Uncoupling Techniques for the Dynamic Characterization of Sub-structures

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ABSTRACT

The characterization of the dynamic behavior of sub-structures may be of great importance in certain complex structures. A possible approach is through the use of uncoupling techniques. This is the subject of the present paper, where various issues are raised and different possibilities to circumvent the encountered problems are discussed. These problems are mostly due to numerical instabilities associated to the calculation procedures, namely to the ill-conditioning that often appears and causes complications in the inversion of the involved matrices. The location of the possible points of measurement may also influence the results. Several variations to the main algorithms are tried in order to reduce the numerical problems. This study will be based on numerical simulations to compare the various approaches and to draw some final conclusions.

1 INTRODUCTION

Recently the subject of structural uncoupling (also referred to by some authors as decoupling) has regained interest among the scientific community [1, 2]. Seeking for the dynamic characterization of sub-structures, there are several techniques combining both experimental and analytical/numerical methods. They can be classified in three different groups, according to the working domain, either modal, mechanical impedance or frequency domain.

In the modal domain one has methods that use modal properties, like the classic Craig-Bampton [3], known as Component-Mode Synthesis (CMS). Meanwhile, the Dual Craig-Bampton method has been presented by Rixen [2].

In what mechanical impedance is concerned, one has the system mass, stiffness and damping matrices, as well as external forces and connection forces. Ahmadian *et al.* and Jalali *et al.* [4, 5] developed a method that allows the identification of a bolted joint using the physical matrices of the system.

In the frequency domain there are a vast number of methods based on the classic one developed by Jetmundesn [6], known as "FRF-based sub-structuring method" (FBS) as classified by de Klerk *et al.* [7]. D'Ambrogio [1] developed two methods based on system impedance and mobility using only the connection co-ordinates. As it will be explained in this work, one of the methods presented here will have the same result. In Jetmundesn's method [6] the inherent numerical problems associated to ill-conditioned matrices are well known; moreover, Frequency Response Function (FRF) measurements have inherent associated noise that make the obtaining of acceptable results a difficult job; several strategies have been appointed to overcome the problems, see for instance Ren and Beard [8]; Maia *et al.* [9] proposed a solution avoiding the connection co-ordinates.

In this work several approaches for the classic uncoupling Jetmundesn's method are developed, analyzed and compared, through numerical examples using beams. Adding local inertial elements, a possible way of circumventing the ill-conditioning problems is proposed.

2 THEORETICAL FORMULATION

2.1 FRF coupling

Coupling techniques are well known (e.g., see [10]); they can be divided into spatial coupling methods, based upon the system matrices (M, $K \in C$), modal coupling methods, using modal properties (natural frequencies, damping ratios and mode shapes) and frequency response function coupling (*FRF* coupling). Let us concentrate in this latter one.

Let us consider two sub-structures A and B, connected though some co-ordinates as represented in figure 1. Together, they form structure C. Let i represent the co-ordinates exclusively belonging to A, k the ones exclusively belonging to B and j the connection co-ordinates.



Fig. 1 Coupling of sub-structures A and B, forming structure C

From the equilibrium of forces and the compatibility conditions of displacements one can write:

$$\boldsymbol{f}_i^A + \boldsymbol{f}_i^B = \boldsymbol{f}_i^C \tag{1}$$

$$\boldsymbol{x}_{j}^{A} = \boldsymbol{x}_{j}^{B} = \boldsymbol{x}_{j}^{C} \tag{2}$$

The receptance matrices \boldsymbol{H} relating the amplitudes of the forces to the amplitudes of the displacements are defined as

$$X = HF \tag{3}$$

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and therefore, the receptance matrices for sub-structures A and B and for structure C are:

$$\boldsymbol{H}^{A} = \begin{bmatrix} \boldsymbol{H}_{ii}^{A} & \boldsymbol{H}_{ij}^{A} \\ \boldsymbol{H}_{ji}^{A} & \boldsymbol{H}_{jj}^{A} \end{bmatrix}; \quad \boldsymbol{H}^{B} = \begin{bmatrix} \boldsymbol{H}_{jj}^{B} & \boldsymbol{H}_{jk}^{B} \\ \boldsymbol{H}_{kj}^{B} & \boldsymbol{H}_{kk}^{B} \end{bmatrix}; \quad \boldsymbol{H}^{C} = \begin{bmatrix} \boldsymbol{H}_{ii}^{C} & \boldsymbol{H}_{ij}^{C} & \boldsymbol{H}_{ik}^{C} \\ \boldsymbol{H}_{ji}^{C} & \boldsymbol{H}_{jj}^{C} & \boldsymbol{H}_{jk}^{C} \\ \boldsymbol{H}_{ki}^{C} & \boldsymbol{H}_{kj}^{C} & \boldsymbol{H}_{kk}^{C} \end{bmatrix}$$
(4)

Using equations (1) and (2), one obtains the receptance matrix for C [11]:

$$\boldsymbol{H}^{C} = \left(\begin{bmatrix} \boldsymbol{H}_{ii}^{A} & \boldsymbol{H}_{ij}^{A} \end{bmatrix}^{-1} & \boldsymbol{0} \\ \boldsymbol{H}_{ji}^{A} & \boldsymbol{H}_{jj}^{A} \end{bmatrix}^{-1} & \boldsymbol{0} \\ \boldsymbol{0} & \boldsymbol{0} & \boldsymbol{0} \end{bmatrix} + \begin{bmatrix} \boldsymbol{0} & \boldsymbol{0} & \boldsymbol{0} \\ \boldsymbol{0} & \begin{bmatrix} \boldsymbol{H}_{jj}^{B} & \boldsymbol{H}_{jk}^{B} \\ \boldsymbol{H}_{kj}^{B} & \boldsymbol{H}_{kk}^{B} \end{bmatrix}^{-1} \\ \boldsymbol{0} & \begin{bmatrix} \boldsymbol{H}_{jj}^{B} & \boldsymbol{H}_{jk}^{B} \\ \boldsymbol{H}_{kj}^{B} & \boldsymbol{H}_{kk}^{B} \end{bmatrix}^{-1} \end{bmatrix} \right)^{-1}$$
(5)

The process of inverting three matrices requires - in general - a high computational effort, implying a strong possibility of encountering ill-conditioned matrices. To try and minimize this problem an alternative formulation is often used [12]:

$$\boldsymbol{H}^{C} = \begin{bmatrix} \boldsymbol{H}_{ii}^{A} & \boldsymbol{H}_{ij}^{A} & \boldsymbol{0} \\ \boldsymbol{H}_{ji}^{A} & \boldsymbol{H}_{jj}^{A} & \boldsymbol{0} \\ \boldsymbol{0} & \boldsymbol{0} & \boldsymbol{H}_{kk}^{B} \end{bmatrix} - \begin{bmatrix} \boldsymbol{H}_{ij}^{A} \boldsymbol{H}_{ji}^{-1} \boldsymbol{H}_{ji}^{A} & \boldsymbol{H}_{ij}^{A} \boldsymbol{H}_{ji}^{-1} \boldsymbol{H}_{ji}^{A} & -\boldsymbol{H}_{ij}^{A} \boldsymbol{H}_{ji}^{-1} \boldsymbol{H}_{jk}^{B} \\ \boldsymbol{H}_{ji}^{A} \boldsymbol{H}_{ji}^{-1} \boldsymbol{H}_{ji}^{A} & \boldsymbol{H}_{ji}^{A} \boldsymbol{H}_{ji}^{-1} \boldsymbol{H}_{ji}^{A} & -\boldsymbol{H}_{ij}^{A} \boldsymbol{H}_{ji}^{-1} \boldsymbol{H}_{jk}^{B} \\ -\boldsymbol{H}_{kj}^{B} \boldsymbol{H}_{ji}^{-1} \boldsymbol{H}_{ji}^{A} & -\boldsymbol{H}_{kj}^{B} \boldsymbol{H}_{ji}^{-1} \boldsymbol{H}_{jk}^{B} \\ -\boldsymbol{H}_{kj}^{B} \boldsymbol{H}_{ji}^{-1} \boldsymbol{H}_{ji}^{A} & -\boldsymbol{H}_{kj}^{B} \boldsymbol{H}_{ji}^{-1} \boldsymbol{H}_{jk}^{B} \end{bmatrix}$$
(6)

where $\boldsymbol{H}_{jj} = \boldsymbol{H}_{jj}^{A} + \boldsymbol{H}_{jj}^{B}$. In a simpler way, one can write:

$$\begin{bmatrix} \boldsymbol{H}_{ii}^{C} & \boldsymbol{H}_{ij}^{C} & \boldsymbol{H}_{ik}^{C} \\ \boldsymbol{H}_{ji}^{C} & \boldsymbol{H}_{jk}^{C} \\ \boldsymbol{H}_{ki}^{C} & \boldsymbol{H}_{jk}^{C} \end{bmatrix} = \begin{bmatrix} \boldsymbol{H}_{ii}^{A} & \boldsymbol{H}_{ij}^{A} & \boldsymbol{0} \\ \boldsymbol{H}_{ji}^{A} & \boldsymbol{H}_{jj}^{A} & \boldsymbol{0} \\ \boldsymbol{0} & \boldsymbol{0} & \boldsymbol{H}_{kk}^{B} \end{bmatrix} - \begin{bmatrix} \boldsymbol{H}_{ij}^{A} \\ \boldsymbol{H}_{jj}^{A} \\ -\boldsymbol{H}_{jj}^{B} \end{bmatrix} \begin{pmatrix} \boldsymbol{H}_{ji}^{A} + \boldsymbol{H}_{jj}^{B} \end{pmatrix}^{-1} \begin{bmatrix} \boldsymbol{H}_{ji}^{A} & \boldsymbol{H}_{jk}^{A} & -\boldsymbol{H}_{jk}^{B} \end{bmatrix}$$
(7)

which replaces the three inversions by only one, restricted to the number of connection co-ordinates j, thereby limiting possible numerical problems.

2.2 FRF uncoupling

Let us suppose that our goal is the dynamic characterization of a joint; joints can be considered as complex components, often difficult to analyze and model. A possible solution relies on obtaining their dynamic behavior by an inverse coupling procedure, i.e., an uncoupling procedure, having eq. (7) as a starting point.

If our joint is defined as sub-structure B, one has co-ordinates i and j, whereas co-ordinates k (internal to B) do not play a role here. Therefore, eq. (7) is re-written as:

$$\begin{bmatrix} \boldsymbol{H}_{ii}^{C} & \boldsymbol{H}_{ij}^{C} \\ \boldsymbol{H}_{ji}^{C} & \boldsymbol{H}_{jj}^{C} \end{bmatrix} = \begin{bmatrix} \boldsymbol{H}_{ii}^{A} & \boldsymbol{H}_{ij}^{A} \\ \boldsymbol{H}_{ji}^{A} & \boldsymbol{H}_{jj}^{A} \end{bmatrix} - \begin{bmatrix} \boldsymbol{H}_{ij}^{A} \\ \boldsymbol{H}_{jj}^{A} \end{bmatrix} \begin{pmatrix} \boldsymbol{H}_{jj}^{A} + \boldsymbol{H}_{jj}^{B} \end{pmatrix}^{-1} \begin{bmatrix} \boldsymbol{H}_{ji}^{A} & \boldsymbol{H}_{jj}^{A} \end{bmatrix}$$
(8)

From eq. (8) it is clear that there are three possibilities for the evaluation of \boldsymbol{H}_{jj}^{B} : using \boldsymbol{H}_{ii}^{C} , \boldsymbol{H}_{ij}^{C} or \boldsymbol{H}_{jj}^{C} .

2.2.1 Without the use of co-ordinates *j*

Sometimes it may be difficult to undertake measurements on co-ordinates j, close to the joint, so ideally one should use measurements on the complete structure C involving only co-ordinates i. From (8), one has:

$$\boldsymbol{H}_{ii}^{C} = \boldsymbol{H}_{ii}^{A} - \boldsymbol{H}_{ij}^{A} \left(\boldsymbol{H}_{jj}^{A} + \boldsymbol{H}_{jj}^{B} \right)^{-1} \boldsymbol{H}_{ji}^{A}$$

$$\tag{9}$$

Rearranging (9),

$$\boldsymbol{H}_{ij}^{A} \left(\boldsymbol{H}_{jj}^{A} + \boldsymbol{H}_{jj}^{B} \right)^{-1} \boldsymbol{H}_{ji}^{A} = \boldsymbol{H}_{ii}^{A} - \boldsymbol{H}_{ii}^{C}$$
(10)

Pre-multiplying (10) by $\left(\boldsymbol{H}_{jj}^{A}\right)^{-1}$ and post-multiplying by $\left(\boldsymbol{H}_{ji}^{A}\right)^{-1}$ leads to:

$$\left(\boldsymbol{H}_{jj}^{A}+\boldsymbol{H}_{jj}^{B}\right)^{-1}=\left(\boldsymbol{H}_{ij}^{A}\right)^{-1}\left(\boldsymbol{H}_{ii}^{A}-\boldsymbol{H}_{ii}^{C}\right)\left(\boldsymbol{H}_{ji}^{A}\right)^{-1}$$
(11)

Note that this operation is mathematically possible only when *i* equals *j*; otherwise, if one of the (pseudo) inversions is possible, the other one is not, and *vice-versa*. Solving for \boldsymbol{H}_{ij}^{B} , it follows that

$$\boldsymbol{H}_{jj}^{B} = \boldsymbol{H}_{ji}^{A} \left(\boldsymbol{H}_{ii}^{A} - \boldsymbol{H}_{ii}^{C} \right)^{-1} \boldsymbol{H}_{ij}^{A} - \boldsymbol{H}_{jj}^{A}$$
(12)

Generalizing to the case where *i* might be different from *j* (in fact, $i \ge j$), one pre-multiplies equation (10) by an arbitrary matrix W_{ji} and post-multiplies it by W_{ij} :

$$\boldsymbol{W}_{ji}\boldsymbol{H}_{ji}^{A}\left(\boldsymbol{H}_{jj}^{A}+\boldsymbol{H}_{jj}^{B}\right)^{-1}\boldsymbol{H}_{ji}^{A}\boldsymbol{W}_{ij}=\boldsymbol{W}_{ji}\left(\boldsymbol{H}_{ii}^{A}-\boldsymbol{H}_{ii}^{C}\right)\boldsymbol{W}_{ij}$$
(13)

Rearranging eq. (13), one can write,

$$\left(\boldsymbol{H}_{jj}^{A}+\boldsymbol{H}_{jj}^{B}\right)^{-1} = \left(\boldsymbol{W}_{ji}\boldsymbol{H}_{ij}^{A}\right)^{-1}\boldsymbol{W}_{ji}\left(\boldsymbol{H}_{ii}^{A}-\boldsymbol{H}_{ii}^{C}\right)\boldsymbol{W}_{ij}\left(\boldsymbol{H}_{ji}^{A}\boldsymbol{W}_{ij}\right)^{-1}$$
(14)

This is only possible if $i \ge j$, not the other way around (which certainly is not a common case). From (14) one has:

$$\boldsymbol{H}_{jj}^{B} = \boldsymbol{H}_{ji}^{A} \boldsymbol{W}_{ij} \left(\boldsymbol{W}_{ji} \left(\boldsymbol{H}_{ii}^{A} - \boldsymbol{H}_{ii}^{C} \right) \boldsymbol{W}_{jj} \right)^{-1} \boldsymbol{W}_{ji} \boldsymbol{H}_{ij}^{A} - \boldsymbol{H}_{jj}^{A}$$
(15)

Now the order of the matrix to invert is equal to the number of co-ordinates *j*. This is advantageous because the size of *j* will be – in general – much smaller than that of *i*. The question is which matrix W_{ij} to use. There is not an easy answer, but probably the most logical one is to use H_{ij}^A :

$$\boldsymbol{H}_{jj}^{B} = \boldsymbol{H}_{ji}^{A} \boldsymbol{H}_{ij}^{A} \left(\boldsymbol{H}_{ji}^{A} \left(\boldsymbol{H}_{ii}^{A} - \boldsymbol{H}_{ii}^{C} \right) \boldsymbol{H}_{ij}^{A} \right)^{-1} \boldsymbol{H}_{ji}^{A} \boldsymbol{H}_{ij}^{A} - \boldsymbol{H}_{jj}^{A}$$
(16)

2.2.2 Using only the coordinates of the joint

Using eq. (7) based on \boldsymbol{H}_{jj}^{C} , one has:

$$\boldsymbol{H}_{jj}^{C} = \boldsymbol{H}_{jj}^{A} - \boldsymbol{H}_{jj}^{A} \left(\boldsymbol{H}_{jj}^{A} + \boldsymbol{H}_{jj}^{B} \right)^{-1} \boldsymbol{H}_{jj}^{A}$$
(17)

Solving eq. (17) with respect to H_{jj}^{B} , it follows that:

$$\boldsymbol{H}_{jj}^{B} = \left(\boldsymbol{H}_{jj}^{A} \left(\boldsymbol{H}_{jj}^{A} - \boldsymbol{H}_{jj}^{C}\right)^{-1} - \boldsymbol{I}_{jj}\right) \boldsymbol{H}_{jj}^{A}$$
(18)

Based on the dynamic stiffness matrix of the structure, Ambrogio [1] obtains an alternative formulation (eq. (19) can be derived from eq. (18)):

$$\boldsymbol{H}_{jj}^{B} = \left(\boldsymbol{I}_{jj} - \boldsymbol{H}_{jj}^{C} \left(\boldsymbol{H}_{jj}^{A}\right)^{-l}\right)^{-1} \boldsymbol{H}_{jj}^{C}$$
(19)

2.2.3 Using coordinates *i* and *j*

Back again to eq. (8), the third option to evaluate H_{jj}^{B} starts from:

$$\boldsymbol{H}_{ij}^{C} = \boldsymbol{H}_{ij}^{A} - \boldsymbol{H}_{ij}^{A} \left(\boldsymbol{H}_{jj}^{A} + \boldsymbol{H}_{jj}^{B} \right)^{-1} \boldsymbol{H}_{jj}^{A}$$
(20)

or

$$\boldsymbol{H}_{ij}^{A} \left(\boldsymbol{H}_{jj}^{A} + \boldsymbol{H}_{jj}^{B} \right)^{-1} \boldsymbol{H}_{jj}^{A} = \boldsymbol{H}_{ij}^{A} - \boldsymbol{H}_{ij}^{C}$$
(21)

Using once again arbitrary matrices, now W_{ji} and W_{jj} , with *i* greater or equal to *j*, one has

$$\boldsymbol{W}_{ji}\boldsymbol{H}_{ij}^{A}\left(\boldsymbol{H}_{jj}^{A}+\boldsymbol{H}_{jj}^{B}\right)^{-1}\boldsymbol{H}_{jj}^{A}\boldsymbol{W}_{jj}=\boldsymbol{W}_{ji}\left(\boldsymbol{H}_{ij}^{A}-\boldsymbol{H}_{ij}^{C}\right)\boldsymbol{W}_{jj}$$
(22)

from which

$$\left(\boldsymbol{H}_{jj}^{A}+\boldsymbol{H}_{jj}^{B}\right)^{-1}=\left(\boldsymbol{W}_{ji}\boldsymbol{H}_{ij}^{A}\right)^{-1}\boldsymbol{W}_{ji}\left(\boldsymbol{H}_{ij}^{A}-\boldsymbol{H}_{ij}^{C}\right)\boldsymbol{W}_{jj}\left(\boldsymbol{H}_{jj}^{A}\boldsymbol{W}_{jj}\right)^{-1}$$
(23)

Solving in order to \boldsymbol{H}_{jj}^{B} , gives:

$$\boldsymbol{H}_{jj}^{B} = \boldsymbol{H}_{jj}^{A} \boldsymbol{W}_{jj} \left(\boldsymbol{W}_{ji} \left(\boldsymbol{H}_{ij}^{A} - \boldsymbol{H}_{ij}^{C} \right) \boldsymbol{W}_{jj} \right)^{-1} \boldsymbol{W}_{ji} \boldsymbol{H}_{ij}^{A} - \boldsymbol{H}_{jj}^{A}$$
(24)

As before, one shall make $W_{ji} = H_{ji}^A$ and $W_{jj} = H_{jj}^A$:

$$\boldsymbol{H}_{jj}^{B} = \boldsymbol{H}_{jj}^{A} \boldsymbol{H}_{jj}^{A} \left(\boldsymbol{H}_{ji}^{A} - \boldsymbol{H}_{ij}^{C} \right) \boldsymbol{H}_{jj}^{A} \right)^{-1} \boldsymbol{H}_{ji}^{A} \boldsymbol{H}_{ij}^{A} - \boldsymbol{H}_{jj}^{A}$$
(25)

2.2.4 Summary

In summary, one has three formulations that allow us to determine matrix H_{ii}^{B} :

First formulation

$$\boldsymbol{H}_{jj}^{B} = \boldsymbol{H}_{ji}^{A} \boldsymbol{H}_{ij}^{A} \left(\boldsymbol{H}_{ji}^{A} \left(\boldsymbol{H}_{ii}^{A} - \boldsymbol{H}_{ii}^{C} \right) \boldsymbol{H}_{ij}^{A} \right)^{-1} \boldsymbol{H}_{ji}^{A} \boldsymbol{H}_{ij}^{A} - \boldsymbol{H}_{jj}^{A}$$
(16)

Second formulation

$$\boldsymbol{H}_{jj}^{B} = \left(\boldsymbol{H}_{jj}^{A} \left(\boldsymbol{H}_{jj}^{A} - \boldsymbol{H}_{jj}^{C}\right)^{-1} - \boldsymbol{I}_{jj}\right) \boldsymbol{H}_{jj}^{A}$$
(18)

Third formulation

$$\boldsymbol{H}_{jj}^{B} = \boldsymbol{H}_{jj}^{A} \boldsymbol{H}_{jj}^{A} \left(\boldsymbol{H}_{ji}^{A} - \boldsymbol{H}_{ij}^{C} \right) \boldsymbol{H}_{jj}^{A} \right)^{-1} \boldsymbol{H}_{ji}^{A} \boldsymbol{H}_{ij}^{A} - \boldsymbol{H}_{jj}^{A}$$
(25)

3 SIMULATION STUDIES

To illustrate the performance of the three formulations one uses the beam of figure 2, constituted by three components, A_1 , B and A_2 , forming structure C; A_1 and A_2 form sub-structure A.



The aim is to characterize component *B* (our "joint"), evaluating H^B , assuming that H^A is calculated analytically and H^C is calculated through experiments (simulated, in this case).

Using the finite element method with beam elements with four degrees of freedom, each component is divided into eight elements and considering only the nodes shown in figure 2 as coordinates i and j where it is possible to measure and excite the structure. The characteristics of each component are displayed in table 1.

Tuble T Characteristics of the components of the count							
Beam	Length	Width	Thickness	ν	E	ρ	
A_1	300 mm	25 mm	3 mm	0.3	210 GPa	7850 Kg/m ³	
В	400 mm	25 mm	6 mm	0.3	210 GPa	7850 Kg/m ³	
A ₂	300 mm	25 mm	3 mm	0.3	210 GPa	7850 Kg/m ³	

Table 1 Characteristics of the components of the beam

3.1 Choice of the formulation

To simulate the experimental errors, one has imposed a 1% perturbation in matrix \boldsymbol{H}^{C} for the three formulations. The elements 11 and 12 (chosen just for comparison) of matrix \boldsymbol{H}_{jj}^{B} obtained with the three formulations are shown in figure 3.

The major disturbances in the responses are due to numerical problems in the matrix inversions in each of the formulations. The variability of the results in figure 3 is quite high. Figure 4 represents the module of the differences between the numerically exact response and the response obtained by the three formulations. From the three results, the second formulation is the one that produces the smallest error throughout the frequency range.



Fig. 4 Errors of *FRF*s H_{11}^B and H_{12}^B for the three formulations

Let us represent H_{11}^{B} and H_{12}^{B} only for the second formulation, superimposed with the numerically exact response (see figure 5).



Fig. 5 H_{11}^{B} and H_{12}^{B} for the second formulation

3.2 Strategies to improve the results

It is clear that one cannot be happy with the results of figure 5. Let us try to improve them. From eq. (18) it is apparent that the problems certainly arise in the inversion of $(H_{ii}^A - H_{ii}^C)$, namely when this difference is small. To try to increase this difference one will change our structure by adding point masses, so to change the behavior of sub-structure A and C, while B remains unchanged.

3.2.1 Adding mass to sub-structure A

A mass of 35 grams has been added to nodes 1, 3, 5 and 15, 17, 19. The new results for H_{11}^B and H_{12}^B are shown in figure 6.



Fig. 6 H_{11}^{B} and H_{12}^{B} with added masses in sub-structure A

It can be observed from figure 6 that the disturbance observed between 1000-1200 Hz in figure 5 moves into the range 800-1000 Hz, generally improving the results.

3.2.2 Adding mass to sub-structure B

An alternative could be to add mass to the sub-structure B itself. Adding 35 grams at co-ordinates j = 9 and 11, leads to the results presented in figure 7.



As sub-structure B is altered, its natural frequency (identified in figure 5) changes to the left. However, the disturbances remain in the area of 1000-1200 Hz and one can conclude that they are caused caused by sub-structure A. So maybe one can improve the results by adding some more mass to sub-structure B, to cause the natural frequency to deviate from the disturbance. Let us now add masses of 70 grams. The results are displayed in figure 8.



The results are clearly better. However, to recover the dynamic response of B, one has to uncouple the added masses. The results of such an operation are shown in figure 9.



Although the results are better than the initial ones (figure 5), they are worse than those of figure 6, when the masses were added to sub-structure *A*. The disturbances come up again in the same frequency range.

3.3 Coupling

One of the main interests of the dynamic characterization of a sub-structure (like a joint) is to be able to predict the dynamic behavior of another structure (or a modified one), possibly a more complex one, inserting (coupling) the identified results from the uncoupling procedure. Whereas the uncoupling procedure tends to be unstable, as shown in this work, the coupling process is usually quite stable.

Based on the results obtained for sub-structure B, a coupling procedure will be undertaken with similar components, two beams A1 and A2 but now with a length of 400 mm. Using the results initially obtained from figure 5, one obtains the behavior illustrated in figure 10.

In figure 10 the disturbance in the area of 1000-1200 Hz remains. One can also observe the results of the coupling of B when the masses were added to A, i.e., when using the responses given in figure 6. Figure 11 illustrates this case. The results have improved and the disturbances have moved down, as expected. Most certainly, better solutions would have been obtained if one had performed a modal analysis identification to the results of figure 6, prior to the coupling procedure.



Fig. 10 H_{11}^{C} and H_{12}^{C} resulting from the coupling of A1 and A2 with the sub-structure B



Fig. 1 H_{II}^{C} and H_{I2}^{C} resulting from the coupling of A1 and A2 with the sub-structure B, with the addition of masses in A

4 CONCLUSIONS

In this paper the authors have presented three formulations for the uncoupling of sub-structures, something that may be of considerable interest, for instance when trying to model complex joints. The formulation that presented the best results requires measurements at the connection points of the structures; unfortunately, this may not always be possible in practice. Any of the three formulations revealed to be numerically unstable due to the inversion of difference matrices. This problem has been investigated in this study and improvements have been obtained when adding point masses to the remaining sub-structures other than the one to be characterized. Those added masses move the natural frequencies, allowing to understand the problems that are happening and as already said, improving the results. Together with the use of modal identification methods, the uncoupling techniques may be of interest in various situations. Experimental implementation still has to be further investigated, as the accurate measurement of rotations is quite difficult to obtain.

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