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Modeling and Measurements of CMUTs with Square Anisotropic Plates

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Abstract—The conventional method of modeling CMUTs use the isotropic plate equation to calculate the deflection, leading to deviations from FEM simulations including anisotropic effects of around 10% in center deflection. In this paper, the deflection is found for square plates using the full anisotropic plate equation and the Galerkin method. Utilizing the symmetry of the silicon crystal, a compact and accurate expression for the deflection can be obtained. The deviation from FEM in center deflection is <0.1%. The deflection was measured on fabricated CMUTs using a white light interferometer. Fitting the anisotropic calculated deflection to the measurement a deviation of 0.5-1.5% is seen for the fitted values. Finally it was also measured how the device behaved under increasing bias voltage and it is observed that the fitted values. It was also measured how the device behaved under increasing bias voltage and it is observed that the model including anisotropic effects is within the uncertainty interval of the measurements.

I. INTRODUCTION

Precise modeling of capacitive micromachined ultrasonic transducers (CMUT) is important for an efficient design process. The deflection \( w(x,y) \) is an important parameter that influences several basic CMUT parameters such as pull-in voltage and capacitance. Most existing analytical approaches use the isotropic plate equation to calculate the deflection [1], [2]. However, when using fusion bonding fabrication technology the plate usually consists of crystalline silicon, which is an anisotropic material. The isotropic approach is then invalidated and this results in deviations in the deflection compared to finite element modeling (FEM) and measurements. Therefore, to get precise modeling of these CMUTs the anisotropy of silicon needs to be taken into account.

For circular plates a simple and exact solution for the deflection exists, but this is not the case for square plates. Existing solutions for the deflection of square plates is based on series expansions with either trigonometric [3] or polynomial basis functions [4]. None of these, however, take the anisotropy of the plate into account.

Previously a model was made for calculating the deflection for an anisotropic plate with circular geometry [5], and in this paper the model is expanded to include square plates as well. The approach used to solve the full anisotropic plate equation is the Galerkin method [6]. Utilizing the symmetry of the silicon crystal, a compact and accurate approximation of the deflection can be obtained. The calculated deflection is compared to the solution for corresponding isotropic cases, a finite element model (FEM) and measurements performed on fabricated devices. Furthermore, the calculated deflection is used to find the stable position of the CMUT plate for a given bias voltage. Equivalent measurements are performed as well and the theory is compared to these.

II. THE ISOTROPIC PLATE EQUATION

Conventionally the deflection \( w(x,y) \) of a CMUT with a thin plate is modeled using the isotropic plate equation [3]

\[
\frac{\partial^4 w}{\partial x^4} + 2 \frac{\partial^2 w}{\partial x^2 \partial y^2} + \frac{\partial^4 w}{\partial y^4} = \frac{p}{D_i},
\]

where \( p \) is the applied pressure difference across the plate. The flexural rigidity is given by

\[
D_i = \frac{E}{12(1-\nu^2)}h^3
\]

with \( E \) being Young’s modulus, \( \nu \) being Poisson’s ratio, and \( h \) being the thickness of the plate. For clamped rectangular and square plates no simple exact solution exists to this equation and approximate methods have to be used. The traditional isotropic approach is based on a series expansion of the deflection and the center deflection for a thin clamped square plate having side length 2\( L \) is [3]

\[
w_{0,\text{isotropic}} = 0.020245 \frac{L^4 p}{D_i}.
\]

However, the plate material is often not isotropic and (1) and (2) are therefore no longer valid. Using the fusion bonding fabrication technique the plate usually consist of silicon which is an anisotropic material with a diamond cubic crystal structure. Having a silicon (001) substrate, which is most often used, Young’s modulus and Poisson’s ratio are strongly anisotropic, and this leads to inaccurate deflection expressions.

III. ANISOTROPIC PLATE EQUATION

To be able to take the anisotropy of the plate into account and avoid the inaccuracy from isotropic modeling, the stiffness of the plate needs to be described through the stiffness matrix of the material instead of Young’s modulus and Poisson’s ratio. The starting point is the relation between stress and strain [7]

\[
\sigma^c = c^c \varepsilon^c, \quad \varepsilon^c = s^c \sigma^c.
\]

Here superscript \( c \) denotes the crystallographic coordinate system, so \( c^c \) is the stiffness matrix and \( s^c = (c^c)^{-1} \) the compliance matrix in this coordinate system. Having a thin plate the stresses in the \( z \) direction can be ignored and plane
The plate coefficients for a uniform load can be used. This is a differential equation for the deflection, $w(x,y)$, of a thin anisotropic plate exposed to a uniform load $p$ given by [10], [9]

$$\frac{\partial^4 w}{\partial x^4} + k_1 \frac{\partial^4 w}{\partial x^2 \partial y^2} + k_2 \frac{\partial^4 w}{\partial y^4} + k_3 \frac{\partial^4 w}{\partial x^2 \partial y^3} + k_4 \frac{\partial^4 w}{\partial x \partial y^4} = \frac{p}{D_a}. \quad (8)$$

The plate coefficients $k_1$-$k_4$ and the anisotropic flexural rigidity, $D_a$, depend on the elastic constants of the plate material

$$k_1 = 4c_{eff}^{11} \frac{c_{11}^{11}}{c_{11}^{11}}, \quad k_2 = 2(c_{eff}^{11} + 2c_{eff}^{22}) \frac{c_{12}^{11}}{c_{11}^{11}}, \quad k_3 = 4c_{eff}^{22} \frac{c_{33}^{11}}{c_{11}^{11}}, \quad k_4 = c_{eff}^{11} \frac{c_{44}^{11}}{c_{11}^{11}}$$

where $c_{eff}^{pq}$ are elements in the effective stiffness matrix. Notice that the stiffness of the plate is no longer expressed through

stress assumed. The relation between strain and stress then becomes

$$\left( \begin{array}{c} \varepsilon_1 \\ \varepsilon_2 \\ \varepsilon_6 \end{array} \right) = \left( \begin{array}{ccc} s_{11} & s_{12} & s_{16} \\ s_{12} & s_{22} & s_{26} \\ s_{16} & s_{26} & s_{66} \end{array} \right) \left( \begin{array}{c} \sigma_1 \\ \sigma_2 \\ \sigma_6 \end{array} \right) = \mathbf{S}_\text{eff} \left( \begin{array}{c} \sigma_1 \\ \sigma_2 \\ \sigma_6 \end{array} \right) \quad (5)$$

and can be defined an effective stiffness matrix, $\mathbf{C}_\text{eff} = (\mathbf{S}_\text{eff})^{-1}$. For silicon the effective compliance matrix becomes

$$\mathbf{S}_\text{eff} = \left( \begin{array}{ccc} s_{11}^c & s_{12}^c & 0 \\ s_{12}^c & s_{22}^c & 0 \\ 0 & 0 & s_{44}^c \end{array} \right). \quad (6)$$

The compliance elements in this matrix are known from measurements and shown in Table I [8]. It is noted that the elements in (6) are known in the crystallographic coordinate system. An equation describing plate deflection on the other hand is valid in the plate coordinate system which is not necessarily the same. To illustrate this further the crystallographic and the plate coordinate systems can be seen in Fig. 1. The solid coordinate system aligned to the (100) direction is where the compliance values for silicon are known and the dashed system shows the rotated coordinate system for the plate where the compliance values needs to be calculated. Having silicon as plate material and performing standard cleanroom fabrication, the plate will usually be on a (001) substrate and aligned to the primary wafer flat. Flat alignment is to the (100) direction and the plate coordinate system will be rotated $\psi = 45^\circ$ and a transformation of the compliance matrix between the two coordinate systems is needed. The resulting effective stiffness matrix for the present case becomes (taking the inverse of the transformed compliance matrix) [9]

$$\mathbf{C}_\text{eff}^{\text{Si(001),(110)}} = \left( \begin{array}{cccc} \frac{1}{s_{11}^c} & \frac{1}{s_{12}^c} & 0 & 0 \\ \frac{1}{s_{12}^c} & \frac{1}{s_{22}^c} & 0 & 0 \\ 0 & 0 & \frac{1}{s_{44}^c} & 0 \\ 0 & 0 & 0 & \frac{1}{s_{11}^c} \end{array} \right) \quad (7)$$

It is seen that the stiffness matrix now has an orthotropic symmetry.

Now having the effective stiffness matrix the generalized plate equation can be used. This is a differential equation for the deflection, $w(x,y)$, of a thin anisotropic plate exposed to a uniform load $p$ given by [10], [9]

$$\frac{\partial^4 w}{\partial x^4} + k_1 \frac{\partial^4 w}{\partial x^2 \partial y^2} + k_2 \frac{\partial^4 w}{\partial y^4} + k_3 \frac{\partial^4 w}{\partial x^2 \partial y^3} + k_4 \frac{\partial^4 w}{\partial x \partial y^4} = \frac{p}{D_a}. \quad (8)$$

The plate coefficients $k_1$-$k_4$ and the anisotropic flexural rigidity, $D_a$, depend on the elastic constants of the plate material

$$k_1 = 4c_{eff}^{11} \frac{c_{11}^{11}}{c_{11}^{11}}, \quad k_2 = 2(c_{eff}^{11} + 2c_{eff}^{22}) \frac{c_{12}^{11}}{c_{11}^{11}}, \quad k_3 = 4c_{eff}^{22} \frac{c_{33}^{11}}{c_{11}^{11}}, \quad k_4 = c_{eff}^{11} \frac{c_{44}^{11}}{c_{11}^{11}}$$

where $c_{eff}^{pq}$ are elements in the effective stiffness matrix. Notice that the stiffness of the plate is no longer expressed through

Young’s modulus and Poisson’s ratio but directly through the stiffness values.

Using the compliance values for silicon in Table I and inserting the stiffness elements in (7) into (9) it follows that $k_1 = k_3 = 0$ and $k_4 = 1$. Thus, aligning the plate to the primary flat simplifies the anisotropic plate equation (8) to

$$\frac{\partial^4 w}{\partial x^4} + k_2 \frac{\partial^4 w}{\partial x^2 \partial y^2} + \frac{\partial^4 w}{\partial y^4} = \frac{p}{D_a}. \quad (10)$$

The same is the case for aligning the plate along the [100] direction where the inverse of (6) is used instead of (7), giving the same values for $k_1$, $k_3$ and $k_4$. For these two special cases the coefficients in the plate equation are summarized in Table II.

### IV. Solving the Plate Equation

Having a rectangular or square plate makes analytical deflection calculations complicated and approximate methods must be used to solve the generalized plate equation. With the anisotropic approach the Galerkin method [6] can be used to find approximate expressions for the deflection of a thin anisotropic square plate. In the most common case for CMUTs the plate is fabricated on a silicon (001) substrate and aligned to the (110) direction. For this orthotropic square plate with side lengths $2L$ the relative deflection is found to [9], [11]

$$\frac{w(x,y)}{w_0} = \left[ 1 - \left( \frac{x}{L} \right)^2 \right] \left[ 1 - \left( \frac{y}{L} \right)^2 \right] \times \left[ 1 + \beta \left( \frac{x}{L} \right)^2 + \beta \left( \frac{y}{L} \right)^2 \right], \quad (11)$$

where the plate parameter is defined as

$$\beta = \frac{182 + 143k_2}{1432 + 91k_2}. \quad (12)$$

---

**TABLE I.** Room temperature (300K) Compliance coefficients for low doped N-type crystalline silicon [8].

<table>
<thead>
<tr>
<th>$s_{11}$</th>
<th>$s_{12}$</th>
<th>$s_{16}$</th>
</tr>
</thead>
<tbody>
<tr>
<td>7.69 x 10^11 Pa^-1</td>
<td>-2.14 x 10^11 Pa^-1</td>
<td>12.58 x 10^11 Pa^-1</td>
</tr>
</tbody>
</table>

**TABLE II.** Selected values for the plate coefficients and anisotropic flexural rigidity for plates on a silicon (001) substrate [9].

<table>
<thead>
<tr>
<th>Orientation</th>
<th>$\psi$</th>
<th>$k_1$</th>
<th>$k_2$</th>
<th>$k_3$</th>
<th>$k_4$</th>
<th>$12D_a/h$ [GPa]</th>
</tr>
</thead>
<tbody>
<tr>
<td>[110]</td>
<td>$\pi/4$</td>
<td>0</td>
<td>1.3241</td>
<td>0</td>
<td>1</td>
<td>140.96</td>
</tr>
</tbody>
</table>

Fig. 1. The two coordinate systems, solid lines are the crystallographic system and the dashed lines the plate system aligned to the (110) direction.
The center deflection can be written

$$w_{0,\text{Si}(001)} = \frac{77(1432 + 91k_2)}{256(16220 + 11k_2(329 + 13k_2))} \frac{L^4 p}{D_a}. \quad (13)$$

Eqn. (11)-(13) are also valid when the plate is aligned to the [100] direction on a silicon (001) substrate. Note that the center deflection depends only on the $k_2$ coefficient. For primary flat alignment it is found by inserting $k_2$ into (12) that $\beta = 0.23920$. This results in a normalized deflection surface for the plate aligned to the (110) direction given by

$$\left. \frac{w(x,y)}{w_0} \right|_{\text{Si}(001),(110)} = \left[ 1 - (x/L)^2 \right]^2 \left[ 1 - (y/L)^2 \right]^2 \quad (14)$$

and the center deflection becomes

$$w_{0,\text{Si}(001),(110)} = 0.02196 \frac{L^4 p}{D_a}. \quad (15)$$

Comparing (3) and (15) it is seen that they are very similar containing the same parameters but different coefficients and the anisotropic instead of the isotropic flexural rigidity.

Fig. 2 shows the deflection cross section through $y = 0$ of a square plate of silicon (001)

$$w_{y=0} = w_0 \left[ 1 - (x/L)^2 \right]^2 \left[ 1 + \beta (x/L)^2 \right]. \quad (16)$$

The deflection calculated with the anisotropic approach uses $k_3 = 1.3241$ in (12) and center deflection (15). This is compared to the isotropic approach using $k_2 = 2$ in (12) and center deflection (3), with Young’s modulus and Poisson’s ratio in the [100] and [110] directions, and to a finite element (FEM) simulation made using the full anisotropic compliance matrix (compliance coefficients from Table 1) in COMSOL. The calculated deflections are normalized to the FEM center deflection. Excellent agreement is shown between the anisotropic curve and FEM with a deviation of less than 0.1 % whereas the isotropic approach leads to deviations in the center deflection of around 10 % for both [100] and [110] directions.

V. CMUT APPLICATION

Many important design parameters for CMUTs depend on the deflection of the plate. By using static analysis it is possible to find the stable position of the plate when applying a certain bias voltage. The stable position is easiest expressed when the strain force balance the electrostatic and pressure forces. The center deflection is found from energy considerations. The total potential energy of the system consists of three terms:

1) Strain energy. Calculated by integrating the strain energy density using (5), (7) and (11) and the result is

$$U_s = \frac{1}{2} \int_{-h/2}^{h/2} \int_{-L}^{L} \left[ \sigma_1 \varepsilon_1 + \sigma_2 \varepsilon_2 + \sigma_6 \varepsilon_6 \right] d\varepsilon_1 d\varepsilon_2 \quad (17)$$

$$U_{s,\text{Si}(001),(110)} = 3.91172 \times 10^{-11} h^2 w_0^2 \frac{L^4}{L}. \quad (18)$$

2) Energy due to applied pressure. This is found from the pressure load on the plate

$$U_p = -\int_{-L}^{L} \int_{-L}^{L} p w(x,y) \ dx \ dy \quad (19)$$

$$U_{p,\text{Si}(001),(110)} = -1.216 pw_0 L^2. \quad (20)$$

3) Electrostatic energy. Expressed through the charge $Q$ or applied voltage $V$, the vacuum permittivity $\varepsilon_0$, gap height $g$ and the total capacitance $C_t$ of the device which for a square plate is found using a Taylor expansion of the integrant with the deflection in (11)

$$U_e = \frac{Q^2}{2C_t} = \frac{1}{2} \int_{-L}^{L} \int_{-L}^{L} \varepsilon_0 \frac{V^2}{g - w(x,y)} d\varepsilon_1 d\varepsilon_2 \quad (21)$$

$$= \frac{1}{2} \int_{-L}^{L} \int_{-L}^{L} \frac{V^2}{g - w(x,y)} d\varepsilon_1 d\varepsilon_2 \quad (22)$$

The total force on the system is then found by differentiating the total potential energy with respect to the center deflection. From this the stable center position of the plate can be found for a given applied voltage as the point where the total force is zero.

VI. COMPARISON TO MEASUREMENTS

CMUTs with square silicon plates have been fabricated using fusion bonding. The fabricated devices have a 65x65 $\mu$m wide and 2.37 $\mu$m thick silicon plate with a gap height of 405 nm and a 198 nm thick oxidizing oxide at the bottom of the cavity. The deflection was measured with a Sensofar PLu Neox 3D Optical Profiler using white light interferometry.

Fig. 3 shows a measured cross section of the normalized deflection for the fabricated device. It is normalized in both center deflection and distance across the plate to compare the shape of the measured deflection with the calculated deflection. The red curve is a fit made to the measurements using the anisotropic model (16). Both the center deflection and the plate parameter $\beta$ is fitted. As it is seen in the figure the fitted value for $\beta$ is 0.243 which matches very well, with a deviation of 1.5%, compared to the calculated value of 0.23920 for this type of plate (silicon (001) substrate aligned to [110] direction). The center deflection found from the fit has a deviation of 0.5% compared to the measurement.
Using isotropic plate theory to calculate the deflection of anisotropic silicon plates results in deviations from FEM or measurements of up to 10%. The full anisotropic plate equation was solved using the Galerkin method. It is seen that the deflection simplifies by utilizing the symmetry of the silicon crystal and a compact solution is obtained for square CMUT plates on a (001) silicon substrate aligned to the [110] direction. The maximum deviation is less than 0.1% compared to FEM. Furthermore, the deflection was measured on fabricated devices and fitting the anisotropic calculated deflection to the measurement a deviation of 0.5-1.5% is observed in the fitted parameters. The stable position for varying bias voltage was also found using the anisotropic theory and comparing this to measurements it is seen that the theory is within the uncertainty interval of the measurements.

REFERENCES


