Budapest University of Technology and Economics Short Course on Topology Optimization of Structures



## Short Course on Topology Optimization of Structures

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### Schedule

DAY 1 Topology optimization of structures: basics

DAY 2

Stress-constrained topology optimization

DAY 3

Mixed finite elements for the optimal design of structures

DAY 4

Analysis and design of no-tension structures, by formulating optimization problems

#### 8 April 2020 Budapest University of Technology and Economics Short Course on Topology Optimization of Structures



## **Topology optimization of structures: basics**

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### Structural and topology optimization



a) sizing optimization:

the areas of the elements of a fixed truss "ground structure" are unknown b) shape optimization:

the parameters describing the geometry of the boundaries are unknown *c) topology optimization:* 

the distribution of material is unknown

### Structural topology optimization: a design tool

"The art of structure is where to put the holes"

Robert Le Ricolais (1894–1977) French/American Engineer, "Father of Spatial Structures"



### Structural topology optimization: a design tool





Gigantic tree-like columns support the overhanging roof canopy of the Qatar National Convention Centre by Japanese Architect Arata Isozaki

### Structural topology optimization: a design tool



A one-of-a-kind project: a conceptual design for the Zendai competition (China) created with topology optimization by Prof. Glaucio Paulino's research group along with Skidmore, Owings & Merrill LLP



Illustration for the concept design of a 288 m tall high-rise in Australia, showing the engineering and architecture expressed together at Skidmore, Owings & Merrill LLP (Beghini, Katz, Baker and Paulino, 2014)

### Outline

## Governing equations:

- > "penalized" elasticity problem and structural compliance
- Problem formulation:
  - conventional volume-constrained minimum compliance formulation
- Solution of the minimization problem:
  - gradient-based algorithms
  - sensitivity computation
  - numerical issues
- Applications:
  - design of stiff structures (with 88-line Matlab code)
  - design of compliant mechanisms
  - design of periodic microstructures

# Structural topology optimization: state equations (linear elastic problem)

 $\sigma_{ij,i} + \mathfrak{X}_i = 0$ 

 $\sigma_{ij} = C_{ijhk} \varepsilon_{hk}$ 



A homogeneous domain  $\Omega \in \mathbb{R}^2$  with a regular boundary  $\Gamma$  is considered, assuming that  $\Gamma = \Gamma_d \cup \Gamma_t$ .

Prescribed displacements with components  $u_{0j}$  and tractions with components  $t_{0j}$  are assigned on  $\Gamma_d$  and  $\Gamma_t$   $g_j$  are the components of the vector of body loads in  $\Omega$  (generally neglected)

 $C_{iihk}$  are the component of the 4th order elasticity tensor

Let be  $\underline{u}$  the unknown displacement field,  $\underline{\sigma}$  the unknown stress field and  $\underline{\varepsilon}$  the unknown strain field. One has:

equilibrium

compatibility

 $arepsilon_{ij} = (u_{i,j} + u_{j,i})/2 ~~$  in  $\Omega$ , along with

$$u_j \mid_{\Gamma_d} = u_{0j}$$

constitutive law

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in  $\Omega$ , along with  $\sigma_{ij}n_i \mid_{\Gamma_t} = t_{0j}$ 

# Structural topology optimization: mathematical formulation

Given a domain with assigned loads and boundary conditions, find the distribution of linear elastic isotropic material that minimizes an assigned scalar function for a fixed set of constraints

The elastic problem depends on  $\psi(\chi)$ : find  $\underline{u} \in H^1$  such that  $\underline{u} \mid_{\Gamma_u} = \underline{u}_0$  and

$$\int_{\Omega} \psi(\chi) C^0_{ijhk} \varepsilon_{hk}(\underline{u}) \varepsilon_{ij}(\underline{v}) \ dx = \int_{\Gamma_t} t_{0j} v_j \ ds, \quad \forall \underline{v} \in H^1$$

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# Structural topology optimization: mathematical formulation

Objective function: "structural compliance", work of the external loads at equilibrium, twice the strain energy stored in the structure (Clapeyron th.), measure of the structural deformability

$$\mathcal{C} = W(\underline{u}) = \int_{\Gamma_t} t_{0j} \, u_j \, ds = \int_{\Omega} \psi(\chi) C^0_{ijhk} \varepsilon_{hk}(\underline{u}) \varepsilon_{ij}(\underline{u}) \, d\Omega$$

Volume constraint: enforcement of the available "volume fraction" of material

$$\Omega_{mat}/\Omega \le V_f$$
  $\int_{\Omega} \chi d\Omega / \int_{\Omega} d\Omega \le V_f$   $V_f \le 1$ 

**Objective function** 

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 $\min_{\boldsymbol{w} \in \boldsymbol{w}} \quad \mathcal{C} = W(\underline{u})$ 

# Structural topology optimization: NAND/SAND

#### Two unknown fields arise in the formulation: density and displacement

In the conventional Nested Analysis and Design Approach (NAND) only the structural design variable is treated as the optimization variable. In the more demanding Symultaneous Analysis and Design Approach (SAND) both variables enter the optimization problem

$$\begin{array}{ll} \min_{\psi \in \psi_{ad}} & \mathcal{C} = W(\underline{u}) & \text{Objective function} \\ \text{s.t.} & \int_{\Omega} \psi(\chi) C_{ijhk}^{0} \varepsilon_{hk}(\underline{u}) \varepsilon_{ij}(\underline{v}) \ dx = \int_{\Gamma_{t}} t_{0j} v_{j} \ ds, \ \forall \underline{v} \in H^{1}, \\ \int_{\Omega} \chi d\Omega \Big/ \int_{\Omega} d\Omega \leq V_{f}, & \text{Volume constraint} \end{array}$$

 $\max_{\psi \in \psi_{ad}} \min_{\underline{v} \in \underline{v}_{ad}} \Pi(\underline{u}) \qquad \Pi(\underline{u}) = U(\underline{u}) - W(\underline{u}) = W(\underline{u})/2 - W(\underline{u}) = -W(\underline{u})/2 = -\mathcal{C}(\underline{u})/2$ 

s.t. 
$$\int_{\Omega} \chi d\Omega \Big/ \int_{\Omega} d\Omega$$

 $\leq V_f$ 

Max-min problem in terms of the total potential energy

# Structural topology optimization: is the continuous problem well-posed?

Minimization problem with discrete values (0-1): the formulation has no feasible solution in the case of isotropic material. The stiffest geometry calls for the largest number of holes, finally achieving "optimal microstructures"



An enlargement of the design domain is needed

"Void" "Full material"

 $0 \le \rho(\chi) \le 1$ 

Function representing a **continuous** material density, i.e. the minimization unknown

1. allowing for microstructures with intermediate densities between 0-1: <u>optimization of composite materials</u>

 alternatively, introducing a penalization of the intermediate densities to achieve pure 0-1 design: <u>optimization by distribution of isotropic material</u>

## Structural topology optimization: topology optimization of composite materials

1. Material density represents the material as a microstructure ("gray" is allowed)

The microstructure is a composite material with an infinite number of infinitely small voids, leading to a porous composite with a density varying between 0 and 1. Macroscopic mechanical features can be derived through homogenization methods.

Reference

domain



Since the macroscopic properties of common types of microstructure are not isotropic an orientation angle is also needed



# Structural topology optimization: topology optimization by distribution of isotropic material

2. Intermediate material density is penalized to achieve 0-1 design (no "gray")



SIMP: Solid Isotropic Microstructure with Penalty (Rozvany et al, 92; Bendsøe and Sigmund, 99; Berke 70)

 $E(\rho) = \rho^p E^0 \qquad \rho_{min} \le \rho(\chi) \le 1 \qquad \rho_{min} = 10^{-3}$ 

- The power p > 1 penalizes intermediate densities to achieve pure 0-1 design (usually p=3).  $\rho_{min}$  is needed against FEM singularities
- It makes even the compliance minimization a nonconvex problem



RAMP: Rationale Approximation of Material Properties (Stolpe and Svanberg, 03)

$$E(\rho) = E_{min} + \frac{\rho}{1 + q(1 - \rho)} (E^0 - E_{min}) \frac{0 \le \rho(\chi) \le 1}{E_{min} = 10^{-9} E^0}$$

- A convex interpolation model implies a convex objective function. Unfortunately, this result is achieved only if  $q \ge E_0/E_{min}$ -1
- It works fine for two-material problems with  $E^0 = E^{01} E_{min} = E^{02}$

# Structural topology optimization: topology optimization by distribution of isotropic material





- ➢ SIMP/RAMP improves the numerical tractability of the continuous setting (topology → sizing), but does not completely solve the problem (mesh dependence of the discrete setting)
- For p ≥ 3, SIMP can be seen as a «material model»

 $C_{ijhk}(\rho)$  corresponds to a composite material constructed from void and the given material at a real density  $\rho$ , since the bulk modulus *k* and the shear modulus  $\mu$  satisfy the Hashin-Shtrikman bounds:

$$\begin{split} 0 &\leq \kappa \leq \frac{\rho \kappa^{0} \mu^{0}}{(1-\rho)\kappa^{0} + \mu^{0}}, \quad 0 \leq \mu \leq \frac{\rho \kappa^{0} \mu^{0}}{(1-\rho)(\kappa^{0} + 2\mu^{0}) + \kappa^{0}}\\ 0 &\leq E \leq E^{*} = \frac{\rho E^{0}}{3 - 2\rho} \end{split}$$

### Governing equations: SIMP-based elasticity problem

Given a domain with assigned loads and boundary conditions, find the distribution of linear elastic isotropic material that minimizes an assigned scalar function for a fixed set of constraints

"Void" "Full material" 
$$\rho_{min} \leq \rho(\chi) \leq 1$$

$$C_{ijhk}(\rho(\chi)) = \rho(\chi)^p C_{ijhk}^0$$

Elasticity tensor of the given isotropic material ("full material") Function representing the **continuous** material density, i.e. the minimization unknown

SIMP: Solid Isotropic Microstructure with Penalty i.e. the material interpolation scheme to represent the constitutive tensor depending on  $\rho$ 

p>1 to penalize intermediate densities and achieve a pure 0-1 design

The elastic problem depends on  $\rho(\chi)$ : find  $\underline{u} \in H^1$  such that  $\underline{u} \mid_{\Gamma_u} = \underline{u}_0$  and

$$\int_{\Omega} \rho^p C^0_{ijkl} \varepsilon_{ij}(\underline{u}) \varepsilon_{kl}(\underline{v}) \ d\Omega = \int_{\Gamma_t} \underline{t}_0 \cdot \underline{v} \ d\Gamma \quad \forall \underline{v} \in H^1$$

### Governing equations: SIMP-based structural compliance

A classical scheme for the discretization of the above problem adopts the same mesh of four-node elements and a two-field interpolation with:

- A piecewise constant density discretization (with unknowns x)
- A bilinear displacement approximation (with unknowns **U**)

$$\mathbf{K}(\mathbf{x}) \mathbf{U} = \sum_{e=1}^{N} x_e^p \mathbf{K}_e^0 \mathbf{U}_e = \mathbf{F}$$



topology optimization  $\rightarrow$  sizing optimization

Compliance: - work of the external loads at equilibrium (Clapeyron th.)
 - measure of the structural deformability

$$\mathcal{C} = \int_{\Omega} \rho_h^p C_{ijkl}^0 \varepsilon_{ij}(\underline{u}_h) \varepsilon_{kl}(\underline{u}_h) \, d\Omega = \mathbf{U}^T \mathbf{K} \mathbf{U} = \sum_{e=1}^N x_e^p \mathbf{U}_e^T \mathbf{K}_e^0 \mathbf{U}_e,$$

**U**<sub>e</sub>: element displacements vector **K**<sub>e</sub><sup>0</sup>: element stiffness matrix for virgin material

## Problem formulation: classical formulation for maximum stiffness

MCW Minimum Compliance with Weight (volume) constraint (Bendsøe and Kikuchi, 88)

Given a domain with assigned loads and boundary conditions, find the distribution of an amount of linear elastic isotropic material that minimizes the compliance

$$\min_{\substack{x_{min} \leq x_e \leq 1 \\ \text{bound/side constraints}}} \mathcal{C} = \mathbf{U}^T \mathbf{K} \mathbf{U} = \sum_{e=1}^N x_e^p \mathbf{U}_e^T \mathbf{K}_e^0 \mathbf{U}_e$$
s.t. 
$$\mathbf{K}(\mathbf{x}) \ \mathbf{U} = \sum_{e=1}^N x_e^p \mathbf{K}_e^0 \mathbf{U}_e = \mathbf{F},$$

$$\sum_{e=1}^N x_e A_e / \sum_{e=1}^N A_e \leq V_f,$$

Structural compliance

Governing eqns. elastic problem

Volume constraint

 $\succ$  For low  $V_f$ , truss-like structures arise / for high  $V_f$ , beam with optimal openings are found



# Solution of the minimization problem: gradient-based minimization

Nested Analysis and Design Approach (NAND): only the structural design variable is treated as the optimization variable



Iterative algorithms:

macrostructural non-gradient approaches, in general heuristic methods: evolutionary approaches, fully stressed design method...

gradient-based approaches:

- optimality criteria
- mathematical programming (MMA, CONLIN...)

sensitivity analysis

### Sensitivity analysis: fundamentals

At each iteration, values and derivatives with respect to x<sub>k</sub> are computed for the objective function and the volume

 $\mathcal{C} = \mathbf{F}^T \mathbf{U}(\mathbf{x}) \quad \text{where } \mathbf{U} \text{ solves } \mathbf{K} \mathbf{U} = \mathbf{F}$ Assuming that  $\mathbf{F}$  does not depend on  $\mathbf{x}_k$  one has:  $\frac{\partial \mathcal{C}}{\partial x_k} = \mathbf{F}^T \frac{\partial \mathbf{U}}{\partial x_k}$ Using the chain rule on the equilibrium equation one has:  $\frac{\partial \mathbf{K}}{\partial x_k} \mathbf{U} + \mathbf{K} \frac{\partial \mathbf{U}}{\partial x_k} = \mathbf{0} \quad \rightarrow \quad \mathbf{K} \frac{\partial \mathbf{U}}{\partial x_k} = -\frac{\partial \mathbf{K}}{\partial x_k} \mathbf{U}$   $\mathcal{V} = \sum_{e=1}^N x_e A_e \quad \frac{\partial \mathcal{V}}{\partial x_k} = A_k$ 

1+N systems of equations must be solved, per step

### Sensitivity analysis: the adjoint method

- - Derivatives with respect to x<sub>k</sub> for the objective function can be computed more efficiently through the adjoint method
    - o.f. is re-written adding a «zero function», with  $\tilde{\mathbf{U}}$  arbitrary but fixed real vector

$$\begin{aligned} \mathcal{C} &= \mathbf{F}^T \mathbf{U} - \tilde{\mathbf{U}}^T (\mathbf{K} \mathbf{U} - \mathbf{F}) \\ \frac{\partial \mathcal{C}}{\partial x_k} &= \mathbf{F}^T \frac{\partial \mathbf{U}}{\partial x_k} - \tilde{\mathbf{U}}^T (\frac{\partial \mathbf{K}}{\partial x_k} \mathbf{U} + \mathbf{K} \frac{\partial \mathbf{U}}{\partial x_k}) \\ \frac{\partial \mathcal{C}}{\partial x_k} &= (\mathbf{F}^T - \tilde{\mathbf{U}}^T \mathbf{K}) \frac{\partial \mathbf{U}}{\partial x_k} - \tilde{\mathbf{U}}^T \frac{\partial \mathbf{K}}{\partial x_k} \mathbf{U} \end{aligned}$$

that can in turn be written as:

$$\frac{\partial C}{\partial x_k} = -\tilde{\mathbf{U}}^T \frac{\partial \mathbf{K}}{\partial x_k} \mathbf{U}$$
 where  $\tilde{\mathbf{U}}$  satisfies the adjoint eqn.  $\mathbf{F}^T - \tilde{\mathbf{U}}^T \mathbf{K} = \mathbf{0}$ 

## Sensitivity analysis: the adjoint method

The adjoint equation is in the form of an equilibrium equation (self-adjoint problem)

$$\mathbf{F}^T - \tilde{\mathbf{U}}^T \mathbf{K} = \mathbf{0} \quad \rightarrow \quad \mathbf{K}\tilde{\mathbf{U}} = \mathbf{F} \quad \rightarrow \quad \tilde{\mathbf{U}} = \mathbf{U}$$

$$\frac{\partial \mathcal{C}}{\partial x_k} = -\mathbf{U}^T \frac{\partial \mathbf{K}}{\partial x_k} \mathbf{U} = -p x_k^{p-1} \mathbf{U}_k^T \mathbf{K}_k^0 \mathbf{U}_k$$

- only 1 system of equations must be solved, per step (very efficient)
- the sensitivity is "localized": only information at the element level is involved (the effect of the other variables is "hidden" in U<sub>k</sub>)
- the sensitivity is negative for any element: according to physical intuition, additional material in any element decreases compliance, i.e. increases stiffness

## Update of the design variables: optimality criterion

Heuristic method based on the Lagrangian function of the optimization problem

the o.j. is augmented by the constraints through a set of non-negative multipliers

$$\mathcal{L} = \mathcal{C} + \lambda (V - V_f V_0) + \lambda_1^T (\mathbf{KU} - \mathbf{F}) + \sum_{e=1}^N \lambda_{2e} (x_{min} - x_e) + \sum_{e=1}^N \lambda_{3e} (x_e - 1)$$
  
o.f. volume con. equilibrium con. lower bounds upper bounds

optimality arises when all derivatives of the Lagrangian with respect to  $x_k$  are zero

$$\frac{\partial \mathcal{L}}{\partial x_k} = \frac{\partial \mathcal{C}}{\partial x_k} + \lambda \frac{\partial V}{\partial x_k} + \lambda_1^T \frac{\partial \mathbf{KU}}{\partial x_k} - \lambda_{2e} + \lambda_{3e} = 0 \qquad k = 1, 2, \dots N$$

# Update of the design variables: optimality criterion

Assuming that side constraints are not active and **F** is design-independent:

$$\frac{\partial \mathcal{L}}{\partial x_k} = \frac{\partial \mathcal{C}}{\partial x_k} + \lambda \frac{\partial V}{\partial x_k} + \lambda_1^T \frac{\partial \mathbf{K} \mathbf{U}}{\partial x_k} - \lambda_{2e} + \lambda_{3e} = -p x_k^{p-1} \mathbf{U}_k^T \mathbf{K}_k^0 \mathbf{U}_k + \lambda A_k = 0$$
  
$$k = 1, 2, \dots N$$

Hence, for intermediate densities, the **strain energy density-like term is constant all over the domain in the optimal solution**. Since areas with high energy are expected to be too low on stiffness, a fix-point type update scheme can be formulated. For intermediate density:

$$x_{k}^{j+1} = x_{k}^{j} \left( \frac{p x_{k}^{p-1} \mathbf{U}_{k}^{T} \mathbf{K}_{k}^{0} \mathbf{U}_{k}}{\lambda A_{k}} \right)^{\eta} = x_{k}^{j} (B_{k}^{j})^{\eta}$$
Add material where  $\mathbf{B}_{k} > 1$   
otherwise remove material updated current current =1 at optimum optimum (0.5 in general)

# Update of the design variables: optimality criterion

A general scheme accounting for side constraints reads:

$$x_{k}^{j+1} = \begin{cases} \max\{(1-\zeta)x_{k}^{j}, x_{min}\} & \text{if } x_{k}^{j}(B_{k}^{j})^{\eta} \leq \max\{(1-\zeta)x_{k}^{j}, x_{min}\} \\ \min\{(1+\zeta)x_{k}^{j}, 1\} & \text{if } \min\{(1+\zeta)x_{k}^{j}, 1\} \leq x_{k}^{j}(B_{k}^{j})^{\eta} \\ x_{k}^{j}(B_{k}^{j})^{\eta} & \text{otherwise} \end{cases}$$

- Add material where B<sub>k</sub> >1, otherwise remove material
- This only takes place if the update does not violate the bounds on x<sub>k</sub>
- A positive move limit ζ (0.2 in general) is introduced to ensure that no big change in relative density arises between two subsequent steps
- The lagrangian multiplier λ is updated in an inner iteration loop using bisection in order to satisfy the active volume constraint

# Update of the design variables: mathematical programming

**Sequential convex programming**: a sequence of explicit sub-problems is used that are convex approximations of the original one

- Sequential linear and quadratic programming techniques attack the problem without accounting for the specific characteristics of the involved functions: they generate sub-problems by linearizing both objective function and constraints in the **direct variables**
- MMA (Svanberg, 1987) and CONLIN (Fleury, 1986) linearize objective function and constraints in the direct variables and in the reciprocal variables, depending on the sign of the gradient. Such an approximation perfectly suits a broad range of structural optimization problems

## Update of the design variables: Method of Moving Asymptotes (MMA)

MMA (Svanberg, 1987) CONLIN (Fleury, 1986) use a sequence of simpler approximated sub-problems of given type. They are **separable**, **convex** and are constructed on the current sensitivity information as well as some history



## Update of the design variables: Method of Moving Asymptotes (MMA)

- separable subproblems → necessary conditions of optimality do not couple the primary variables (the design variables)
- convex approximations  $\rightarrow$  efficient dual methods can be used

For the volume-constrained minimum compliance problem, the sensitivity is negative and an MMA **convex** approximation of the o.f. after the j-th step reads:

$$\mathcal{C}(\mathbf{x}^j) - \sum_{e=1}^N \frac{(x_e^j - L_e)^2}{x_e - L_e} \frac{\partial \mathcal{C}}{\partial x_e}(\mathbf{x}^j)$$

A dual method can be used:

1) the Lagrange functional is minimized **element by element** with respect to all  $x_e$ 

2) then, the resulting functional is maximized with respect to  $\lambda$ 

$$\mathcal{L} = \mathcal{C}(\mathbf{x}^j) - \sum_{e=1}^N \frac{(x_e^j - L_e)^2}{x_e - L_e} \frac{\partial \mathcal{C}}{\partial x_e}(\mathbf{x}^j) + \lambda \left(\sum_{e=1}^N x_e A_e - V_f V_0\right)$$

Flexibility + Excellent performance in case of a limited number of constraints

### **Numerical issues**

Checkerboard patterns and mesh dependence (Sigmund and Petersson, 98)



Checkboarded layouts are optimal solutions from a mathematical point of view, but they are not feasible from a physical point of view (this depends on the adopted FEM and density discretization)





Mesh dependence: different solutions arise for different meshes (this is the discrete counterpart of not well-posedness)



# Numerical issues: filters

 Filtered sensitivities of the objective function can be successfully used to prevent numerical instabilities, i.e. mesh dependence and checkerboard patterns

 $r_{min}$  = 1.5 d<sub>m</sub> is the minimum value to avoid the arising of undesired checkerboard patterns

 $r_{min}$  > 1.5 d<sub>m</sub> provides control on the minimum thickness of any member of the design



e

### Numerical issues: filters

Alternatively, filtered densities can be easily implemented for increased robustness 



The chain rule is needed to compute the derivatives of o.f. and constraints:  $\frac{\partial \psi}{\partial x_k} = \sum_{i=1}^{n} \frac{\partial \psi}{\partial \widetilde{x}_l} \frac{\partial x_l}{\partial x_k}$ 

Heaviside projection filters can be implemented to reduce blurred edges in case of big r<sub>min</sub> 



*e*-th intermediate variable *l*-th design variable

 $\beta \rightarrow \infty$  Heaviside step function

A continuation approch is needed starting the first optimization from large values of  $\beta$ : high CPU cost

# Numerical issues: filters



Optimal layouts using the same filtering radius  $r_{min}$  and finer meshes: density filter vs. HS projection filter



Optimal layouts using increasing values of the filtering radius *r<sub>min</sub>* on the same mesh: HS projection filter

#### Other numerical issues



• What about uniqueness of the solution?

Most problems are non-convex  $\rightarrow$  multiple solutions arise (one big bar/many small bars under uni-axial tension)

What about local minima?

Most problems are non-convex  $\rightarrow$  many local minima arise (multi-start procedures along with globally convergent algorithms can be used, but there is no guarantee of global optimality)

• How to choose the starting guess?

 $x_e = V_f$  or  $x_e = 1$  are generally used to start the minimization all over the domain, no matter for their feasibility with respect to the volume constraint

How to choose the convergence criterion?

In general, check on the maximum change in density / o.f. between two subsequent steps

Efficient topology optimization in MATLAB using 88 lines of code (Andreassen et al, 2011)

It exploits a mesh of *nelx* x *nely* four-node finite elements with  $d_m=1$ . For each element, one density unknown. It allows for two filtering techniques (ft=1/2 sensitivity/density filter)

```
%%%% AN 88 LINE TOPOLOGY OPTIMIZATION CODE %%%%
1
   function top88(nelx,nely,volfrac,penal,rmin,ft)
   %% MATERIAL PROPERTIES
                                                                       E(\rho) = E_{min} + \rho^p (E^0 - E_{min})
3
   E0 = 1;
   Emin = 1e-9;
                                                                                0 < \rho(\chi) \leq 1
5
6 nu = 0.3;
7 %% PREPARE FINITE ELEMENT ANALYSIS
8 A11 = [12 \ 3 \ -6 \ -3; \ 3 \ 12 \ 3 \ 0; \ -6 \ 3 \ 12 \ -3; \ -3 \ 0 \ -3 \ 12];
                                                                      A12 = [-6 -3 \ 0 \ 3; \ -3 -6 -3 -6; \ 0 \ -3 -6 \ 3; \ 3 -6 \ 3 -6];
9
  B11 = \begin{bmatrix} -4 & 3 & -2 & 9; \\ 3 & -4 & -9 & 4; \\ -2 & -9 & -4 & -3; \\ 9 & 4 & -3 & -4 \end{bmatrix};
10
 B12 = [2 -3 4 -9; -3 2 9 -2; 4 9 2 3; -9 -2 3 2];
11
  KE = 1/(1-nu^2)/24*([A11 A12;A12' A11]+nu*[B11 B12;B12' B11]);
                                                                              12
   nodenrs = reshape(1:(1+nelx)*(1+nely),1+nely,1+nelx);
13
   edofVec = reshape(2*nodenrs(1:end-1,1:end-1)+1,nelx*nely,1);
14
   edofMat = repmat(edofVec,1,8)+repmat([0 1 2*nely+[2 3 0 1] -2 -1],nelx*nely,1);
15
   iK = reshape(kron(edofMat,ones(8,1))',64*nelx*nely,1);
16
   jK = reshape(kron(edofMat,ones(1,8))',64*nelx*nely,1);
17
```

### **Efficient implementation in Matlab**

- 18 % DEFINE LOADS AND SUPPORTS (HALF MBB-BEAM)
- 19 F = sparse(2,1,-1,2\*(nely+1)\*(nelx+1),1);
- 20 U = zeros(2\*(nely+1)\*(nelx+1),1);
- 21 fixeddofs = union([1:2:2\*(nely+1)],[2\*(nelx+1)\*(nely+1)]);
- 22 alldofs = [1:2\*(nely+1)\*(nelx+1)];
- 23 freedofs = setdiff(alldofs,fixeddofs);




## **Efficient implementation in Matlab**

```
%% PREPARE FILTER
24
   iH = ones(nelx*nely*(2*(ceil(rmin)-1)+1)^2,1);
25
   jH = ones(size(iH));
26
   sH = zeros(size(iH));
27
   k = 0;
28
   for i1 = 1:nelx
29
      for j1 = 1:nely
30
        e1 = (i1-1)*nely+j1;
31
        for i2 = max(i1-(ceil(rmin)-1),1):min(i1+(ceil(rmin)-1),nelx)
32
          for j2 = max(j1-(ceil(rmin)-1), 1):min(j1+(ceil(rmin)-1), nely)
33
            e2 = (i2-1)*nely+j2;
34
            k = k+1;
35
            iH(k) = e1;
36
            jH(k) = e2;
37
            sH(k) = max(0, rmin-sqrt((i1-i2)^2+(j1-j2)^2));
38
          end
39
        end
40
      end
41
    end
42
   H = sparse(iH, jH, sH);
43
   Hs = sum(H,2);
44
   %% INITIALIZE ITERATION
45
   x = repmat(volfrac,nely,nelx);
46
                                           x design unknown
   xPhys = x;
47
                                           xPhys physical unknown
   loop = 0;
48
   change = 1;
49
```



## **Efficient implementation in Matlab**

50 %% START ITERATION  
51 while change > 0.01 convergence criterion  
52 loop = loop + 1;  
53 %% FE-ANALYSIS  
54 SK = reshape(KE(:)\*(Emin+xPhys(:)'.^penal\*(EO-Emin)),64\*nelx\*nely,1); K(x) U = 
$$\sum_{e=1}^{N} x_e^p K_e^0 U_e = F$$
  
55 K = sparse(iK,jK,sK); K = (K+K')/2;  
56 U(freedofs) = K(freedofs,freedofs)\F(freedofs);  
57 %% OBJECTIVE FUNCTION AND SENSITIVITY ANALYSIS  
58 ce = reshape(sum((U(edofMat)\*KE).\*U(edofMat),2),nely,nelx);  $C = \mathbf{U}^T \mathbf{K} \mathbf{U} = \sum_{e=1}^{N} x_e^p \mathbf{U}_e^T \mathbf{K}_e^0 U_e$   
59 c = sum(sum((Emin+xPhys.^penal\*(EO-Emin)).\*ce));  
60 dc = -penal\*(EO-Emin)\*xPhys.^(penal-1).\*ce;  $\frac{\partial C}{\partial x_k} = -px_k^{p-1}\mathbf{U}_k^T \mathbf{K}_k^0 \mathbf{U}_k \qquad \frac{\partial \mathcal{V}}{\partial x_k} = A_k$   
61 dv = ones(nely,nelx);  $\frac{\partial C}{\partial x_k} = -px_k^{p-1}\mathbf{U}_k^T \mathbf{K}_k^0 \mathbf{U}_k \qquad \frac{\partial \mathcal{V}}{\partial x_k} = A_k$   
63 if ft == 1  
64 dc(:) = H\*(x(:).\*dc(:))./Hs./max(1e-3,x(:));  
65 elseif ft == 2  
66 dc(:) = H\*(dc(:)./Hs);  
67 dv(:) = H\*(dv(:)./Hs);

## **Efficient implementation in Matlab**

```
%% OPTIMALITY CRITERIA UPDATE OF DESIGN VARIABLES AND PHYSICAL DENSITIES
69
      11 = 0; 12 = 1e9; move = 0.2;
70
                                                                                          inner loop in the o.c. update to
      while (12-11)/(11+12) > 1e-3
71
        lmid = 0.5*(l2+l1);
                                                                                          enforce the volume constraint
72
        xnew = max(0,max(x-move,min(1,min(x+move,x.*sqrt(-dc./dv/lmid)))));
                                                                                             through the multiplier \lambda
73
        if ft == 1
74
          xPhys = xnew;
75
                                                                                         \tilde{x}_e = \frac{1}{\sum_N H_{el}} \sum_N H_{el} x_l,
        elseif ft == 2
76
          xPhys(:) = (H*xnew(:))./Hs;
77
        end
78
        if sum(xPhys(:)) > volfrac*nelx*nely, l1 = lmid; else l2 = lmid; end
79
      end
80
      change = max(abs(xnew(:)-x(:)));
81
      x = xnew;
82
                                                                                              xPhys(passive==1) = 0;
      %% PRINT RESULTS
83
      fprintf(' It.:%5i Obj.:%11.4f Vol.:%7.3f ch.:%7.3f\n',loop,c, ...
                                                                                              xPhys(passive==2) = 1;
84
        mean(xPhys(:)),change);
                                                                                             to enforce black/white regions
85
      %% PLOT DENSITIES
86
      colormap(gray); imagesc(1-xPhys); caxis([0 1]); axis equal; axis off; drawnow;
87
    end
88
```

## A variation of the theme: multiple loads

Optimal design for more than one load case (M) can be achieved working with the sum of the relevant strain energies

$$\min_{x_{min} \le x_e \le 1} \quad \mathcal{C} = \sum_{i=1}^{M} \mathbf{U}_i^T \mathbf{K} \mathbf{U}_i = \sum_{i=1}^{M} \sum_{e=1}^{N} x_e^p \mathbf{U}_{i,e}^T \mathbf{K}_e^0 \mathbf{U}_{i,e}$$
s.t. 
$$\mathbf{K}(\mathbf{x}) \quad \mathbf{U}_i = \sum_{e=1}^{N} x_e^p \mathbf{K}_e^0 \mathbf{U}_{i,e} = \mathbf{F}_i, \quad i = 1...M$$

$$\sum_{e=1}^{N} x_e A_e / \sum_{e=1}^{N} A_e \le V_f,$$







1 + 2 (1 load case)



1 and 2 (2 load cases)

# A variation of the theme: multiple loads

- - Multiple load cases can be implemented to improve robustness of the layouts



statically determinate and completely constrained

Load condition can be probabilistic to account for several uncertainties (Lógó 2013)

# A variation of the theme: compliant mechanisms

Design of flexible (micro-)mechanisms that transfer an input force or displacement to another point through elastic body deformation

Given a domain with assigned loads and boundary conditions, find the distribution of a prescribed amount of linear elastic isotropic material that maximizes the output displacement:



## Sensitivity analysis: the adjoint method

- > Derivatives with respect to  $x_k$  for the o.f. are computed via the adjoint method
  - o.f. is re-written adding a «zero function», with  $\lambda$  arbitrary but fixed real vector

$$\begin{aligned} u_{out} &= \mathbf{L}^T \mathbf{U} - \boldsymbol{\lambda}^T \left( [\mathbf{K}(\mathbf{x}) + \mathbf{K}_s] \ \mathbf{U} - \mathbf{F} \right) \\ \frac{\partial u_{out}}{\partial x_k} &= \mathbf{L}^T \frac{\partial \mathbf{U}}{\partial x_k} - \ \boldsymbol{\lambda}^T \left( \frac{\partial \mathbf{K}}{\partial x_k} \mathbf{U} + [\mathbf{K}(\mathbf{x}) + \mathbf{K}_s] \frac{\partial \mathbf{U}}{\partial x_k} \right) \\ \frac{\partial u_{out}}{\partial x_k} &= \left( \mathbf{L}^T - \boldsymbol{\lambda}^T [\mathbf{K}(\mathbf{x}) + \mathbf{K}_s] \right) \frac{\partial \mathbf{U}}{\partial x_k} - \mathbf{\lambda}^T \frac{\partial \mathbf{K}}{\partial x_k} \mathbf{U} \end{aligned}$$

that can in turn be written as:

$$\frac{\partial u_{out}}{\partial x_k} = -\boldsymbol{\lambda}^T \frac{\partial \mathbf{K}}{\partial x_k} \mathbf{U} = -p x_k^{p-1} \boldsymbol{\lambda}_k^T \mathbf{K}_k^0 \mathbf{U}_k \qquad \text{(not a self-adjoint problem)}$$

where  $\lambda$  satisfies the adjoint eqn.  $[\mathbf{K}(\mathbf{x}) + \mathbf{K}_s] \lambda = \mathbf{L}$ 

# A variation of the theme: compliant mechanisms





- Control of "hinge" dimension in «lumped» mechanisms
- Need for geometric non-linear models

# A variation of the theme: compliant mechanisms

- - > Design of «distributed» compliant mechanism (structures with no internal hinge)





constitutive tensor

- Nonconization and pariodic boundary conditions
- Homogenization and periodic boundary conditions



Unit strains are enforced at the boundaries of the cell to compute the so-called SIMP-based mutual energies:

$$(11 \rightarrow 1), \boldsymbol{\varepsilon}^{0(1)}$$

$$(22 \rightarrow 2) \boldsymbol{\varepsilon}^{0(2)}$$

$$(12 = 21 \rightarrow 3), \boldsymbol{\varepsilon}^{0(3)}$$
Unit strains in 2D

$$E_{ijkl}^{H} = \frac{1}{|Y|} \int_{Y} C_{pqrs} \varepsilon_{pq}^{A(ij)} \varepsilon_{rs}^{A(kl)} dY$$
$$E_{ij}^{H} = \frac{1}{|Y|} \sum_{e}^{N} x_{e}^{p} (\mathbf{U}_{e}^{A(i)})^{T} \mathbf{K}_{0e} \mathbf{U}_{e}^{A(j)}$$

 $\max E_{11}^H + E_{22}^H$  s.t.  $\frac{1}{V} \sum_N x_e V_e \le V_f$ 

Max bulk modulus

$$\min_{x_{min} < x_e < 1} f\left(E_{ij}^H(\mathbf{U}^{A(i)}, \mathbf{U}^{A(j)})\right)$$
  
s.t. 
$$\left[\sum_e^N x_e^p \mathbf{K}_{0e}\right] \mathbf{U}^{A(i)} = \mathbf{F}^{A(i)} \quad i = 1, 2, 3$$
$$f(E_{ij}^H) \le \overline{f}$$
$$\frac{1}{V} \sum_N x_e V_e \le V_f .$$

Max bulk modulus with limited shear modulus



$$\max E_{11}^H + E_{22}^H \text{ s.t. } E_{33}^H \le \overline{f}$$



 $\blacktriangleright$  Auxetic microstructures  $\rightarrow$  o.f. - $\epsilon_2/\epsilon_1$  , - $\epsilon_1/\epsilon_2$ 

$$\begin{cases} \min_{x_{min} < x_e < 1} E_{12}^H - \beta^{iter} (E_{11}^H + E_{22}^H) \\ \text{s.t.} \quad \frac{1}{V} \sum_N x_e V_e \le V_f . \qquad 0 < \beta < 1 \end{cases}$$

Highly nonconvex: many (local) solutions!



v\*=-0.33

v\*=-0.71



### Some remarks

- Topology optimization by distribution of isotropic material is a powerful numerical tool to perform conceptual design of structures, structural components and materials
- Key ingredients are: adoption of a "material model", formulation of a constrained minimization problem, implementation of ad hoc/general algorithms for the solution, iterative computation of obj. fun., constraints and sensitivities
- Numerical issues: instabilities, mesh dependence, non uniqueness of the solution, local/global minima
- Very easy implementation for basic problems (design for stiff structures / compliant mechanisms / periodic microstructures)
- Many advanced issues can be dealt with...



### Suggested readings

- Bendsøe MP, Sigmund O (2003) Topology optimization Theory, methods and applications. Springer, Berlin Beghini LL, Beghini A, Katz N, Baker WF, Paulino GH (2014) Connecting architecture and engineering through structural topology optimization. Eng Struct;59:716-726
- Bendsøe MP, Sigmund O (1999) Material interpolation schemes in topology optimization. Arch Appl Mech 69:635–654
- Andreassen E, Clausen A, Schevenels M, Lazarov BS, Sigmund O (2011) Efficient topology optimization in MATLAB using 88 lines of code. Struct Multidiscip Optim 43:1–16
- Sigmund O, Petersson J (1998) Numerical instabilities in topology optimization: a survey on procedures dealing with checkerboards, mesh-dependencies and local minima. Struct Optim 16(1): 68–75
- Svanberg K (1987) Method of moving asymptotes-a new method for structural optimization. Int J Numer Methods Eng 24(2):359–373
- Lógó J (2013) On the optimal layout of structures subjected to probabilistic or multiply loading. Struct Multidiscip Opt;48(6):1207-1212.
  - Pictures and numerical examples from authored and co-authored papers

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