}\tilde{\mathbf{x}}$ and $\dot{\mathbf{x}}(t) = \mathcal{F}^{-1}\nabla \tilde{\mathbf{x}}$ where ∇ is the frequency-domain derivation operator, nonlinear forces are evaluated in the time domain on physical sectors according to the ex-
- 3. nonlinear forces are evaluated in the time domain on physical sectors according to the expression of the nonlinearity $f_{nl}(\mathbf{x}(t), \dot{\mathbf{x}}(t)),$
- 4. the time-dependant nonlinear forces are transformed back to the frequency domain by using the DFT matrix such as $f_{\text{nl}} = \mathcal{F}f_{\text{nl}}(t),$

5. the frequency domain physical nonlinear forces are then used to compute the spectral components following $\widehat{\tilde{\mathbf{f}}}_{\text{nl}} = \widetilde{\mathbf{G}} \widetilde{\mathbf{f}}_{\text{nl}}.$

This type of alternation actually nests the classical AFT from steps 2 to 4. Steps 1 and 5 are added in order to perform the back and forth transformations between cyclic and physical domains. The whole procedure is summed up in Fig. 2. For the computation of the jacobian matrices of the nonlinear forces for this alternating procedure, it is necessary to apply the following formula

$$
\frac{\partial \widetilde{\mathbf{f}}_{\rm nl}}{\partial \widetilde{\mathbf{x}}} = \frac{\partial \widetilde{\mathbf{f}}_{\rm nl}}{\partial \widetilde{\mathbf{f}}_{\rm nl}} \frac{\partial \mathbf{f}_{\rm nl}}{\partial \mathbf{x}} \frac{\partial \mathbf{x}}{\partial \widetilde{\mathbf{x}}} \frac{\partial \widetilde{\mathbf{x}}}{\partial \widetilde{\mathbf{x}}} + \frac{\partial \widetilde{\mathbf{x}}}{\partial \widetilde{\mathbf{x}}} \frac{\partial \widetilde{\mathbf{f}}_{\rm nl}}{\partial \widetilde{\mathbf{x}}} \frac{\partial \mathbf{f}_{\rm nl}}{\partial \widetilde{\mathbf{x}}} \frac{\partial \widetilde{\mathbf{x}}}{\partial \widetilde{\mathbf{x}}} \frac{\partial \widetilde{\mathbf{x}}}{\partial \widetilde{\mathbf{x}}} \tag{68}
$$

which finally yields

$$
\frac{\partial \hat{\tilde{\mathbf{f}}}_{\text{nl}}}{\partial \hat{\tilde{\mathbf{x}}}} = \tilde{\mathbf{G}} \mathcal{F} \frac{\partial \mathbf{f}_{\text{nl}}}{\partial \mathbf{x}} \mathcal{F}^{-1} \overline{\tilde{\mathbf{G}}} + \tilde{\mathbf{G}} \mathcal{F} \frac{\partial \mathbf{f}_{\text{nl}}}{\partial \dot{\mathbf{x}}} \mathcal{F}^{-1} \nabla \overline{\tilde{\mathbf{G}}}
$$
(69)

Figure 2. Alternating Frequency/Time-Cyclic/Physical procedure, $c_j^* = \sqrt{(a_j^*)^2 + (b_j^*)^2}$.

3.4.2 Regularized penalty law (R-HBM)

The use of a regularized penalty law with the HBM for the characterization of nonlinear mechanical systems featuring vibro-impact was recently shown to provide excellent results by Colaïtis and Batailly [20]. Due to its regularized nature, it allows to perform a large variety of analyses that are usually restricted to smooth problems. Even if the regularization parameter affects the simulated dynamics of the system, a good agreement with state-of-the-art strategies is oftentimes reached $[11]$. For each nonlinear node j, the nonlinear force in the normal direction is given by

$$
f_{\mathrm{nl},j}^{\mathrm{N}}(t) = -\kappa \frac{g_j(t)}{2} + \sqrt{\left(\kappa \frac{g_j(t)}{2}\right)^2 + \gamma^2} \tag{70}
$$

where $g_j(t) = g_0(t) - x_j^N(t)$ is the gap function with $g_0(t)$ a periodic function that defines the evolution of the initial gap along a period. $x_j^N(t)$ is the displacement of the jth node in the normal direction. This function allows to model casings of specific shapes when considering rotating systems and rotor/stator interactions. Since this formula is purely analytical, the use of the AFT-CP as presented in Sec. 3.4.1 is directly possible. For details about the computation of the jacobian matrix of the nonlinear forces in the frequency domain, one should refer to the work of Colaïtis [20].

3.4.3 Augmented lagrangians (DLFT-HBM)

A second strategy for the evaluation of normal contact forces is used throughout this article. The previously published DLFT-HBM [27, 28] relies on a prediction-correction scheme that is presented in this section. The underlying mathematical background of this method is outside of the scope of this paper but it can be found in the literature [27, 28].

The prediction of nonlinear forces, superscripted \bullet^p , relies on the HBM equation formulated in the cyclic domain Eq. (43) and it reads

$$
\widehat{\widetilde{\mathbf{f}}}_{n l,r}^{p,N} = \left[\mathbf{Z}_r \widehat{\widetilde{\mathbf{x}}}_r - \widehat{\widetilde{\mathbf{f}}}_{ex,r} - \widehat{\widetilde{\mathbf{f}}}_{0,r} \right]^N - \varepsilon \widehat{\widetilde{\mathbf{g}}}_{\mathbf{x}} \tag{71}
$$

where \tilde{g}_x is a multiharmonic cyclic vector standing for the normal gap function $g_j(t) = g_{0,j} - x_j^N(t)$ of each contact
node $j \in S$ weighting parameter responsible for the respect of the unilateral contact conditions. T node j, ε a weighting parameter responsible for the respect of the unilateral contact conditions. The right term of Eq. (71) is restricted to dof of the normal direction only, indicated by the notation $\lbrack \bullet \rbrack^N$. Like for the AFT-CP β procedure, β p,N $n_{\text{nl,r}}$ should be firstly transformed to the frequency-physical domain and then to the time-physical domain, such as

$$
\mathbf{f}_{\mathrm{nl},\mathrm{r}}^{\mathrm{p},\mathrm{N}} = \mathcal{F}^{-1} \overline{\widetilde{\mathbf{G}}}\widetilde{\mathbf{f}}_{\mathrm{nl},\mathrm{r}}^{\mathrm{p},\mathrm{N}}.
$$

This vector encompasses the predicted physical normal contact forces applied to the contact nodes belonging to all sectors in S in the time domain. This prediction is then corrected by looping over all node numbers j and time instants i. Two cases arise:

Separation if $f_{\text{nl},j}^{\text{p},\text{N}}(t_i) > 0$, the *j*th node is separated at time t_i and $f_{\text{nl},j}^{\text{N}}(t_i) = 0$,

Contact if $f_{\text{nl},j}^{\text{p},\text{N}}(t_i) < 0$, the *j*th node is in contact at time t_i and $f_{\text{nl},j}^{\text{N}}(t_i) = f_{\text{nl},j}^{\text{p},\text{N}}(t_i)$.

After looping on all nodes and all instants, the contact forces are aggregated into a single vector $f_{nl,r}^N$ and transformed back to cyclic-frequency variables such as

$$
\widehat{\widetilde{\mathbf{f}}}_{\mathrm{nl,r}}^{\mathrm{N}} = \widetilde{\mathbf{G}} \mathcal{F} \mathbf{f}_{\mathrm{nl,r}}^{\mathrm{N}}.
$$
\n(73)

At this stage, the nonlinear forces are finally used to compute the residual of Eq. (43). The jacobian matrix of the nonlinear forces should also be provided to the nonlinear solver for the resolution to advance, such developments are not included in the article for the sake of brevity but can be found in the literature [28, 32].

The DLFT-HBM strategy is, unlike the R-HBM, a nonsmooth strategy which makes it more numerically sensitive but also more accurate. The use of this second method to evaluate the contact forces allows to put to the test the criterions derived in Sec. 2.4 and Sec. 3.2.1. Then, this article aims to demonstrates that the criterions not only work for a regularized contact law, but that they can also be extended to the nonsmooth and non-differentiable contact laws that approximate unilateral contact.

4 Results on a simplified bladed disk model

This section presents the results obtained on a simplified model composed of a cyclic chain of oscillators that undergo unilateral contacts. This model is approximating a bladed disk that experiences blade-tip/casing contacts. The demonstration of the nodal diameters selection method proposed in this article is made on several blade numbers N as well as several loading scenarios for both the regularized and the nonsmooth contact law. In all cases, the arc-length continuation procedure [29] is employed to compute the forced responses that feature numerous turning points.

4.1 Presentation of the model

The model of simplified bladed disk is adapted from the literature [33]. Each sector is composed of 5 masses linked with stiffnesses. A single sector of this model is illustrated in Fig. 3. The links between the different sectors are made through the stiffnesses between m_5^j and m_4^{j+1} where j denotes the sector number. All the stiffnesses and

Figure 3. Simplified model of a bladed disk (sector j), adapted from Ref. [33].

masses of a sector are given in Tab. 1.

m (kg)	J.2	0.3			
$k(10^6 N \cdot m^{-1})$			40	U.U	50

Table 1. Structural parameters of a sector.

The excitation force on each sector is applied to the mass m_1 that models the tip of the blade. Modal damping is used such that all modes of the system are damped with the same modal damping ratio of $\xi = 1 \cdot 10^{-3}$. For this model, no pre-loading is applied to the system: $f_0 = 0$. The number of sectors of the system is parameterized by the variable N allowing to study the dynamics of the system with different number of sectors. This allows to validate the rule derived in Sec. 2.4 for multiple values of N.

For the HBM simulations, the number of harmonics is set to $N_h = 10$ and the number of time instants is $N_t = 10000$. For R-HBM simulations, $\kappa = 1 \cdot 10^8 \,\mathrm{N \cdot m^{-1}}$ and $\gamma = 2.5 \,\mathrm{N}$. For DLFT-HBM simulations, a wall of stiffness $\kappa = 1 \cdot 10^8 \,\mathrm{N \cdot m^{-1}}$ is added (in order to be consistent with R-HBM) and $\varepsilon = 1 \cdot 10^9 \,\mathrm{N \cdot m^{-1}}$. Both configurations are displayed in Fig. 3a and Fig. 3b respectively.

4.2 Contact scenario

The nonlinear dof are all the masses m_1^j with $j \in [1, N]$ that are impacting a wall at a given distance $g_{0,j}(t)$. Unless
otherwise mentioned, the gap function is constant along the period with $g_{\alpha,j}(t) = g_{\alpha} = 1$, 10 otherwise mentioned, the gap function is constant along the period with $g_{0,i}(t) = c_i = 1 \cdot 10^{-3}$ m for all sectors, making the casing purely circular like displayed in Fig. 4a. In some contact scenarios, the casing is chosen to be a rigid diametral shape, such that

$$
g_{0,j}(t) = c_j + p_j \cos\left(h_c \Omega \left(t - j\frac{2\pi}{N\Omega}\right)\right) \tag{74}
$$

where c_j is the mean value of the gap function of node j, p_j is the amplitude of oscillation of the gap function around c_j , Ω is the rotation speed (synchronous with the excitation frequency ω) and h_c is the the nodal diameter index of the casing. An example of diametral casing is displayed in Fig. 4b as an illustration for $h_c = 5$. In order

Figure 4. Casing shapes and contact scenario.

to initiate contact interactions, the system is also excited for several engine orders by traveling or standing waves excitations. The definition of a traveling wave excitation of order h_{ex} is Eq. (48) and a standing wave excitation is the sum of two traveling wave excitations of the same order traveling in the opposite directions, such as

$$
\mathbf{f}_{\text{ex},j}^{\text{SW}}(t) = f_{\text{ex}} \left[\cos \left(t + \frac{\alpha (j-1) h_{\text{ex}}}{\omega} \right) + \cos \left(t - \frac{\alpha (j-1) h_{\text{ex}}}{\omega} \right) \right] \quad \forall j \in [\![1, N]\!]. \tag{75}
$$

4.3 Even number of sectors: $N = 24$

The first application chosen for the presented methodologies is the case study of Fig. 3 with $N = 24$ sectors. Since N is even, this case is suitable for the presentation of the parity criterion for unilateral nonlinearities in HBM (see Sec. 3.2.1). Following this criterion, the expected reduction in terms of total size and number of sectors to evaluate inside the AFT-CP procedure are presented in Tab. 2.

engine order	$\overline{0}$		$\overline{2}$	3	4	5	6	7	8	9	10	11	12
value of K		12	6	4	3	12	$\overline{2}$	12	3	$\overline{4}$	6	12	
reduced size	21	252	126	84	63	252	42	252	63	84	126	252	21
sectors for AFT-CP		12	6	4	3	12	$\overline{2}$	12	3	$\overline{4}$	6	12	
reduction $(\%)$	95.8	50.0	75.0	83.3	87.5	50.0	91.7	50.0	87.5	83.3	75.0	50.0	95.8

Table 2. Reduction for the case of the simplified system with $N = 24$ sectors for 10 harmonics (full size: $N \times n_{\text{nl}} \times (2N_{\text{h}}+1) = 504$ unknowns).

For this case, almost all engine orders benefit from the the criterion on the parity of Sec. 3.2.1. Only $h_{\rm ex} = 8$ does not fall into this category: $K = 3$ and $Kh_{ex} \equiv 0$ [N], thus, no reduction is achieved on the parity of the harmonics. Overall, the mean over all engine orders of the expected reduction is 75%. This number is significantly impacted by the parity criterion and would be of only 51.3% without it. Note that reductions presented in Tab. 2 do not take into account the type of excitation that is applied to the system. Indeed, by using the traveling wave hypothesis formulated in Sec. 3.3, the achieved reduction is higher. As a reminder, when considering the traveling wave hypothesis, the size of the system (and the AFT-CP) is brought down to the size of a single sector, yielding a reduction of 95.8% for any engine orders. Note that the value of K is systematically equal to the numbers of sectors needed for the calculation. This means that K has a physical meaning related to a relation of symmetry of along the blades of the cyclic system.

For this case study, as a way of demonstrating all the reduction capacities of the HBM method for different loading scenarios, several types of calculations are made. Computations are carried out for traveling wave excitations in Sec. 4.3.1, for standing wave excitations in Sec. 4.3.2 and for a mixed contact scenario involving a traveling wave excitation and a casing deformation in Sec. 4.3.3. For all cases, the result obtained with the proposed reduction is always compared to the full calculations (with all sectors) that serve as reference. The latter is directly built by assembling all sectors without the use of cyclic symmetry.

4.3.1 Traveling wave excitation

In this section, traveling wave excitations are applied to the system. Two spatial excitation orders are retained for this example: $h_{\text{ex}} = 1$ and $h_{\text{ex}} = 3$. The so-called traveling wave hypothesis (see Sec. 3.3) is used to perform all the calculations in cyclic symmetry in order to speed up the calculations. The calculations are carried out with the basis \mathcal{D}^k that are chosen by reading the zig-zag diagram. It comes that the lists of \mathcal{D}^k used for computations in cyclic symmetry is

$$
[\mathcal{D}^k]_{k \in [0, N_h]} = [\{0\}, \{-1\}, \{-2\}, \{-3\}, \{-4\}, \{-5\}, \{-6\}, \{-7\}, \{-8\}, \{-9\}, \{-10\}] \tag{76}
$$

for $h_{\text{ex}} = 1$ and

$$
[\mathcal{D}^k]_{k \in [\![0,N_h]\!]} = [\{0\}, \{-3\}, \{-6\}, \{-9\}, \{12\}, \{9\}, \{6\}, \{3\}, \{0\}, \{-3\}, \{-6\}] \tag{77}
$$

for $h_{\rm ex} = 3$. The results for both engine orders are displayed in Fig. 5, for the regularized contact law as well as the lagrangian contact law. The forcing amplitude is $f_{\text{ex}} = 4 N$ for $h_{\text{ex}} = 1$ and $f_{\text{ex}} = 1.75 N$ for $h_{\text{ex}} = 3$.

The forced responses in cyclic symmetry with the traveling wave hypothesis match perfectly with the full calculations for both contact laws and both engine orders. This shows that the reduced calculation allows to retrieve

Figure 5. Traveling wave excitation on the simplified bladed disk with $N = 24$ sectors for various engine orders, linear (\ldots ,), R-HBM: cyclic symmetry $(-)$, full $(--)$, DLFT-HBM: cyclic symmetry $(-)$, full $(--)$.

the same dynamics as the full model at the expense of 4.2% of the computational effort. This type of reduction was already demonstrated by Petrov for friction nonlinearities [6], but the application to systems featuring unilateral contact nonlinearities allows to extend the scope of the traveling wave hypothesis. As a way of demonstrating that the choice of the diametral basis Eq. (76) and Eq. (77) is appropriate, the spatial decomposition of the resonance point of the full calculations are analyzed and displayed in Fig. 6. As a way of analyzing the direction of propagation of the waves, the forward components are displayed in Fig. 6a and the backward components are displayed in Fig. 6b. Note that non-degenerated nodal diameters (here 0 and 12) are displayed as half forward and half backward components since they correspond to standing waves. The spatial content of each harmonic is analyzed independently by a spatial Fourier analysis and the magnitude of the components are displayed as a bar plot in logarithmic scale. The analysis is made on the stiffer contact law, *i.e.* the full DLFT-HBM calculation in Fig. 5b for $h_{ex} = 1$.

The cyclic contributions displayed in Fig. 6 are totally inline with what was predicted through the zig-zag diagram in Eq. (76). This justifies *a posteriori* the choice of the cyclic decomposition for the reduced system. A slight discrepancy appears on harmonic 0 where a standing ND10 appears. However, it is of a negligible magnitude compared to the predicted ND0 magnitude. Though, this contribution falls into the more general criterion of Sec. 2.4 and appears due to the coupling in the nonlinearity. As a secondary validation, the same type of analysis is made in Fig. 7 on the resonance of Fig. 5d for the traveling wave of order $h_{\text{ex}} = 3$.

The cyclic decomposition of the solution at the resonance reveals the same behavior as for $h_{\text{ex}} = 1$ in Fig. 6: the basis predicted through the zig-zag diagram in Eq. (77) is perfectly recovered on the full calculation. The

(b) Backward traveling diameters

Figure 6. Cyclic components over all harmonics at the resonance point of the full calculation in Fig. 5b, traveling wave with $h_{\rm ex} = 1$. Nodal diameter: 0 (**iii)**, 1 (**iii)**, 2 (**iii)**, 3 (**iii)**, 4 (**iii)**, 5 (**iii)**, 6 (**iiii)**, 7 (**iiii)**, 8 (**iiii)**, 9 (**iiii)**.

(b) Backward traveling diameters

Figure 7. Cyclic components over all harmonics at the resonance point of the full calculation in Fig. 5d, traveling wave with $h_{\rm ex} = 3.$ Nodal diameter: 0 (\Box), 3 (\Box), 6 (\Box), 9 (\Box), 12 (\Box).

same behavior as for the harmonic 0 of Fig. 6 happens but is explained the same way. Indeed, the standing ND6 contribution can be attributed to a coupling in the nonlinearity, but remains negligible with respect to the ND0 contribution.

4.3.2 Standing wave excitation

Now that the behavior of the system was demonstrated under traveling wave excitations, simulations are carried out on the nonlinear dynamics of the system under a standing wave excitation. This type of excitation allows to observe how the nonlinearity couples the different nodal diameters together. Hence, it is possible to verify that the analytical developments made in Sec. 2.4 and Sec. 3.2.1 are relevant and actually are exact reductions of the problem. Standing wave excitations are applied to the system for $h_{\text{ex}} = 1$ with $f_{\text{ex}} = 2$ N, and for $h_{\text{ex}} = 4$ with $f_{\text{ex}} = 1$ N.

Following the rules of determination of the nodal diameters to account for in each harmonic, the basis chosen are

$$
\mathcal{D}^{2k}|_{k \in [0,5]} = \{0, 2, 4, 6, 8, 10, 12\} \text{ and } \mathcal{D}^{2k-1}|_{k \in [1,5]} = \{1, 3, 5, 7, 9, 11\}
$$
\n
$$
(78)
$$

for $h_{\text{ex}} = 1$ ($K = 12 \Rightarrow Kh_{\text{ex}} = 12 \equiv \frac{N}{2}$ [N]) and

$$
\mathcal{D}^{2k}|_{k \in [0,5]} = \{0,8\} \text{ and } \mathcal{D}^{2k-1}|_{k \in [1,5]} = \{4,12\} \tag{79}
$$

for $h_{\text{ex}} = 4$ ($K = 3 \Rightarrow Kh_{\text{ex}} = 12 \equiv \frac{N}{2}$ [N]). The results of the different HBM computations are then displayed for all sectors in Fig. 8 for both contact laws and values of h_{ex} .

Once again, a perfect match is observed between the full calculations and the reduced calculations is cyclic symmetry. The standing wave nature of the response of the system can be seen to the presence of curves of different amplitude (linear and nonlinear), related to the level of forcing that a given sector is actually experiencing in the standing wave excitation. Similarly as for the case of traveling wave excitations, the cyclic contributions to the displacement field at the resonance for the full DLFT calculation is displayed in Fig. 9 for $h_{\text{ex}} = 1$ and in Fig. 10 for $h_{\text{ex}}=4.$

Figure 8. Standing wave excitation on the simplified bladed disk with $N = 24$ sectors for various engine orders, linear (......), R-HBM: cyclic symmetry $(-)$, full $(--)$, DLFT-HBM: cyclic symmetry $(-)$, full $(--)$.

Figure 9. Cyclic components over all harmonics at the resonance point of the full calculation in Fig. 8b, standing wave with $h_{\rm ex} = 1.$ Nodal diameter: 0 (**m**), 1 (**m**), 2 (**m**), 3 (**m**), 4 (**m**), 5 (**m**), 6 (**m**), 7 (**m**), 8 (**m**), 9 (**m**), 10 (**m**), 11 (**m**), $12 \left(\Box \right)$

For the case of $h_{\text{ex}} = 1$ of Fig. 9, the basis was correctly identified *a priori* in Eq. (78). This case demonstrates that even if no multiplicity is observed between h_{ex} and N, the HBM can take advantage of the criterion on the parity of the harmonics to still reduce the computational burden of 50%. This can be observed in Fig. 9 where all ND intervene in the whole solution, however one half only responds on even harmonics and the other half on odd harmonics. The sampling strategy of AFT-CP is then able to perform the computations on only half of the sectors, yielding a significant gain in terms of computational time. This type of strategy would not be possible for time-integration strategies since all cyclic components would have been needed to compute the response.

(b) Backward traveling diameters

Figure 10. Cyclic components over all harmonics at the resonance point of the full calculation in Fig. 8d, standing wave with $h_{\rm ex} = 4.$ Nodal diameter: 0 (\Box), 4 (\Box), 8 (\Box), 12 (\Box)

Similarly, the cyclic contributions responding in Fig. 10 for the case $h_{\text{ex}} = 4$ were correctly identified in Eq. (79) and the split among even and odd harmonics is also highlighted. For this couple (N, h_{ex}) a reduction of 87.5% is reached, and the results are strictly equivalent to the full model. This section has demonstrated the relevance of the method used to determine the nodal diameter couplings as a way of reducing the computational effort.

For both Fig. 10 and Fig. 9, one can note that the response is purely stationary, meaning that forward and backward contributions are strictly equal. Hence, when such responses emerge (standing wave excitations) in the rest of the article, only the first graph is displayed for the sake of brevity.

4.3.3 Mixed contact scenario

Finally, the system's response is analyzed under a loading scenario that involves both the excitation through a traveling wave and also a nodal diameter shape of the surrounding casing, as presented in Fig. 4b. When considering them independently, the excitation and the casing both induce a traveling wave response in cyclic system. However, if $h_{\text{ex}} \neq h_{\text{c}}$ a coupling between the two types of loading happens and a partly standing wave response is expected to emerge. This scenario could not have been studied without a specific criterion in cyclic symmetry.Indeed, Petrov's method [6] is restricted to a single traveling wave loading.

The chosen configuration is a case where the basis associated to h_{ex} and h_c are different. Indeed, trivial cases occur when one of the two basis is included in the other $(\mathcal{D}_{ex}^k \subset \mathcal{D}_{c}^k$ or $\mathcal{D}_{c}^k \subset \mathcal{D}_{ex}^k)$. In the non trivial case, one should consider the basis that is formed by the greatest common divisor (GCD) of $h_{\rm ex}$ and $h_{\rm c}$. In order to demonstrate this property, a case with $h_{\rm ex} = 6$ and $h_{\rm c} = 8$ is studied. The basis for each of the two are respectively

$$
\mathcal{D}^{2k}|_{k \in [0,5]} = \{0,12\} \text{ and } \mathcal{D}^{2k-1}|_{k \in [1,5]} = \{6\}
$$
\n
$$
(80)
$$

for the excitation

$$
\mathcal{D}^{k}|_{k \in [0,10]} = \{0,8\}.
$$
\n(81)

and for the casing, no parity criterion exists for an excitation over 8ND. In order to cover all the possible couplings that can happen, one should chose the basis that would emerge from a 2ND excitation, such as

$$
\mathcal{D}^{k}|_{k\in\llbracket 0,10\rrbracket} = \{0,2,4,6,8,10,12\}.
$$
\n(82)

Note that this basis not only contains the nodal diameters in Eq. (80) and Eq. (81) but also supplementary nodal diameters that are expected to appear through the coupling of both basis (4ND and 10ND). Also, the parity criterion has not been used for this expression. The rule of parity is expected to play a role in the response for this case but more developments are needed in order to generalize it when two loadings are applied to the system. In the current case, the casing has a diametral shape of order $h_c = 8$ and the rotation is considered to be synchronous with the excitation, hence the casing excitates the 8th harmonic. Then, in order to derive if the parity criterion applied, one should compute the couplings between a loading with a fundamental on the 2ND first harmonic $(e^{i\omega t})$ and a loading with a fundamental on 8ND on the 8th harmonic $(e^{8i\omega t})$ but it exceeds the scope of this paper. Since the detection of the parity criterion can be prone to error when two loadings are applied to the system, the conservative basis Eq. (82) is used. The results of the simulations are displayed in Fig. 11 for both DLFT-HBM and R-HBM on two sectors.

Figure 11. Traveling wave excitation on the simplified bladed disk with $N = 24$ sectors for $h_{\text{ex}} = 6$ and a casing deformed with $h_c = 8$, linear (.......), R-HBM: cyclic symmetry (-), full (---), DLFT-HBM: cyclic symmetry (-), full (---).

Firstly, the forced responses obtained in cyclic symmetry match the full calculations for both methods. The responses of DLFT-HBM and R-HBM are not exactly identical because of the difference in contact treatment. This configuration being a highly nonlinear very sensitive case study, it highlights the discrepancies. The standing wave nature of the displacement field can be seen by the fact that all sectors do not experience the same response. This fact can be seen more clearly in Fig. 12 where the cyclic components are displayed.

The decomposition of the solution at the resonance in terms of cyclic components reveals that the basis Eq. (82), emerging from the coupling of the excitation and the casing, was correctly identified *a priori*. Indeed, the only nodal diameters that are involved in the response are in the basis. Moreover, as discussed beforehand, a parity criterion is observed on the response: nodal diameters $\{0, 4, 8, 12\}$ respond on even harmonics while nodal diameters $\{2, 6, 10\}$ intervene only on odd harmonics. Detecting the coupling of the casing and the excitation in terms of parity would have then allowed to divide the computational effort by two in this particular case. However, the general developments in order to detect such couplings are outside the scope of this article. Moreover, it is observed that

(b) Backward traveling diameters

Figure 12. Cyclic components over all harmonics at the resonance point of the full calculation in Fig. 11, traveling wave with hex = 6 and casing with h^c = 8. Nodal diameter: 0 (), 2 (), 4 (), 6 (), 8 (), 10 (), 12 ()

the couplings of Eq. (80) and Eq. (81) make appear ND4 and ND10 contributions in the simulations even though it is in neither of the two basis, justifying that the basis should be built according to the GCD of $h_{\rm ex}$ and $h_{\rm c}$ in order to cover all possible couplings.

The test case studied in depth here was featuring $N = 24$ sectors and was prone to large reductions in all types of loading scenarios. Indeed, N being even ensures that all even values of h_{ex} will provide a significant reduction of around at least 50%.

This section compared full calculations with the reduced methodology proposed in this article. The results are in perfect agreement and thus these numerical simulations give further credibility to the theoretical demonstration. Notice that the numerical simulations showed that the reduction strategy also works for non-regularized contact law. The arbitrary case of $N = 24$ was dealt with here, but the proposed methodology can be applied for a wide number of cases. In order to cover all configurations, an odd number of sectors of $N = 45$ is also dealt with in C. The latter is dealt with for the sake of completeness as the reduction rates are quite lower for systems with odd numbers of sectors.

5 Results on an industrial model

Now that the reduction capabilities of the proposed strategy was extensively demonstrated on academic examples, the study is taken further to tackle an industrial size realist configuration. The same range of loading scenarios is applied to the system in order to show the applicability of the method to large scale systems featuring blade-tip/casing contacts.

5.1 Presentation of the model

The model chosen to demonstrate the capabilities of the strategy is the open test case fan stage ECL5/Catana designed at Ecole Centrale de Lyon [34] to be a step towards more representative open numerical models of ultrahigh bypass ratio architectures. The stage is composed of 16 blades built out of unidirectional carbon fibers and epoxy composite plies. The disk is made of titanium. This geometry was recently characterized in terms of aerodynamics [35], acoustics [36], aeroelastics [37], and structural dynamics [38], from both experimental and numerical perspectives. The full finite-element mesh is composed of 87,310 nodes for a total of 19,642 elements. It is displayed in Fig. 13a for information. Both the material data and the boundary conditions are those given in Ref. [34]. The system is considered in cyclic symmetry and the cyclic boundaries, as long as other boundary conditions, are displayed in Fig. 13b. In this work, the contacts between the blades and the disk are linearized and no relative motion is possible.

In order to decrease the computational costs of simulating the dynamics of this system, a Craig-Bampton reduced-order modeling technique [39] is applied to the system for each nodal diameter. A total of 11 boundary nodes are retained in the reduced-order model: $n_{nl} = 9$ boundary nodes are kept at the tip of the blade at the intrados side for contact treatment purposes, one node is kept at the blade-tip at mid chord in order to apply the excitation and one observation node is retained in order to monitor the bending amplitude away from the nonlinear

Figure 13. ECL5/Catana fan stage model, cyclic symmetry boundary (**iii**), clamped (**iii**), axial blockage (**iii**), nonlinear nodes (\Box) , observation node for NFRC (\Box) , excitation node (\Box) .

interface. The contact nodes are evenly spaced along the blade-tip. The displacement of the observation node in the circumferential direction is systematically used throughout the article in order to plot the forced response curves. 20 internal modes are retained for the Craig-Bampton procedure. Since the system is formulated using the cyclic symmetry, the number of unknowns associated to non-degenerated nodal diameters (ND0 and ND8) is 53 and 106 for all other diameters. Modal damping is used to build the damping matrix. The first five modes are damped with a ratio $\xi_{1-5} = 1 \cdot 10^{-3}$ and the remaining modes with a ratio of $\xi_{5+} = 5 \cdot 10^{-3}$. The mode of interest in this study is the first bending mode. The eigenfrequencies of the first bending mode along all nodal diameters are given in Tab. 3.

nodal diameter			∼						
(Hz) frequency	263.2	284.1	285.5	285.6	286.1	286.8	287.4	287.7	287.9

Table 3. Eigenfrequencies of the first bending mode along all nodal diameters.

Given the number of sectors of the ECL5 fan stage of $N = 16$, the expected reductions with the proposed methodology are provided in Tab. 4. Since N is even, it is subject to the parity criterion, and parity reduction is predicted along all nodal diameters.

5.2 Contact scenario

In this article, the dynamics of the casing at the contact interface is neglected. The casing is considered to be a rigid mathematical profile. The casing is located a distance $c_j = 4 \cdot 10^{-4}$ m of each boundary node after accounting for the pre-loading forces f_0 . The latter are centrifugal forces that deform the blade into its working conditions. The system is rotating at an angular speed of $\Omega = 1,151.9$ rad·s⁻¹ = 11,000 rpm. These characteristics are chosen to reflect the same order of magnitude that was experimentally measured on the PHARE#2 test rig [35]. However, blade-tip/casing contacts were not observed during the experimental work. Thus, the blade-tip contact is numerically

engine order	θ		2	3	4	5	6		
value of K		8	4	8	$\overline{2}$	8	4	8	
reduced size	567	4536	2268	4536	1134	4536	2268	4536	567
sectors for AFT-CP		8	4	8	$\overline{2}$	8	4	8	
reduction $(\%)$	93.8	50.0	75.0	50.0	87.5	50.0	75.0	50.0	93.8

Table 4. Reduction for the case of the industrial system with $N = 16$ sectors for 10 harmonics (full size: $N \times 3n_{\text{nl}} \times (2N_{\text{h}} + 1)$ 9072 unknowns).

triggered by decreasing the nominal gap and applying a sufficient level of excitation to the system.

The excitation is applied to a node at the tip of the blade on each of the 16 blades in the circumferential direction in order to favor the response of bending modes. The excitation is considered to be synchronous with the rotation of the blade, the excitation frequency is $\omega = \Omega$. The contact scenario considered for this study is similar as the one that was considered on the academic model in Sec. 4.2.

When contact interactions occur, dry friction forces are applied to the blade. Given the high relative speeds at stake in such interactions, sliding is assumed at all times. The friction forces are expressed on the jth node at time t_i , such as

$$
f_{\mathrm{nl},j}^{\theta}(t_i) = -\mu \frac{v_j^{\theta}(t_i) + \rho_j \Omega}{\sqrt{(v_j^{\theta}(t_i) + \rho_j \Omega)^2 + v_j^z(t_i)^2}} f_{\mathrm{nl},j}^{\mathrm{N}}(t_i)
$$
\n(83)

for the circumferential direction and

$$
f_{\mathrm{nl},j}^{z}(t_{i}) = -\mu \frac{v_{j}^{z}(t_{i})}{\sqrt{(v_{j}^{\theta}(t_{i}) + \rho_{j}\Omega)^{2} + v_{j}^{z}(t_{i})^{2}}} f_{\mathrm{nl},j}^{\mathrm{N}}(t_{i})
$$
(84)

for the axial direction. v_j^{θ} and v_j^z are the circumferential and axial components of the vibrational velocity, $\mu = 0.15$ is the dry friction coefficient and ρ_i are the radial distances between each of boundary node and the rotation axis. Note that, since R-HBM uses a regularized contact law, the residual normal force associated with the regularization induces a residual friction force as well.

For all contact laws, since permanent sliding is assumed, the tangent force expressed in the friction direction is just the product of the normal contact forces f_{nl}^{N} and the friction coefficient μ . In Eq. (83) and Eq. (84), a projection of the tangent force in the cylindrical frame is performed depending on the rotation speed and the vibrational velocities of the nodes. Even though the expressions are nonlinear with respect to v_j^z and v_j^{θ} , the friction direction is highly driven by the rotation speed of the whole system. Indeed, for the given application $\rho_j \Omega \gg v_j^z$ and $\rho_j \Omega \gg v_j^{\theta}$. Under this hypothesis, it comes that $f_{\text{nl},j}^{\theta} \simeq \mu f_{\text{nl},j}^{N}$ and $f_{\text{nl},j}^{z} \simeq 0$ and thus the friction forces being proportional to normal forces it should not create any supplementary couplings of nodal diameters. Though, this hypothesis is only made for the determination of \mathcal{D}^k and the full expressions (Eq. (83) and Eq. (84)) are kept for the nonlinear solving process. It then comes that the nonlinear factors associated with the projections in the directions θ and z are assumed to create no further couplings than what was already identified in Sec. 2.4 and the same procedure is used for the industrial model than what was used in Sec. 4. This assumption will be verified numerically in the next section through the comparison between the reduced model and the full system.

For this study, $N_h = 10$ harmonics and $N_t = 2000$ instants are used for the HBM. For DLFT-HBM, the value $\varepsilon = 1 \cdot 10^7 \,\mathrm{N \cdot m^{-1}}$ is used. The friction coefficient is set to $\mu = 0.15$. In this section, only the simulations run with DLFT-HBM are shown. The simulations were also run with the R-HBM and provided near identical results. They are not presented here for the sake of brevity.

5.3 Traveling wave excitations

The first case study for this industrial model is the configuration where the casing is purely circular and the system is excited by a traveling wave excitation. In order to demonstrate the applicability of the traveling wave hypothesis formulated in Sec. 3.3, the cases of $h_{\text{ex}} = 1$ and $h_{\text{ex}} = 4$ (odd and even) are dealt with. Like for the studies in Sec. 4, the results are obtained through DLFT-HBM for the cyclic symmetric model as well as the full model. For these values of h_{ex} , the forcing amplitude is $f_{\text{ex}} = 2$ N and the results are displayed in Fig. 14.

Figure 14. Traveling wave excitation on the Catana industrial blade model for various engine orders, linear (......), DLFT-HBM: cyclic symmetry $(-)$, full $(--)$.

First of all, the same response is obtained through both methodologies, thus consolidating the relevance of both methodologies to tackle blade-tip/casing contact problems. Moreover, as for the academic case in Sec. 4, there is an exact match between the full computations and the reduced calculations made with the traveling wave hypothesis formulation. The latter yields a 93.75% reduction, allowing to perform calculations in a few minutes, where the full calculations usually last for between 5 to 15 hours, depending on the choice of continuation parameters, contact treatment algorithm and other HBM parameters. Indeed, the scaling of the reduction rate with the computational time is not linear: numerous operations made during the solving process scale in $\mathcal{O}(n \log n)$ or even $\mathcal{O}(n^2)$ and higher order. Then, a reduction of the size of the system of a certain percentage, say $R_{\%}^{\text{size}}$, can be translated into an effective reduction in terms of time $R_{\%}^{\text{time}}$ that is significantly higher. Such that generally speaking $R_{\%}^{\text{time}} > R_{\%}^{\text{size}}$, depending on the implementation. While this effect is not systematically highlighted on small systems, it underlines that relatively small reductions in terms of size can yield significant speed-ups of the overall execution of a calculation.

For instance, for the case $h_{\rm ex} = 4$ for R-HBM, the full calculation took 7 hours and 14 minutes while the cyclic symmetric one only lasted for 1 minute a 31 seconds, making the reduction of computing time of over $R_{\%}^{\text{time}} = 99.6\%$. This is almost 6 points greater than the expected size reduction of $R_{\%}^{\text{size}} = 93.75\%$. The same phenomenon is observed for DLFT-HBM and the effective reduction is of $R_{\%}^{\text{time}} = 98.9\%$. Obviously, this reduction is highly dependent on implementation and hardware. Throughout the article, only effective size reductions are presented but it must be kept in mind that the announced reductions are most likely underestimated.

For $h_{\text{ex}} = 1$, the decomposition used for computations in cyclic symmetry is

$$
[\mathcal{D}^k]_{k \in [0, N_h]} = [\{0\}, \{-1\}, \{-2\}, \{-3\}, \{-4\}, \{-5\}, \{-6\}, \{-7\}, \{8\}, \{7\}, \{6\}].
$$
\n
$$
(85)
$$

It can be seen in Fig. 15 that the cyclic components emerging from the full DLFT-HBM calculation match the exact basis of Eq. (85). The match is also observed for $h_{\text{ex}} = 4$ but not shown here for the sake of brevity. This section has then demonstrated the usefulness of the traveling wave hypothesis on industrial systems featuring blade-tip/casing contacts.

Figure 15. Cyclic components over all harmonics at the resonance point of the full calculation in Fig. 14a, traveling wave with $h_{\rm ex} = 1.$ Nodal diameter: 0 (**m**), 1 (**m**), 2 (**m**), 3 (**m**), 4 (**m**), 5 (**m**), 6 (**m**), 7 (**m**), 8 (**m**).

5.4 Partly standing wave excitation

Now that the capabilities of the overall strategy is demonstrated on the case of traveling wave excitations, a new scenario is envisioned. Pure standing wave excitations are not likely to be encountered in turbomachinery applications since the aerodynamic loadings are oftentimes synchronous with the rotation of the shaft. The order of the traveling excitations is generally determined by the number of blades in the immediately adjacent stages of the studied stage. However, it is possible for a stage to be excited by stages further than its immediate neighbors. In this section, the loading scenario that is studied is one where the blade is under a backward excitation for 90% (adjacent stage) and under a forward excitation for 10% (further stage) for the same engine order. This means that the excitation is a sum of a backward traveling wave (80%) and of a standing wave (20%). For this case, the engine order considered is $h_{\text{ex}} = 4$ with $f_{\text{ex}} = 1$ N and the excitation is such as

$$
\mathbf{f}_{\text{ex},j}(t) = f_{\text{ex}} \left[0.9 \cos \left(t + \frac{\alpha (j-1) h_{\text{ex}}}{\omega} \right) + 0.1 \cos \left(t - \frac{\alpha (j-1) h_{\text{ex}}}{\omega} \right) \right] \quad \forall j \in [\![1, N]\!].
$$
 (86)

Since the excitation includes stationary components, the traveling wave hypothesis is not suitable for this case. Hence, it is necessary to resort to the criterion of Sec. 2.4. It is then predicted that the following basis is needed for such a problem:

$$
\mathcal{D}^{2k}|_{k\in[0,5]} = \{0,8\} \text{ and } \mathcal{D}^{2k-1}|_{k\in[1,5]} = \{4\}. \tag{87}
$$

Indeed, this case is subject to the parity criterion since $K = 2$ and $Kh_{\text{ex}} = 8 \equiv \frac{N}{2}$ [N]. Using this basis, the results of the simulations are displayed in Fig. 16 for DLFT-HBM. Firstly, it is observed that the match is once again excellent between the full and reduced calculations even if the contact law is nonsmooth. The standing characteristics of the response are clearly identified by comparing the responses of sectors 1 (Fig. 16a) and 2 (Fig. 16b). An interesting behavior even occurs at the resonance of sector 1. Indeed, the cyclic decomposition given in Fig. 17 shows that a purely standing wave displacement field emerges from the excitation that is only 20% stationary. Actually, this behavior is caused by the fact that the dynamic motion of the system reaches a point where only one in two blades experiences contact along a period at this resonance. On the frequency response curves of Fig. 16, all even numbered blades (sector 2 in Fig. 16b) experience a linear behavior while odd sectors feature resonances around 293 Hz. The associated branch for the even sectors also experience a turning point at the same frequency. These points can be visualized in the zooms of Fig. 16. One can see that it is superimposed with the classical linear response of the system.

This type of response underlines the need of investigating the dynamics of the system using general nodal diameter coupling criteria such as the one presented in this article. Indeed, even if the system is under an almost purely backward traveling wave, the small standing components introduced for this scenario creates new couplings with the nonlinearity and makes a pure standing wave motion emerge. Hence, it highlights that no part of the excitation should be neglected in a nonlinear framework because small contributions can still induce specific dynamics in the system.

Figure 16. Partly standing wave excitation on the Catana industrial blade model for $h_{ex} = 4$, linear (\cdots) , DLFT-HBM: cyclic symmetry $(-)$, full $(-)$, equivalent backward traveling wave excitation $(-)$.

(b) Backward traveling diameters

Figure 17. Cyclic components over all harmonics at the resonance point of the full calculation in Fig. 16a (sector 1), 90% forward and 10% backward traveling wave with $h_{\text{ex}} = 4$. Nodal diameter: 0 (...), 8 (...).

Besides this observation, the diametral basis chosen in Eq. (87) was correctly identified beforehand and matches perfectly with the cyclic components emerging in the full response in Fig. 17. The predicted parity criterion is also respected like expected. This shows that on a fairly complex configuration, the rules derived in Sec. 2.4 and Sec. 3.2.1 are still suitable in partly standing wave excitations scenarios on large scale industrial models.

In order to demonstrate the usefulness of the proposed strategy once again, a similar example is studied with $h_{\rm ex} = 7$. The percentages of forward and backward excitations are kept the same and $f_{\rm ex} = 0.3$ N. $h_{\rm ex} = 7$ is actually prime with $N = 16$ and no reduction is expected in terms of overall diameters in the response. However, the parity criterion yields that only half of the diameters are considered on odd and even harmonics. This means that even if h_{ex} and N are prime with one another, a reduction of 50% is reached (as showed in Tab. 4). The basis considered for these computation is thus

$$
\mathcal{D}^{2k}|_{k\in[0,5]} = \{0,2,4,6,8\} \text{ and } \mathcal{D}^{2k-1}|_{k\in[1,5]} = \{1,3,5,7\}. \tag{88}
$$

The responses obtained for this configuration are displayed in Fig. 18. The response of the system to a pure backward traveling excitation of amplitude $f_{\rm ex} = 0.3$ N is also displayed for comparison purposes. For this case, the reduction method also produces an exact approximation of the solution obtained on the full model. This further supports the reduction capabilities of the method exposed in this article for the reduction of blade-tip/casing problems using cyclic symmetry. The forced responses displayed in Fig. 18 highlight that the purely traveling wave excitation response is very different from the response on the case where 10% of forward traveling excitation is introduced. This shows the level of sensitivity that blade-tip/casing problems experience when it comes to the loading scenario. Actually, when looking at the cyclic components of the response at the resonance of Fig. 18a in Fig. 19, one can see that the response is actually a standing wave (both forward components in Fig. 19a and backwards components in

Figure 18. Partly standing wave excitation on the Catana industrial blade model for $h_{\text{ex}} = 7$, linear (......), DLFT-HBM: cyclic symmetry $(-)$, full $(-)$, equivalent backward traveling wave excitation $(-)$.

Fig. 19b are equal). For instance, at this point only two sectors out of the $N = 16$ sectors actually experience contact along a period of vibration. This means that even if the excitation is mainly composed of a backward traveling wave, the small forward traveling perturbation yields a localized nonlinear response on a few sectors. These localizations are actually symptoms of the appearance of standing wave displacement fields. One should also note that the parity

(b) Backward traveling diameters

Figure 19. Cyclic components over all harmonics at the resonance point of the full calculation in Fig. 18a (sector 1), 90% forward and 10% backward traveling wave with $h_{\text{ex}} = 7$. Nodal diameter: 0 (**iii**), 1 (**iii**), 2 (**iii**), 3 (**iii**), 4 (**iii**), 5 (**iii**). $6 \ (\blacksquare), 7 \ (\blacksquare), 8 \ (\blacksquare).$

criterion (that yields all the reduction properties for this example) is strictly respected as predicted by the analysis of the loading that allowed to find the basis in Eq. (88).

As a way of providing more insights about the phenomenon of localization observed in Fig. 19, the cyclic components of the first harmonic are displayed in Fig. 20 for the whole response for the full calculation carried out with DLFT-HBM in Fig. 18. The results displayed in Fig. 18 show that in the linear domain, *i.e.* at low and high frequencies, the backward 7ND component is significantly higher than the forward 7ND component and this reflects the excitation applied to the system. However, once the nonlinearity is activated, all other nodal diameters have both their cyclic components increased of the same amount, meaning that theses nodal diameters are actually responding on a standing wave. The behavior is quite different with ND7 (associated to the excitation $h_{\rm ex} = 7$) since the backward contribution suddenly drops and the forward contribution raises rapidly. This continues until the resonance where both contributions are almost equal, as already highlighted in Fig. 19. This demonstrates how the response of the system goes from an almost pure traveling wave displacement field in the linear domain (90%) to a standing wave response at its nonlinear resonance.

Overall, the reduction capabilities of the strategy was here demonstrated on the industrial test case for various loading scenarios: traveling wave excitations and partly standing wave excitations. A localization behavior was shown

Figure 20. Cyclic components of the first harmonic for the whole response of the full calculation in Fig. 18a, 90% forward and 10% backward traveling wave with $h_{\text{ex}} = 7$. Nodal diameter: 1 (\Box), 3 (\Box), 5 (\Box), 7 (\Box), backward components (--), forward components $(-)$.

to emerge in the cases where the traveling excitation was not pure. Though, this behavior was correctly captured by the reduction strategy. The emergence of such solutions vibrational solutions was not previously highlighted and this contributions demonstrates the importance of modeling accurately the excitation in order not to simplify the response of the bladed disk.

6 Conclusion

This paper demonstrates the reduction capabilities of the cyclic symmetry in order to tackle vibro-impact applications, as much on small size systems than on industrial models. Firstly, the nodal diameter coupling criterion is derived by analyzing a regularized contact law. Its implementation is quite straightforward and allows to drastically decrease the size of the system to be solved. The criterion relies on the analysis of the Taylor expansion of the contact law. The reduction is dependent on both the number of sectors of the system and the nodal diameters that are excited by the external loadings.

A method to further reduce the computing times is also proposed by analyzing how the different harmonics of the harmonic balance method interact with the nodal diameters. It results in a criterion that is able to reduce even more the size of the problem by a factor 2 for certain combinations of numbers of sectors and nodal diameters of the excitation.

These two original contributions are then put to the test on numerous forced response calculations. In order to cover all possible configurations, academic models with both even and odd numbers of sectors are studied through different loading scenarios: traveling wave excitations, standing wave excitations and loadings along multiple nodal diameters (traveling wave excitation and casing deformation). Firstly, the traveling wave hypothesis, oftentimes used on friction damping applications, was proven to be suitable for highly nonlinear vibro-impact applications to reduce significantly the size of the system. Secondly, the two criterion used to build the basis of diameters that serve to reduce the size of the problem are shown to produce results that match exactly with the full reference calculations. Indeed, the reduction is exact and yields no loss in precision whatsoever.

These observations of the response are also true when analyzing the results obtained on the state-of-the-art fan stage ECL5/Catana under blade-tip/casing contacts. In fact, even though the friction forces are added in the modeling of blade-tip/casing interactions, the criterion holds true and the results on the full model are matching exactly with the reduced ones, even on the most complex loading scenarios.

In both cases, the analysis of the cyclic components of the response validate *a posteriori* the choice of the different basis of nodal diameters for the reduced calculations. As a whole, the analytical derivation about the nodal diameter couplings were consolidated by numerous examples and the criterion appears to be industry-ready for the simulation of complex blade-tip/casing interactions on full bladed disks at a low computational cost. Moreover, it appears that these derivation made on the regularized contact law are still valid when considering a truly nonsmooth contact law. This means that the reductions still hold when the contact laws are non differentiable.

As a way of concluding this study, a wide view of the reduction rates achieved by the proposed criterion is given hereafter. In the case of a traveling wave excitation, the traveling wave hypothesis systematically yields a drastic reduction of a rate

$$
R_{\%}^{\text{TW}} = 100 \left(1 - \frac{1}{N} \right),\tag{89}
$$

that is over 90% as soon as the system is composed of more than $N = 10$ blades.

For the more general case where the loading is not necessarily a traveling wave, the mean reductions achieved along all non-zero engine orders for numbers of sectors up to 150 are displayed in Fig. 21. First of all, the reduction

Figure 21. Expected reduction as a function of N until $N = 150$, even number of sectors with the parity rule (\blacksquare) , odd number of sectors (\blacksquare) , even number of sectors (\blacksquare) .

strategy is unfortunately not suitable when the number of sectors of the system is a prime number: no reduction can be achieved. Moreover, it can be seen that even numbers of sectors are much more subject to reduction that odd numbers of sectors. Considering a resolution through harmonic balance method that allows to use the parity rule, the mean reduction of even numbers of sector reaches 62.4% while non-prime odd numbers of sectors are reduced of 23.4% on average and it drops down to 12.5% when considering the prime numbers. Though, the overall mean reduction is 37.6% which is quite large when considering the fact that the reduction is exact. If a resolution through other methods than the harmonic balance method is preferred, the mean reduction among even number of sectors drops down to 37.7% and the overall reduction falls to an average of 25.2%.

Overall, this work has participated in making a major step forward in the efficient simulation of the nonlinear vibrations of bladed disk featuring blade-tip/casing contacts. Indeed, such simulations were conducted on a single blade or on the full bladed disk to date, making them unrealistic or extremely cumbersome. Hence, the development of a reduction strategy such as the one presented here enables to carry out increasingly realistic calculations at an affordable computational burden.

Among the perspectives of this work, the stability analysis of the response of these cyclic systems featuring vibro-impact is one of the main avenues of research. Indeed, this study has focused on solving the equation of motion through the harmonic balance method to validate the reduction procedure in terms of nodal diameters. It has shown that reduction is exact since the same results are obtained with the cyclic symmetric formulation as well as with the full system. However, this study does not give any information about the stability of the presented solutions and a stability analysis would be required to state on the nature of these solutions. Moreover, a cross-verification of the stability of these nonlinear solutions was not carried out in this article because it exceeded its scope. Nevertheless, it would be necessary in order to validate the physical nature of these solutions. The characterization of eventual branches in terms of cyclic components would also allow to understand more finely the dynamics response of such cyclic systems.

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A Real spatial fourier matrices

This appendix provide the expression used to build the spatial Fourier matrix used in the real cyclic symmetry formalism for both even and odd numbers of sectors N . If the number of sectors of the system is even, the Fourier matrix is expressed such as

$$
\mathbf{S} = \sqrt{\frac{2}{N}} \begin{bmatrix} \frac{1}{\sqrt{2}} & \cdots & \frac{1}{\sqrt{2}} & \cdots & \frac{1}{\sqrt{2}} \\ \cos(\alpha) & \cdots & \cos(j\alpha) & \cdots & \cos(N\alpha) \\ \sin(\alpha) & \cdots & \sin(j\alpha) & \cdots & \sin(N\alpha) \\ \vdots & \vdots & \vdots & & \vdots \\ \cos(h\alpha) & \cdots & \cos(hj\alpha) & \cdots & \cos(hN\alpha) \\ \sin(h\alpha) & \cdots & \sin(hj\alpha) & \cdots & \sin(hN\alpha) \\ \vdots & \vdots & & \vdots & \vdots \\ \frac{(-1)^1}{\sqrt{2}} & \cdots & \frac{(-1)^j}{\sqrt{2}} & \cdots & \frac{(-1)^N}{\sqrt{2}} \end{bmatrix}
$$
(90)

whereas its expression becomes the following when considering an odd number of sectors

$$
\mathbf{S} = \sqrt{\frac{2}{N}} \begin{bmatrix} \frac{1}{\sqrt{2}} & \cdots & \frac{1}{\sqrt{2}} & \cdots & \frac{1}{\sqrt{2}} \\ \cos(\alpha) & \cdots & \cos(j\alpha) & \cdots & \cos(N\alpha) \\ \sin(\alpha) & \cdots & \sin(j\alpha) & \cdots & \sin(N\alpha) \\ \vdots & \vdots & \vdots & \vdots \\ \cos(h\alpha) & \cdots & \cos(hj\alpha) & \cdots & \cos(hN\alpha) \\ \sin(h\alpha) & \cdots & \sin(hj\alpha) & \cdots & \sin(hN\alpha) \\ \vdots & \vdots & \vdots & \vdots \\ \cos((\frac{N-1}{2})\alpha) & \cdots & \cos((\frac{N-1}{2})j\alpha) & \cdots & \cos((\frac{N-1}{2})N\alpha) \\ \sin((\frac{N-1}{2})\alpha) & \cdots & \sin((\frac{N-1}{2})N\alpha) \end{bmatrix} .
$$
 (91)

Note that these Fourier matrices are orthogonal, yielding that $S^{-1} = S^{\top}$.

B Developments associated with the parity criterion for the selection of nodal diameters in the case of unilateral nonlinearities

When using the HBM, the choice of the sets \mathcal{D}^k can lead to further reduction by evaluating which harmonic will respond on which nodal diameter. This section aims to demonstrate the existence of such a relation for unilateral nonlinearities. For simplicity, only a single nodal diameter will be considered in the initial basis $\mathcal{D}_{q,0}$. Extension to multiple initial nodal diameters is then straightforward.

Consider h_{ex} , the initial nodal diameter associated to K the smallest integer which respects the following

$$
2Kh_{\text{ex}} \equiv 0 \,[N].\tag{92}
$$

This definition is derived in Sec. 2.4. As a reminder, to determine the full nodal diameter basis, one has to proceed with the following steps

- Create a vector $\mathbf{v}_K = [0, 1, 2, \cdots, q, \cdots, K]$ (a scalar of this vector is noted q)
- Multiply it by h_{ex} leading to a vector $\mathbf{v}_{K,h_{\text{ex}}} = [0, h_{\text{ex}}, 2h_{\text{ex}}, \cdots, qh_{\text{ex}}, \cdots, Kh_{\text{ex}}].$
- for each component $q h_{\text{ex}}$, its reminder through the Euclidian division by N is evaluated, noted r_q . Then, if $0 < r_q < m$ with m the maximum number of nodal diameters for the given number of sectors, the excited diameter is $d_q = r_q$. Otherwise, if $m < r_q$ then the excited diameter is $d_q = N - r_q$. A final vector is obtained with the list of the interacting nodal diameter $\mathcal{D} = [d_0, d_1, d_2 \cdots, d_q, \cdots d_K]$ (where $d_0 = 0$, and $d_1 = h_{\text{ex}}$).

Two scenarios exist:

- if $Kh_{\text{ex}} \equiv 0$ [N], then no reduction is obtained with the HBM,
- if $Kh_{\text{ex}} \equiv \frac{N}{2}$ [N], then the nodal diameters d_q with q even only respond on even harmonics, and the nodal diameters \bar{d}_q with q odd only respond on odd harmonics.

This second scenario allows to divide the number of unknowns by 2. The following sections aim to demonstrate this property: each scenario is dealt with separately in B.1 and B.2 respectively.

B.1 First scenario

For the first scenario, one assumes that

$$
Kh_{\text{ex}} \equiv 0 \,[N].\tag{93}
$$

The number K is then necessarily an odd number. Indeed, if K were to be even, then there would exist an integer K_2 which satisfies $K = 2K_2$. Substituting in Eq. (93) gives

$$
2K_2 h_{\text{ex}} \equiv 0 \,[N].\tag{94}
$$

This breaks the definition of K , *i.e.* K is defined as the smallest integer verifying Eq. (92). Therefore, K is necessarily an odd number in the following. Then, it comes that

$$
\mathbf{v}_K = \left[0, 1, \cdots \frac{K-1}{2}, \frac{K+1}{2}, \cdots, K-1, K\right].
$$
 (95)

Following the expression of \mathbf{v}_K in this case, a symmetry then occurs in the interacting nodal diameters $\mathcal D$ between both ends of the list: $d_K = d_0 = 0$, then $d_{K-1} = d_1 = h_{ex}$, etc. As a consequence, each nodal diameter appears twice in $\mathcal D$ and it is such that

$$
\mathcal{D} = \left[0, h_{\text{ex}}, \cdots d_{\frac{K-1}{2}}, d_{\frac{K-1}{2}}, \cdots, h_{\text{ex}}, 0\right].
$$
\n(96)

Notice that both indices 1 and $K - 1$ (respectively an even and an odd number) are associated to the the same nodal diameter h_{ex} .

Consider now the HBM in an iterative procedure as described in Sec. 2.4. Initially, only the nodal diameter h_{ex} constitutes the basis and responds only on the first harmonic $e^{\pm i\omega t}$. Then through Eq. (25), the second diameter $2h_{\text{ex}}$ responds on its second harmonic $e^{\pm 2i\omega t}$, and so on. Notice that $K-1$ is even and thus the nodal diameter $r_{K-1} = h_{\text{ex}}$ will also have an even harmonic component. Due to the nodal diameter interaction, the harmonics gets coupled and eventually all harmonics respond on all nodal diameters.

B.2 Second scenario

For this second scenario, one assumes that

$$
Kh_{\text{ex}} \equiv r\,[N],\ r \neq 0. \tag{97}
$$

Since $2Kh_{\text{ex}} \equiv 0$ [N], Eq. (97) implies that $2Kh_{\text{ex}}$ is proportional to N with an odd number. Moreover, since $2Kh_{\text{ex}}$ is an even number then N must be even, and K must be odd. Therefore, it comes that

$$
2Kh_{\text{ex}} = (p+1)N
$$

\n
$$
Kh_{\text{ex}} = \frac{N}{2} + 2pN,
$$
\n(98)

hence $r = \frac{N}{2}$.

Now, assume that there exists an integer $1 < q < K$ such that

$$
q h_{\text{ex}} \equiv h_{\text{ex}} \,[N].\tag{99}
$$

This relation is equivalent to

$$
(q \pm 1) h_{\text{ex}} \equiv 0 \,[N]. \tag{100}
$$

To ensure that K is the smallest integer verifying Equation Eq. (92), the relation $q \pm 1 > K$ must be respected. Therefore, either of $q > K + 1$ or $q > K - 1$ is true. Neither of the two relations is not compatible with $q < K$. As a consequence, there exists no integer $1 < q < K$ such that Eq. (99) holds true. Therefore, in this scenario, each nodal diameter that is contained in $\mathcal D$ is only present for a single value q unlike for the first scenario in B.1. Hence it comes that

$$
\mathbf{v}_K = [0, 1, \cdots q, \cdots, K] \tag{101}
$$

and

$$
\mathcal{D} = \left[0, h_{\text{ex}}, \cdots d_q, \cdots \frac{N}{2}\right].\tag{102}
$$

To determine which harmonic will respond on which nodal diameter, the HBM can be employed with an iterative process. Without loss of generality, the excited diameter $d_1 = h_{ex}$ is assumed to be excited along its first harmonic $e^{i\omega t}$. Through the first iteration (determination of the coupling with only d_1), the nodal diameter d_q will systematically have a component along the qth harmonic $e^{\pm i q \omega t}$. After this first iteration process, the entire set of interacting nodal diameters is determined. Due to the aforementionned property, after this first iteration, each nodal diameter is only associated to a single harmonic component. Notice also that if q is even (resp. odd) then d_q is associated with an even (resp. odd) harmonic.

The next iterations of the procedure to determine the harmonic couplings in HBM consist in identifying how the nodal diameter will combine with each other and hence how their temporal harmonic content will be combined. To determine this, Eq. (27) is recalled

$$
\widehat{f}_{\rm nl,3}^q = \frac{\kappa}{\sqrt{N}\gamma^3} \sum_{l=0}^{\infty} \left(\beta_l \sum_{k \in \mathcal{T}_l} \left(\alpha_k \prod_{p_k=1}^{Q_k} \left(\widehat{u}_{h_{p_k}} \right)^{\beta_{p_k}} \left(\sum_{j=1}^N e^{i\alpha(j-1)\left(\sum_{p_k=1}^{Q_k} h_{p_k} \beta_{p_k} - q\right)} \right) \right) \right). \tag{103}
$$

The second iteration concerns $l = 1$, hence \sum Q_k $p_k=1$ $\beta_{p_k} = 2$. So, here, the interaction between only two nodal

diameters is studied. The nodal diameter d_q can be obtained through the combination of two other nodal diameter d_a and d_b if $a \pm b \equiv q \mod N$. Two scenarios arise

- If q is even, then necessarily a and b must both be either odd or even. In either case, the combination of their temporal harmonic $e^{i\omega a t}$ and $e^{i\omega b t}$ will be an even harmonic $e^{i\omega(a \pm b)t}$ associated to d_q .
- If q is odd, then necessarily a and b one is even and the other odd, leading to an odd harmonic $e^{i\omega(a\pm b)t}$ associated to d_q .

Following these elements, the second iteration shows that a nodal diameter d_q can be generated through another interaction (involving two nodal diameter) but the parity of its associated harmonic does not change.

By induction, considering the $(k+1)$ th iteration concerns $l = k$, the coupling relation is conditioned by Q_k

 $\sum \beta_{p_k} = 2k$. Here, the interaction between 2k nodal diameters is studied. A similar result is obtained: nodal $p_k=1$

diameter d_q can be obtained through this interaction but the parity property must still hold. This is caused by the fact that $2k$ is an even number.

In the end, if q is even, only even harmonics will respond on the nodal diameter d_q and conversely if q is odd.

C Academic case with an odd number of sectors : $N = 45$

The case study of this appendix is the same model as in Sec. 4.3 but with $N = 45$ sectors. It is dealt with for completeness: the reductions between even and odd numbers of sectors are significantly different. The fact that N is odd allows to showcase the reduction properties of the proposed method. This time, no calculation is able to take advantage of the parity criterion on the harmonics as mentioned in Sec. 3.2.1. With this number of sectors, the expected reductions are displayed in Tab. 5 for all engine orders. It can be seen that the reductions are lower

Table 5. Reduction for the case of the simplified system with $N = 45$ sectors for 10 harmonics (full size: $N \times n_{\text{nl}} \times (2N_{\text{h}} + 1) = 945$ unknowns).

reduction (%) 66.7 0.0 0.0 93.3 0.0 0.0 88.9 0.0 80.0 66.7 0.0

that for the previous case in Sec. 4.3: the mean reduction only reaches 38.1% (mean reduction of 75% for $N = 24$). Moreover, Tab. 5 reveals that 12 values of $h_{\rm ex}$ are not subject to any reduction at all.

For the simulations, the approach is similar as in Sec. 4.3: firstly the response of the system under traveling wave excitations is analyzed, then standing wave are studied in order to validate the choice of the nodal diameter basis, and finally a loading scenario involving both a traveling wave excitation and a casing deformation is studied.

C.1 Traveling wave excitations

For the traveling wave excitations, two cases are analyzed: $h_{\text{ex}} = 2$ with $f_{\text{ex}} = 6$ N and $h_{\text{ex}} = 17$ with $f_{\text{ex}} = 4$ N. The casing is circular and the reduced problem is solved by means of the traveling wave hypothesis. For the case

with $h_{\text{ex}} = 2$, the nodal diameter basis retained is

 \lceil

$$
[\mathcal{D}^k]_{k \in [\![0,N_h]\!]} = [\{0\}, \{-2\}, \{-4\}, \{-4\}, \{-6\}, \{-8\}, \{-10\}, \{-12\}, \{-14\}, \{-16\}, \{-18\}, \{-20\}] \tag{104}
$$

and

$$
\mathcal{D}^{k}\big]_{k\in\llbracket 0,N_{\rm h}\rrbracket} = [\{0\},\{-17\},\{11\},\{-6\},\{22\},\{5\},\{-12\},\{16\},\{-1\},\{-18\},\{10\}] \tag{105}
$$

for $h_{\text{ex}} = 17$. The results of the different simulations are displayed in Fig. 22.

Figure 22. Traveling wave excitation on the simplified bladed disk with $N = 45$ sectors for various engine orders, linear (\ldots ,), R-HBM: cyclic symmetry $(-)$, full $(--)$, DLFT-HBM: cyclic symmetry $(-)$, full $(--)$.

The forced responses displayed in Fig. 22 indicate that the traveling wave hypothesis is also valid when considering an odd number of sectors. There is an exact match between computations led in cyclic symmetry and full calculations. For the sake of brevity, the cyclic components of these calculations are not displayed, but the nodal diameters predicted by the zig-zag diagram are recovered. The reduction rate achieved in the case of traveling wave excitations is $100 \cdot (1 - \frac{1}{N}) = 97.7\%.$

C.2 Standing wave excitations

In this section, more complex excitations are applied to the system: standing wave excitations. In order to demonstrate the reduction capabilities of the methodology, two values of h_{ex} where a reduction is predicted are picked from Tab. 5. Simulations are led for $h_{\text{ex}} = 9$ with $f_{\text{ex}} = 1$ N and $h_{\text{ex}} = 15$ with $f_{\text{ex}} = 1$ N. The two basis associated with these values of h_{ex} are

$$
\mathcal{D}^{k}|_{k\in[0,10]} = \{0,9,18\} \text{ for } h_{\text{ex}} = 9 \quad \text{ and } \quad \mathcal{D}^{k}|_{k\in[0,10]} = \{0,15\} \text{ for } h_{\text{ex}} = 15. \tag{106}
$$

For $h_{\text{ex}} = 9$, the procedure of Sec. 2.4 yields $K = 5$ and $K = 3$ for $h_{\text{ex}} = 15$. With these basis, the results are displayed in Fig. 23 for both R-HBM and DLFT-HBM.

Figure 23. Standing wave excitation on the simplified bladed disk with $N = 45$ sectors for various engine orders, linear (......), R-HBM: cyclic symmetry $(-)$, full $(--)$, DLFT-HBM: cyclic symmetry $(-)$, full $(--)$.

Again, the reduction criterion on the nodal diameters produces an exactly matching response for both the regularized and the nonsmooth contact laws. As a way of verifying the choice of the basis of Eq. (106), the cyclic contributions are displayed in Fig. 24a for $h_{\text{ex}} = 9$ and in Fig. 24b for $h_{\text{ex}} = 15$. For this case, since the response is purely stationary (forward contributions are strictly equal to backward contributions), no distinction is made on the direction of propagation and only one of the two graphs is displayed for each h_{ex} . For both cases, the basis

Figure 24. Cyclic components over all harmonics at the resonance point of the full calculation in Fig. 23b and in Fig. 23d, standing wave excitations. Nodal diameter: $0 \ (\blacksquare), 9 \ (\blacksquare), 15 \ (\blacksquare), 18 \ (\blacksquare).$

identified in Eq. (106) are exactly matching with what the response of the full model features. This shows that the coupling rule derived in Sec. 2.4 is also relevant for cases with an odd number of sectors. Contrarily with the case in Sec. 4.3.2, there is no parity rule among the cyclic components and the whole basis responds on all harmonics. This property is also correctly predicted by the parity rule of Sec. 3.2.1: the criterion is only possible for even number of sectors, which is not the case here.

C.3 Mixed contact scenario

Finally, the academic test case is characterized through an even more complex contact scenario: it is excited through a traveling wave and the casing it interacts with is deformed with a diametral shape. The system is excited along a traveling wave excitation on the 12th engine order $(h_{\rm ex} = 12)$ with an amplitude of $f_{\rm ex} = 2$ N and the casing is deformed with a 3ND shape $(h_c = 3)$ with $p_j = 0.05c_j$. Both excitations are traveling waves, however both orders are expected to couple themselves and make a displacement field of standing wave shape emerge. In order to build the basis of the nodal diameters that intervene in the response, one should take a decomposition that is related to the GCD of h_c and h_{ex} , 3 in this case, and the basis is

$$
\mathcal{D}^{k}|_{k\in\llbracket 0,10\rrbracket} = \{0,3,6,9,12,15,18,21\}.
$$
\n(107)

Even if the GCD of h_{ex} and h_c is equal to h_{ex} , the nodal diameter basis is extended. Indeed, if no deformation of the casing were present, a simple traveling wave basis could have been chosen. The computations are then led for both DLFT-HBM and R-HBM and displayed in Fig. 25 for two sectors (1 and 4).

Figure 25. Traveling wave excitation on the simplified bladed d isk with $N = 24$ sectors for $h_{ex} = 12$ and a casing deformed with $h_c = 3$, linear (......), R-HBM: cyclic symmetry (-), full (---), DLFT-HBM: cyclic symmetry (-), full (---).

In this scenario and with the chosen diametral basis, there is again a perfect match between the computations led in cyclic symmetry and the full calculations. This agreement for such of a sophisticated contact scenario demonstrates the usefulness of the proposed strategy. It offers reduction capabilities in numerous configurations. For the present case, a reduction of 66.7% is achieved. As a way of demonstrating that the diametral basis is correctly identified,

Figure 26. Cyclic components over all harmonics at the resonance point of the full calculation in Fig. 25, traveling wave with $h_{\text{ex}} = 12$ and a casing deformed with $h_{\text{c}} = 3$. Nodal diameter: 0 (...), 3 (...), 6 (...), 9 (...), 12 (...), 15 (...), 18 (...), $21 \, (\blacksquare).$

the cyclic contributions are displayed in Fig. 26. Contrarily with standing wave excitations, the displacement field is not purely stationary: the system experiences a vibrational motion composed of multiple traveling waves of different orders as well as stationary components that can not be easily predicted by analyzing the loadings (excitation and casing) that are applied. Nevertheless, the nodal diameters responding in Fig. 26 correspond to the basis identified in Eq. (107) justifying once again the relevance of the nodal diameter coupling criterion for unilateral nonlinearities.