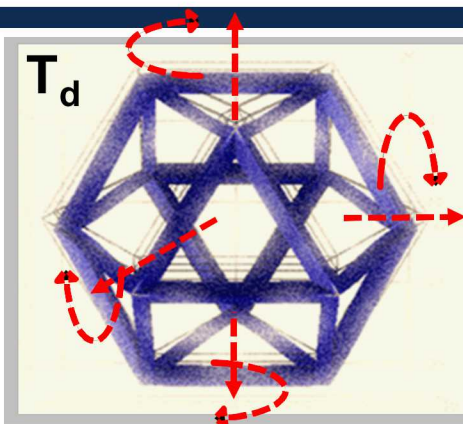
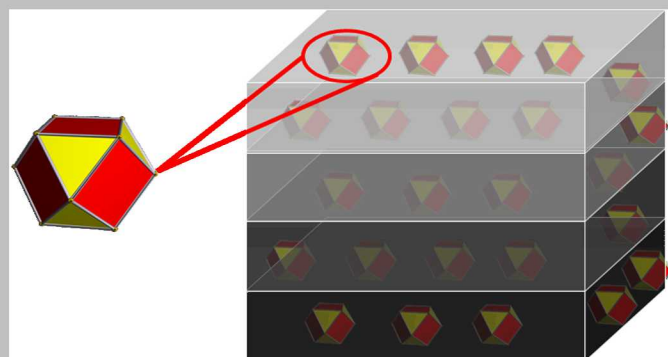


Group Theory Approach to Pentamode-Like Acoustic Metamaterials for Underwater Cloaking and Impedance Matching to Fluids

SAND2019-4414C



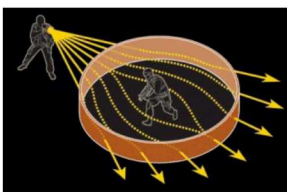
Character table for T_d point group							
	E	$8C_3$	$3C_2$	$6S_4$	$6\sigma_d$	linear, rotations	quadratic
A_1	1	1	1	1	1		$x^2+y^2+z^2$
A_2	1	1	1	-1	-1		
E	2	-1	2	0	0		$(2z^2-x^2-y^2, x^2-y^2)$
T_1	3	0	-1	1	-1	(R_x, R_y, R_z)	
T_2	3	0	-1	-1	1	(x, y, z)	(xy, xz, yz)

Ihab El-Kady & Charles M. Reinke

Applied Photonic Microsystems, Sandia National Laboratories, Albuquerque, NM, USA

M. Ghasemi Baboly & C.Z. Leseman

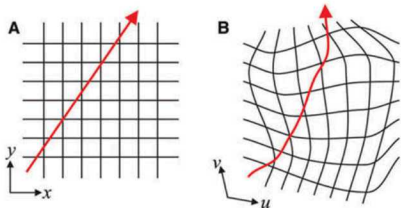
Department of Mechanical and Nuclear Engineering, Kansas State University, Manhattan KS, USA



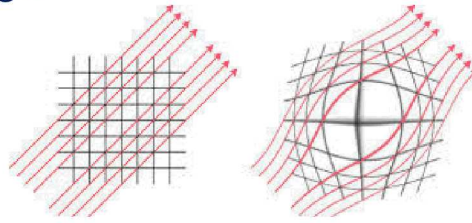
Cloaking in the EM Domain



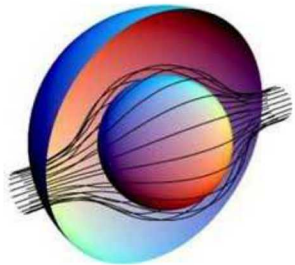
Transformation Optics^{*,1}



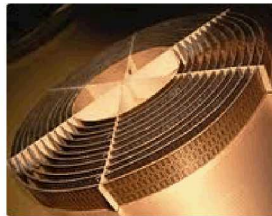
Coordinate Transformation



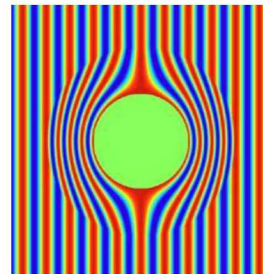
Transformation selected to open up a hole in space



3D Cloak Concept

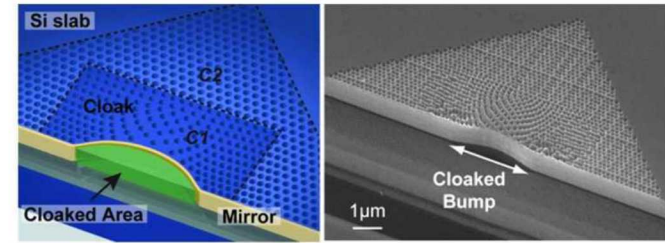


Duke's 2D RF Cloak

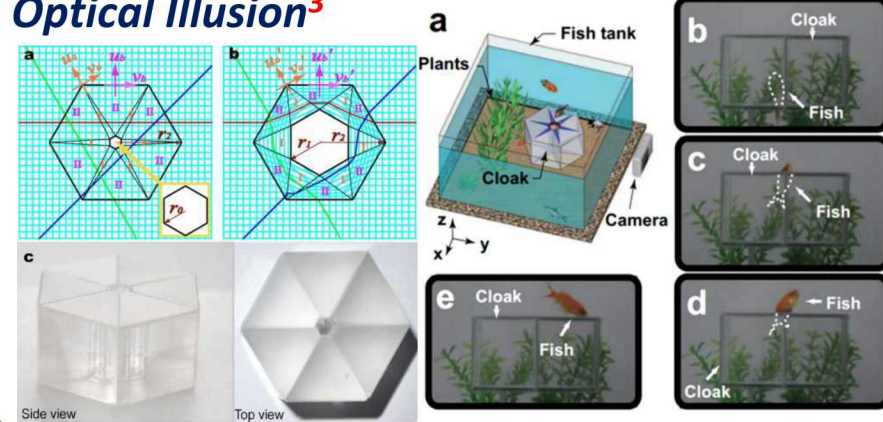


Cloak EM Signature

Diffraction Optics²

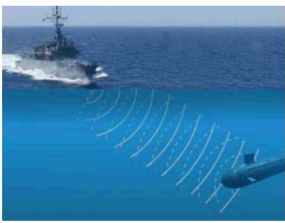


Optical Illusion³



Disadvantages:

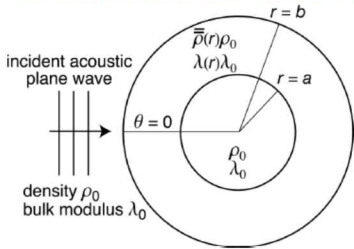
- Cloak's size
- Cloaked area $\sim \lambda^2$
- Blind inside out & outside in!



Cloaking in the Acoustic Domain

Acoustic Cloaking

Transformation Acoustics^{*,4}



3D Acoustic Cloaking Shell

$$\frac{r^2}{\rho_r} = \frac{(r-a)^2}{k_1},$$

$$\rho_\phi = k_1,$$

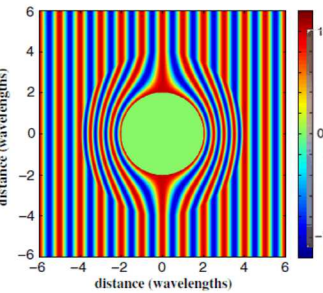
$$\frac{\rho_\phi}{\lambda} k_0^2 r^2 = k_{sh}^2 (r-a)^2,$$

$$\rho_\phi = \rho_\theta = \frac{b-a}{b},$$

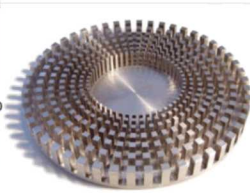
$$\rho_r = \frac{b-a}{b} \frac{r^2}{(r-a)^2},$$

$$\lambda = \frac{(b-a)^3}{b^3} \frac{r^2}{(r-a)^2}.$$

Coordinate Transformation

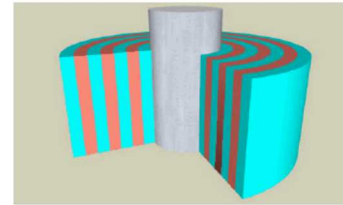


Pressure Field



MIT 2D Ultrasonic Cloak

Fluidic Acoustic Cloak⁵

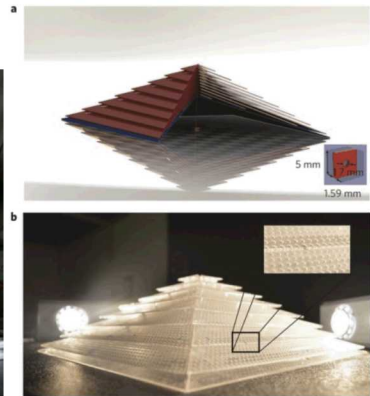


Layers with varying fluid density



3D rendering using fluid filled pipes!

Ground Acoustic Cloak⁶



Disadvantages:

- Cloak's size
- Cloaked area $\sim \lambda^2$
- Blind inside out & outside in!

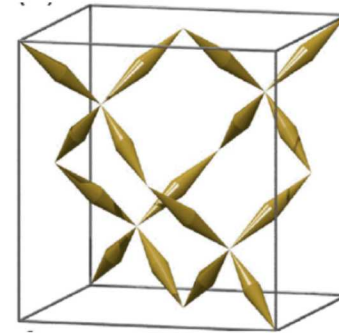
Pentamode Approach to Cloaking

Why is Acoustic Underwater cloaking difficult?

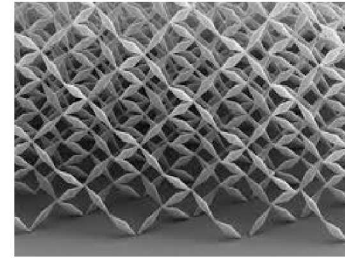
- Fluids are Shearless
- Solids support both shear and bulk waves
- The impedance mismatch will always create reflections!

Pentamode: The Sufficient but **NOT** Necessary Condition:

- 3D solid that behaves like a fluid.
- Finite bulk but vanishing shear modulus
- Hard to compress yet easy to deform
- Elasticity tensor with only one non-zero eigenvalue and five (penta) vanishing eigenvalues.



Pentamode Milton^{7,8}.



$$\tilde{\mathbf{C}}_{\text{Pentamode}} = \begin{bmatrix} C_{11} & 0 & 0 & \hat{0} \\ 0 & 0 & 0 & \\ 0 & 0 & 0 & \\ \hat{0} & & & \hat{0} \end{bmatrix}$$

The Necessary **AND** Sufficient Condition:

- Modal energy distribution is based on ratio of Bulk (**K**) to Shear (**G**) moduli
- Need to minimize **G/K**⁸ or ideally have **G/K** → 0

$$\tilde{\mathbf{C}}_{\text{Needed}} = \begin{bmatrix} K_{ij} & KG_{ij} \\ GK_{ij} & G_{ij} \end{bmatrix}$$

$$KG_{ij} \rightarrow 0$$

$$GK_{ij} \rightarrow 0$$

$$\& G_{ij}/K_{ij} \rightarrow 0$$

An Elastic Tensor by Design:

- **#1 Sigmund Numerical Search:** Space = boundary of hypercube ~10⁹ possibilities⁹.
- **#2 Milton logical deduction:** 1D → 2D → 3D; limited number of starting points → Severely restrict outcome → Requiring an infinitely small connection

A Metamaterial Approach

❖ **What is a Metamaterial?**

Engineered artificial materials that exhibit properties different and likely unattainable by constituent components

❖ **Context:**

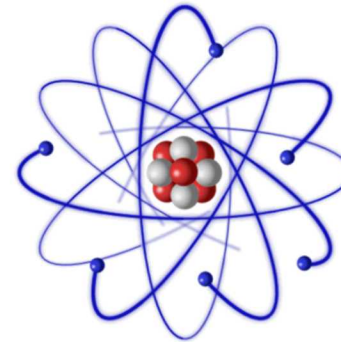
- Interested in the mechanical properties e.g. stiffness, and elastic modulus
- Think of two atoms connected by a spring (often nonlinear)

❖ **Stiffness:**

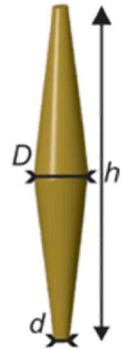
- Depends on the type of bond

❖ **Elastic Modulus:**

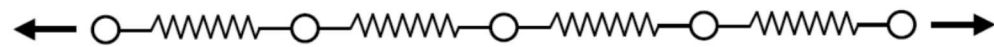
- Depends on the arrangement/packing of atoms



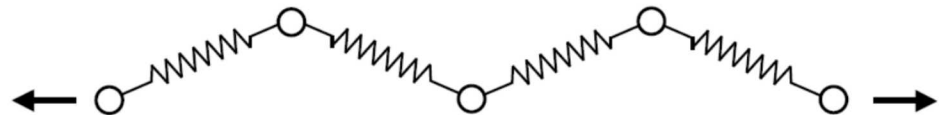
**Actual
Natural Atom**



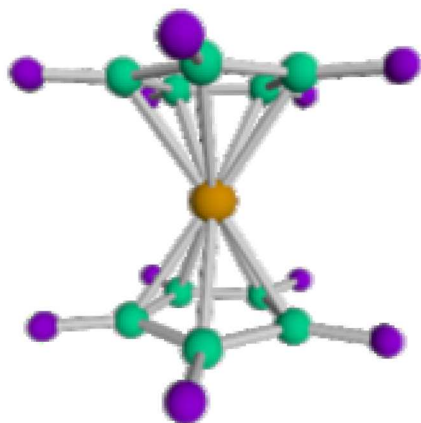
**Artificial
Metamaterial Atom**



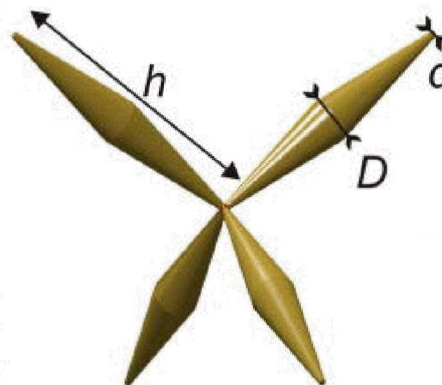
versus



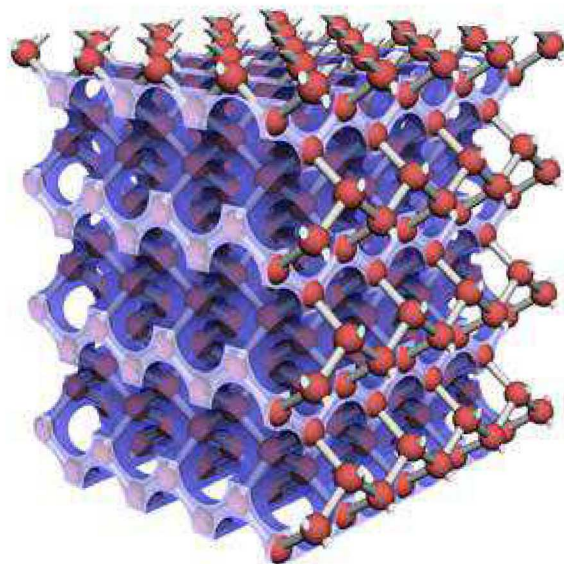
What is a Metamaterial?



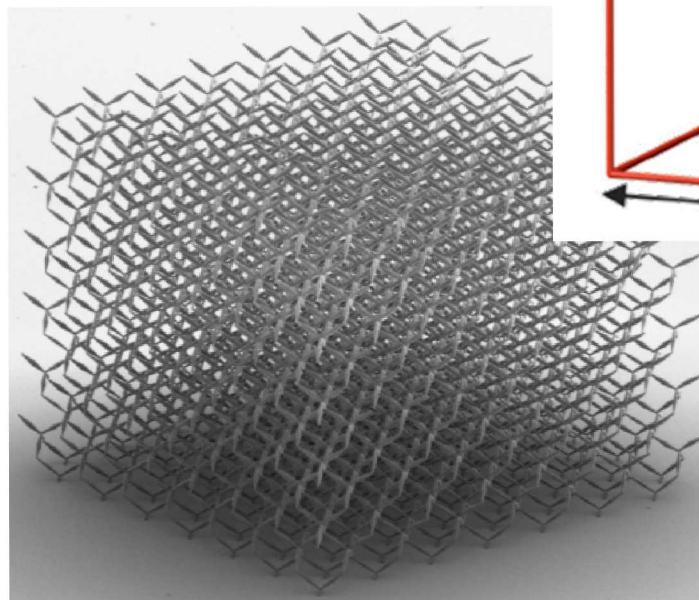
Natural Molecule



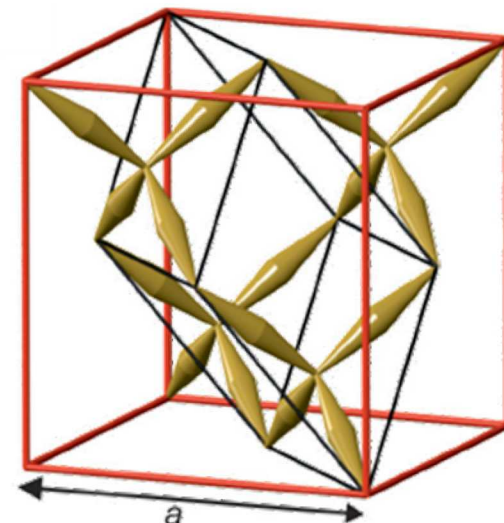
Metamaterial Molecule



Molecular Lattice



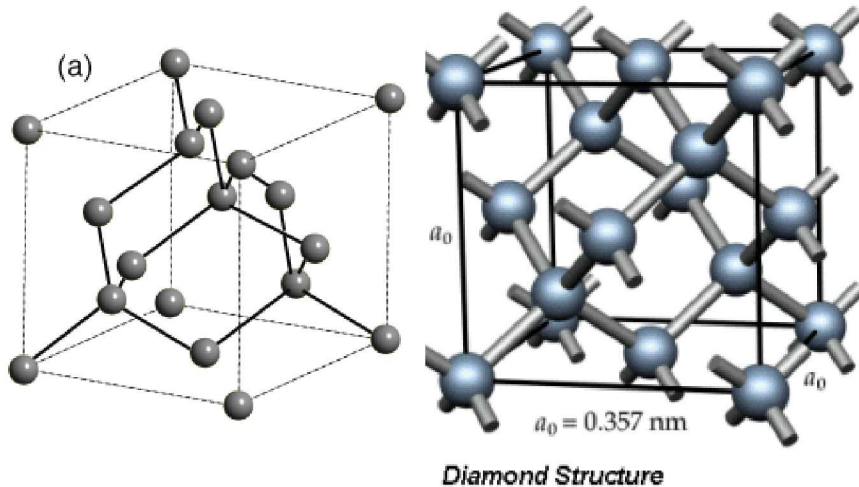
Metamaterial Lattice



Arrangement Matters!

Take Carbon atom as an example

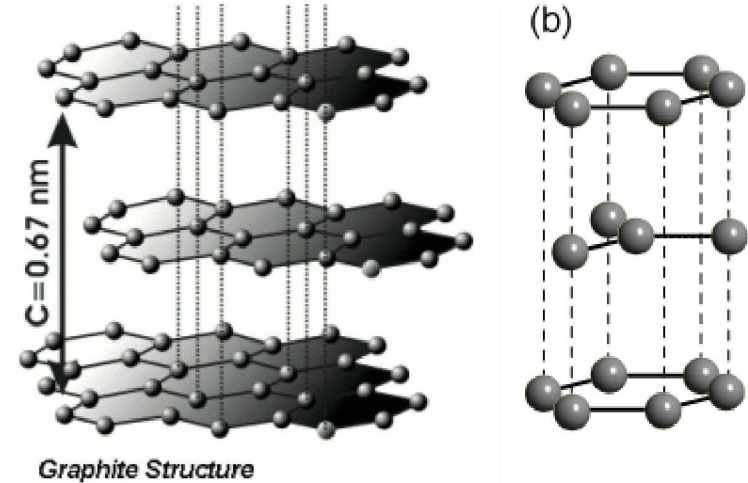
Diamond



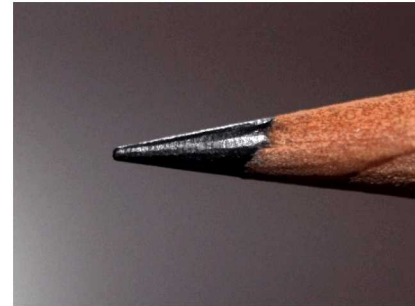
One of the **hardest** yet optically **transparent** materials known to man!



Graphite



One of the **softest** yet optically **opaque** materials known to man!



*In the language of metamaterials: “Not only does the **local resonance** (atom) matter, but also the **global resonance** and the **interactions** of the **adjacent cells**!”*

How do we arrange Metamaterial “atoms” to achieve a desired functionality?

❖ Why is this such a hard problem?

➤ Take the simplest MM element; Split Ring Resonator:

- Trial and Error?

- For a simple SRR on the 6 faces of a cube:

- In generally: $4^6 = 4096$ ways!
 - Invoking reciprocity: $4^3 = 256$ ways!
 - Invoking the Quasi-static limit: $4^{3/2} = 128$ ways!

➤ No Design rules of thumb! Physical intuition can only get you so far!

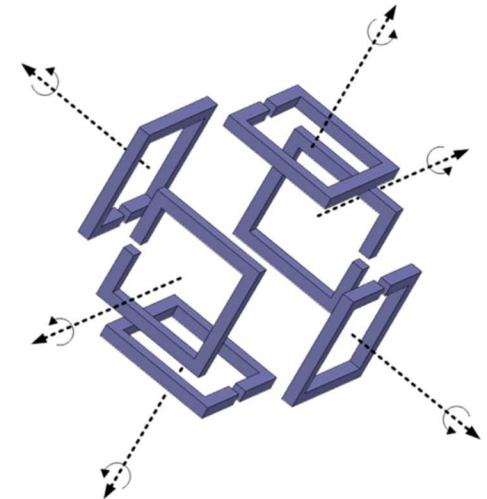
- Element-Element Interactions:

- Brute force EM simulations can only tell us the net result with very limited insight in to element-element interactions.

- Extrapolation:

- Starting with a well studied MM element (e.g. SRR) can we predict the outcome of a chosen topological arrangement?

❖ Turn back to Natural atoms, and ask the same question!



Question: How are natural atoms arranged to achieve a given property?

❖ *Bare in mind a few issues:*

- *Natural atoms assemble to give the **minimum energy configuration**, NOT necessarily our objective here!*
- ***NOT all properties** are awarded by naturally occurring materials! Entire field of MM arose in an attempt to achieve **unusual functionalities**!*
- ***Manybody Problem:***
 - *Equations of motion are almost impossible to solve.*
- *We are only primarily interested in **Mechanical properties***
 - *Basically **vibrations** and **rotations** in our artificial atoms*
- *Quantum Mechanics tricks → come to selection rules, “**solutionless approach**”:*
 - Symmetry of Wavefunction
 - Symmetry of the interaction Hamiltonian/potential
 - Symmetry of the boundary value problem

Code Word → Symmetry

Molecular Spectroscopy: Determining the properties of Natural Molecules

➤ Symmetry elements of a molecule constitute a complete group, “Point Group”

➤ Catalogue molecules into “58 Character” tables:

- Despite sharing **same symmetry**, molecules may have vastly **different** motional resonances!

➤ Generalized rotational-vibrational motion of a molecule:

- Can be viewed as a superposition of elemental/fundamental motions called “**normal modes**”.
- Use normal modes as basis → Symmetry elements are diagonal Matrices

➤ Good news:

- 1-to-1 analogy with a molecule’s spectroscopic activity
 - Active **Vibrational** modes → **Raman** active
 - Active **Rotational** modes → **IR** active

Nonaxial groups	<u>C₁</u>	<u>C_s</u>	<u>C_i</u>	-	-	-	-
C _n groups	<u>C₂</u>	<u>C₃</u>	<u>C₄</u>	<u>C₅</u>	<u>C₆</u>	<u>C₇</u>	<u>C₈</u>
D _n groups	<u>D₂</u>	<u>D₃</u>	<u>D₄</u>	<u>D₅</u>	<u>D₆</u>	<u>D₇</u>	<u>D₈</u>
C _{nv} groups	<u>C_{2v}</u>	<u>C_{3v}</u>	<u>C_{4v}</u>	<u>C_{5v}</u>	<u>C_{6v}</u>	<u>C_{7v}</u>	<u>C_{8v}</u>
C _{nh} groups	<u>C_{2h}</u>	<u>C_{3h}</u>	<u>C_{4h}</u>	<u>C_{5h}</u>	<u>C_{6h}</u>	-	-
D _{nh} groups	<u>D_{2h}</u>	<u>D_{3h}</u>	<u>D_{4h}</u>	<u>D_{5h}</u>	<u>D_{6h}</u>	<u>D_{7h}</u>	<u>D_{8h}</u>
D _{nd} groups	<u>D_{2d}</u>	<u>D_{3d}</u>	<u>D_{4d}</u>	<u>D_{5d}</u>	<u>D_{6d}</u>	<u>D_{7d}</u>	<u>D_{8d}</u>
S _n groups	<u>S₂</u>	<u>S₄</u>	<u>S₆</u>	<u>S₈</u>	<u>S₁₀</u>	<u>S₁₂</u>	-
Cubic groups	<u>T</u>	<u>T_h</u>	<u>T_d</u>	<u>O</u>	<u>O_h</u>	<u>I</u>	<u>I_h</u>
Linear groups	<u>C_{∞v}</u>	<u>D_{∞h}</u>	-	-	-	-	-

Application of Group Theory to Molecular Spectroscopy:

❖ Group Theory Terminology:

- **Reducible Representation Γ** : Most general form of vibration (current, mechanical, ..etc).
- **Irreducible Representations (IrredRep)**
 $A_1, A_2, B_1, B_2, \dots$: The normal modes classified and named based upon their symmetry
- **Characters χ** : The trace of the matrix representation of a symmetry op in the normal mode basis
- **L&Q Functions (X, Y, Z) and (R_x, R_y, R_z)**
Represent the behavior of the Irred. Rep. in a Cartesian basis.

Character Table:

Most compact form of symmetry Representation

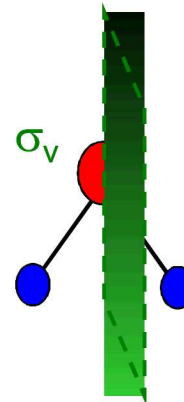
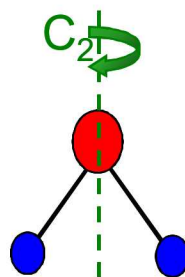
Point Group

Symmetry Elements/Operations

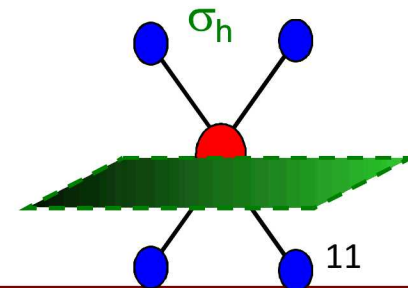
functions that transform as the various irreps of the group.

C_{2v}	E	C_2	$\sigma_v(xz)$	$\sigma_v(yz)$	Linear	Quadratic
A_1	1	1	1	1	z	x^2, y^2, z^2
A_2	1	1	-1	-1	R_z	xy
B_1	1	-1	1	-1	x, R_y	xz
B_2	1	-1	-1	1	y, R_x	yz

Irreducible Representations



Characters (trace of the matrix representation) of the group elements



Group Theory Mapping

❖ Mapping:

Bulk modes \leftrightarrow Generalized vibrations about structural joints
 \Rightarrow Transform like $\vec{r} = (x, y, z)$

Shear modes \leftrightarrow Generalized rotations about structural joints
 \Rightarrow Transform like axial vector $\vec{R} = (R_x, R_y, R_z)$

$$\begin{bmatrix} \sigma_1 \\ \sigma_2 \\ \sigma_3 \\ \sigma_4 \\ \sigma_5 \\ \sigma_6 \end{bmatrix} = \begin{bmatrix} \boxed{C_{11} \ C_{12} \ C_{13}} & C_{14} \ C_{15} \ C_{16} \\ C_{12} \ C_{22} \ C_{23} & C_{24} \ C_{25} \ C_{26} \\ C_{13} \ C_{23} \ C_{33} & C_{34} \ C_{35} \ C_{36} \\ C_{14} \ C_{24} \ C_{34} & \boxed{C_{44} \ C_{45} \ C_{46}} \\ C_{15} \ C_{25} \ C_{35} & C_{45} \ C_{55} \ C_{56} \\ C_{16} \ C_{26} \ C_{36} & C_{46} \ C_{56} \ C_{66} \end{bmatrix} \begin{bmatrix} \epsilon_1 \\ \epsilon_2 \\ \epsilon_3 \\ \epsilon_4 \\ \epsilon_5 \\ \epsilon_6 \end{bmatrix}$$

Stress **Strain**

❖ Quadrant Access:

Coupled Linear Oscillations
 (r_i, r_j)

$\bar{C} =$

C_{11}	C_{12}	C_{13}	C_{14}	C_{15}	C_{16}
C_{21}	C_{22}	C_{23}	C_{24}	C_{25}	C_{26}
C_{31}	C_{32}	C_{33}	C_{34}	C_{35}	C_{36}
C_{41}	C_{42}	C_{43}	C_{44}	C_{45}	C_{46}
C_{51}	C_{52}	C_{53}	C_{54}	C_{55}	C_{56}
C_{61}	C_{62}	C_{63}	C_{64}	C_{65}	C_{66}

r_i, R_j (R_i, R_j)

Coupled Linear/axial Motions

Independent Linear Oscillations $r_i's$

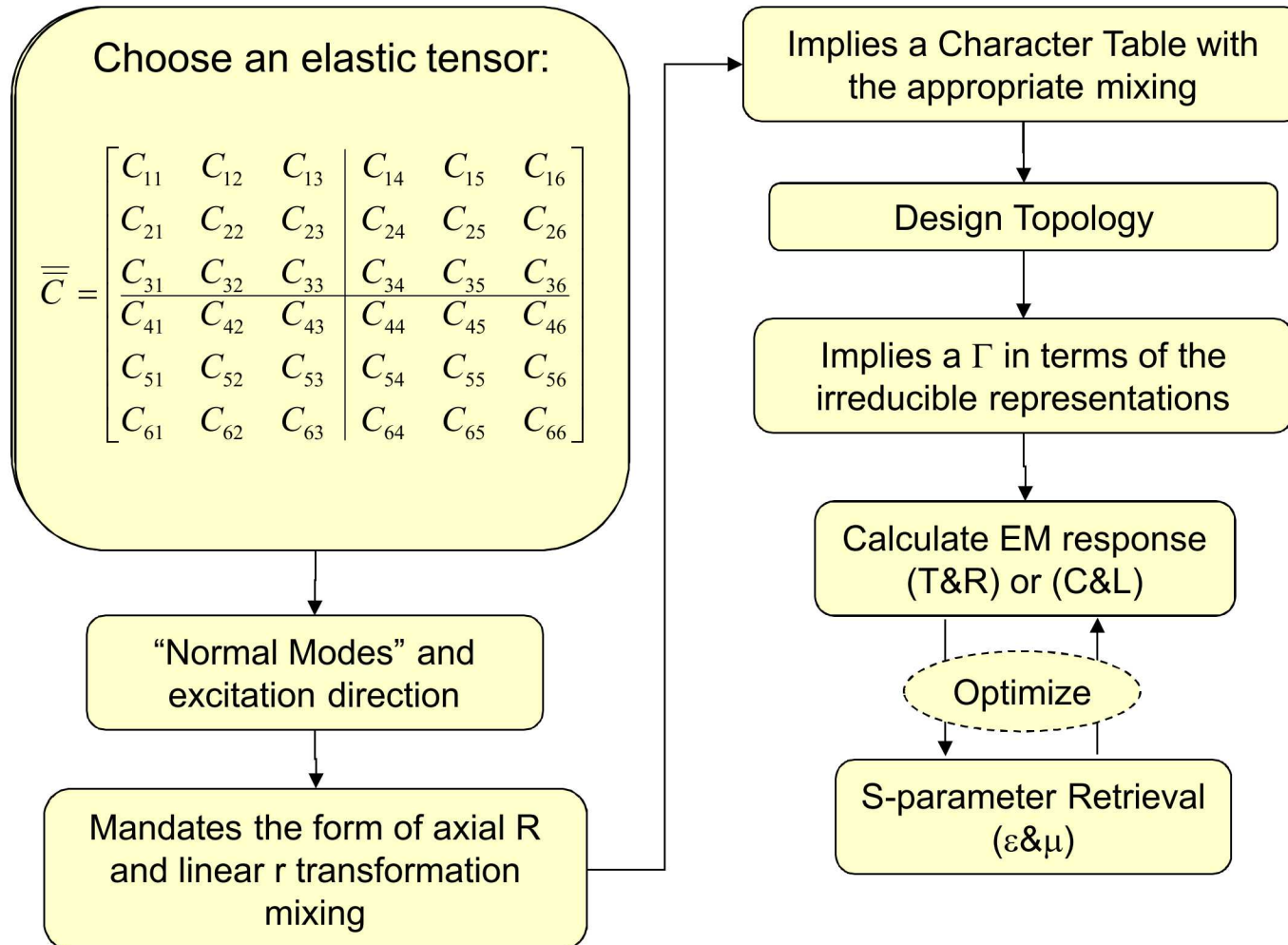
$\bar{C} =$

C_{11}	0	0	0	0	0
0	C_{22}	0	0	0	0
0	0	C_{33}	0	0	0
0	0	0	C_{44}	0	0
0	0	0	0	C_{55}	0
0	0	0	0	0	C_{66}

Independent Axial motion $R_i's$

Group Theory Inverse Problem Approach

Engineering the MM Constitutive Tensor



Group Theory Approach Pentamode

Practically what we require is:

- Diagonal elasticity tensor
- Optimize such that all diagonal elements are small relative to C_{11}

Linear and axial mixing:

- Need a block diagonal tensor with no shear/bulk coupling
- (x, y, z) & (R_x, R_y, R_z)

Cavite:

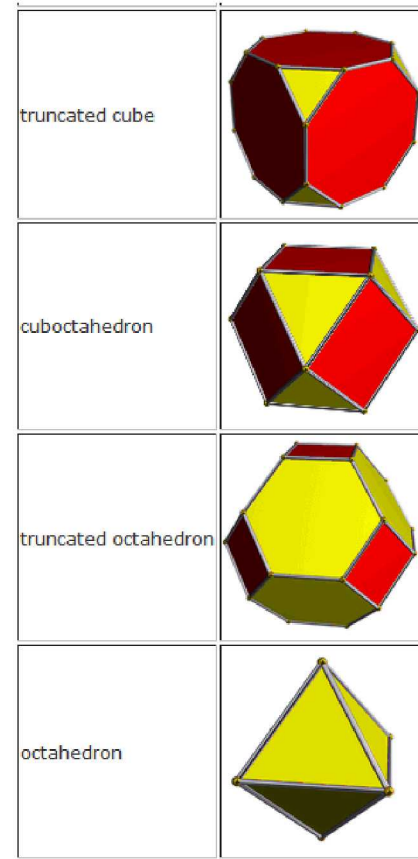
GpTh only tells us about the symmetry but nothing about the structural or material parameters!

Character table for T_d point group

	E	$8C_3$	$3C_2$	$6S_4$	$6\sigma_d$	linear, rotations	quadratic
A_1	1	1	1	1	1		$x^2+y^2+z^2$
A_2	1	1	1	-1	-1		
E	2	-1	2	0	0		$(2z^2-x^2-y^2, x^2-y^2)$
T_1	3	0	-1	1	-1	(R_x, R_y, R_z)	
T_2	3	0	-1	-1	1	(x, y, z)	(xy, xz, yz)

Character Table for O_h point group

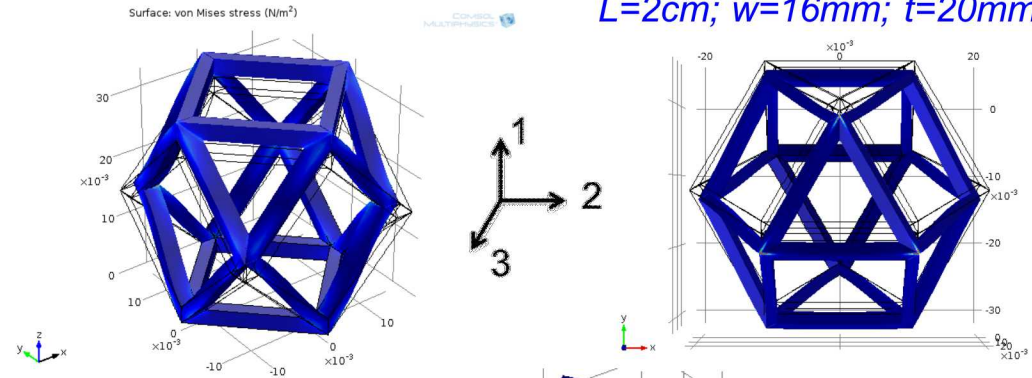
O_h	E	$8C_3$	$6C_2$	$6C_4$	$3C_2=(C_4)^2$	i	$6S_4$	$8S_6$	$3\sigma_h$	$6\sigma_d$	linear functions, rotations
A_{1g}	+1	+1	+1	+1	+1	+1	+1	+1	+1	+1	-
A_{2g}	+1	+1	-1	-1	+1	+1	-1	+1	+1	-1	-
E_g	+2	-1	0	0	+2	+2	0	-1	+2	0	-
T_{1g}	+3	0	-1	+1	-1	+3	+1	0	-1	-1	(R_x, R_y, R_z)
T_{2g}	+3	0	+1	-1	-1	+3	-1	0	-1	+1	-
A_{1u}	+1	+1	+1	+1	+1	-1	-1	-1	-1	-1	-
A_{2u}	+1	+1	-1	-1	+1	-1	+1	-1	-1	+1	-
E_u	+2	-1	0	0	+2	-2	0	+1	-2	0	-
T_{1u}	+3	0	-1	+1	-1	-3	-1	0	+1	+1	(x, y, z)
T_{2u}	+3	0	+1	-1	-1	-3	+1	0	+1	-1	-



Modeling & Results

Results:

$$\tilde{C} = \begin{bmatrix} 1 & 0.0821 & 0.003 & \sim 0 & \sim 0 & \sim 0 \\ 0.01 & 0.1 & 0.007 & \sim 0 & \sim 0 & \sim 0 \\ 0.067 & 0.006 & 0.008 & \sim 0 & \sim 0 & \sim 0 \\ \sim 0 & \sim 0 & \sim 0 & 0.009 & \sim 0 & \sim 0 \\ \sim 0 & \sim 0 & \sim 0 & \sim 0 & 0.0098 & \sim 0 \\ \sim 0 & \sim 0 & \sim 0 & \sim 0 & \sim 0 & 0.002 \end{bmatrix}$$



Calculating G/K:

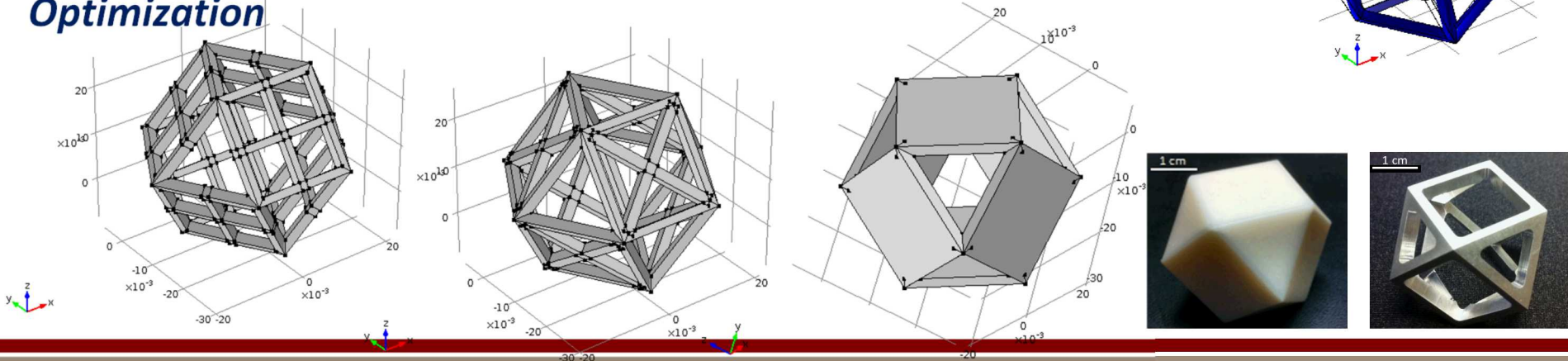
1. Calculated the volumetric strain on each side. ($\Delta V/V$)
2. Fixed the same boundary on the opposite side and calculated the force and displacement on those boundaries.

$$K = \frac{\frac{1}{3}(f_{11} + f_{22} + f_{33})}{\frac{\Delta V}{V}}$$

$$G/K \sim 7.14 \times 10^{-4}$$

$$\epsilon_{13} = \frac{\text{displacement}}{L} = \frac{\Delta L}{L}; \quad G = \frac{f_{13}}{\epsilon_{13}}$$

Optimization



Summary and Conclusions

- a. Introduced the Group theory as an inverse design methodology to engineered metamaterials¹⁰
- b. Used symmetry arguments to engineer an elastic tensor with the desired format
- c. Used modeling to optimize the bulk to shear modulus ratio
- d. Were able to achieve Pentamode-like behavior with $G/K \sim 10^{-4}$

