

# Multi-stability of the hexagonal origami hypar based on group theory and symmetry breaking

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## A B S T R A C T

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The origami hyperbolic paraboloid, or the hypar, is widely known for its characteristic non-zero Gaussian curvature and multi-stable states. Previous investigations have mainly considered two particular cases of the origami hypar patterns, namely the square and the circular hypars. As a representative example of general polygonal hypar patterns, the hexagonal origami hypar displays more desirable energy properties and subtler multi-stable configurations, which is however, rarely studied. In this paper, we investigate the multi-stability of the hexagonal origami hypar by combining a group-theoretic approach, a symmetry-breaking method, and a bar-and-hinge structural model, to simplify the kinematic analysis of this highly symmetric structure. Notably, the kinematic path of the hexagonal origami hypar is divided into three bifurcation branches by symmetry breaking. Each branch corresponds to two symmetric stable states according to the equilibrium loading and potential energy simulated using the bar-and-hinge model. The non-rigid deformation of the hexagonal origami hypar is mainly controlled by the folding of creases and the bending of facets. Moreover, the energy barrier among the stable states becomes increasingly stronger with higher symmetry orders, thicker sheets of material, and longer creases. This work provides a new strategy for analyzing multi-stable origami structures with high symmetry orders, which can be used in the design and development of novel adaptive or deployable engineering structures.

## 1. Introduction

Multi-stable systems possess non-trivial energy landscapes that lead to multiple distinct stable configurations. As a result of various desirable characteristics of such systems, in recent years, they have been increasingly adopted in bionics [1,2], metamaterials [3,4], robotics [5–8], engineering structures [9–11,85] and other fields. From a structural design viewpoint, many multi-stable systems are designed by combining bistable units [12]. In general, the local properties and method of combination of these units have a critical impact on the global properties of a multi-stable system. For example, a well-assembled double-layer Kresling tube can achieve the same total height in two different stable states [13]. The rich mechanics of origami structures provide many candidates for such bistable units and have been a main source of inspiration for the design of multi-stable structures, such as the

Kresling configurations with different stable-state numbers [14], the Kresling assembly [15,16], and kirigami-based metastructures [17].

A number of multi-stable origami structures are derived from traditional origami patterns, such as the Miura-ori [18–31], Tachi-Miura polyhedron [32,33], Waterbomb tubes [34,35], and Kresling tubes [36,37]. By combining multi-stable origami units such as the square twist [38,39], leaf-like origami [40], origami-inspired cellular [41,42], and origami hypar [43], we can produce higher numbers of stable states and richer levels of programmability in designing origami structures. To ensure smooth deployment and transition processes, kinematic analysis is a key consideration in the design and analysis of multi-stable origami structures [44]. The existing methods for the kinematic analysis of rigid or non-rigid origami can be divided into two types: (1) mechanism analysis, and (2) structural analysis. In general, the conditions of rigid folding are more complex than those of non-rigid folding; therefore, the

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kinematic analysis of rigid folding is simpler and more compatible with a wide range of computational models [45]. Kinematics of rigid origami can be simulated by means of mechanism analysis methods such as polygonal surfaces [46] and spherical polygonal linkages [47], as well as by structural analysis methods such as the bar-and-hinge [48] and shell [49] models. However, non-rigid folding involves the deformations of origami facets, which greatly increase the variables to be considered [50]. Thus, in general, for non-rigid origami, mechanism analysis methods are not applicable to the study of non-rigid origami structures. The most common analytical methods for analyzing non-rigid origami structures are as follows: (1) the bar-and-hinge model with elastic bars for the tension and compression of facets and rotating springs for the bending of them [51,52]; (2) the membrane-and-hinge model with elastic membranes for the tension and compression of facets and rotating springs for the bending of them [53]; and (3) the thin plate elements model [54].

The origami hypar, as a non-rigid and multi-stable origami unit, is composed of concentrically pleated polygons with a minimum of four sides [55]. Among this family of patterns, two extreme cases have been widely studied, namely, the square origami hypar [56] and the circular origami hypar (note that a circle can be considered as a polygon with an infinite number of sides) [57,58]. The square origami hypar has two stable states that are macroscopically symmetric with a rotation of 90° [43]. When transitioning from one stable state to the other, it first unfolds to a flat state and then refolds up, going through a significant energy barrier caused by the competing folding, bending, and stretching energies [51]. Moreover, the folding angle of the square origami hypar can be regulated by controlling the defects of creases [59] to make the structure reach the appropriate stable state. On the other hand, the circular origami hypar can reach different stable states under certain approximate conditions [60]; in the simplest condition, it forms a saddle-like shape. Nevertheless, except for these two extreme cases, the multi-stability of the polygonal origami hypar is still not studied in detail. Therefore, in this paper, we will present a thorough investigation of the multi-stable behavior of the hexagonal origami hypar.

The identification of all stable states is essential for investigating the multi-stability of the hexagonal origami hypar, as it has been experimentally observed to possess more than two stable states. Since the tangent stiffness matrix may become ill-conditioned, simulating the global properties of the hexagonal origami hypar by the same method as used for solving the problem of bistability may be infeasible. Therefore, the simulation of the global behavior needs to be simplified as the kinematic bifurcation analysis of deployable structures [61]. Bifurcations can be determined by the singular value decomposition [62]. A wide range of numerical algorithms has been applied to the geometric, kinematic, or mechanical analysis of origami structures, including the geometric-graph-theoretic method [63], particle swarm optimization method [64], and mixed-integer linear programming method [65]. Another part of studies has simplified multi-stability analysis of systems by bifurcation branches, such as assuming the bifurcation parameters [66,67] and the group-theoretic approach [68,69]. The latter is particularly effective in solving bifurcation branches. More specifically, for structures with high orders of symmetry, it can be simplified according to group theory [70,71]. Based on these theories, Chen et al. [72,73] developed a group-theoretic approach to studying the kinematic behavior of deployable bar structures with cyclic symmetry. However, the applications of group theory and symmetry in the field of origami structures have been largely focused on designing innovative crease patterns (see, e.g. [74–76]), whereas there has been little attention to their folding kinematics and stability problems.

Here, using group theory and its mathematical representation capabilities, we propose a new strategy to study the multi-stability of the hexagonal origami hypar. The group-theoretic approach is introduced to simplify the whole folding process of origami structures. The bifurcation branches of origami structures with high orders of symmetry as the hexagonal origami hypar can be solved by symmetry breaking. The

corresponding stable states can be obtained by kinematic analysis with the bar-and-hinge models of different bifurcation branches. The bifurcation branches and the corresponding stable states will manifest multi-stability during the entire deployment process. Moreover, our work will study the effect of material properties on the multi-stability of this structure. This article is organized as following: Section 2 introduces the group-theoretic method for multi-stability analysis; Section 3 demonstrates the multi-stability of the hexagonal origami hypar in different bifurcation branches; Section 4 discusses possible stable states and the effect of crease parameters; finally, Section 5 provides the main concluding remarks of this research.

## 2. Methodology

We utilize the equivalent bar-and-hinge model to analyze the deformation of the hexagonal origami hypar. Using group theory, the analysis process is simplified to studying bifurcation branches which can be obtained by symmetry breaking.

### 2.1. Bar-and-hinge model of non-rigid origami

The bar-and-hinge model is a simplified model in the finite element analysis of origami [77,78]. On this basis, the total potential energy  $\Pi$  can be expressed as

$$\Pi = U - V \quad (1)$$

where  $V$  is the external work and  $U$  is the strain energy. According to the principle of the minimum total potential energy, the finite element equation can be expressed as

$$\mathbf{F} = \mathbf{K}\Delta \quad (2)$$

where  $\mathbf{F}$ ,  $\mathbf{K}$ , and  $\Delta$  donate the force vector, tangent stiffness matrix, and generalized displacement vector, respectively. The internal force vector and the tangent stiffness matrix can be assembled from equivalent nodal loads and elementary stiffness matrices.

In comparison with rigid origami, the potential energy  $U$  of non-rigid origami contains not only folding energy, but also facet deformation energy [50].

$$U = U_F + U_S + U_B \quad (3)$$

where  $U_F$ ,  $U_S$ , and  $U_B$  represent the crease folding energy, bar stretching energy, and facet bending or twisting energy, respectively. Then, the tangent stiffness matrix can be calculated by finding displacement variations of Eq. (3) twice [50].

$$\mathbf{K} = \mathbf{K}_F + \mathbf{K}_S + \mathbf{K}_B = \mathbf{K}_{spr} + \mathbf{K}_{bar} \quad (4)$$

where the stiffness of rotational spring elements  $\mathbf{K}_{spr}$  only affects the folding stiffness of the crease  $\mathbf{K}_F$ . The stiffness of bar elements  $\mathbf{K}_{bar}$  contains in-plane stretching and shearing stiffness of the bar  $\mathbf{K}_S$  and the out-plane bending stiffness of the facet  $\mathbf{K}_B$ . Thus, the elementary stiffness of bar elements and rotational spring elements should be calculated. The hyperplastic model is selected to simulate the constitutive relationship of bar elements [77,78]. The elementary stiffness matrices are obtained by

$$\begin{cases} \mathbf{K}_{spr}^e &= \mathbf{K}_F^e = \partial^2 U_F^e / \partial \mathbf{u}^2 = k_{spr}^e d\theta / d\mathbf{u} \otimes d\theta / d\mathbf{u} + M_F d^2\theta / d\mathbf{u}^2 \\ \mathbf{K}_B^e &= \partial^2 U_B^e / \partial \mathbf{u}^2 = k_B^e d\theta / d\mathbf{u} \otimes d\theta / d\mathbf{u} + M_B d^2\theta / d\mathbf{u}^2 \\ \mathbf{K}_S^e &= \partial^2 U_S^e / \partial \mathbf{u}^2 = k_S^e (\mathbf{B}_1^T + \mathbf{B}_2 \mathbf{u}) (\mathbf{B}_1^T + \mathbf{B}_2 \mathbf{u})^T + f_s \mathbf{B}_2 \end{cases} \quad (5)$$

During the simulation of the kinematic path, the tangent stiffness matrix near the bifurcations will be severely ill-conditioned. However, it can be effectively avoided by utilizing group theory [79,80].

## 2.2. A group-theoretic approach to the global bifurcation analysis

Central to performing the global bifurcation analysis is to solve the vector equilibrium equation

$$\mathbf{f}(\mathbf{u}, \lambda) = \mathbf{0} \quad (6)$$

where  $\mathbf{u}$  belongs to an  $N$ -dimensional real vector space and  $\lambda$  is a real number. Then, using the symmetry-adapted coordinate system, the Newton-Rapson procedure can be implemented by the block diagonalization method as follows

$$\begin{bmatrix} \mathbf{K}_{v(1)} & 0 & \cdots & 0 \\ 0 & \mathbf{K}_{v(2)} & \cdots & 0 \\ \vdots & \vdots & \cdots & \vdots \\ 0 & 0 & \cdots & \mathbf{K}_{v(p)} \end{bmatrix} \begin{bmatrix} \Delta \mathbf{u}^{(1)} \\ \Delta \mathbf{u}^{(2)} \\ \vdots \\ \Delta \mathbf{u}^{(p)} \end{bmatrix} = \begin{bmatrix} -\mathbf{f}_\lambda \\ 0 \\ \vdots \\ 0 \end{bmatrix} \quad (7)$$

where  $\mathbf{K}$  represents the tangent stiffness matrix  $\mathbf{K} = d\mathbf{f}/d\mathbf{u}$ . Decomposed by many block matrices, the original problem can be reduced to a series of independent problems with specific symmetry subspaces.

$$\mathbf{K}_{v(i)} \Delta \mathbf{u}^{(i)} = -\mathbf{f}_\lambda \quad (8)$$

In other words, the solution of Eq. (6) can be obtained by solving Eq. (8) associated with different subspaces. In particular, for a symmetric structure, the symmetry subgroups are associated with different symmetry subspaces. Therefore, the bifurcation branches of the original structure will consist of the bifurcation branches of the corresponding symmetry subgroups [81].

## 2.3. Symmetry breaking behavior

The transformation from the primary path to bifurcation paths of deployable structures is often accompanied by symmetry breaking behavior [72,73]. In other words, different bifurcation paths can be obtained by reducing the symmetry order of structures.

The  $C_{nv}$  symmetry group contains a total of  $2n$  independent symmetry operations [82], and is represented as

$$C_{nv} = \{C_n^j, \sigma_j | j = 1, 2, \dots, n\} \quad (9)$$

where rotation  $C_n^j$  represents the rotation of the original structure by an angle  $2\pi j/n$  around the symmetry axis,  $j \in [1, n]$ , and  $\sigma_j$  denotes the  $j$ -th reflection along the vertical symmetry plane. Among these symmetry operations, the one-dimensional and two-dimensional irreducible representations are selected to describe the partial symmetry of the original structure. The irreducible representations are expressed as

$$n = \text{even} \begin{cases} \Gamma^{(1)} = A_1, \Gamma^{(2)} = A_2, \Gamma^{(3)} = B_1, \Gamma^{(4)} = B_2 & \text{One - dimensional} \\ \Gamma^{(5)} = E_1, \dots, \Gamma^{(3+0.5n)} = E_{0.5n-1} & \text{Two - dimensional} \end{cases} \quad (10)$$

$$n = \text{odd} \begin{cases} \Gamma^{(1)} = A_1, \Gamma^{(2)} = A_2 & \text{One - dimensional} \\ \Gamma^{(3)} = E_1, \dots, \Gamma^{(1.5+0.5n)} = E_{0.5n-0.5} & \text{Two - dimensional} \end{cases} \quad (11)$$

where  $\Gamma^{(i)}$  denotes an irreducible representation of the structure, and  $A_1, \dots, B_1, \dots, E_1, \dots$  represent the basic symmetry representations. The number of one-dimensional irreducible representations of  $C_{nv}$  is  $2\alpha + 2$ , whereas that of two-dimensional irreducible representations is  $(n-1-\alpha)/2$  [83]. The value of  $\alpha$  equals 0 (or 1) when  $n$  is odd (or even). The first and last items indicate the highest and lowest orders of symmetry respectively. Then, the subgroups can be obtained from these

irreducible representations.

$$C_1, \dots, C_i, \dots, C_m, C_n, C_v, \dots, C_{iv}, \dots, C_{mv} \subset C_{nv} \quad (12)$$

where  $m$  represents the maximum proper divisor of  $n$ , and  $i$  includes all the proper divisors of  $n$ . Accordingly, the affiliation and symmetry order of these subgroups of  $C_{nv}$  can be determined. For instance,  $C_1$  is the lowest-order symmetry group.

A kinematically indeterminate structure is mobile if the symmetry order of the internal mechanism mode is higher than that of the self-stress state [71]. In other words, the symmetry order of the self-stress state has to be reduced to ensure foldability. If the symmetry order of the mechanism mode fails to satisfy the full symmetry, the structure will transform into a bifurcation branch with lower-order symmetry. The bifurcation branch matches the symmetry order of the associated constraint modes. That is, we can effectively program the symmetry order of the structure by controlling the constraint mode.

## 3. Multi-stability of the hexagonal origami hyper

Fig. 1 shows the crease pattern of a typical hexagonal origami hyper. The blue dashed lines and red solid lines represent the valley and mountain creases, respectively, while the black solid dots represent the  $z$ -direction constraints. The geometric and material parameters of the initial pattern are defined according to the square origami hyper in [51]. The fold spacing  $d$  and inner fold spacing  $D$  are  $d = D = 8$  mm. The number of the crease layers is  $n = 8$ , the thickness  $t$  is  $127 \mu\text{m}$ , the elastic modulus  $E$  for the material is 5 GPa, and the Poisson's ratio is 0.35. In contrast to the square origami hyper, there are more possibilities for the 3D geometry of the hexagonal origami hyper. In this section, we will investigate the multi-stability of the hexagonal origami hyper by studying its stable states with different symmetry orders.

The initial planar pattern of the hexagonal origami hyper belongs to symmetry group  $C_{6v}$ . To allow the structure to be foldable, the symmetry order of the self-stress state is required to be decreased to a value less than that of the initial pattern [71,72]. The subgroups of  $C_{6v}$  group can be computed by

$$C_1, C_2, C_3, C_4, C_5, C_6, C_v, C_{2v}, C_{3v} \subset C_{6v} \quad (13)$$

According to established theorems for assigning mountain and valley creases [84], crease assignments in the initial patterns must be symmetric with respect to the diagonals. In other words, the initial pattern cannot belong to a symmetry subgroup containing only cyclic rotations, i.e.  $C_6, C_3, C_2$ , and  $C_1$ . Therefore, we only need to study the cases with mirror symmetries, namely  $C_{3v}, C_{2v}$ , and  $C_v$ .

According to Section 2, a specific symmetry-breaking of the structure

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is achieved by adjusting the symmetry order of the constraint mode. Fig. 1(a) shows the three-dimensional view of the pattern, and Figs. 1(b), (c), and (d) depict different constraint modes leading to origami configurations with different symmetry groups. The mountain and valley assignments of the innermost creases depend on the symmetry group. Figs. 1(b), (c), and (d) corresponds to the  $C_{3v}, C_{2v}$ , and  $C_v$  symmetries, respectively.

### 3.1. Hexagonal origami hyper with $C_{3v}$ symmetry

Among the subgroups, the highest-symmetry mechanism mode is the  $C_{3v}$  symmetry. The initial 3D configuration can be obtained by

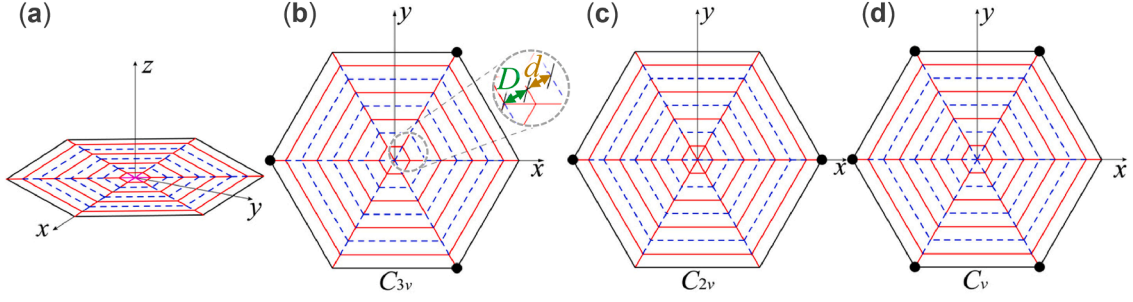


Fig. 1. Crease patterns of the hexagonal origami hyar: (a) Three-dimensional view of the crease pattern. (b–d) Different constraint modes and crease patterns corresponding to symmetry groups  $C_{3v}$ ,  $C_{2v}$ , and  $C_v$ .

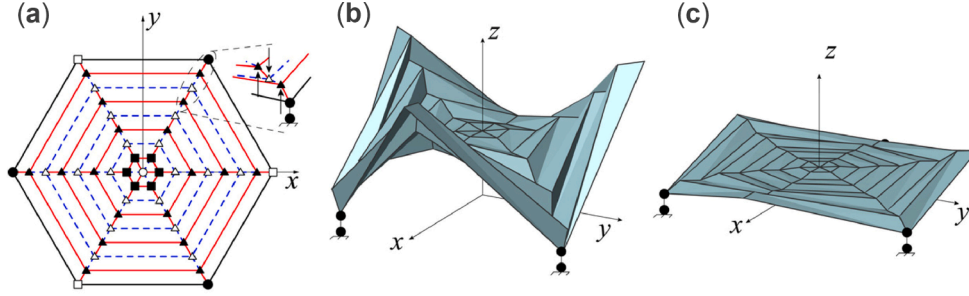


Fig. 2. 3D origami configuration with  $C_{3v}$  symmetry. (a) Applied forces in the  $z$ -direction (the hollow circle  $\circ$  represents downward force  $F_0$ , the solid square  $\blacksquare$  represents upward force  $2.25F_0$ , the hollow triangle  $\triangle$  represents downward force  $4F_0$ , the solid triangle  $\blacktriangle$  represents upward force  $4F_0$ , the hollow square  $\square$  represents upward force  $2F_0$ , and the solid circle  $\bullet$  represents constraints in the  $z$ -direction) (b) Initial 3D configuration. (c) 3D configuration in a stable state.

loading the equivalent bar-and-hinge model in the  $z$ -direction. The applied forces to fold the initial pattern are adjusted based on the folding of the square origami hyar [51]. The loading in the  $z$ -direction used in this study is shown in Fig. 2(a), which provides us with the initial 3D configuration (see Fig. 2(b), obtained using MERLIN2\_v1.1 based on the constraints and forces depicted in Fig. 2(a)). However, it is a random state rather than a stable state. When the external loads are relieved, the system will recover into one of its stable states, as shown in Fig. 2(c).

In order to find other stable states with  $C_{3v}$  symmetry, we further analyze the stable 3D configurations by using the bar-and-hinge model. In Fig. 3(a1), the three lower outermost vertices are constrained, whereas downward displacements are applied to the remaining three vertices. As displacements increase, the structure first unfolds from the first stable state (State 1) to an approximately fully flat state (State 2). Then, it refolds along the creases to a new equilibrium (State 3). Importantly, the total potential energy reaches another minimum at State 3, as shown in Fig. 3(e). According to the energy principle, the new equilibrium is the second stable state of the hexagonal origami hyar with  $C_{3v}$  symmetry.

The stored energy-displacement relation clearly manifests the bistability of the structure. In Fig. 3(e), the total potential energy  $E_T$  contains the tensile and shear energy  $E_S$ , the bending energy  $E_B$ , and the folding energy  $E_F$ . During the transition from the stable State 1 to the stable State 3,  $E_F$  increases with the decrease of the folding angle and accounts for the largest proportion of  $E_T$ . This property indicates that this structure folds preferentially at the creases. The facet deformation energy  $E_B + E_S$  accounts for a certain proportion of the total potential energy, especially when the structure is close to the stable states. Thus, in terms of the folding energy, the facet deformation energy cannot be ignored. In other words, the hexagonal origami hyar is a non-rigid origami structure the facet deformations of which need to be considered during the folding process.

Concerning the specific composition of the deformation energy,  $E_S$  increases while  $E_B$  decreases with the flattening of the facets.  $E_B$  accounts for the most of the facet deformation energy around the stable states

whereas  $E_S$  accounts for the most of the facet deformation energy near the fully-expanded state. That is, the deformation of the facets is dominated by bending near the stable states, and by tension, compression, and shear near the fully-expanded state. Moreover,  $E_B$  changes substantially during the entire process while  $E_S$  only changes considerably near the fully flattened state. This implies that the bending deformation of the facets exists in the entire process while the tensile deformation of the facets only occurs near the fully flattened state. In addition, the proportion of  $E_S$  in  $E_T$  is still small even near the fully expanded state, whereas the proportion of  $E_B$  in  $E_T$  is large near the stable states. Therefore, it can be concluded that the facets deformation of the hexagonal origami hyar is mainly controlled by the bending deformation.

The equilibrium load-displacement curves further verify these two stable states. In Fig. 3(d), these two stable states correspond exactly to the two zero points that cross the reference line  $F = 0$  in the positive direction. State P1 and State P2 both reach the peak value of the equilibrium load during folding and unfolding. Moreover, according to Fig. 3(d), they are roughly symmetric about the center point corresponding to State 2. Similarly, the two stable states (i.e., State 1 and State 3) are symmetric about State 2 as the center point. Fig. 3(c) shows the changing profile of the  $y = 0$  section during folding. The two stable states (i.e., State 1 and State 3) and the two peak force states (i.e., State P1 and State P2) are mirror symmetric along  $x = 0$ . The intermediate State 2 is mirror symmetric about the  $x$ -axis, but not fully flattened. Moreover, except for the innermost folds, the mountain and valley properties of the creases remain constant during the transition between the two stable states.

### 3.2. Hexagonal origami hyar with $C_{2v}$ symmetry

The first stable state with  $C_{2v}$  symmetry is obtained by the same method described at the beginning of Section 3.1. The main difference is that this case requires the origami pattern and constraints shown in Fig. 1(c). The unconstrained outermost vertices are all load points in this



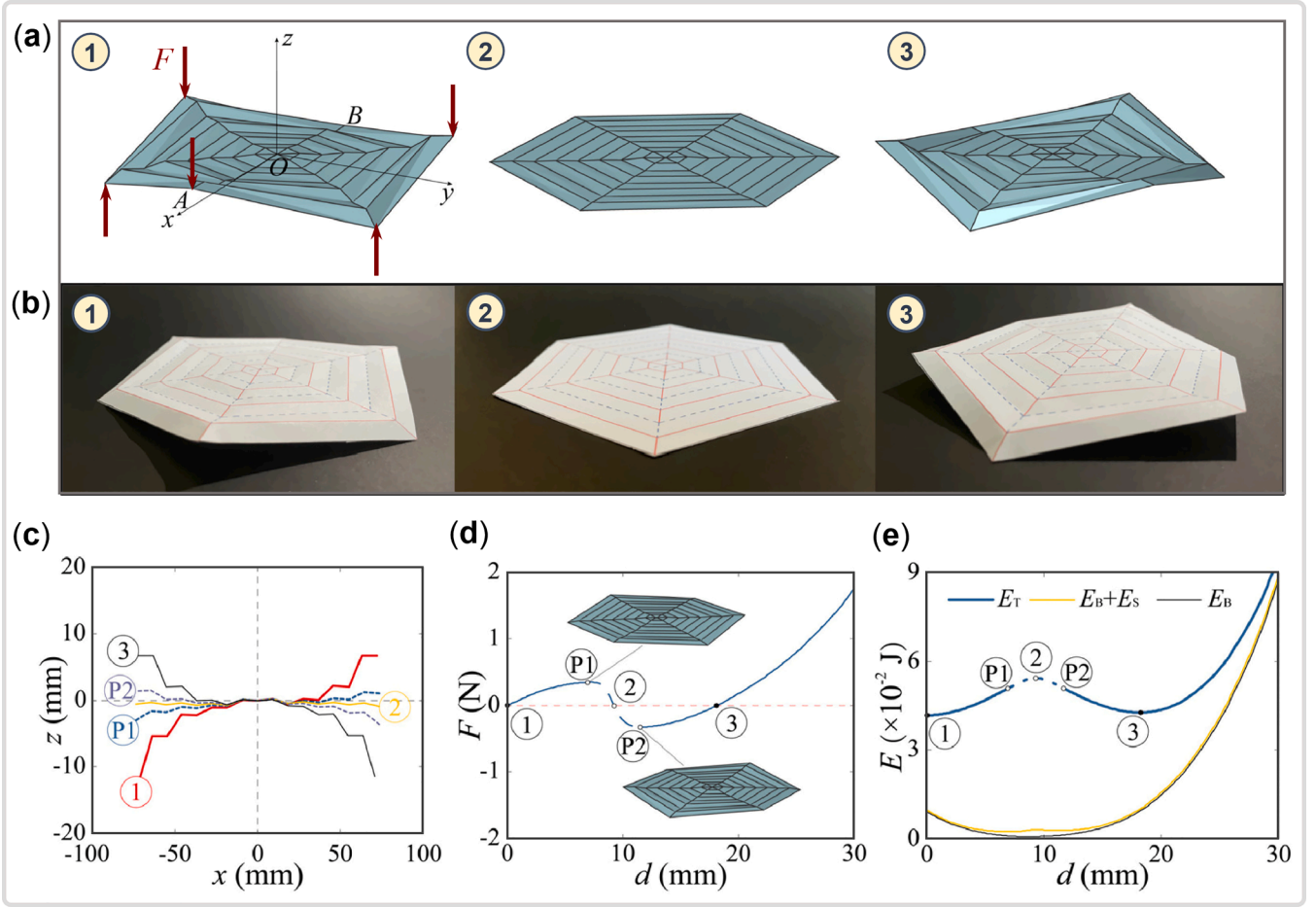


Fig. 3. Multi-stability of the hexagonal origami hyper with  $C_{3v}$  symmetry. (a) Three key states of the shape transition process. (b) A physical models in the three key states. (c) Changing profile of the set of diagonal creases on the  $x$ -axis (i.e., diagonal AOB). (d) Equilibrium load versus displacement curves of single load point. (e) Stored energy versus displacement curves.

case because the constraint conditions have been altered. The load size and direction of the new load points are the same as the load size and direction represented by the hollow square  $\square$  in Fig. 2(a). The loading conditions for the remaining nodes are the same as those shown in Fig. 2(a).

Here, we analyze the stable 3D configurations with  $C_{2v}$  symmetry by a similar method as in Section 3.1. The two outermost vertices on an adjacent diagonal are constrained, and upward loads are applied to the remaining four outermost vertices on two adjacent diagonals. Then, two lower outermost vertices are constrained, whereas the downward displacements are applied to the remaining four outermost vertices, as shown in Fig. 4(a1). As displacements increase, the structure first unfolds from the initial stable State 1 to approximately fully flattened State 2. Then, it refolds along the creases to a new equilibrium State 3. The total potential energy at State 3 reaches another minimum, as shown in Fig. 4(e). According to the energy principle, the new equilibrium is the second stable state of the hexagonal origami hyper with  $C_{2v}$  symmetry.

Similar to the hexagonal origami hyper with  $C_{3v}$  symmetry, in Fig. 4(e), the bistability of the hexagonal origami hyper with  $C_{2v}$  symmetry can be directly seen from the stored energy versus displacement curves. As can be seen from Fig. 4(e), the hexagonal origami hyper with  $C_{2v}$  symmetry is non-rigid, and the deformations of its facets are dominantly controlled by the bending deformation. Moreover, the variation tendencies of equilibrium load and potential energy during the transition between the two stable states illustrated in Fig. 4(d) and 4(e) are similar to those shown in Fig. 3(d) and 3(e), respectively. Fig. 4(c) reflects the same mountain and valley properties of the creases and symmetry

between corresponding states during the entire process. In contrast to the hexagonal origami hyper with  $C_{3v}$  symmetry, the key states during the entire transition of this structure are all mirror symmetric about  $x = 0$ .

### 3.3. Hexagonal origami hyper with $C_v$ symmetry

The 3D configuration with  $C_v$  symmetry can be achieved by the further reduction of the symmetry order. The first stable state with  $C_v$  symmetry is obtained by the same method as explained in Section 3.1. The origami pattern and constraints of this structure are shown in Fig. 1(c). The only outermost free vertex is subjected to an upward load of  $3F_0$  in the  $z$ -direction. The loading conditions for the remaining nodes are the same as those shown in Fig. 2(a).

For this origami hyper with  $C_v$  symmetry, the five outermost vertices are constrained, and upward loads are applied to the remaining outermost vertex. Then, the five lower outermost vertices are constrained and an downward displacement is applied to the remaining vertex, as shown in Fig. 5(a1). As the displacement increases, the structure first unfolds from the first stable State 1 to an approximately fully flattened State 2. Then, it refolds along the creases to a new stable State 3. The total potential energy reaches another minimum at State 3, as shown in Fig. 5(e). According to the energy principle, the new equilibrium is the second stable state of the hexagonal origami hyper with  $C_v$  symmetry.

Similar to the two cases studies earlier in this paper, the bistability of the hexagonal origami hyper with  $C_v$  symmetry can be demonstrated by Fig. 5(e). It is a non-rigid origami structure with facet deformations

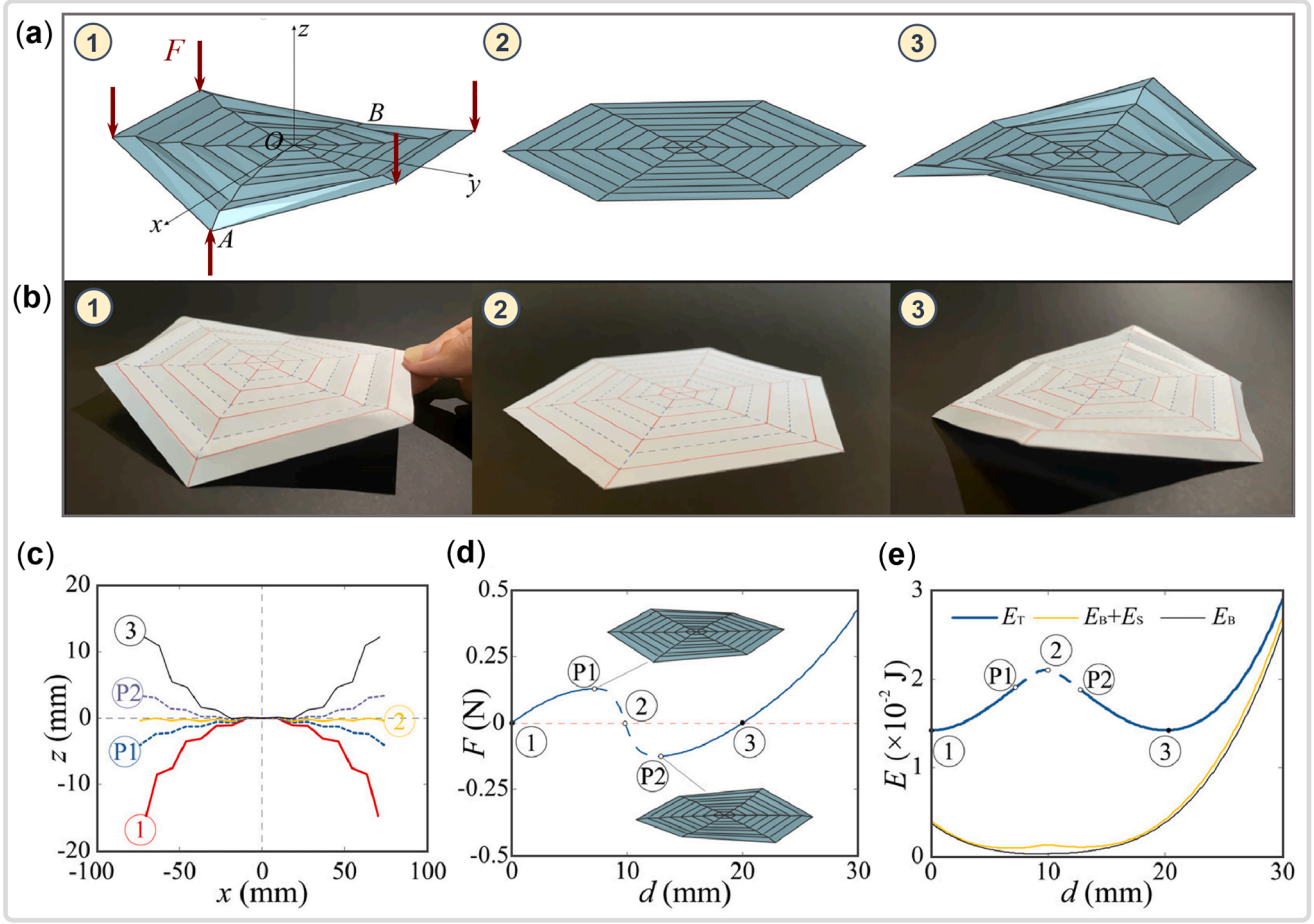


Fig. 4. Multi-stability of the hexagonal origami hyper with  $C_{2v}$  symmetry. (a) Three key states of the shape transition process. (b) A physical model in the three key states. (c) Changing profile of the set of diagonal creases on the  $x$ -axis (i.e., diagonal AOB). (d) Equilibrium load versus displacement curves of a single load point. (e) Stored energy versus displacement curves.

which are mainly regulated by bending. Moreover, Fig. 5(d) and Fig. 5(e) indicate similar variation tendencies of the equilibrium load and potential energy as those of the other bifurcation paths. Fig. 5(c) depicts the same mountain and valley properties of the creases during the entire process as in Fig. 3(c) and Fig. 4(c). Compared to the hexagonal origami hyper with  $C_{3v}$  and  $C_{2v}$  symmetries, the key states of this structure show no symmetry in the  $y = 0$  section.

In summary, the bistability of the hexagonal origami hyper in each bifurcation branch is the same as the square origami hyper [43]. There are two stable states along each bifurcation branch. The two stable states in the same bifurcation branch are approximately symmetric along the horizontal plane. Except for the innermost folds, the mountain and valley properties of the creases remain unchanged between these two stable states.

#### 4. Discussion

In this section, we discuss the differences among these three bifurcation branches. Several metrics are chosen to represent the energy barrier of multistable transitions in Section 4.1, and some relevant influencing factors are studied in Section 4.2.

##### 4.1. Multi-stability of the hexagonal origami hyper with different symmetries

To compare and discuss different bifurcation branches, we transform them into a unified coordinate system, as shown in Fig. 6. Points  $I_a$ ,  $P1_a$ ,

$P2_a$ ,  $III_a$ , points  $I_b$ ,  $P1_b$ ,  $P2_b$ ,  $III_b$ , and points  $I_c$ ,  $P1_c$ ,  $P2_c$ ,  $III_c$  correspond to the key States 1, P1, P2, and 3 on the bifurcation branches with  $C_v$ ,  $C_{2v}$  and  $C_{3v}$  symmetries described in Section 3, respectively. To ensure the fairness of the comparison among the different numbers of loading points for different bifurcation branches, the total equilibrium load is chosen as the  $y$ -axis. In Section 3, we obtained the equilibrium load curves at a single loading point for each bifurcation branch. Therefore, the total equilibrium load should be the sum of the equilibrium loads at all loading points in each branch. Then, based on the symmetry of each bifurcation branch, the equilibrium loads of all the loading points will be identical. That is, we can obtain the total equilibrium load by multiplying the equilibrium load at a single loading point by the number of loading points. For the bifurcation branch with  $C_{3v}$  symmetry, we applied downward displacements on the three outmost free vertices. Thus, the total equilibrium load bifurcation branch with  $C_{3v}$  symmetry will be equal to the equilibrium load at a single loading point multiplied by three. For the bifurcation branch with  $C_{2v}$  symmetry, there are four loading points. Therefore, its total equilibrium load will be equal to the equilibrium load at a single loading point multiplied by four. For the bifurcation branch with  $C_v$  symmetry, we only applied downward displacement on one vertex. As a result, the total equilibrium load of the bifurcation branch with  $C_v$  symmetry will be equal to the equilibrium load at a single loading point itself.

According to Fig. 6(a), seven possible critical points can be found in total, including point X, Y, Z, X', Y', Z', and II. Fig. 6(c) shows the profiles of the  $y = 0$  section of the structure in each bifurcation branch at State 2. Since the displacements are calculated by finite element

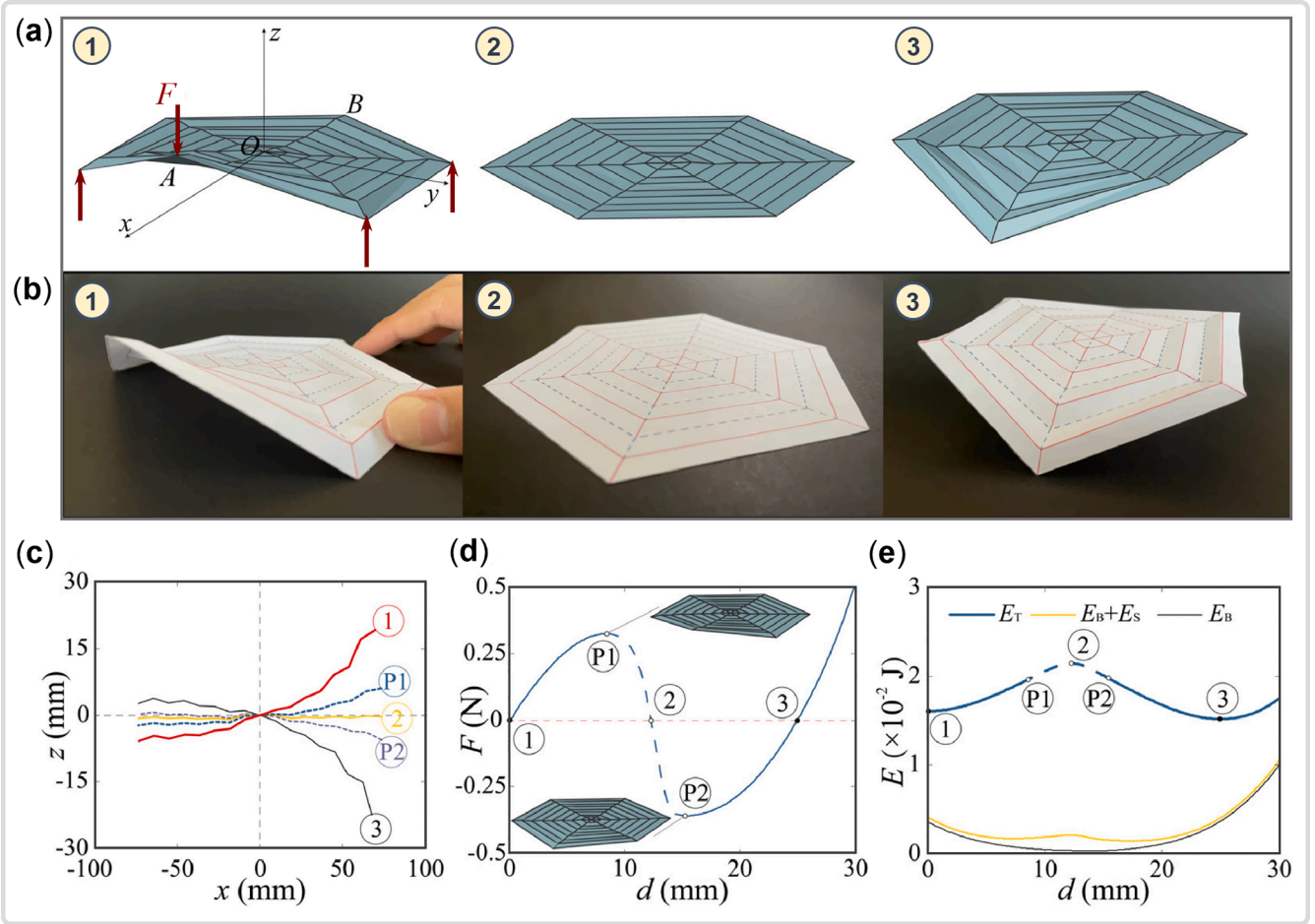


Fig. 5. Multi-stability of the hexagonal origami hypar with  $C_v$  symmetry. (a) Three key states of the shape transition process. (b) A physical model in the three key states. (c) Changing profile of the set of diagonal creases on the  $x$ -axis (i.e., diagonal AOB). (d) Equilibrium load versus displacement curves of a single load point. (e) Stored energy versus displacement curves.

analysis, the State 2 shown in Fig. 6(c) is an approximate solution. However, it can be judged from the folding tendency that the hexagonal origami hypar is fully expanded in State 2 as point II in Fig. 6(d). Importantly, point II is one of the critical points of these three bifurcation branches. As for the other possible critical points, the states corresponding to various bifurcation branches at the intersections are different. In Fig. 6(d), these three bifurcation branches have no intersection except point II. In other words, point II is the only critical point of these three bifurcation branches.

We select the state corresponding to the critical point II as the zero potential energy state. Fig. 6(e) shows the total energy ( $E_T$ ) variation of the hexagonal origami hypar from the stable State 1 to the stable State 3. Critical point II represents the fully flattened state in the three bifurcation branches. The energy barrier of conversion among these stable states can be reflected by the difference between the extreme values of the stored energy curves. In comparison with the stable states with  $C_{2v}$  and  $C_v$  symmetries, unfolding from the stable state with  $C_{3v}$  symmetry to the fully unfolded state is the most difficult transformation to achieve. Unfolding from the stable states with  $C_{2v}$  symmetry is relatively easier, while this is the easiest for the stable states with  $C_v$  symmetry. In other words, the stable states with  $C_{3v}$  symmetry have the greatest energy barrier against transformations to the other stable states, followed by the stable states with  $C_{2v}$  symmetry; the stable states with  $C_v$  symmetry have the slightest energy barrier.

## 4.2. Influence of geometric parameters

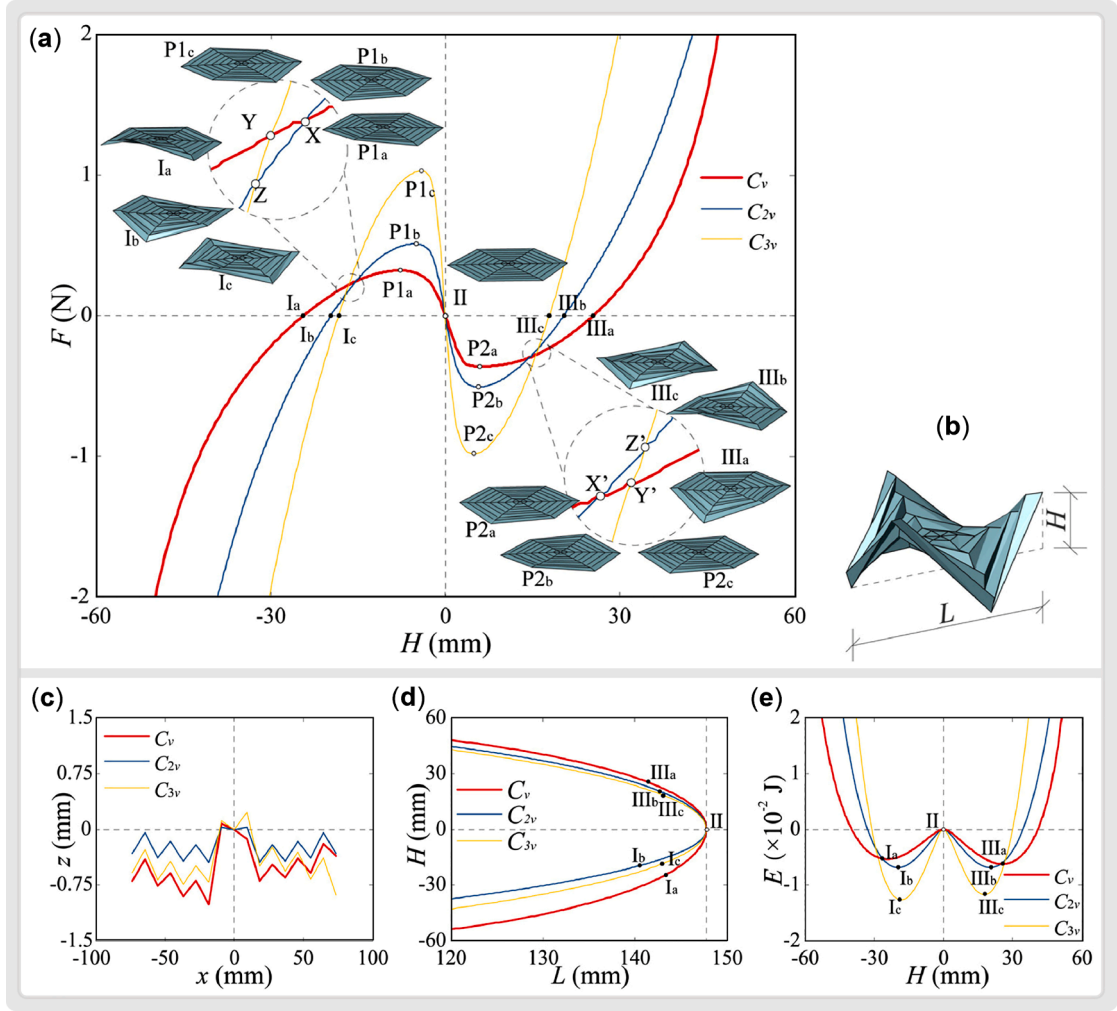
To obtain more general conclusions, here we examine and discuss the effects of different geometric parameters on the stable states of the hexagonal origami hypar, including the thickness of material  $t$  and the fold spacing  $d$ . The geometric parameters of different numerical examples are summarized in Table 1.

### 4.2.1. Influence of material thickness

Various geometric parameters associated with each model are listed in Table 1. We utilize the same method, as used for the reference model T0, to simulate the multi-stability of models T1 to T4. Fig. 7(a), (c), and (e) show the results corresponding to  $C_{3v}$ ,  $C_{2v}$ , and  $C_v$  symmetries, respectively. Point T0<sub>I</sub> represents the stable State 1 in model T0, and the point T0<sub>III</sub> represents the stable State 3 in model T0.

Fig. 7(a), (c), and (e) show that the peak values of the equilibrium load increase with the material thickness during the transitions of two stable states. The absolute value of  $F$  indicates its structural height, and the sign (positive or negative) of  $F$  indicates whether it increases along the positive  $z$ -axis. If  $F$  increases along the positive  $z$ -axis, it is negative; otherwise,  $F$  is positive. To make the results clearer, we have extracted the peaks of the balanced load and marked them in Fig. 7(b), (d), and (f). The equilibrium load versus height curves during the transition between two stable states consists of two peaks, and the tendencies of them are centrosymmetric along the origin point. The entire process can be divided into two phases with the origin as the cut-off point. The first phase, called the loading phase, is the transition from stable State 1 to a





**Fig. 6.** Comparison among various bifurcation branches. (a) Total equilibrium load-displacement curves of the three bifurcation branches. (b) Physical representation of span ( $L$ ) and height ( $H$ ). (c) Profile of the  $y = 0$  section in State 2. (d) Span versus height curves. (e) Total energy versus height curves.

**Table 1**

Geometric parameters of numerical examples.

Model	T0	T1	T2	T3	T4	F1	F2	F3	F4	F5
$L$ (mm)	64	64	64	64	64	60	60	60	60	60
$D$ (mm)	8	8	8	8	8	12	12	12	12	12
$d$ (mm)	8	8	8	8	8	4	6	8	12	16
$t$ ( $\mu\text{m}$ )	127	40	76	200	240	127	127	127	127	127

fully flattened state under external loads. Similarly, the second phase, called the recovery phase, is the transition from the fully flattened state to the other stable State 3. The peak equilibrium load in the loading phase is  $F_L$ , and the peak equilibrium load in the recovery phase is  $F_R$ . Fig. 7(b), (d), (f) show that  $F_L$  and  $F_R$  are proportional to the third power of the material thickness (i.e.,  $t^3$ ). The fitting plots  $f_L$  and  $f_R$  in Fig. 7(b), (d), and (f) are obtained by a linear fitting with  $t^3$  as the independent variable in MATLAB.

In the simulation process, the stiffness of the creases and the bending stiffness of the facets are related to  $t^3$ , whereas the tension-compression stiffness of the facets is related to  $t$  [77]. Similar to the results presented in Section 3, we can conclude that the folding of the structure is mainly realized by the folding of creases and the bending of facets, and the tension and compression deformations of facets only account for a small part.

As for the height of the structure, we refer to the height at which the

structure reaches a stable state as the stable height  $H$ . Fig. 7(a), (c), and (e) show that the stable height increases with material thickness. Similar to the peak equilibrium load, each equilibrium load-height curve has two stable heights corresponding to the two stable states. The stable height corresponding to stable State 1 is called  $H_I$  and the stable height corresponding to stable State 3 is called  $H_{III}$ . The absolute value of  $H$  is its height and the plus or minus sign indicates whether  $H$  increases along the positive  $z$ -axis. If  $H$  increases along the positive  $z$ -axis, it is positive. Otherwise,  $H$  is negative. Thus, for the equilibrium load versus height curve, the farther the stable state from  $H = 0$ , the higher the stable height will be.

In summary, the peak equilibrium loads and the stable heights with the same symmetry, all increase with material thickness. In particular, the magnitude of the peak equilibrium load is positively related to the third power of the material thickness.



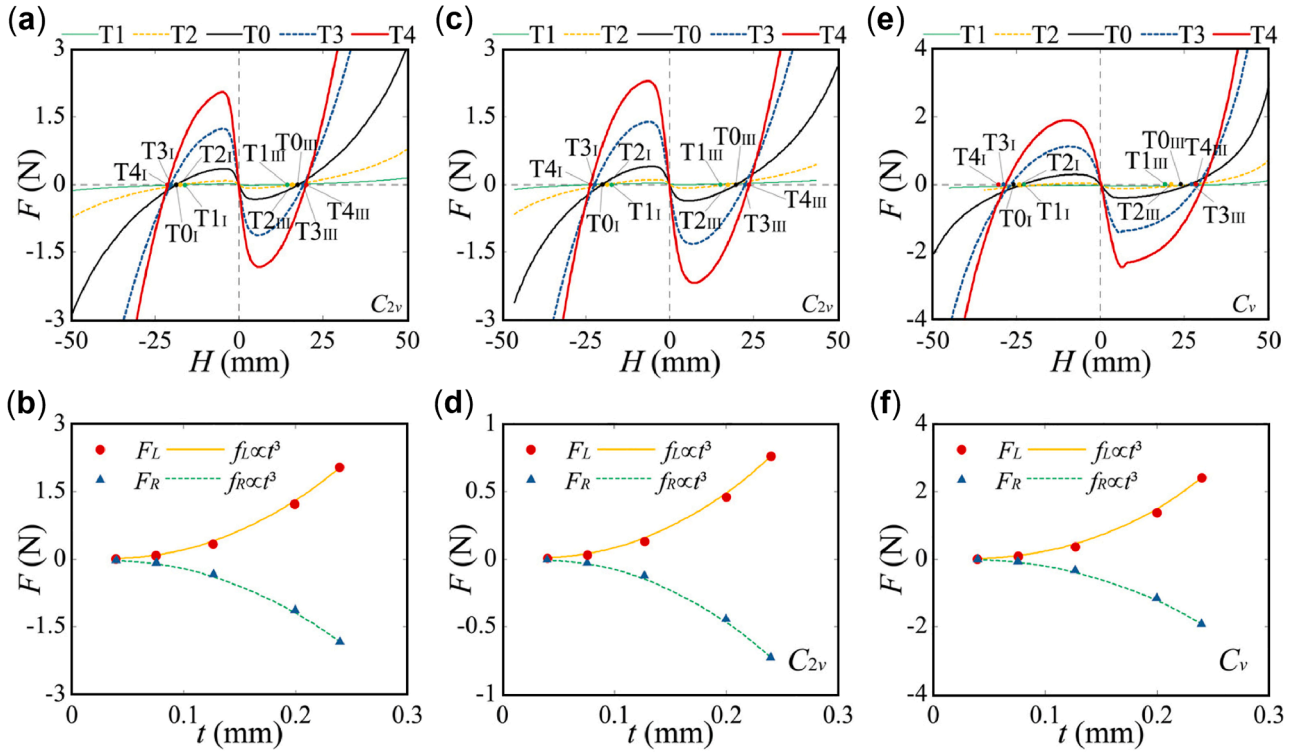


Fig. 7. Effects of material thickness. (a), (c), and (e): Equilibrium load versus height curves of a single load point. (b), (d), and (f): Relationship between the material thickness  $t$  and the peak equilibrium load  $F$ .

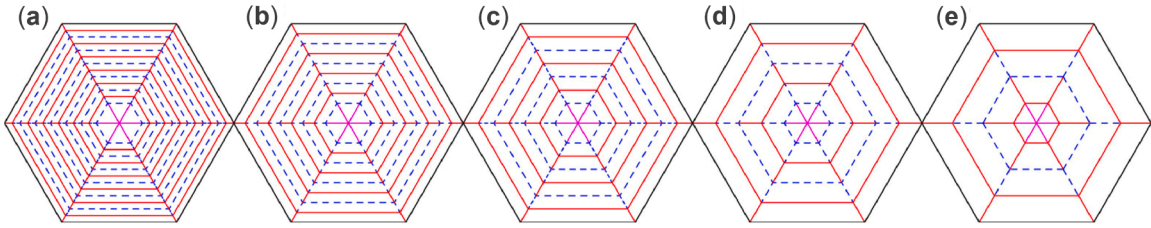


Fig. 8. Crease patterns with different values for parameter  $d$ . Sub-figures (a) to (e) correspond to models F1 to F5, respectively.

#### 4.2.2. Influence of crease length

We adjust the geometric parameters of the reference model to the variation of  $d$  while keeping  $L$  constant. Parameter  $n$  varies with  $d$  as models F1 to F5 (see Table 1). The outermost creases are kept at peak and various patterns are generated as shown in Fig. 8. Then, the bistability of models F1 to F5 can be simulated. Fig. 9(a) to (c) show the results corresponding to  $C_{3v}$ ,  $C_{2v}$ , and  $C_v$  symmetries, respectively. Point  $F1_{P1}$  represents State P1 in model F1, and point  $F1_{P2}$  represents State P2 in model F1.

In Fig. 9, the peak equilibrium loads increase with creases spacing  $d$  during the transitions of two stable states with the same symmetry. Since  $L$  does not change, the change of  $d$  will directly lead to the change of  $n$ . For simplicity, we unify these two variables as the total length of all creases, called crease length (denoted by  $D$ ). The relationships between the crease length and the peak equilibrium loads are shown in Fig. 9(d). For the three different symmetry cases, the peak equilibrium loads in both the loading and recovery phases increase with crease length. In other words, hexagonal origami hypars with longer crease lengths are more difficult to be folded. As for the stable height, it is related to the crease spacing, crease layers number and crease length. The details of their influence need to be further studied.

In summary, the magnitudes of peak balanced loads increase with increasing the crease length. In other words, the stable states of

hexagonal origami hypars with longer crease lengths have more significant energy barriers. The effects of different spans  $L$  and the number of initial polygon sides on the stable states of the hexagonal origami hypar can be further studied. The obtained data can be analyzed using deep learning [85] methods to develop a more precise numerical model, and could facilitate the design of novel programmable metamaterials [86, 87] based on the hexagonal origami hypar.

## 5. Conclusions

In this work, we utilized a bar-and-hinge model, a group-theoretic approach, and a symmetry breaking process for exploring the multi-stability of the hexagonal origami hypar. The bar-and-hinge model was applied to simulating the entire folding and unfolding process of the structure. Concurrently, the group-theoretic approach was exploited to avoid singularities in the stiffness matrix near the bifurcations. The symmetry subgroups obtained by symmetry breaking were introduced to trace the corresponding bifurcation branches. The multi-stability of the structure was demonstrated by analyzing the multi-stability of all the bifurcation branches. The main conclusions of this study can be summarized as follows:

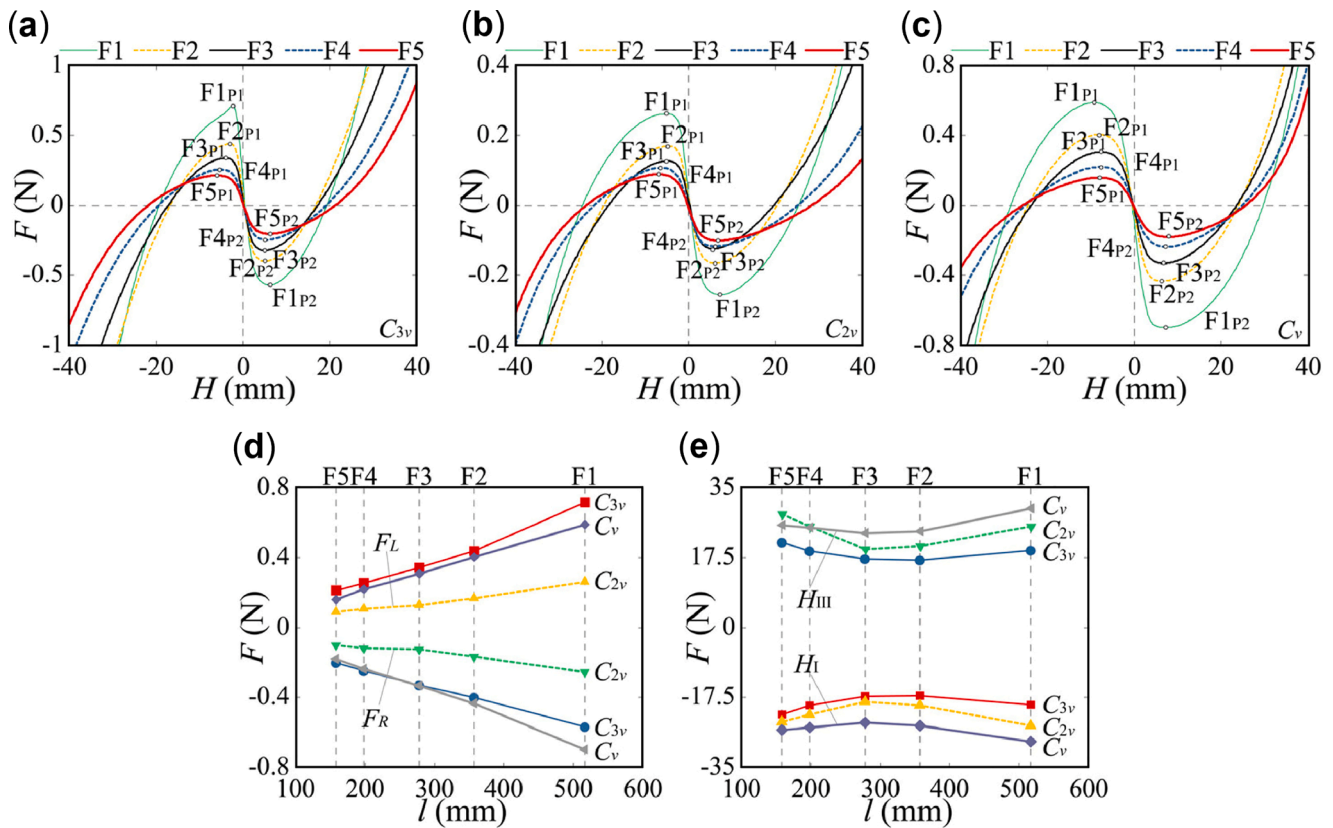


Fig. 9. Effect of crease spacing. (a) to (c) Equilibrium load versus height curves of a single load point. (d) Relationship between the creases length  $l$  and the peak equilibrium loads  $F_L$  and  $F_R$ . (e) Relationship between the creases length  $l$  and the stable heights  $H_I$  and  $H_{III}$ .

- 1) The hexagonal origami hypar is always non-rigid during folding. Its non-rigid deformation is mainly realized by the bending deformation of each facet.
- 2) The exagonal origami hypar has three distinct types of bifurcation paths, including  $C_{3v}$ ,  $C_{2v}$ , and  $C_v$  symmetries. All the bifurcation branches exhibit bistability. Thus, the hexagonal origami hypar has six independent stable states.
- 3) These stable states can be converted to each other only through the fully flattened state. The initial state with the greatest energy barrier is the stable states with  $C_{3v}$  symmetry, and the slightest one is the stable states with  $C_v$  symmetry. The energy barrier increases with increasing the material thickness or the crease length.

Furthermore, the method proposed in this study can be extended for the multi-stability analysis of origami structures with higher-order symmetries. It can systematically describe the transitions among their stable states by converting the original folding analysis into a collection of multi-stability investigations in different bifurcation branches.

#### CRediT authorship contribution statement

**Yao Chen:** Methodology, Resources, Formal analysis, Validation, Investigation, Writing – original draft, Writing – review & editing, Supervision, Project administration, Funding acquisition. **Ruizhi Xu:** Conceptualization, Investigation, Data curation, Formal analysis, Validation, Methodology, Software, Writing – original draft. **Chenhao Lu:** Formal analysis, Software, Writing – review & editing. **Ke Liu:** Supervision, Methodology, Validation, Writing – review & editing. **Jian Feng:** Supervision, Validation, Writing – review & editing. **Pooya Sareh:** Supervision, Validation, Writing – review & editing, Funding acquisition.

#### Declaration of Competing Interest

The authors declare that they have no known competing financial interests or personal relationships that could have appeared to influence the work reported in this paper.

#### Data availability

Data will be made available on request.

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