

# **TOPOLOGY OPTIMIZATION**

**MAE 598 DESIGN OPTIMIZATION**

**MONISH DEV SUDHAKHAR**

**1220859588**

## Topology Optimization

### Problem Formulation:

A MBB is a classical problem in which The design domain, the boundary conditions, and the external load for the MBB beam are shown in Fig. 1 problem is to find the optimal material distribution, in terms of minimum compliance, with a constraint on the total amount of material.

### Approach:

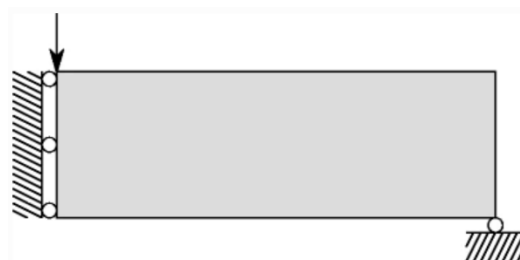
The design domain is discretized by square finite elements and a “density-based approach to topology optimization” is followed (Bendsøe [1989](#); Zhou and Rozvany [1991](#)); i.e. each element  $e$  is assigned a density  $x_e$  that determines its Young’s modulus  $E_e$ :

$$E_e(x_e) = E_{\min} + x_e p (E_0 - E_{\min}), x_e \in [0,1]$$

where  $E_0$  is the stiffness of the material,  $E_{\min}$  is a very small stiffness assigned to void regions in order to prevent the stiffness matrix from becoming singular, and  $p$  is a penalization factor (typically  $p = 3$ ) introduced to ensure black-and-white solutions.

### The compliance minimization problem

Topology optimization has been commonly used to design structures and materials with optimal mechanical, thermal, electromagnetic, acoustical, or other properties. The structure under design is segmented into  $n$  finite elements, and a density value  $x_i$  is assigned to each element  $i \in \{1, 2, \dots, n\}$ : A higher density corresponds to a less porous material element and higher Young's modulus. Reducing the density to zero is equivalent to creating a hole in the structure. Thus, the set of densities  $\mathbf{x} = \{x_i\}$  can be used to represent the topology of the structure and is considered as the variables to be optimized. A common topology optimization problem is compliance minimization, where we seek the "stiffest" structure within a certain volume limit to withstand a particular load:



$$\text{subject to: } \mathbf{h} := \mathbf{K}(\mathbf{x})\mathbf{d} = \mathbf{u},$$

$$\mathbf{g} := V(\mathbf{x}) \leq v,$$

$$\mathbf{x} \in [0, 1].$$

Here  $V(\mathbf{x})$  is the total volume;  $v$  is an upper bound on volume;  $\mathbf{d} \in \mathbb{R}^{n_d \times 1}$  is the displacement of the structure under the load  $\mathbf{u}$ , where  $n_d$  is the degrees of freedom (DOF)

of the system (i.e., the number of x- and y-coordinates of nodes from the finite element model of the structure);  $\mathbf{K}(\mathbf{x})$  is the global stiffness matrix for the structure.

$\mathbf{K}(\mathbf{x})$  is indirectly influenced by the topology  $\mathbf{x}$ , through the element-wise stiffness matrix

$$\mathbf{K}_i = \bar{\mathbf{K}}_e E(x_i),$$

$$\mathbf{K}(\mathbf{x}) = \mathbf{G}[\mathbf{K}_1, \mathbf{K}_2, \dots, \mathbf{K}_n],$$

where the matrix  $\bar{\mathbf{K}}_e$  is predefined according to the finite element type (we use first-order quadrilateral elements throughout this tutorial) and the nodal displacements of the element (we use square elements with unit lengths throughout this tutorial),  $\mathbf{G}$  is an assembly matrix,  $E(x_i)$  is the element-wise Young's modulus defined as a function of the density  $x_i$ :  $E(x_i) := \Delta E x_i^p + E_{\min}$ , where  $p$  (the penalty parameter) is usually set to 3. This cubic relationship between the topology and the modulus is determined by the material constitutive models, and numerically, it also helps binarize the topologies, i.e., to push the optimal  $x_i$  to 1 or 0 (why?). The term  $E_{\min}$  is added to provide numerical stability.

## Design sensitivity analysis

This problem has both inequality and equality constraints. However, the inequality ones are only related to the topology  $\mathbf{x}$ , and are either linear ( $V(\mathbf{x}) \leq v$ ) or simple bounds ( $x \in [0, 1]$ ). We will show that these constraints can be easily handled. The problem thus involves two sets of variables: We can consider  $\mathbf{x}$  as the **decision variables** and  $\mathbf{u}$  as the state variables that are governed by  $\mathbf{x}$  through the equality constraint  $\mathbf{K}(\mathbf{x})\mathbf{d} = \mathbf{u}$ .

The reduced gradient (often called design sensitivity) can be calculated as

$$\frac{df}{d\mathbf{x}} = \frac{\partial f}{\partial \mathbf{x}} - \frac{\partial f}{\partial \mathbf{u}} \left( \frac{\partial \mathbf{h}}{\partial \mathbf{u}} \right)^{-1} \frac{\partial \mathbf{h}}{\partial \mathbf{x}},$$

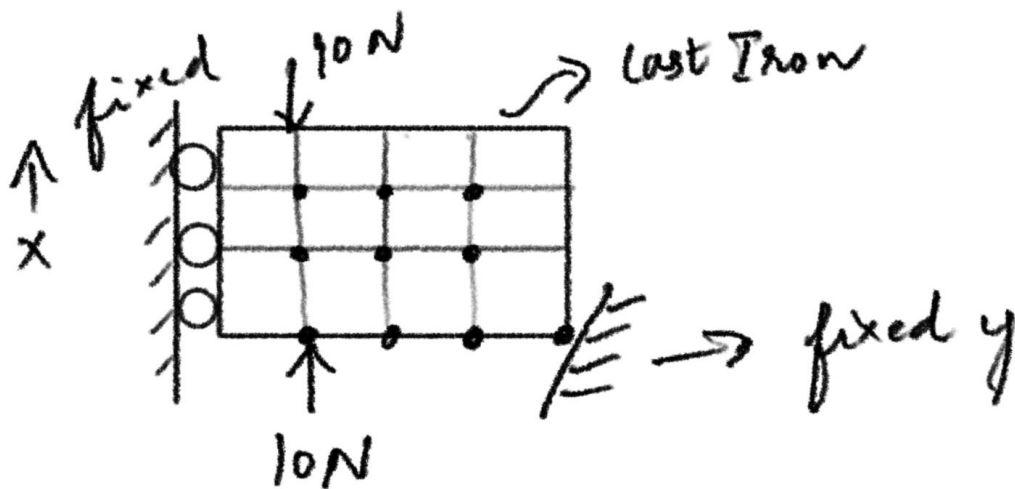
$$\frac{df}{d\mathbf{x}} = -\mathbf{u}^T \frac{\partial \mathbf{K}}{\partial \mathbf{x}} \mathbf{u}.$$

Recall the relation between  $\mathbf{K}$  and  $\mathbf{x}$ , and notice that

$$\mathbf{u}^T \mathbf{K} \mathbf{u} = \sum_{i=1}^n \mathbf{u}_i^T \mathbf{K}_i \mathbf{u}_i,$$

Problem Formulation:

The system is a beam with Cast iron as the material. It is subjected to a compressive load of 10 N. The material properties are changed with respect to the material and the boundary



## CODE IN PYTHON

```
# Monish Dev Sudhakhar
from __future__ import division
import numpy as np

from scipy.sparse import coo_matrix
from scipy.sparse.linalg import spsolve
from matplotlib import colors
import matplotlib.pyplot as plt
import cvxopt;
import cvxopt.cholmod

def main(nelx, nely, volfrac, penal, rmin, ft):
    print("Minimum compliance problem with OC")
    print("ndes: " + str(nelx) + " x " + str(nely))
    print("volfrac: " + str(volfrac) + ", rmin: " + str(rmin) + ", penal: " + str(penal))
    print("Filter method: " + ["Sensitivity based", "Density based"][ft])

    # Max and min stiffness (Cast iron is being considered)
    Emin = 1e-9
    Emax = 1.0
```

```

# dofs:
ndof = 2 * (nelx + 1) * (nely + 1)

# Allocate design variables (as array), initialize and allocate sens.
x = volfrac * np.ones(nely * nelx, dtype=float)
xold = x.copy()
xPhys = x.copy()

g = 0 # must be initialized to use the NGuyen/Paulino OC approach
dc = np.zeros((nely, nelx), dtype=float)

# FE: Build the index vectors for the for coo matrix format.
KE = lk()
edofMat = np.zeros((nelx * nely, 8), dtype=int)
for elx in range(nelx):
    for ely in range(nely):
        el = ely + elx * nely
        n1 = (nely + 1) * elx + ely
        n2 = (nely + 1) * (elx + 1) + ely
        edofMat[el, :] = np.array(
            [2 * n1 + 2, 2 * n1 + 3, 2 * n2 + 2, 2 * n2 + 3, 2 * n2, 2 * n2 + 1, 2 * n1, 2 * n1 + 1])
# Construct the index pointers for the coo format
iK = np.kron(edofMat, np.ones((8, 1))).flatten()
jK = np.kron(edofMat, np.ones((1, 8))).flatten()

# Filter: Build (and assemble) the index+data vectors for the coo matrix format
nfilter = int(nelx * nely * ((2 * (np.ceil(rmin) - 1) + 1) ** 2))
iH = np.zeros(nfilter)
jH = np.zeros(nfilter)
sH = np.zeros(nfilter)
cc = 0
for i in range(nelx):
    for j in range(nely):
        row = i * nely + j
        kk1 = int(np.maximum(i - (np.ceil(rmin) - 1), 0))
        kk2 = int(np.minimum(i + np.ceil(rmin), nelx))
        ll1 = int(np.maximum(j - (np.ceil(rmin) - 1), 0))
        ll2 = int(np.minimum(j + np.ceil(rmin), nely))
        for k in range(kk1, kk2):
            for l in range(ll1, ll2):
                col = k * nely + l
                fac = rmin - np.sqrt(((i - k) * (i - k) + (j - l) * (j - l)))
                iH[cc] = row
                jH[cc] = col
                sH[cc] = np.maximum(0.0, fac)
                cc = cc + 1
# Finalize assembly and convert to csc format
H = coo_matrix((sH, (iH, jH)), shape=(nelx * nely, nelx * nely)).tocsc()
Hs = H.sum(1)

# BC's and support
dofs = np.arange(2 * (nelx + 1) * (nely + 1))
fixed = np.union1d(dofs[0:2 * (nely + 1):2], np.array([2 * (nelx + 1) * (nely + 1) - 1]))
free = np.setdiff1d(dofs, fixed)

# Solution and RHS vectors
f = np.zeros((ndof, 1))
u = np.zeros((ndof, 1))

```

```

# Set load
f[1, 0] = - 10
# Initialize plot and plot the initial design
plt.ion() # Ensure that redrawing is possible
fig, ax = plt.subplots()
im = ax.imshow(-xPhys.reshape((nelx, nely)).T, cmap='gray', \
                interpolation='none', norm=colors.Normalize(vmin=-1, vmax=0))
fig.show()

loop = 0
change = 1
dv = np.ones(nely * nelx)
dc = np.ones(nely * nelx)
ce = np.ones(nely * nelx)
while change > 0.0001 and loop < 100:
    loop = loop + 1
    # Setup and solve FE problem
    sK = ((KE.flatten()[np.newaxis]).T * (Emin + (xPhys) ** penal * (Emax - Emin))).flatten(order='F')
    K = coo_matrix((sK, (iK, jK)), shape=(ndof, ndof)).tocsc()
    # Remove constrained dofs from matrix and convert to coo
    K = deleterowcol(K, fixed, fixed).tocoo()
    # Solve system
    K = cvxopt.spmatrix(K.data, K.row.astype(int), K.col.astype(int))
    B = cvxopt.matrix(f[free, 0])
    cvxopt.cholmod.linsolve(K, B)
    u[free, 0] = np.array(B)[:, 0]

    # Objective and sensitivity
    ce[:] = (np.dot(u[edofMat].reshape(nelx * nely, 8), KE) * u[edofMat].reshape(nelx * nely, 8)).sum(1)
    obj = ((Emin + xPhys ** penal * (Emax - Emin)) * ce).sum()
    dc[:] = (-penal * xPhys ** (penal - 1) * (Emax - Emin)) * ce

    dv[:] = np.ones(nely * nelx)
    # Sensitivity filtering:
    if ft == 0:
        dc[:] = np.asarray((H * (x * dc))[np.newaxis].T / Hs)[:, 0] / np.maximum(0.001, x)
    elif ft == 1:
        dc[:] = np.asarray(H * (dc[np.newaxis].T / Hs))[:, 0]
        dv[:] = np.asarray(H * (dv[np.newaxis].T / Hs))[:, 0]

    # Optimality criteria
    xold[:] = x
    (x[:, g] = oc(nelx, nely, x, volfrac, dc, dv, g)

    # Filter design variables
    if ft == 0:
        xPhys[:] = x
    elif ft == 1:
        xPhys[:] = np.asarray(H * x[np.newaxis].T / Hs)[:, 0]

    # Compute the change by the inf. norm
    change = np.linalg.norm(x.reshape(nelx * nely, 1) - xold.reshape(nelx * nely, 1), np.inf)

    # Plot to screen
    im.set_array(-xPhys.reshape((nelx, nely)).T)
    fig.canvas.draw()

    # Write iteration history to screen (req. Python 2.6 or newer)
    print("it.: {0} , obj.: {1:.3f} Vol.: {2:.3f}, ch.: {3:.3f}".format( \
        loop, obj, (g + volfrac * nelx * nely) / (nelx * nely), change))

```

```

# Make sure the plot stays and that the shell remains
plt.show()
input("Press any key...")

# element stiffness matrix (Titanium is being considered)
def lk():
    E = 9.6
    nu = 0.21
    k = np.array(
        [1 / 2 - nu / 6, 1 / 8 + nu / 8, -1 / 4 - nu / 12, -1 / 8 + 3 * nu / 8, -1 / 4 + nu / 12, -1 / 8 - nu / 8,
         nu / 6, 1 / 8 - 3 * nu / 8])
    KE = E / (1 - nu ** 2) * np.array([[k[0], k[1], k[2], k[3], k[4], k[5], k[6], k[7]],
                                       [k[1], k[0], k[7], k[6], k[5], k[4], k[3], k[2]],
                                       [k[2], k[7], k[0], k[5], k[6], k[3], k[4], k[1]],
                                       [k[3], k[6], k[5], k[0], k[7], k[2], k[1], k[4]],
                                       [k[4], k[5], k[6], k[7], k[0], k[1], k[2], k[3]],
                                       [k[5], k[4], k[3], k[2], k[1], k[0], k[7], k[6]],
                                       [k[6], k[3], k[4], k[1], k[2], k[7], k[0], k[5]],
                                       [k[7], k[2], k[1], k[4], k[3], k[6], k[5], k[0]]]);

    return (KE)

def oc(nelx, nely, x, volfrac, dc, dv, g):
    l1 = 0
    l2 = 1e9
    move = 0.2
    # reshape to perform vector operations
    xnew = np.zeros(nelx * nely)

    while (l2 - l1) / (l1 + l2) > 1e-3:
        lmid = 0.5 * (l2 + l1)
        xnew[:] = np.maximum(0.0,
                             np.maximum(x - move, np.minimum(1.0, np.minimum(x + move, x * np.sqrt(-dc / dv / lmid))))))
        gt = g + np.sum((dv * (xnew - x)))
        if gt > 0:
            l1 = lmid
        else:
            l2 = lmid
    return (xnew, gt)

def deleterowcol(A, delrow, delcol):
    # Assumes that matrix is in symmetric csc form !
    m = A.shape[0]
    keep = np.delete(np.arange(0, m), delrow)
    A = A[keep, :]
    keep = np.delete(np.arange(0, m), delcol)
    A = A[:, keep]
    return A

if __name__ == "__main__":
    # Default input parameters
    nelx = 190
    nely = 35
    volfrac = 0.39
    rmin = 3.5
    penal = 2.9

```

```
ft = 1 # ft==0 -> sens, ft==1 -> dens

import sys

if len(sys.argv) > 1: nelx = int(sys.argv[1])
if len(sys.argv) > 2: nely = int(sys.argv[2])
if len(sys.argv) > 3: volfrac = float(sys.argv[3])
if len(sys.argv) > 4: rmin = float(sys.argv[4])
if len(sys.argv) > 5: penal = float(sys.argv[5])
if len(sys.argv) > 6: ft = int(sys.argv[6])

main(nelx, nely, volfrac,
```

## **OPTIMIZATION RESULTS:**

C:\Anaconda\python.exe "C:/Users/Monish Dev  
Sudhakar/AppData/Roaming/JetBrains/PyCharmCE2021.2/scratches/scratch\_4.p  
y"

**Minimum compliance problem with OC**

**ndes: 190 x 35**

**volfrac: 0.39, rmin: 3.5, penal: 2.9**

**Filter method: Density based**

**it.: 1 , obj.: 106344.188 Vol.: 0.390, ch.: 0.200**

**it.: 2 , obj.: 48799.330 Vol.: 0.390, ch.: 0.200**

**it.: 3 , obj.: 31953.044 Vol.: 0.390, ch.: 0.200**

**it.: 4 , obj.: 24227.188 Vol.: 0.390, ch.: 0.200**

**it.: 5 , obj.: 21879.368 Vol.: 0.390, ch.: 0.200**

**it.: 6 , obj.: 21191.599 Vol.: 0.390, ch.: 0.200**

**it.: 7 , obj.: 20801.539 Vol.: 0.390, ch.: 0.124**

**it.: 8 , obj.: 20530.824 Vol.: 0.390, ch.: 0.195**

**it.: 9 , obj.: 20301.349 Vol.: 0.390, ch.: 0.115**

**it.: 10 , obj.: 20068.117 Vol.: 0.390, ch.: 0.168**



it.: 11 , obj.: 19818.626 Vol.: 0.390, ch.: 0.168  
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it.: 97 , obj.: 14127.980 Vol.: 0.390, ch.: 0.078

**it.: 98 , obj.: 14125.158 Vol.: 0.390, ch.: 0.089**

**it.: 99 , obj.: 14122.487 Vol.: 0.390, ch.: 0.100**

**it.: 100 , obj.: 14119.269 Vol.: 0.390, ch.: 0.111**

**Press any key...**

### **OBERVATION/ CONCLUSION :**

Hence the optimization for the problem formulation is done and the result are obtained.