

Acoustic displacement triangle based on the individual element test

Abstract

A three node -displacement based- acoustic element is developed. In order to avoid spurious rotational modes, a higher order stiffness is introduced. This higher order stiffness is developed from an incompatible strain field which computes element volume changes under nodal rotational displacements fields. The higher order strain resulting from the incompatible strain field satisfies the Individual Element Test (IET) requirements without affecting convergence. The higher order stiffness is modulated, element by element, with a factor. As a result, the displacement based formulation presented on this paper is capable of placing the spurious rotational modes above the range of the physical compressional modes that can be accurately calculated by the mesh.

Keywords

Fluid-structure interaction, Displacement-based formulation, Spurious modes, Individual Element Test, Parameterized Variational Principle

S. Correa*, C. Militello** and M. Recuero***

*Department of Product Design Engineering. EAFIT University, Medellín-Colombia

**Department of Fundamental Physics, Experimental, Electronics and Systems. Universidad de La Laguna, S/C de Tenerife, Spain

***Department of Mechanics. School of Industrial Engineers. Universidad Politécnica de Madrid. Madrid, Spain

Received 27 Mar 2012;
In revised form 16 Apr 2012

* Author email: scorrea5@eafit.edu.co

1 INTRODUCTION

Acoustic propagation in an inviscid media is generally studied using the pressure p as primitive variable. Consequently, after the finite element approximation, only one unknown variable per node is obtained. This is a drastic reduction on the number of unknown variables as compared to a displacement based formulation where two unknown displacements u, v at each node are necessary to describe the problem. The only drawback in using a pressure formulation can be found when the fluid is interfaced with an elastic solid because both of them do not share the same variables. To overcome this issue, an equilibrium constraint must be imposed at the fluid-solid boundary. Therefore, an acoustic fluid element cannot be handled by a finite element code as any other structural element. Displacement, pressure, displacement-pressure and velocity potential elements have been developed in the past. References [2–4, 7, 9, 10, 12–14] are, in the authors' opinion, the most relevant.

Displacement formulations have been reported by many authors. The formulation of these elements following finite element standard procedures produces a rank deficient stiffness matrix. This rank deficiency is not an error because the strain energy in an acoustic fluid is only

computed from volume changes. The displacement field inside the element computes element shape changes with no volume changes. Any displacement field that describes shape changes without volume variation expands the null space of the stiffness matrix. This displacement field is called spurious because, despite being consistent with the formulation, it is not expected in a true irrotational formulation. Moreover, when an eigenfrequency problem is solved, the spurious displacement field produces low frequency rotational modes. Many authors [2, 10, 14] compute the rotational of the displacement field and add a fictitious rotational energy. There is not a constitutive equation for such behavior and a fictitious elastic coefficient must be introduced, appearing in the formulation as a penalty factor. Generally, the value suggested for the factor ranges from 1 to 1,000 times the compressibility modulus. This variability of the factor is a serious drawback because the results strongly depend on this *a priori* tuning.

In this paper, a displacement-based triangular element is proposed. In order to obtain an irrotational displacement field in weak form, an incompatible mode is added. From the added incompatible mode, a higher order strain is computed. After filtering the higher order strain to satisfy the IET from Bergan [5, 11], a higher order stiffness matrix is assembled. No rotational measure is introduced and no fictitious elastic coefficient is needed. However, it is possible to introduce a coefficient in front of the resulting higher order stiffness. The coefficient changes from element to element without affecting convergence. A stabilization strategy is proposed and the coefficient is computed in a closed form as a function of the element size. In this way, the spurious modes are placed over the higher compressional mode that still retains a physical meaning. This formulation is not based in the Raviart-Thomas polynomial [6] nor in the displacement/pressure formulation [2, 14], hence no centerface nor midside degrees of freedom are necessary, resulting in lower computational cost.

A 2D acoustic cavity and a 2D fluid interaction problem are presented to show convergence, non uniform meshes behavior and boundary normal definition issues [2, 14]. The 2D acoustic cavity is also used to check the appearance of spurious modes.

2 THE STIFFNESS MATRIX ASSEMBLY

2.1 Displacement field and incompatible modes

To describe the displacement field, the same orientation of the coordinate system is assumed for all the elements. Each element uses a system with origin at its center of gravity. In a linear triangle the general form for the displacement field is:

$$u_c = \begin{pmatrix} u_c \\ v_c \end{pmatrix} = \begin{pmatrix} a_1 + a_2x + a_3y \\ b_1 + b_2x + b_3y \end{pmatrix} \quad (1)$$

Where \mathbf{u} is the displacement field and the subscript c stands for the compatible part. The field is linear and the a and b factors can be obtained as a function of the nodal displacements values u_j, v_j with $j = 1, 2, 3$ and the nodal coordinates [15].

The following incompatible modes are introduced,

$$u_i = \begin{pmatrix} u_i \\ v_i \end{pmatrix} = \begin{pmatrix} \varsigma g(x, y) \\ \eta g(x, y) \end{pmatrix} \quad (2)$$

Where \mathbf{u} is the displacement field and the subscript i stands for the incompatible part and

$$g(x, y) = xy^2 + yx^2 \quad (3)$$

Constants ς and η must be determined element by element and take values different from zero when the fluid tends to rotate.

A trial displacement field, $u^* = u_c + u_i$ is proposed in order to obtain ς and η as a function of the a and b coefficients.

The rotor of the trial field is computed as:

$$\nabla \times u^* = (u_{,y}^* - v_{,x}^*) = ((a_3 + \varsigma g_{,y}) - (b_2 + \eta g_{,x})) \quad (4)$$

From (4), it becomes clear that asking the rotor to cancel at each point inside the element will not produce the sufficient conditions. Instead, the following weak form is tried:

$$\int_{\Omega_e} (x^2 + y^2)(u_{,y}^* - v_{,x}^*) d\Omega = 0 \quad (5)$$

By integrating and rearranging terms, (5) results in the following matrix relationship:

$$\begin{pmatrix} f_{11} & f_{12} \\ f_{21} & f_{22} \end{pmatrix} \begin{pmatrix} \varsigma \\ \eta \end{pmatrix} = \begin{pmatrix} p_{11} & p_{12} \\ p_{21} & p_{22} \end{pmatrix} \begin{pmatrix} a_3 \\ b_2 \end{pmatrix} \quad (6)$$

Which, in compact form, becomes:

$$\mathbf{F}\psi = \mathbf{P}\mathbf{d} \quad (7)$$

Where ψ and \mathbf{d} are vectors and the terms of matrix \mathbf{F} are:

$$\begin{aligned} f_{11} &= \int_{\Omega_e} y^2(2yx + x^2) d\Omega \\ f_{12} &= - \int_{\Omega_e} y^2(y^2 + 2yx) d\Omega \\ f_{21} &= \int_{\Omega_e} x^2(2yx + x^2) d\Omega \\ f_{22} &= - \int_{\Omega_e} x^2(y^2 + 2yx) d\Omega \end{aligned} \quad (8)$$

And the terms of matrix \mathbf{P} are:

$$\begin{aligned} p_{11} &= \int_{A_e} x^2 dA \\ p_{12} &= - \int_{A_e} x^2 dA \\ p_{13} &= \int_{A_e} y^2 dA \\ p_{14} &= - \int_{A_e} y^2 dA \end{aligned} \quad (9)$$

By nodal collocation, a linear relationship can be obtained between vector \mathbf{d} and the nodal displacements \mathbf{v} :

$$\mathbf{d} = \mathbf{Q} \mathbf{v} \quad (10)$$

Where matrix \mathbf{Q} is:

$$\mathbf{Q} = \frac{1}{\det[\mathbf{J}]} \begin{bmatrix} x_{32} & 0 & x_{13} & 0 & x_{21} & 0 \\ 0 & y_{23} & 0 & y_{31} & 0 & y_{12} \end{bmatrix} \quad (11)$$

Being x_i, y_i the nodal coordinates and $\mathbf{v}^t = (u_1 \ v_1 \ u_2 \ v_2 \ u_3 \ v_3)$

Assuming \mathbf{F}^{-1} exists, equation (6) shows that ψ is a null vector for nodal displacements that produces volume changes because it is independent from a_2 and b_3 , and $\psi \neq 0$ for rotational displacements fields. Thus, rotational fields activate the incompatible modes.

For simplicity, replacing (10) in (7):

$$\psi = \mathbf{F}^{-1} \mathbf{P} \mathbf{Q} \mathbf{v} = \mathbf{R} \mathbf{v} \quad (12)$$

2.2 The strain measure.

In a displacement based acoustic element the only strain measure is the unitary change of volume:

$$e = u_{,x} + v_{,y} \quad (13)$$

The pressure inside the element is computed as:

$$p = \beta e \quad (14)$$

where β is the compressibility modulus of the fluid. In (14) the continuous mechanics convention is used, i.e., a negative change of volume is associated with a negative pressure (stress).

For the element strain two contributions are considered, one from the compatible part of the displacement, e_c , and a higher order one, e_h , computed from the incompatible modes u_i . The computation of e_h is not simple because it follows the rules presented in [11], in order to satisfy the IET.

First, a unitary volume change is computed from the incompatible modes:

$$e_i = \varsigma g_{,x} + \eta g_{,y} = \begin{pmatrix} g_{,x} & g_{,y} \end{pmatrix} \begin{pmatrix} \varsigma \\ \eta \end{pmatrix} = \mathbf{G} \psi \quad (15)$$

The mean volume averaged strain is computed as:

$$\bar{e}_i = \frac{1}{\Omega_e} \int_{\Omega_e} \begin{pmatrix} g_{,x} & g_{,y} \end{pmatrix} d\Omega \psi = \bar{\mathbf{G}} \psi \quad (16)$$

By subtracting (16) from (15) and replacing (12), the higher order strain field is obtained:

$$e_h = e_i - \bar{e}_i = (\mathbf{G} - \bar{\mathbf{G}}) \psi = (\mathbf{G} - \bar{\mathbf{G}}) \mathbf{R} \mathbf{v} \quad (17)$$

This, in a more standard notation, becomes:

$$e_h = (G - \bar{G})Rv = B_h v \quad (18)$$

From which the higher order stiffness can be computed as:

$$K_h = \int_{\Omega_e} B_h^t \beta B_h d\Omega \quad (19)$$

This higher order stiffness matrix maintains the irrotationality of the fluid. Additionally, the basic stiffness computes the constant strain state, in this case a change of volume, and produces zero energy under rigid body motions, i.e., translation and rotation. This is achieved using \mathbf{u}_c in the more usual shape function expansion:

$$\mathbf{u}_c = \mathbf{N}^t \mathbf{v} \quad (20)$$

From this displacement field, the basic strain field is defined as:

$$e_b = \begin{pmatrix} N_{1,x} & N_{1,y} & \dots & N_{3,x} & N_{3,y} \end{pmatrix} \mathbf{v} = B_b \mathbf{v}, \quad (21)$$

and the basic stiffness is computed,

$$K_b = \int_{\Omega_e} B_b^t \beta B_b d\Omega \quad (22)$$

The total element stiffness is calculated by adding the basic and higher order contributions:

$$K_e = K_b + \alpha K_h \quad (23)$$

The value of α can be changed from element to element without affecting the capability of the assembly to correctly define a constant strain state and rigid body modes [8].

3 STABILIZATION PROCEDURE

As mentioned in the introduction, one of the disadvantages of the previous formulations is the selection of the penalty factor α . A low factor will contaminate the correct modes, whereas a high factor will conceal the contribution of the basic stiffness.

Considering that the basic stiffness computes the irrotational modes from a linear displacement field, it seems reasonable to assume that three elements in a line is the limit to correctly capture half the shortest wavelength. Under these circumstances, it is proposed that the energy computed for a rotational mode of the same wavelength should be of the same order as the energy computed for an irrotational mode. In order to make both energies comparable the eigenfrequencies are asked to match. The displacement fields that produce rotational and irrotational modes are defined as:

$$\mathbf{u}_{ir}^t = \begin{pmatrix} u_{ir} & v_{ir} \end{pmatrix} = \begin{pmatrix} \sin\left(\frac{\pi x}{\lambda}\right) & 0 \end{pmatrix} \quad (24)$$

$$\mathbf{u}_r = \begin{pmatrix} u_r \\ v_r \end{pmatrix} = \begin{pmatrix} \sin\left(\frac{\pi x}{\lambda}\right) \cos\left(\frac{\pi y}{\lambda}\right) \\ -\sin\left(\frac{\pi y}{\lambda}\right) \cos\left(\frac{\pi x}{\lambda}\right) \end{pmatrix} \quad (25)$$

Where subscript *ir* refers to the irrotational displacement field and subscript *r* refers to the rotational displacement field. These fields satisfy $\nabla \times \mathbf{u}_{ir} = 0$ and $\nabla \cdot \mathbf{u}_r = 0$.

To compute the value of α in (23) a regular mesh, as shown in Figure 1, is used. The displacement fields are computed at the mesh nodes and the values are arranged in vectors \mathbf{U}_{ir} and \mathbf{U}_r .

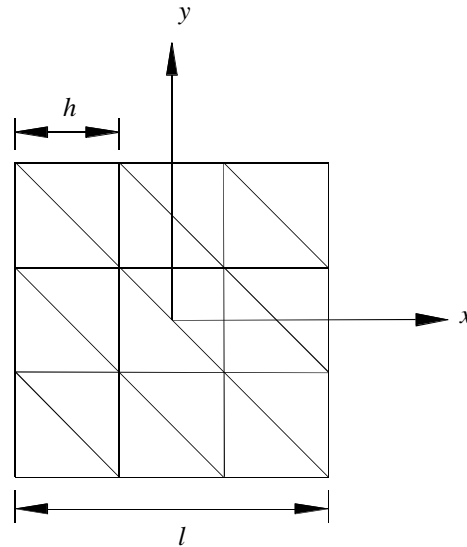


Figure 1 Sample surface mesh used to compute stabilization coefficient α .

The approximation to the eigenvalue (ω) is obtained from the Rayleigh quotient. In both cases, a lumped diagonal mass matrix (\mathbf{M}) is used.

$$\omega_{ir}^2 = \frac{\mathbf{U}_{ir}^T \mathbf{K}_b^a \mathbf{U}_{ir}}{\mathbf{U}_{ir}^T \mathbf{M}^a \mathbf{U}_{ir}} \quad (26)$$

$$\omega_r^2 = \alpha \frac{\mathbf{U}_r^T \mathbf{K}_h^a \mathbf{U}_r}{\mathbf{U}_r^T \mathbf{M}^a \mathbf{U}_r} \quad (27)$$

The superscript *a* indicates the use of the assembled element stiffness matrix from Figure 1. It should be noted that the irrotational field expands the null space of the higher order stiffness matrix, therefore α does not contribute to equation (26). On the other hand, the rotational field expands the null space of the basic stiffness matrix without contributing to equation (27). By equating both frequencies the following expression for α is obtained:

$$\alpha = \frac{\mathbf{U}_{ir}^t \mathbf{K}_b^a \mathbf{U}_{ir}}{\mathbf{U}_r^t \mathbf{K}_h^a \mathbf{U}_r} \frac{\mathbf{U}_r^t \mathbf{M}^a \mathbf{U}_r}{\mathbf{U}_{ir}^t \mathbf{M}^a \mathbf{U}_{ir}} \quad (28)$$

The value of α is independent from physical fluid properties. Equation (28) is computed from the mesh shown in Figure 1. The quadrilateral side is $\lambda/2$, being λ the wavelength, and the relation between quadrilateral side and element side (h) satisfies:

$$\frac{\lambda}{2} = 3h \quad (29)$$

Quadrilaterals are constructed for values of h ranging from 0.01 to 10 meters. Rotational and irrotational fields for the corresponding λ are computed at the mesh nodes. Finally, equation (28) is evaluated. Hence, a value of α is computed for each value of h .

The adjusted function for the pairs(α, h) , for h in meters is:

$$\alpha = \frac{12.9}{h^2} \quad (30)$$

Equation (30) is implemented in the finite element code and is calculated for each element. To compute the element size h , the diameter of the circle inscribed in the triangle is used. This method for computing α is conservative because the diameter of the circle is always smaller than the smallest triangle side.

4 NUMERICAL RESULTS

The new formulated element is used to solve two 2D problems in order to assess its performance. One consists of a closed rectangular cavity and the other consists of a closed cavity with a skewed corner and a rigid moving piston. In both cases, convergence of the solution varying the element size of the mesh is presented. Since the definition of the normal direction in the solid – fluid interface is critical when imposing impenetrability conditions, the problem of a skewed cavity with moveable piston is also solved for a random variation of that normal direction. To test convergence, four finite element meshes are tried. The first three have one predominant element size and the fourth one has three predominant element sizes in order to demonstrate that the results do not depend on maintaining a constant element size.

5 CLOSED RECTANGULAR CAVITY.

When a fluid-structure interaction problem is solved using the displacement based proposed element, it is necessary to impose the impenetrability condition in the interface. For the closed rectangular cavity problem discussed here, the condition is imposed directly in the fluid element restraining the displacement perpendicular to the rigid wall.

The fluid properties and dimensions are shown in Figure 2. The eigenfrequencies can be computed from

$$f = c\pi\sqrt{\left(\left(\frac{l}{a}\right)^2 + \left(\frac{m}{b}\right)^2\right)} \quad (31)$$

Where f is the frequency in [Hz], c is the speed of sound in water [m/s], a is the width in [m], b is the height in [m], l and $m = 0,1,2,\dots$

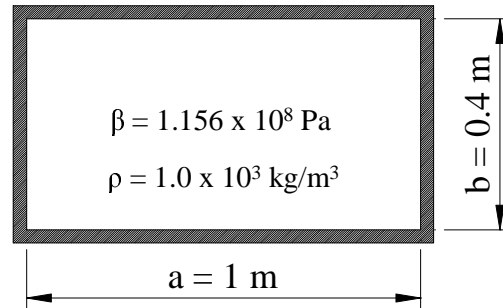


Figure 2 Rectangular cavity. Dimensions and fluid properties.

Four meshes are presented in Figure 3 (left). The first three have one predominant element size (coarse, medium and fine) and the fourth one has three predominant element sizes in the same mesh. The modal pressure distribution for mode 1 is shown at the right side. In Table 1, the convergence to the first four modes is shown.

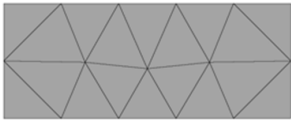
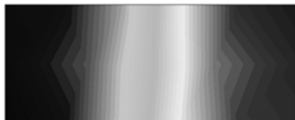
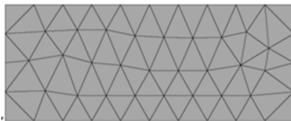
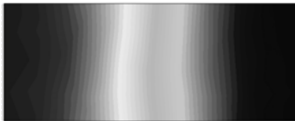
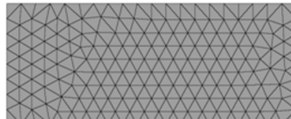

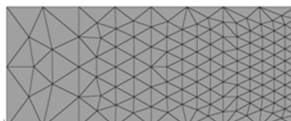
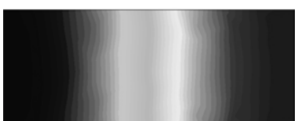
Mesh	Mode
 17 nodes - 18 elements	 169.1 Hz
 54 nodes - 78 elements	 170 Hz
 202 nodes - 346 elements	 170 Hz
 137 nodes - 244 elements	 169.8 Hz

Figure 3 Rectangular cavity results. Meshes at left. Pressure distribution for the 1th mode at right.

5.1 Spurious mode placement

According to the stabilization procedure proposed and noting that the coarse mesh from Figure 3 has only two elements in the vertical direction, the first spurious mode must appear after

Table 1 Resonant frequencies for a room with rigid walls. Convergence analysis

Mesh nodes	1st mode $f = 170$ Hz	2nd mode $f = 340$ Hz	3rd mode $f = 425$ Hz	4th mode $f = 457.7$ Hz
17	169.1	350.9	384.1	478.1
54	170	341.3	421	470.5
202	170	340.3	423.3	463.7
137	169.8	340.2	430.3	459.5

the first half wavelength in that direction. Figures 4a and 4b show the first correct modes in the longitudinal and vertical direction respectively. Figures 4c and 4d show modes with a high tendency to rotate, i.e., these are incorrect (spurious) modes. Also note that the mode in 4b is the last one to be correctly calculated using three elements in a half wavelength. In this way, the spurious modes can be located at a convenient place, i.e., with an eigenfrequency higher than the last one that can be adequately calculated by the mesh.

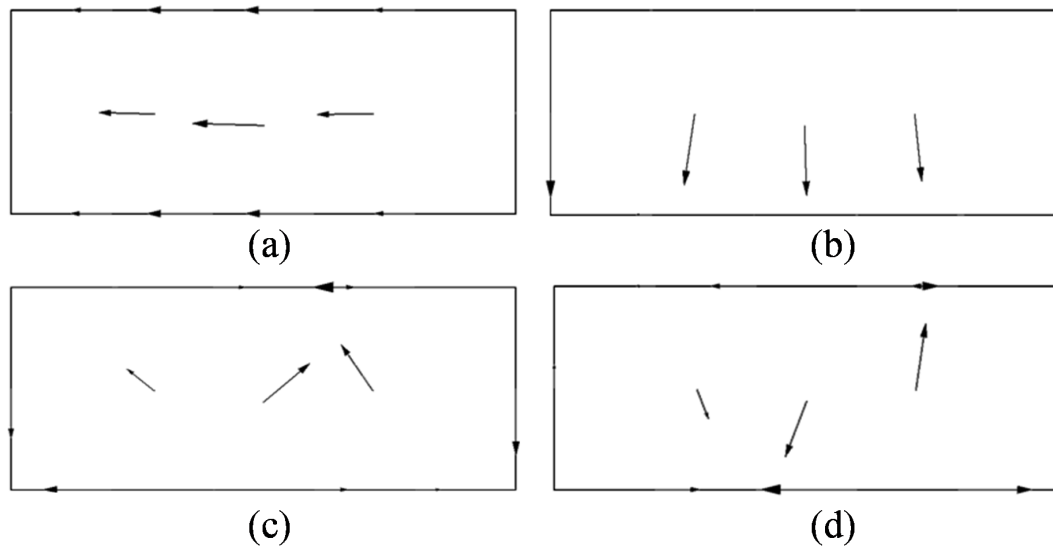


Figure 4 Spurious modes appearance in a rectangular cavity. (a) First longitudinal mode. (b) First vertical mode (3^{rd} mode). (c-d) 5^{th} and 6^{th} modes with a high rotational tendency.

5.2 Element size sensitivity

It is necessary to test the sensitivity of the proposed formulation for element sizes (h) out of the range of equation 30 (0.01 to 10 m.). In consequence, two meshes with predominant element sizes of 0.001 m. and 100 m. respectively and having the same number of nodes and elements are tried. The results are summarized in Table 2.

Table 2 Resonant frequencies for a room with rigid walls. Element size sensitivity

Mesh nodes	h [m]	α	1st mode $f = 170$ Hz	2nd mode $f = 340$ Hz	3rd mode $f = 425$ Hz	4th mode $f = 457.7$ Hz
202	0.001	12.9×10^6	170	340.8	423.9	464.1
	100	12.9×10^{-4}	170	340	423.1	463.4

6 SKEWED CAVITY WITH RIGID MOVABLE PISTON.

A common fluid-structure interaction problem [2], consisting of a skewed rigid piston capable of moving back and forth is presented in Figure 5. For such problem, the impenetrability condition is imposed directly in the fluid element restraining the displacement perpendicular to the rigid wall. Furthermore, the nodes in the piston-fluid interface are forced to move together via Lagrange multipliers.

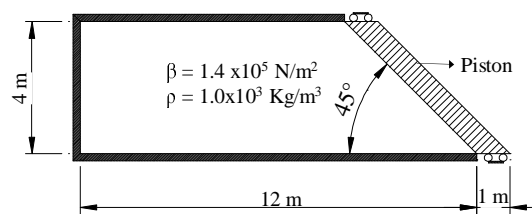


Figure 5 Skewed cavity coupled with movable piston.

Figure 6 shows the meshes and the results obtained for the 3rd and 4th modes.

In Table 3, the convergence of the first four modes is shown. A coarse mesh captures a vertical and a longitudinal mode respectively due to a lack of convergence. A fine mesh captures the correct modes [2] that result from a linear combination of the previous ones. The eigensolver is Eispack through the Matlab interface [1].

Table 3 Resonant frequencies for the skew cavity with rigid piston. Convergence analysis

Mesh nodes	1st mode $f = 0.29$ Hz	2nd mode $f = 0.88$ Hz	3rd mode $f = 1.45$ Hz	4th mode $f = 1.48$ Hz
114	0.29	0.86	1.36	1.41
498	0.29	0.87	1.44	1.48
2362	0.29	0.88	1.45	1.48
1358	0.29	0.90	1.46	1.48

For the piston-fluid interface, the normal direction is randomly varied ± 5 degrees. The effect in the convergence is shown in Table 4.

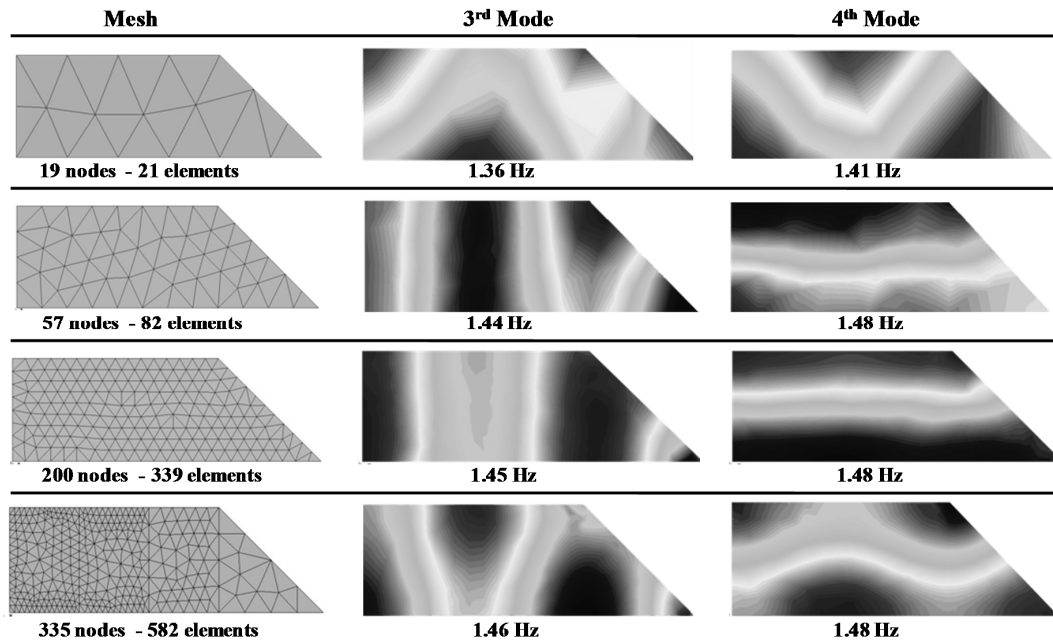


Figure 6 Skew cavity convergence analysis.

Table 4 Resonant frequencies for the skew cavity with rigid piston. Convergence analysis with ± 5 randomly changed interface normal directions.

Mesh nodes	1st mode $f = 0.29$ Hz	2nd mode $f = 0.88$ Hz	3rd mode $f = 1.45$ Hz	4th mode $f = 1.48$ Hz
114	0.29	0.86	1.37	1.41
498	0.29	0.87	1.44	1.48
2362	0.29	0.88	1.45	1.47
1358	0.29	0.89	1.45	1.5

7 DISCUSSION AND CONCLUSIONS.

The element proposed is a 2D linear triangle with degrees of freedom that can be easily coupled to 2D solid elements. Although it has a penalty factor, the proposed energy balance formulation can be implemented directly in the finite element code without user intervention. The penalty factor depends on the element size and does not influence the element convergence. This results in a clear advantage with respect to other fluid-structure interaction elements in which the factor must be selected by the user [2, 10, 14] and can vary between 1 and 1,000 times the compressibility modulus of the fluid.

The convergence is not altered by a reasonable error in the normal direction definition. The spurious modes do not disappear, as in previous formulations [2, 6, 10, 14]. Instead, the penalty factor places spurious modes in frequencies higher than the ones corresponding to the last compression mode accurately calculated by the mesh.

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