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Layout optimization for multi-bi-modulus materials system under multiple load cases

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Abstract An optimization model is presented for obtaining optimal layout of multiple bi-modulus materials systems under multiple load cases (MLC). In the optimization model, the objective function is the linearly weighted structural compliance under MLC. The bi-modulus materials in a finite element are replaced by isotropic materials according to the stress state of that element. The equivalent mechanical properties of an element are expressed as the power–law function of the volume fractions (design variables) and moduli of the solid phases. Numerical experiments are presented to verify the validity and efficiency of the present algorithm. The effects of factors including the bi-modulus behavior of materials, the load directions and the weighting schemes of MLC are also investigated numerically.

Keywords Topology optimization · Bi-modulus materials · Multi-material structures · Multiple load cases · SIMP

1 Introduction

During recent decades, topological optimization has gained considerable attention in both theoretical research and

Qing-Hua Qin Qinghua.qin@anu.edu.au practical applications [1–3]. Due to their complexity, topology optimization problems with large numbers of design variables are still the most challenge task in the structural optimization field. Several optimization schemes have been reported, including the homogenization-based method [4], the SIMP method (solid isotropic material with penalization) [5, 6], the ESO (evolutionary structural optimization) method [7, 8], and the level set method [9, 10]. Topology optimization has been used widely in the design of engineering structures such as MEMS [11], acoustics [12], crashworthiness [13], fluidics [14], bone structures [15], and heat conduction [16].

In practical engineering, a structure is often under multiple loads (MLs). To apply topology optimization on such structures, Díaz and Bendsøe [17] developed an optimization model with a single objective function using a linear weighting scheme for MLs. Similarly, Bendsøe et al. [18] addressed the design of material properties and material distribution in structures under MLC. Luo et al. [19] presented a hybrid fuzzy-goal multi-objective programming scheme for topology optimization that considered both static and dynamic loadings. Sui et al. [20] suggested an independent continuous mapping method for solving the topology optimization problems of a continuum under MLs. Balamurugan et al. [21] used a genetic algorithm for the topology design of structures under MLs.

Most of the work above focused on a continuum with single material only. Topological design with multi-phase materials is more complex than traditional 0–1 design of structures with only one solid phase. For solving topology optimization problems with multi-phase materials, the level set method was developed [22–25]. Other methods including the SIMP method [26–28], the ESO method [29], phase-field method [30], and pseudo-sensitivities scheme [31] have also been proposed.

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Few of the works mentioned above have considered optimization of the layout of bi-modulus materials. Nevertheless, bi-modulus materials that have different tensile and compressive moduli along the same direction are very common in engineering. For example, materials such as rubber, concrete, cast iron, graphite, foam materials, masonry, bone, alloys, and ropes/membranes exhibit bi-modulus behavior. Due to the stress dependency of bi-modulus materials, deformation analysis of bi-modulus structures is more complex than that of structures with isotropic materials [32]. Achtziger [33] considered the bi-modulus effect on the final topology of a truss and found that the final structure under tension was obviously different from that under compression. Chang et al. [34] approximated the original piecewise linear stress-strain curve of a bi-modulus material with a derivable nonlinear curve for the topological design of a tension-only or compression-only material. Cai [35] solved the tension-only or compression-only design using a modified SIMP method, in which the tension-only or compression-only material is replaced with an isotropic material. Querin et al. [36] used orthotropic materials to replace the original bi-modulus material according to the local stress state in topology optimization of truss-like structures. Cai et al. [37] suggested a sampler scheme for finding the optimal topology of a continuum structure with one bi-modulus material.

In the present work, topology optimization of multiphase bi-modulus materials under MLs is studied. Numerical examples are presented showing the applicability and efficiency of the proposed algorithm.

2 Methodology

2.1 Statement of linear elasticity problem

For a linear elastic structure, the basic equations read

$$-\nabla \cdot \sigma(u) = f \text{ in } \Omega,$$

$$\varepsilon(u) = \frac{1}{2} [\nabla u + (\nabla u)^{\mathrm{T}}],$$

$$\sigma(u) = D : \varepsilon(u),$$
(1)

and the corresponding boundary conditions are:

$$u = u_0 \qquad \text{on } \Gamma_1, \sigma(u) \cdot n = T \qquad \text{on } \Gamma_2.$$
(2)

where Γ_1 and Γ_2 are Dirichlet condition and Neumann condition, respectively, and $\Gamma_1 \cup \Gamma_2 = \partial \Omega \ \Omega \subset R^2 \text{ or } R^3$ is the design domain. ε and σ are the strain and stress tensors, respectively. *f* is the body force vector, *u* is the displacement field, u_0 is the prescribed displacement on Γ_1 and *T* is boundary force on Γ_2 . *D* is the elasticity tensor.

2.2 SIMP approach for optimization of layout of multiple materials

Topology optimization based on the density-like method of the SIMP approach [5] is adopted. In the density-like method, the equivalent modulus of a composite material is calculated using the material interpolation scheme of the moduli and volume fractions of the components in the composite. For manufacturing, the final design should have only one component material in one element, and interfaces between component materials should be on common boundaries between elements.

For a [0, 1] design of only one component material using the SIMP method, the interpolation for material modulus in an element is defined as:

$$\mathbf{E}^{1}(\rho_{1,j}^{p}) = \rho_{1,j}^{p} \cdot \mathbf{E}_{1} + (1 - \rho_{1,j}^{p}) \cdot \mathbf{E}_{\text{void}} = \rho_{1,j}^{p} \cdot \mathbf{E}_{1}$$
(3)

where $E^{(1)}$ means the equivalent modulus of a composite material with one solid material and void. $\rho_{1,j} \in [0, 1]$ are the volume fraction, a design variable of the *j*th element j = 1, 2, ..., n and *n* is the total number of elements in design variable, E_1 is the elastic modulus of the solid phase, and *p* is the penalization parameter, typically p = 3

when the composite has *m* types of solid with the moduli of E_1, E_2, \ldots, E_m ($E_1 > E_2, \ldots, > E_m$). E^m can be given based on the interpolation scheme:

$$E^{(m)} \rho_{1,j}, \rho_{2,j}, \dots \rho_{m-1,j} = \rho_{m-1,j}^{p}, E^{(m-1)} + 1 - \rho_{m-1,j}^{p} \cdot E_{m}.$$
(4)

In this equation, $\rho_{i,j}$ is the summation of the volume fractions of the first *i* types of solid in the *j*th composite element, and $\rho_{m,j} = 1.0$. This interpolation scheme is used in the present study.

2.3 Optimization model

For the stiffness design of a continuum with multiple bimodulus materials under MLC, the optimization model reads

Find
$$\{\rho_{i,j} | i \in \{1, 2, ..., m\}, j \in \{1, 2, ..., n\}\}$$

min $c = \sum_{l=1}^{N_{LC}} w_l \cdot \overline{c_l},$
s.t. $\sum_{j=1}^{n} \rho_{r,j} \cdot v_j = V_0 \cdot \sum_{i=1}^{r} f_i,$
 $\sum_{l=1}^{N_{LC}} w_l = 1.0,$
 $\overline{K}_l \cdot \overline{U}_l = F_l, \quad (l = 1, 2, ..., N_{LC}),$
 $0 < \rho_{\min} \le \rho_{1,j} \le \rho_{2,j} \le \cdots \le \rho_{m,j} = 1,$
 $0 < w_l, \quad (l = 1, 2, ..., N_{LC}).$
(5)

where the design variable $\rho_{i,j}$ is the volume fraction of the *i*th component material in the *j*th element. The objective function, c, is the linear weighted mean compliance of structure under the multiple load cases $(\overline{c_1}, l = 1, 2, ..., N_{LC} N_{LC}$ is the number of loading cases. F_l and \overline{U}_l denote the global nodal force and displacement vector in the *l*th load case , respectively. \overline{K}_l is the final global stiffness matrix of a structure with many bi-modulus materials, and can be calculated using the well-established finite element formulation [38, 39], "n" and "m" are, respectively, the total number of finite elements and the number of solid materials in the design domain. v_i is the *i*th element volume, f_i is the *i*th material volume fraction, and V_0 denotes the volume of the design domain. ρ_{\min} is the minimum value of relative densities. To avoid singularity of the stiffness matrix K_l , here we set $\rho_{\min} = 0.001$.

2.4 Material replacement scheme for bi-modulus material

Figure 1a shows the stress–strain curve of a bi-modulus material with tensile modulus of $E^{T} = \tan \alpha$ (*Green* line) and compressive modulus of $E^{c} = \tan \beta$ (*Orange* line). The stress–strain curve is piecewise linear if $\alpha \neq \beta$. $\alpha = \beta$ means that the material is degenerated into an isotropic material. If $\beta = 0$, the material is a tensile-only material (Fig. 1b). If $\alpha = 0$, the material is a compressive-only material (Fig. 1c). σ_{T} and σ_{c} are the allowable stresses of the material under tension and compression, respectively. σ_{T} and σ_{C} are usually different.

To represent the difference between tension and compression performance, the ratio between E^{T} and E^{C} is defined as:

$$R_{\rm TCE} = \frac{E^{\rm T}}{E^{\rm C}}.$$
 (6)

In the deformation analysis of a structure with bi-modulus materials, many structural reanalysis could be required due to the nonlinearity of materials. However, the nonlinearity here has some particular features. For example, the bi-modulus material appears isotropic when it is under pure compression or under pure tension. When the material is in a complex stress state, e.g., the first principal stress is positive but the third is negative, the eigen pairs of the elasticity tensor depend on the direction and value of the second principal stress. Hence, in a structure with an optimal loadtransmission path (LTP), most of the structure is under a simple stress state. At the junction of adjacent parts, however, the material may be transverse isotropic because of its complex stress state. This condition implies that we can use an appropriate isotropic material to replace the bi-modulus material during structural deformation analysis. The difference in structural stiffness caused by such replacement can be modified step by step. The merit of material replacement is that the structural deformation analysis becomes a linear analysis after replacement. The modification of the local stiffness difference can be performed during the topology optimization process. Two aspects must, therefore, be given in detail for the material replacement scheme. The first is selection of the isotropic material that will replace the bimodulus material in structural deformation analysis. The second is modification of the local stiffness due to material replacement.

2.4.1 Selection of isotropic material for local replacement

For a given element in a design domain, σ_s and ε_s (s = 1, 2, 3) denote the principal stresses and principal strains.

- 1. If $0 \ge \sigma_1 \ge \sigma_2 \ge \sigma_3$, the compressive modulus of the bi-modulus material should be the same as that of the isotropic material;
- 2. If $\sigma_1 \ge \sigma_2 \ge \sigma_3 \ge 0$, the tensile modulus of the bimodulus material should be the same as that of the isotropic material;
- 3. If the element is under a complex stress state, i.e., $\sigma_1 \cdot \sigma_3 < 0$, the elastic modulus depends on the comparison between the values of the tension strain energy density (SED) and compression SED, which considers the influence of the second principal stress.

Mathematically, the modulus of the isotropic material to replace the *i*th bi-modulus material can be obtained from the equation:

$$E_{i,j} = \begin{cases} E_{i,j}^{\mathrm{T}}, & \text{if SED}_{\mathrm{T}} > \mathrm{SED}_{\mathrm{C}}, \\ E_{i,j}^{\mathrm{C}}, & \text{if SED}_{\mathrm{T}} < \mathrm{SED}_{\mathrm{C}}, \\ \max(E_{i,j}^{\mathrm{T}}, E_{i,j}^{\mathrm{C}}), & \text{others.} \end{cases}$$
(7)

where the tension SED (SED_T) and compression SED (SED_C) are determined by the equations:

$$\text{SED}_{\text{T}} = \sum_{l=1}^{N_{LC}} w_l \cdot \left(\sum_{G=1}^{N_G} \sum_{s=1}^3 \left(\frac{1}{4N_G} (|\sigma_s| + \sigma_s) \cdot \varepsilon_s \right)_G \right), \quad (8)$$

$$\operatorname{SED}_{\mathrm{C}} = \sum_{u=1}^{t} w_{u} \cdot \left(\sum_{G=1}^{N_{G}} \sum_{s=1}^{3} \left(\frac{-1}{4N_{G}} (|\sigma_{s}| - \sigma_{s}) \cdot \varepsilon_{s} \right)_{G} \right).$$
(9)

where N_G is the number of Gaussian integral points in the element. From Eq. (8), one can find that a negative principal stress (i.e., $\sigma_s < 0$) contributes zero to the value of SED_T. Only positive principal stress can result in positive value of SED_T. On the contrary, Eq. (9) indicates that only negative principal stress can produce positive value of SED_C. Hence,

the third formulation in Eq. (7) indicates that the stronger stress will determine the modulus of the new isotropic material which is to be used to replace the original bi-modulus material. For example, if $SED_C > SED_T$, the compressive principal (negative) stress is stronger than the tensile principal (positive) stress, and E^C should be used.

2.4.2 Modification of local stiffness

Accurate deformation of a structure depends on an accurate global stiffness matrix, which is formed using the local (element) stiffness matrix. When a difference appears after material replacement, the local stiffness matrix of the element with the isotropic material is different from that of the same element with the original bi-modulus material. To eliminate the difference, under the same stress state, the same element should have the same strain energy density before and after replacement. According to this principle, we can calculate the modification factor by comparing the SEDs before and after replacement.

Under a complex stress state at the *k*th iteration, the SED of the element with the new isotropic material is

$$\operatorname{SED}_{j,k} = \sum_{l=1}^{N_{LC}} \left(w_l \cdot \sum_{G=1}^{N_G} \sum_{s=1}^3 \left(\frac{1}{2N_G} \sigma_s \cdot \varepsilon_s \right)_G \right), \tag{10}$$

which is the summation of SED_{T} and SED_{C} of the *j*th element at the *k*th iteration. And the effective SED of the element with the original bi-modulus materials is

$$\begin{split} & \text{SED}_{j,k}^{\text{effective}} \\ &= \sum_{i=1}^{m} \left\{ \gamma_i \cdot \sum_{l=1}^{N_{LC}} \left(w_l \cdot \sum_{G=1}^{N_G} \sum_{s=1}^{3} \left(\frac{1}{2N_G} \text{sign}_i(\sigma_s) \cdot \sigma_s \cdot \varepsilon_s \right)_G \right) \right\}_{j,k} \end{split}$$
(11)

where $\gamma_i = \rho_i - \rho_{i-1}$ is the volume fraction of the *i*th material in the element.

The value of the $sign_i(\cdot)$ can be calculated using either Eq. (11) or Eq. (12).

(a) If the element has *compressive moduli* at the (k-1)th iteration and *tensile moduli* should be used at the current (*k*th) iteration, the value of the sign_i(·) is

$$\operatorname{sign}_{i}(\sigma_{s}) = \begin{cases} 1 & \text{if } \sigma_{s} \ge 0, \\ R_{\operatorname{TCE}}^{(i)} & \text{if } \sigma_{s} < 0. \end{cases}$$
(12)

where $R_{\text{TCE}}^{(i)}$ is the moduli ratio of the *i*th bi-modulus material.

(b) If the element has *tensile moduli* at the (k-1)th iteration and *compressive moduli* should be used at the current kth iteration, the value of the sign_i(·) is

$$\operatorname{sign}_{i}(\sigma_{s}) = \begin{cases} 1 & \text{if } \sigma_{s} \leq 0, \\ (R_{\operatorname{TCE}}^{(i)})^{-1} & \text{if } \sigma_{s} > 0. \end{cases}$$
(13)

By comparing Eqs. (10) and (11), we find that the two SEDs are identical when the element is under pure tension or pure compression. If the element is under a complex stress state, the two SEDs are usually different. The difference in the local effective stiffness of the *j*th element at the *k*th iteration is defined as:

$$M_f = \max\left(10^{-6}, \frac{\text{SED}_{j,k}^{effective}}{\max(\text{SED}_{j,k}, 10^{-10})}\right)$$
(14)

Only when the *j*th element is under pure compression or pure tension state at both of (k-1)th and *k*th iteration, $M_f = 1.0$, i.e., there is unnecessary to modify the local stiffness.

The stiffness matrix of the *j*th element with the "new" isotropic materials can be given as [40, 41]:

$$k_j = \int_{\nu_j} B_j^{\mathrm{T}} \cdot D_j \cdot B_j d\nu \tag{15}$$

where B_j is the displacement-strain matrix, and D_j is the elasticity matrix of the *j*th element.

The modified stiffness matrix of the element is defined as:

$$\bar{k}_j = M_f k_j = \int_{v_j} B_j^{\mathrm{T}} (M_f \cdot D_j) B_j dv.$$
(16)

For the *l*th loading case, the mean compliance of the structure with bi-modulus materials (in Eq. (5)) can be obtained, i.e.,

$$\overline{c_l} = \sum_{j=1}^n u_j^{\mathrm{T}}(M_f k_j) u_j = \sum_{j=1}^n M_f \cdot u_j^{\mathrm{T}} \cdot k_j \cdot u_j.$$
(17)

2.5 Optimization procedure

In the present study, the method of moving asymptotes (MMA) [42] is adopted to solve the optimization problem defined in Eq. (5). The partial differential equations for structural deformation expressed in Eq. (1) with boundaries in Eq. (2) are solved by the commercial software ANSYS 12.0 [43]. In the following numerical examples, the initial elastic moduli of the materials in the elements are the same as the tensile modulus of the first solid material (Material 1). All the initial design variables are considered to be equal.

The MMA procedure contains the following steps:

Step 1 Build a finite element model of the structure, initiate parameters in optimization, and let i = 1.

- Step 2 Find the deformation fields of the structure under MLC by finite element analysis.
- Step 3 Calculate the SED, tension SED, compression SED and the local effective SED of each element in the design domain.
- Step 4 Chose the moduli of the materials in each element by comparing tension SED and compression SED; calculate the value of M_f for each element under a complex stress state.
- Step 5 Compute the values of the objective and constraint functions and their sensitivities.
- Step 6 Update the design variables for each element by the MMA optimizer.
- Step 7 Check the convergence: if the termination criterion is not satisfied, return to Step 2, else go to Step 8.
- Step 8 Judge, if i < m 1, then, i = i + 1, return to Step 2, otherwise, go to Step 9.

Step 9 Save and stop.

The termination criterion is either that the iteration number is greater than 100 or that the change of compliance of the structure satisfies the condition:

$$\left|\frac{c_{k-t}-c_k}{c_k}\right| \le \eta, \ t \in \{1, \ 2, \ 3, \ 4, \ 5\},\tag{18}$$

where η is the algorithm tolerance.

3 Examples and discussion

In this section, numerical examples are considered and assessed by the present algorithm. The code is compiled by combining software MATLAB and ANSYS. In all examples, four-node quadrilateral plane stress elements are employed in the finite element analysis. In optimization, the objective is to minimize the compliance of the structure. The Poisson's ratios of materials in all examples are set to be 0.2. In the following examples, Material 1 is in *Red*, Material 2 is in *Green*. Material 3 is void which is in *White*.

3.1 Example 1: validity assessment

The structure shown in Fig. 2a is a cantilever beam with dimensions 0.7 m × 0.4 m × 0.02 m, and is meshed with 120 × 60 elements. The left side of the structure is fixed. There are two solids and a void phase in the structure. The material tensile moduli of the two solids are 80 and 40 GPa, respectively. $R_{\text{TCE}}^{(1)} = 2 > 1$ and $R_{\text{TCE}}^{(2)} = 0.5 < 1$. Volume fractions of the two solids and void phase are 0.12, 0.12, and 0.76, respectively.

The structure is under two loading cases, i.e., $P_1 = 2000$ N in the first case, $P_2 = 2000$ N in the second

case. P_1 is applied on the center of the right side of the structure. P_2 is applied on the center of the design domain (Fig. 2a). The weighting scheme is $w_1 = 0.2$ and $w_2 = 0.8$. From Fig. 2b, we find that Material 1 (Red) is mainly under tension and Material 2 (Green) is under compression. If we consider the components in the final structure as LTP, Material 1 is mainly on tensile LTPs and Material 2 on compressive LTPs. From the R_{TCE} values of the two materials, we know that the two materials have higher stiffness under the current loading states with $w_1 = 0.2$ and $w_2 = 0.8$. Briefly, the moduli of the two materials are the same, i.e., 80 GPa, in the final structural topology, a result that can also be obtained using single phase topology optimization with an approach such as the SIMP method. With a different weighting scheme, we believe that the amount of material supporting the two forces would be different. But the optimal materials layout must have the property stated above, that the majority of the materials in the final structure should have higher moduli to decrease structural compliance. It is concluded that the correctness of the algorithm is verified.

Figure 3 demonstrates that the structural compliance approaches 0.19782 N m after 101 iterations for the structure under MLC with the weighting scheme of $w_1 = 0.2$ and $w_2 = 0.8$. Because no internal iteration occurs within each update of the design variables, only 101 structural deformation analyses are needed. Hence, the present algorithm has acceptable efficiency during optimization.

3.2 Example 2: effect of R_{TCE} on final materials layouts

The design domain is shown in Fig. 4 with the dimensions 1.6 m × 0.5 m × 0.01 m. The structure is modeled with 160 × 50 elements. Two loading cases are considered. In the first case, $P_1 = 2000$ N is applied at the point " K_1 ". In the second case, two concentrated forces $P_2 = 2000$ N are applied at two points " K_2 ". The design domain contains two solid materials and one void phase, with the volume fractions 0.16, 0.16, and 0.68, respectively. The tensile moduli of the two solids (Material 1 and Material 2) are 100 and 50 GPa, respectively. If the two materials are bimodulus materials, the values of R_{TCE} of the two solids are 2 and 0.5, respectively. For comparison, the optimal layouts of the two isotropic solids are also given.

Three weighting schemes are considered: $w_1 = 1, w_2 = 0$ (only P_1 is active), $w_1 = 0, w_2 = 1$ (only P_2 is active),

 $w_1 = 0.5, w_2 = 0.5$

When the structure is only under the load P_1 , the final layout of the two solid bi-modulus materials (Fig. 5a) is different from the traditional design, that is, the isotropic materials layout. In Fig. 5a, Material 1 is under pure tension Table 1 Final iterations and values of the objective function (Obj) (or mean compliance) of the structure with different load direction schemes

Weighting	Scheme 1 ($\eta = 0.01$)		Scheme 2 ($\eta = 0.01$)		Scheme 3 ($\eta = 0.001$)	
	Iterations	Obj (N m)	Iterations	Obj (N m)	Iterations	Obj (N m)
$w_1 = 0.2, w_2 = 0.8$	77	0.51137	75	0.52606	186	0.43058
$w_1 = 0.5, w_2 = 0.5$	68	0.51368	68	0.51055	168	0.73453
$w_1 = 0.8, w_2 = 0.2$	67	0.50587	79	0.50121	121	0.52510

Fig. 1 Stress-strain curves for a common bi-modulus material and two special cases







(a) Bi-modulus material

(b) Tension-only material

(c) Compression-only material







(a) Initial design



 $\begin{array}{c|c} K_1 & K_2 \\ \hline & P_1 & P_2 \end{array}$

0.5 m



Fig. 3 Iteration histories of the mean compliance of structure under two loading cases with three different weighting schemes ($\eta = 0.01$)

because its tensile modulus is greater than the compressive modulus ($R_{\text{TCE}}^{(1)} = 2 > 1$). Almost all of Material 2 is under compression due to its compressive modulus being greater than the tensile modulus ($R_{\text{TCE}}^{(2)} = 0.5 < 1$). In Fig. 5b, the interfaces between the two isotropic solids are more complex than in Fig. 5a

Fig. 4 Final materials distribution in the structure for different cases (Material 1: Red, Material 2: Green, void: White)

0.4 m

0.4 m

When the structure is subjected only to P_2 , the bi-modulus materials layout is also different from that of isotropic materials. The interface of the two bi-modulus solids (Fig. 5c) is also simpler than that between two isotropic solids (Fig. 5d). Hence, when the moduli of the two solids are clearly different, the complex bi-modulus behavior of the materials does not imply that they have complex interface, which would be difficult for manufacturing.

Under the two loading cases, the final materials layouts (Fig. 5e, f) are different from those in the structure under a single load. Material 1 (bi-modulus) is still mainly under tension, and most of Material 2 is still under compression (Fig. 5e).

Fig. 5 Final materials distribution in the structure for different cases (Material 1: Red, Material 2: Green, void: White)



Fig. 6 The final materials layouts in the structure under different schemes (Material 1: Red, Material 2: Green, void: White)

3.3 Example 3: effect of load directions

The dimensions of the structure used in this example are $1.6 \text{ m} \times 0.4 \text{ m} \times 0.01 \text{ m}$. The structure is simply supported and the finite element mesh is 160×40 . There are two solids (Material 1 and Material 2) and one void phase in the structure. The two solids have the tensile moduli of 100 and 50 GPa, respectively. They are bi-modulus materials, and the values of the R_{TCE} of the two solids are 2 and 0.5, respectively. The volume fractions of the three phases are 0.2, 0.2, and 0.6, respectively.

The structure is under two loading cases with two vertical concentrated forces, P_1 and P_2 , applied separately on the centers of the upper and lower sides of structure. Both



Fig. 7 Iteration histories of the mean compliance of the structure under MLC with different load directions/schemes

forces have a magnitude of 2000 N. In different schemes, the directions of the two forces may be different. Three schemes are considered (see the uppermost layer of Fig. 6).

The results for scheme 1 (the left column) demonstrate that Material 1 (*Red*) is mainly under tension for two loading cases, and Material 2 is under compression when the directions of the two forces are vertical upward. The reason is that the tensile modulus of Material 1 is greater than the compressive modulus ($R_{TCE}^{(1)} = 2 > 1$), whereas Material 2 has higher compressive stiffness ($R_{TCE}^{(2)} = 0.5 < 1$).

This situation changes when the directions of the two forces are vertical downward (scheme 2). In the central column (scheme 2), Material 1 layouts are near the lower side of the beam rather than near the upper side as in scheme 1.

If the directions of the two forces are different, as in scheme 3, the layouts of the two bi-modulus materials are clearly different from those in schemes 1 and 2. Moreover, the locations of Material 1 in the final structure depend on the weighting coefficients of the two load cases. For example, Material 1 layouts are near the upper side when $w_1 = 0.2$, $w_2 = 0.8$ (P_2 has higher influence on structural compliance than P_1), in the central area of the structure when $w_1 = 0.5$, $w_2 = 0.5$, or near the bottom when $w_1 = 0.8$, $w_2 = 0.2$.

This phenomenon does not occur in either scheme 1 or scheme 2. The weighting coefficients of the two loading cases have only a slight influence on the positions of the material interfaces in scheme 1. Only when $w_1 = 0.8$, $w_2 = 0.2$ is the topology of the structure different from the other topologies in scheme 2.

In Fig. 7 and Table 1, we find that the number of iterations does not exceed 200 in the 9 different cases. With the same weighting coefficients, namely $w_1 = 0.2$, $w_2 = 0.8$, the optimal value of structural compliance is different for different load directions. Simultaneously, the layout differences among the three schemes are attributable to the bi-modulus behavior of the two materials under loads with different directions [44].

4 Conclusions

Using the algorithm presented in this study to achieve optimal layout of multiple bi-modulus materials in a continuum under MLC, three numerical tests are considered. From the numerical results, some remarkable conclusions are drawn.

- 1. The computational cost of the present algorithm is very close to that of simple single material layout optimization by the SIMP method
- 2. In a stiffness design, materials with higher modulus should be laid out on the main LTPs. When the differences among the moduli of bimodulus materials are not great, on the tensile LTPs, materials with $R_{\text{TCE}} > 1$ are usually laid out. Materials with $R_{\text{TCE}} < 1$ are usually laid out on the compressive LTPs.
- 3. The final layouts of bi-modulus materials are sensitive to the values of R_{TCE} and the load conditions. Under the same loading conditions, the interfaces between bi-modulus materials may be clearer than those between isotropic materials
- 4. The optimal layout of bi-modulus materials depends on the force directions (forward and reverse).

Hence, the present algorithm is applicable and effective for analyzing the performance of structures with many bimodulus materials and under MLC. **Acknowledgments** Financial support from the National Natural Science Foundation of China (Grant No. 51179164) and the Australian Research Council (Grant No. DP140103137) is acknowledged.

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