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# CS 519: Scientific Visualization

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## Working with Data: Representation and Interpolation

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# Today...

- ▣ Visualization of the day
  - ▣ <https://www.youtube.com/watch?v=EgumU0Ns1YI>
- ▣ Data representation
- ▣ (Linear) Interpolation

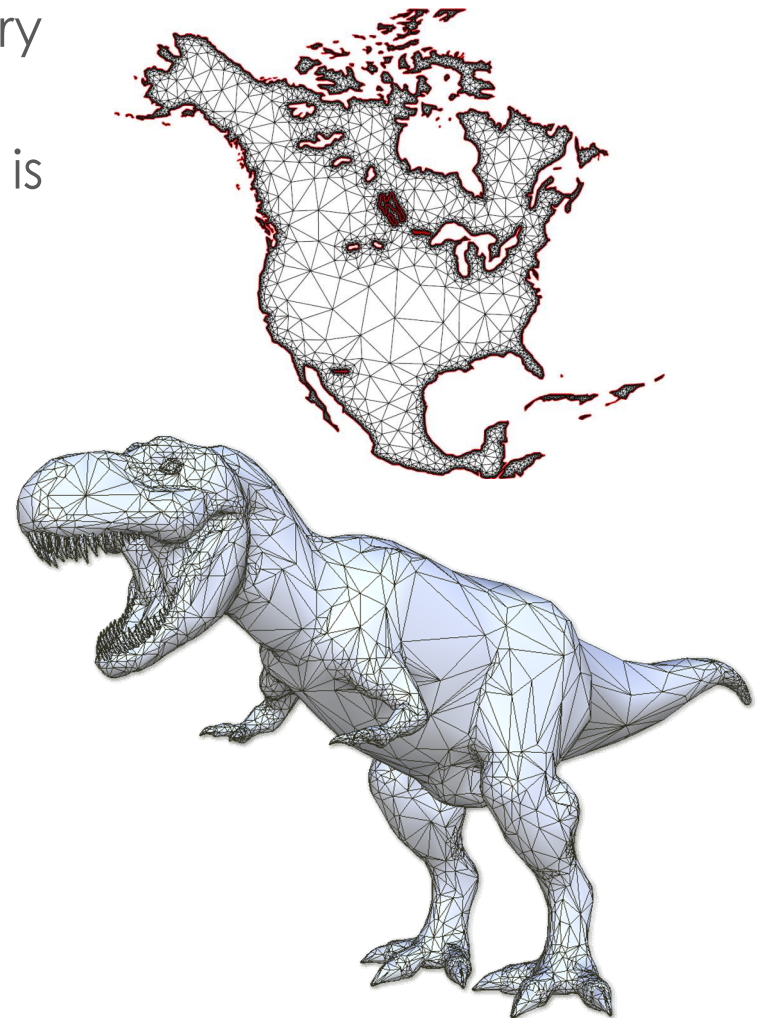
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# Things for later

- ▣ These are mentioned in Chapter 3...but we'll get to them later
    - ▣ Color
    - ▣ Advanced spatial data structures
    - ▣ Tensors
    - ▣ Calculating derivatives
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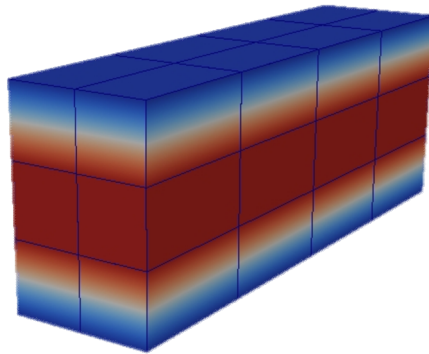
# Data Representation

- ❑ Disclaimer: add “typically” before every sentence
- ❑ “Typically” Scientific Visualization data is discretely sampled
  - ▣ The original function is continuous
- ❑ The Domain over which the data is sampled is discretized
  - ▣ In 2D using a polygonal mesh of cells
    - ▣ This includes data sampled on a 2D surface embedded in 3-space
  - ▣ In 3D using a polyhedral mesh of cells
- ❑ Data values are either vertex or cell-centered



# Interpolation

- ❑ Interpolation is a mathematical process for filling in missing data
- ❑ An interpolating function approximates some sampled function
  - ❑ Approximation matches original function value at the sample points
- ❑ Useful in visualization
  - ❑ We usually don't have original function values at each pixel.



# How Can We Organize Data Sampled in Euclidean Space ?

- ▣ Chop the space up (discretize it) into **cells**
- ▣ A structure of cells is called a **grid** or **mesh**
- ▣ Lots of cell types are possible
  - ▣ **0D** point
  - ▣ **1D** line
  - ▣ **2D** triangle, quad, rectangle
  - ▣ **3D** tetrahedron, parallelepiped, box, pyramid, prism,

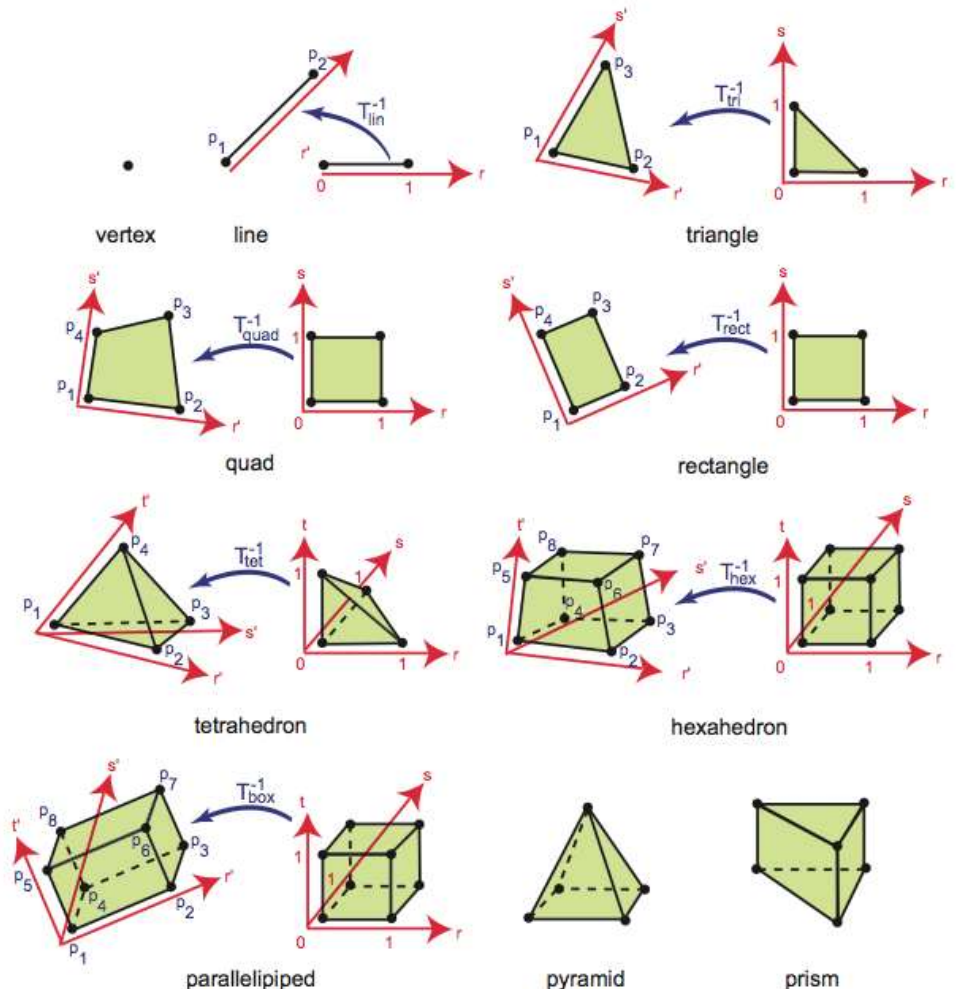


Figure 3.5. Cell types in world and reference coordinate systems.

# Grids (or Meshes)

## Cells

- provide interpolation over a small, simple-shaped spatial region

## Grids (or meshes)

- partition our complex data domain  $D$  into cells
- allow applying per-cell interpolation (as described so far)

Given a domain  $D$ ...

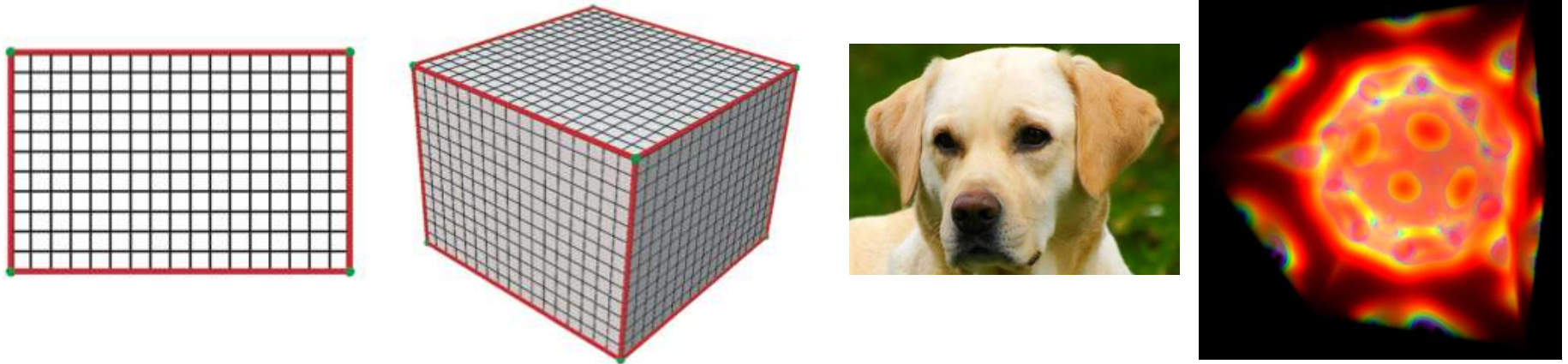
A **grid**  $G = \{c_i\}$  is a set of cells such that

$c_i \cap c_j = \emptyset, \forall i \neq j$       no two cells overlap

$\bigcup_i c_i = D$       the cells cover all our domain

The dimension of the domain  $D$  constrains which cell types we can use

# Uniform Grids



**Figure 3.7.** Uniform grids. 2D rectangular domain (left) and 3D box domain (right).

image

volume

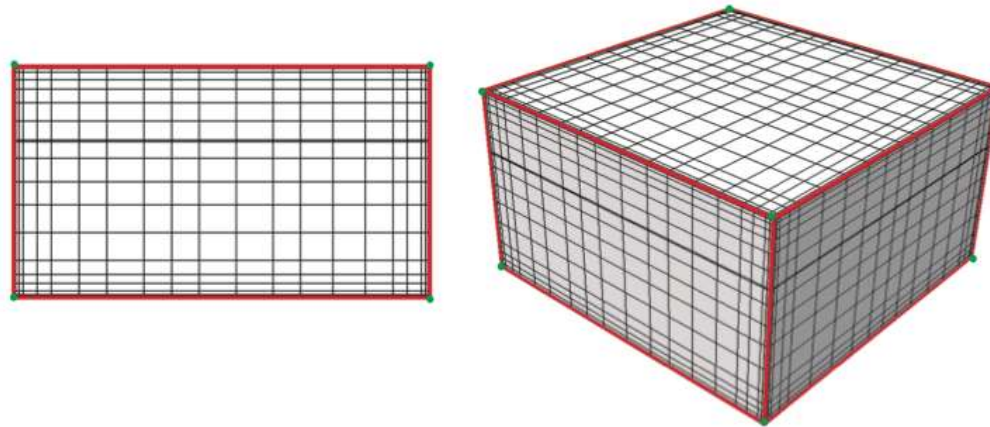
- all cells have identical size and type (typically, square or cubic)

## Storage requirements for the structure (not the data)

- $m$  integers for the #cells along each of the  $m$  dimensions of  $D$  (e.g.  $m=2$  or  $3$ )
- two corner points



# Rectilinear Grids



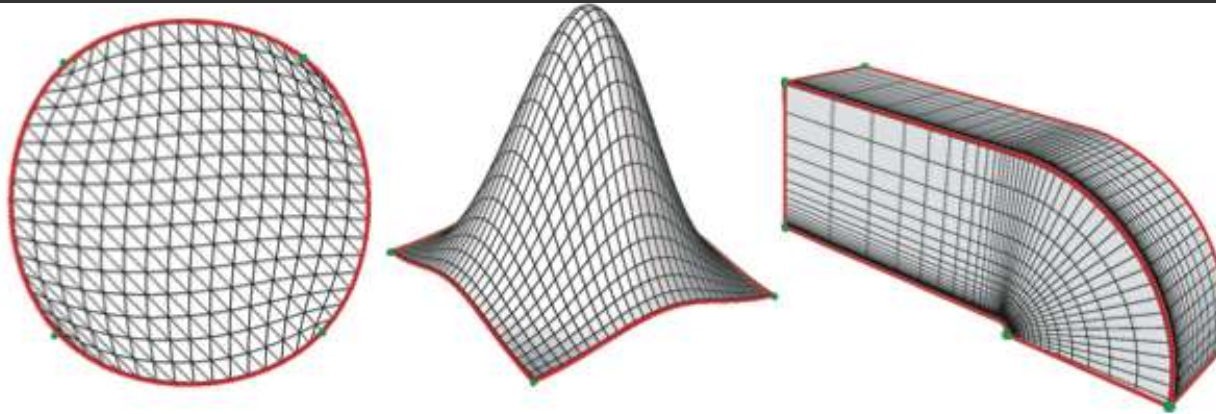
**Figure 3.8.** Rectilinear grids. 2D rectangular domain (left) and 3D box domain (right).

- all cells have same type
- cells can have different sizes but share them along axes
- Book says → “cannot model non-axis-aligned domains”. Is that true?

## Storage requirements for the structure

$\sum_{i=1}^m d_i$  floats (coordinates of vertices along each of the  $m$  axes of  $D$ )  
And what else?

# Structured Grids



**Figure 3.9.** Structured grids. Circular domain (left), curved surface (middle), and 3D volume (right). Structured grid edges and corners are drawn in red and green, respectively.

- all cells have same type
- cell vertex coordinates are freely (explicitly) specifiable...
- ...as long as cells assemble in a matrix-like structure
- can approximate more complex shapes than rectilinear/uniform grids

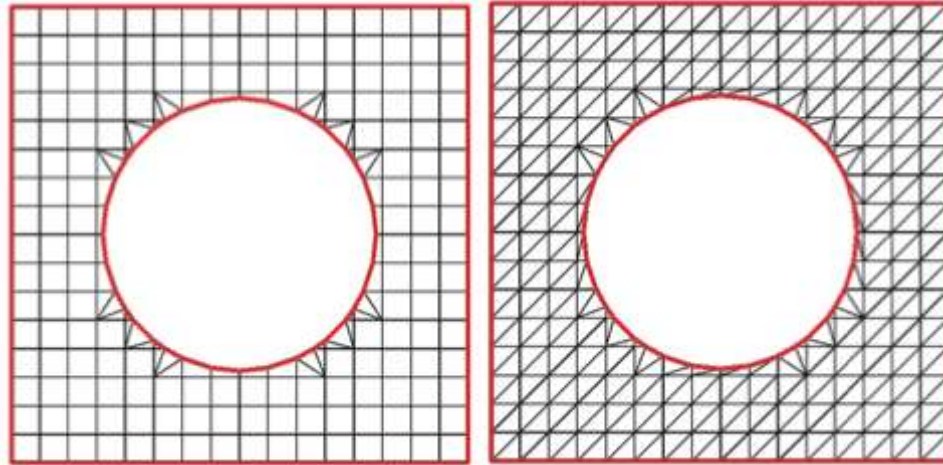
## Storage requirements

$\prod_{i=1}^m d_i$  floats (coordinates of all vertices)

And what else?

# Unstructured Grids

Consider the domain  $D$ : a square with a hole in the middle

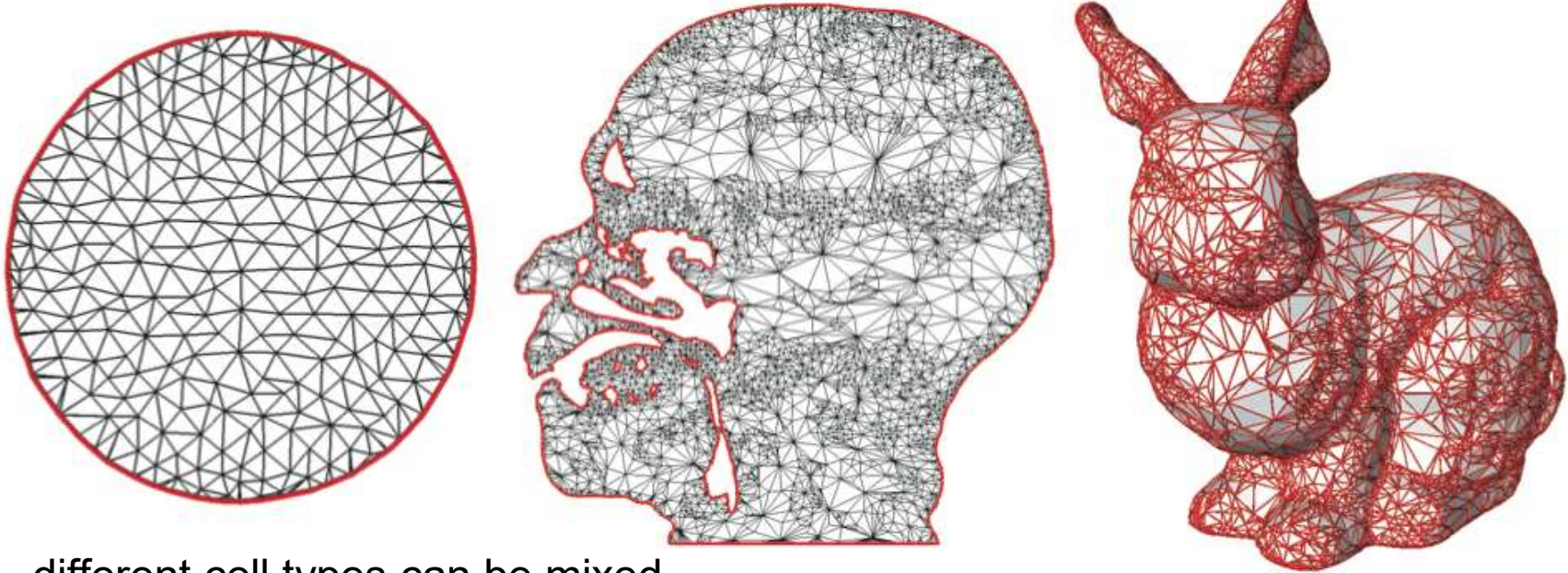


**According to book:** We cannot cover such a domain with a structured grid (why?)

- it's not of genus 0, so cannot be covered with a matrix-like distribution of cells
- Or could it?
- What about genus 2?
- BTW, genus means how many holes there are in the domain.

For more generality, we need **unstructured grids**

# Unstructured Grids



- different cell types can be mixed
- both vertex coordinates and cell themselves are freely (explicitly) specifiable
- implementation
  - vertex set  $V = \{v_i\}$
  - cell set  $C = \{c_i = (\text{indices of vertices in } V)\}$
- most flexible, but most complex/expensive grid type

## Storage requirements

$m\|V\| + s\|C\|$  for a  $m$ -dimensional grid with cells having  $s$  vertices each  
What operation is hard to do with just this information?

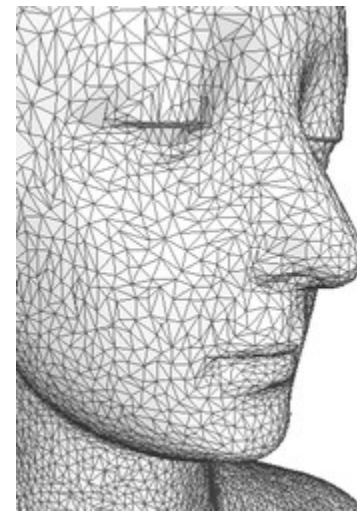
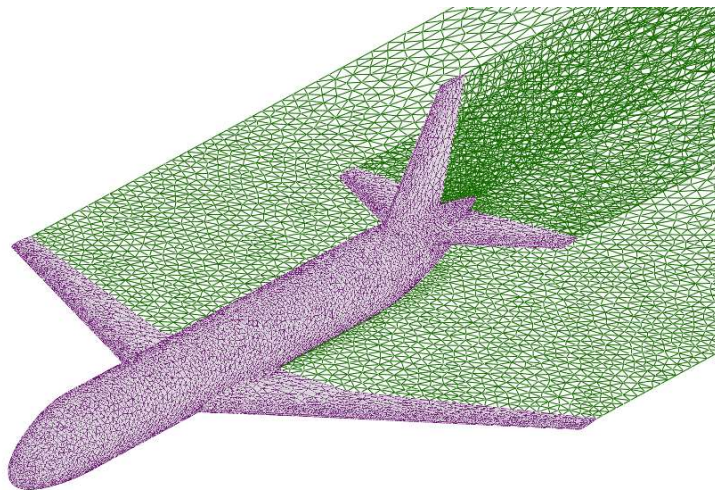


# Example: Unstructured Surface Mesh

This Wavefront/OBJ file describes a single triangle.

Could add more vertices and faces for a bigger mesh

```
# Simple Wavefront file
v 0.0 0.0 0.0
v 0.0 1.0 0.0
v 1.0 0.0 0.0
f 1 2 3
```



# Data Attributes

$$f: \mathbf{R}^m \rightarrow \mathbf{R}^n$$

- $n=0$  no attributes (we model a shape only e.g. a surface)
- $n=1$  scalars (e.g. temperature, pressure, curvature, density)
- $n=2$  2D vectors
- $n=3$  3D vectors (e.g. velocity, gradients, normals, colors)
- $n=6$  symmetric tensors (e.g. diffusion, stress/strain)
- $n=9$  asymmetric general tensors (not very common)

## Remarks

- an attribute is usually specified for **all** sample points in a dataset
- each attribute is interpolated separately
- different visualization methods for each  $n$

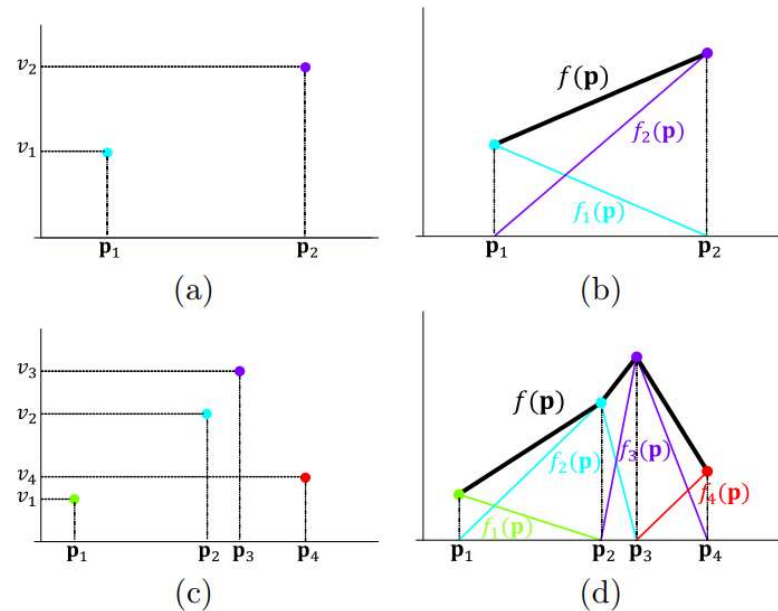
# A Simple Approach to Interpolation

- Suppose we have sampled values at two points  $P_1$  and  $P_2$  and want to guess at the function values on the line between the points.

- We can perform linear interpolation or LERP

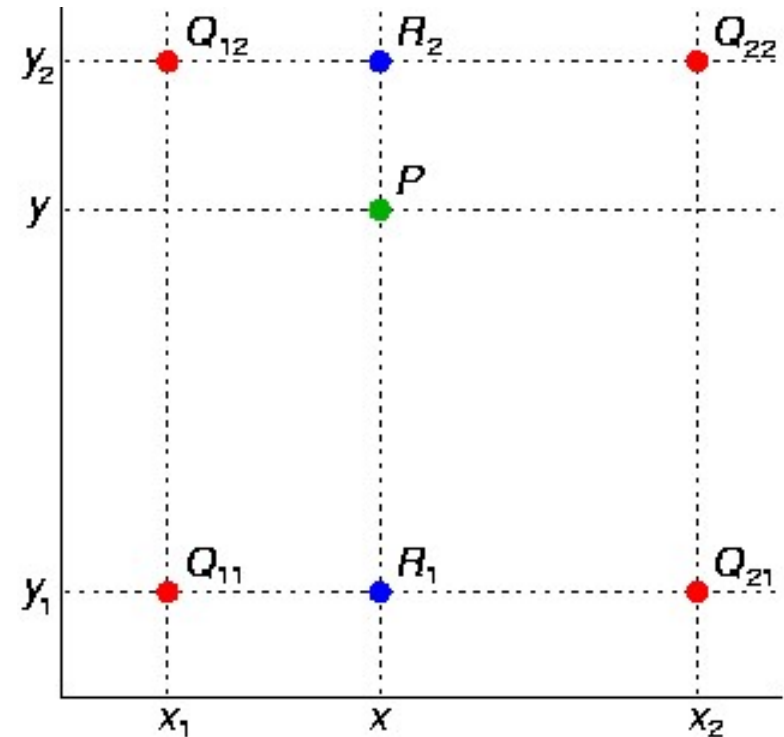
$$f(\mathbf{p}) = \frac{\mathbf{p}_2 - \mathbf{p}}{\mathbf{p}_2 - \mathbf{p}_1} v_1 + \frac{\mathbf{p} - \mathbf{p}_1}{\mathbf{p}_2 - \mathbf{p}_1} v_2$$

- When might this be a poor choice?
- Also, what dimension is this domain?



# Bilinear Interpolation

- If we have a function defined on a 2D domain we need to do more
- You've already seen it ...but repetition can be helpful
- Assume we know a function value at the four points  
 $Q_{11} = (x_1, y_1)$ ,  $Q_{12} = (x_1, y_2)$ ,  
 $Q_{21} = (x_2, y_1)$ ,  $Q_{22} = (x_2, y_2)$
- We first do linear interpolation in the x-direction
- ...and then in the y direction





# Bilinear Interpolation

$$f(R_1) \approx \frac{x_2 - x}{x_2 - x_1} f(Q_{11}) + \frac{x - x_1}{x_2 - x_1} f(Q_{21})$$

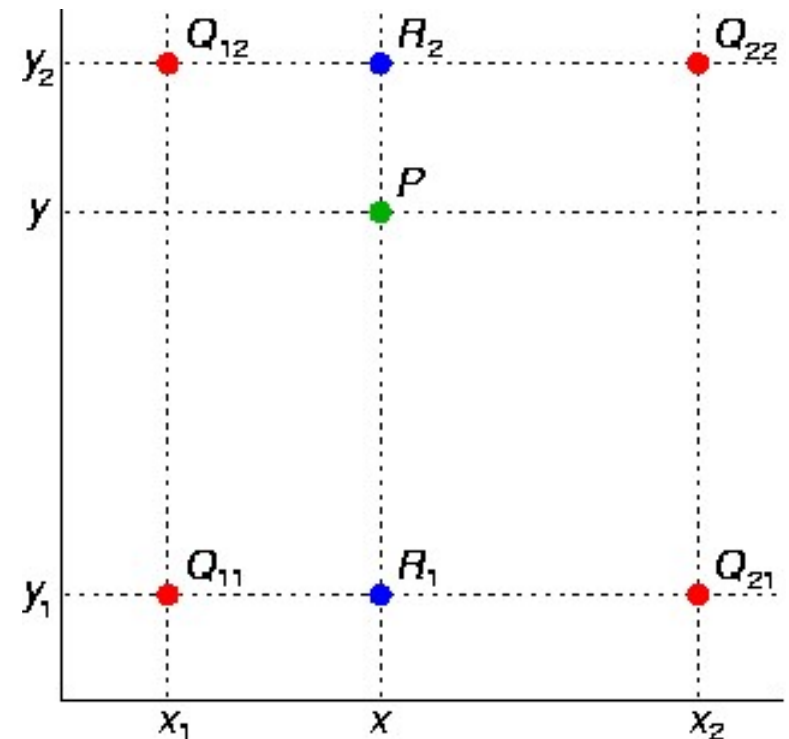
where  $R_1 = (x, y_1)$ ,

$$f(R_2) \approx \frac{x_2 - x}{x_2 - x_1} f(Q_{12}) + \frac{x - x_1}{x_2 - x_1} f(Q_{22})$$

where  $R_2 = (x, y_2)$ .

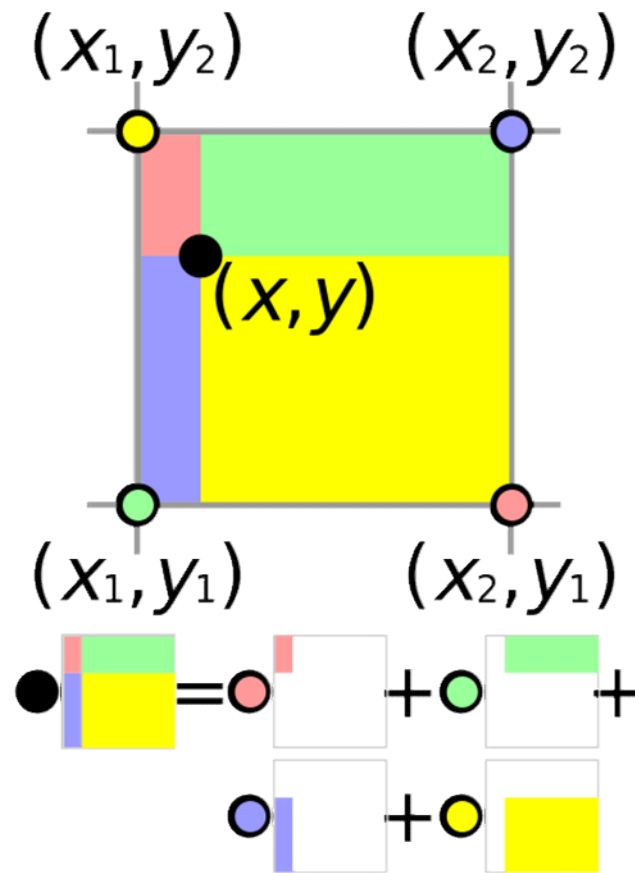
We proceed by interpolating in the  $y$ -direction.

$$f(P) \approx \frac{y_2 - y}{y_2 - y_1} f(R_1) + \frac{y - y_1}{y_2 - y_1} f(R_2).$$



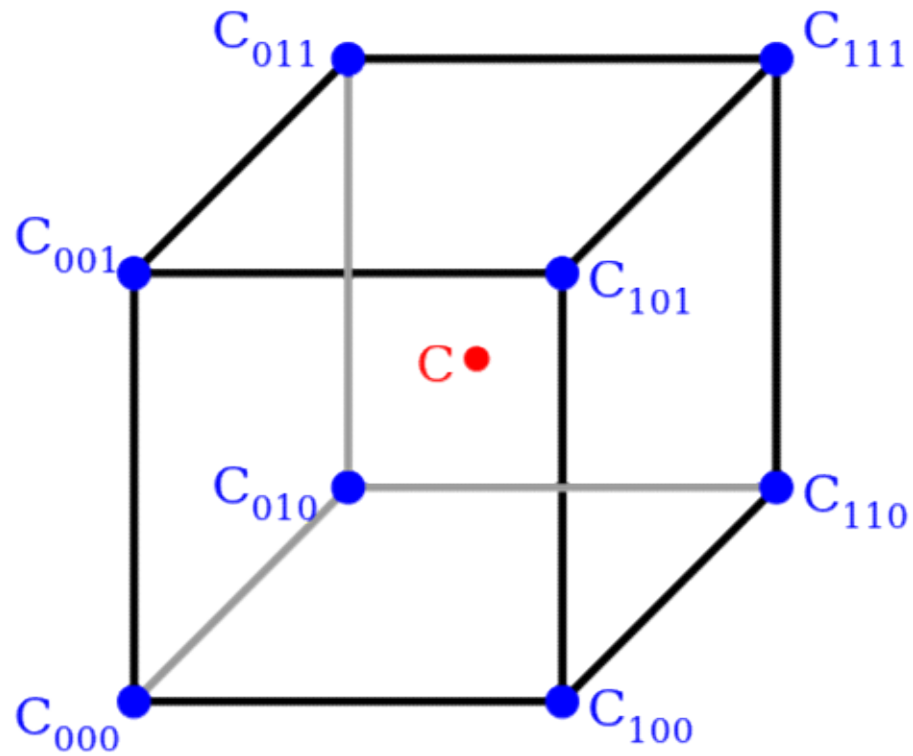
from Wikipedia

# Bilinear Interpolation



What is the image telling us?

# What about 3D?

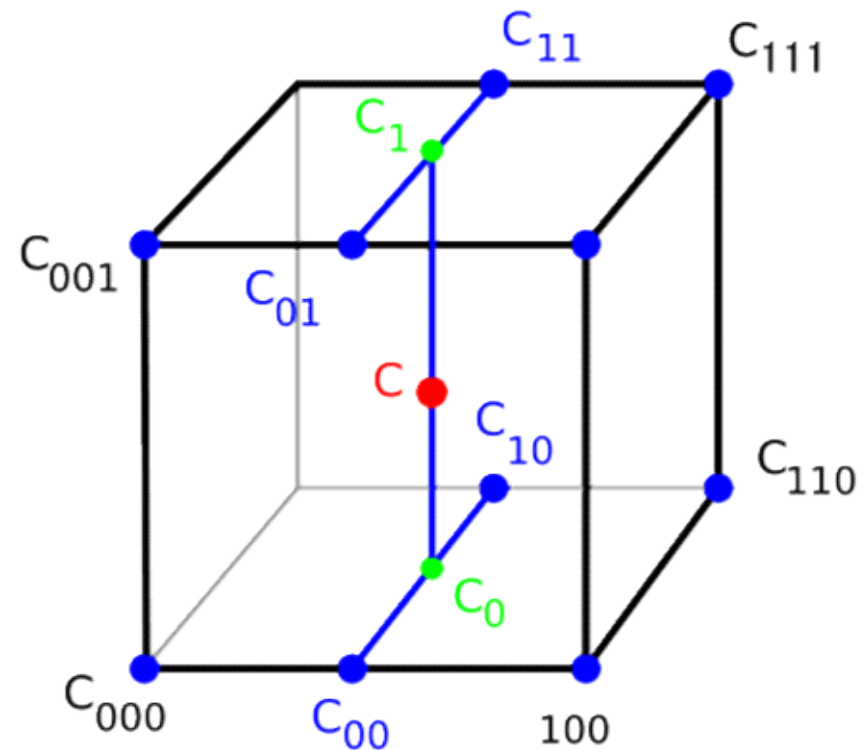


# Trilinear Interpolation

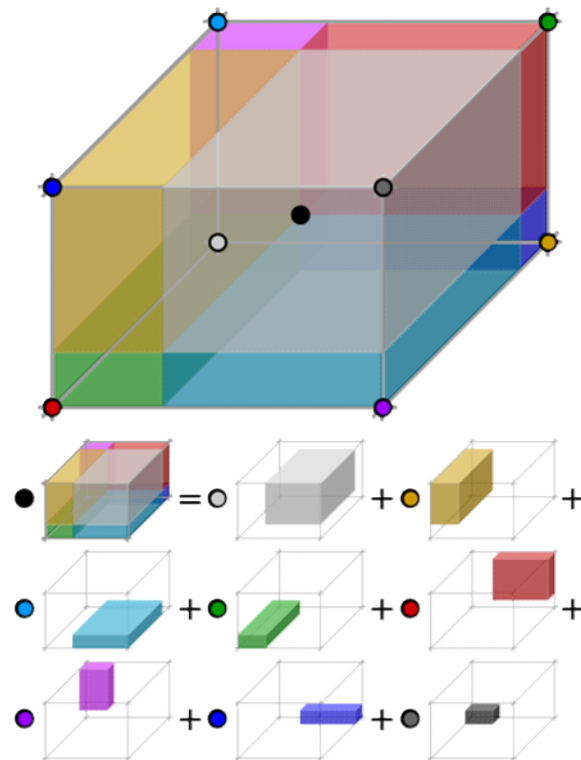
First interpolate in x to find  $C_{00}$ ,  $C_{01}$ ,  $C_{10}$ , and  $C_{11}$

Then in y to find  $C_0$  and  $C_1$

And then in z to find  $C$

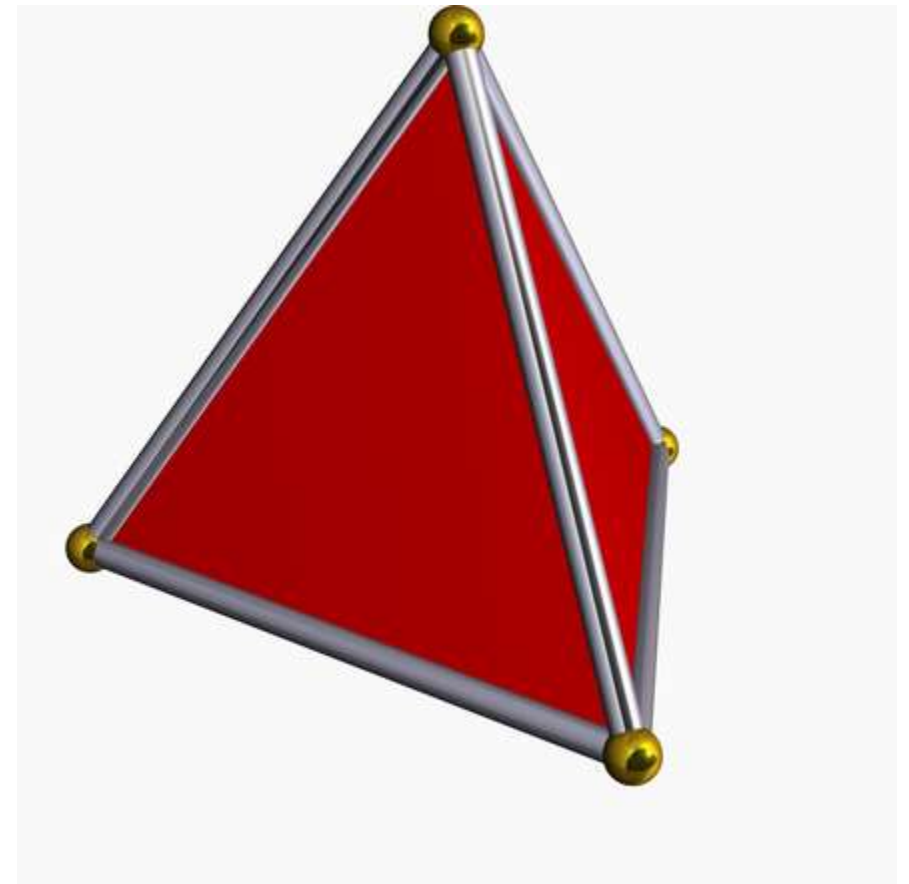


# Trilinear Interpolation



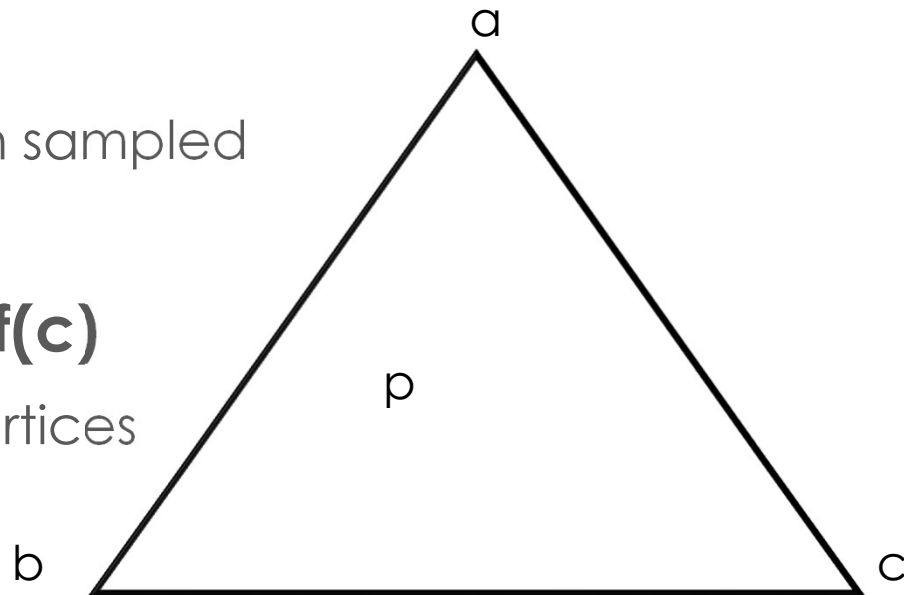
# Barycentric Coordinates

- What about non-quad-like things?
- A simplex is...
  - Convex hull of  $k+1$  points in a  $k$ -dimensional space
  - Simplest convex “polygon” in a  $k$ -dimensional space
  - A 3-simplex is a....
- Barycentric coordinates provide a simple way to interpolate over simplices



# Barycentric Coordinates for Triangles

- Describe location of point in a triangle in relation to the vertices
- $p = (\lambda_1, \lambda_2, \lambda_3)$  where the following are true
  - $p = \lambda_1 a + \lambda_2 b + \lambda_3 c$
  - $\lambda_1 + \lambda_2 + \lambda_3 = 1$
- To interpolate a function sampled at the vertices we just do:  
 **$f(p) = \lambda_1 f(a) + \lambda_2 f(b) + \lambda_3 f(c)$**
- Does the order of the vertices need to be specified?



# Barycentric Coordinates for Triangles

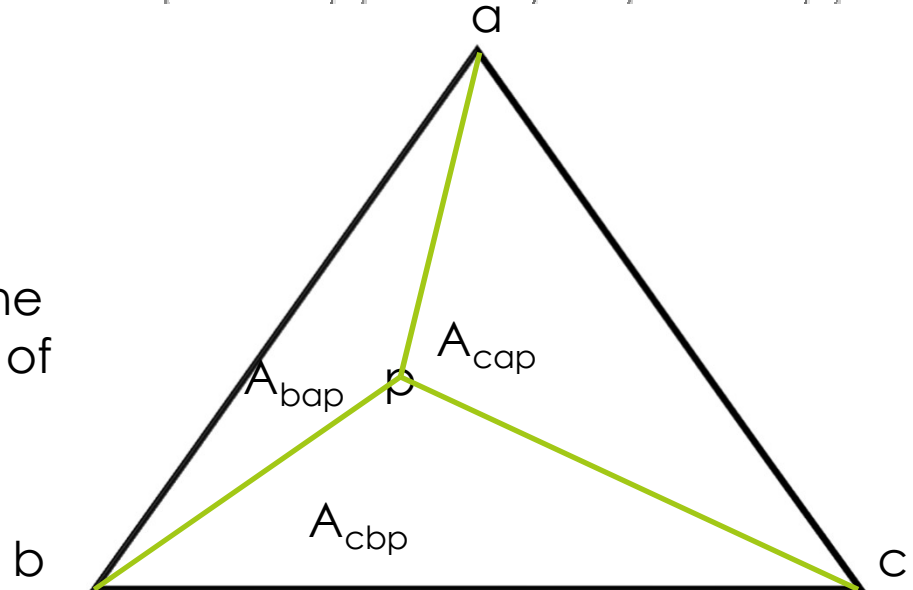
$$\lambda_1 = \frac{(y_2 - y_3)(x - x_3) + (x_3 - x_2)(y - y_3)}{\det(T)} = \frac{(y_2 - y_3)(x - x_3) + (x_3 - x_2)(y - y_3)}{(y_2 - y_3)(x_1 - x_3) + (x_3 - x_2)(y_1 - y_3)},$$

$$\lambda_2 = \frac{(y_3 - y_1)(x - x_3) + (x_1 - x_3)(y - y_3)}{\det(T)} = \frac{(y_3 - y_1)(x - x_3) + (x_1 - x_3)(y - y_3)}{(y_2 - y_3)(x_1 - x_3) + (x_3 - x_2)(y_1 - y_3)},$$

$$\lambda_3 = 1 - \lambda_1 - \lambda_2.$$

Coordinates are the signed area of the opposite subtriangle divided by area of the triangle

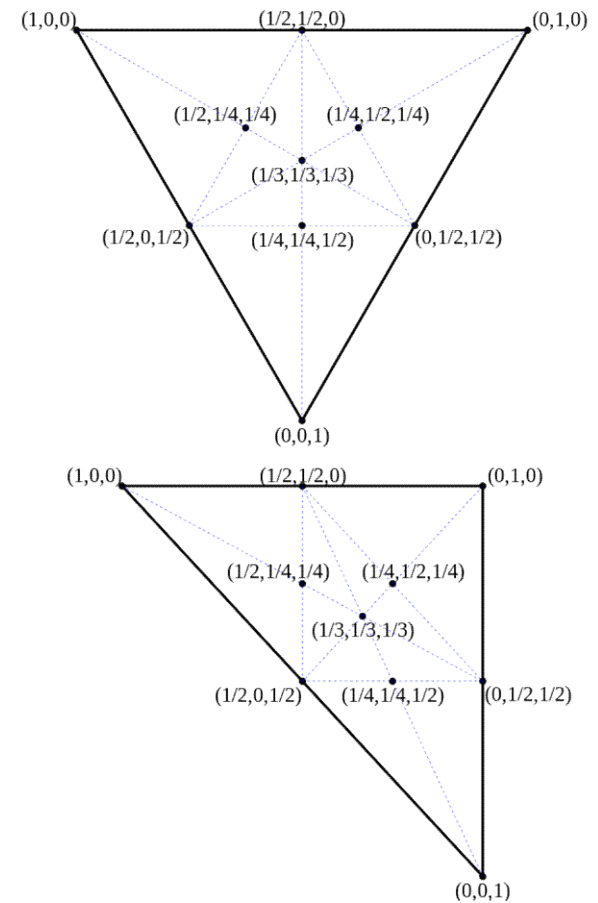
What about triangles in  $\mathbf{R}^3$  ?





# Barycentric Coordinates for Triangles

- Can barycentric coordinates be negative?
- What do you know about a point if it has a coordinate not in  $[0,1]$ ?



# Barycentric Coordinates for Tetrahedra

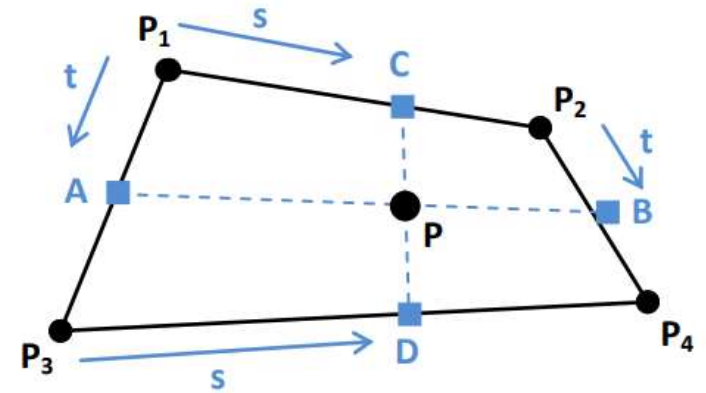
$$\begin{pmatrix} \lambda_1 \\ \lambda_2 \\ \lambda_3 \end{pmatrix} = \mathbf{T}^{-1}(\mathbf{r} - \mathbf{r}_4)$$
$$\mathbf{T} = \begin{pmatrix} x_1 - x_4 & x_2 - x_4 & x_3 - x_4 \\ y_1 - y_4 & y_2 - y_4 & y_3 - y_4 \\ z_1 - z_4 & z_2 - z_4 & z_3 - z_4 \end{pmatrix}$$

We can solve a linear system to find the coordinates.  
Here,  $\mathbf{r}$  is a point in  $\mathbf{R}^3$  and  $\mathbf{r}_4$  is the 4<sup>th</sup> corner of the tetrahedron.

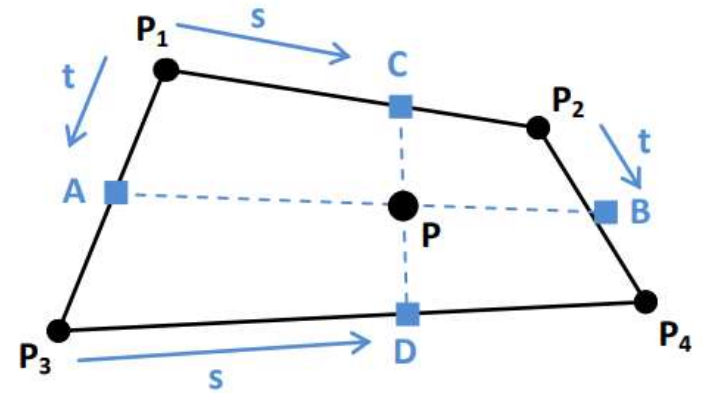
How do we find  $\lambda_4$ ?

# Scattered Data Interpolation

- What happens if you have data sampled irregularly?
  - Can you use bilinear interpolation?
- There are interpolation methods for unstructured data
  - i.e. for data sampled in any pattern
- Could we use barycentric coordinates?
  - How?
  - What would the main drawbacks be?



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# Scattered Data Interpolation

- ▣ There are interpolation methods for unstructured data
    - ▣ i.e. for data sampled in any pattern
  - ▣ We will look at 2 types:
    - ▣ Shepard's interpolation
    - ▣ Radial Basis Functions
  - ▣ Both allow us to interpolate without meshing
  - ▣ Functions are defined in terms of distance from a point
    - ▣ Hence the the term *radial*
-

# Shepard's Interpolation

$$f_i(\mathbf{p}) = \frac{(\|\mathbf{p} - \mathbf{p}_i\|^{-\alpha})}{\sum_{j=1}^n (\|\mathbf{p} - \mathbf{p}_j\|^{-\alpha})} v_i$$

$$f(\mathbf{p}) = \sum_{i=1}^n f_i(\mathbf{p})$$

- ▣ Point  $\mathbf{p}$  is the location at which we are interpolating
  - ▣  $f(\mathbf{p})$  is the interpolated value
- ▣ The  $\mathbf{p}_i$  are the sample points
- ▣ The  $v_i$  are the function values at  $\mathbf{p}_i$
- ▣ Alpha is a positive real number
  - ▣ What does it control?

# Shepard's Interpolation

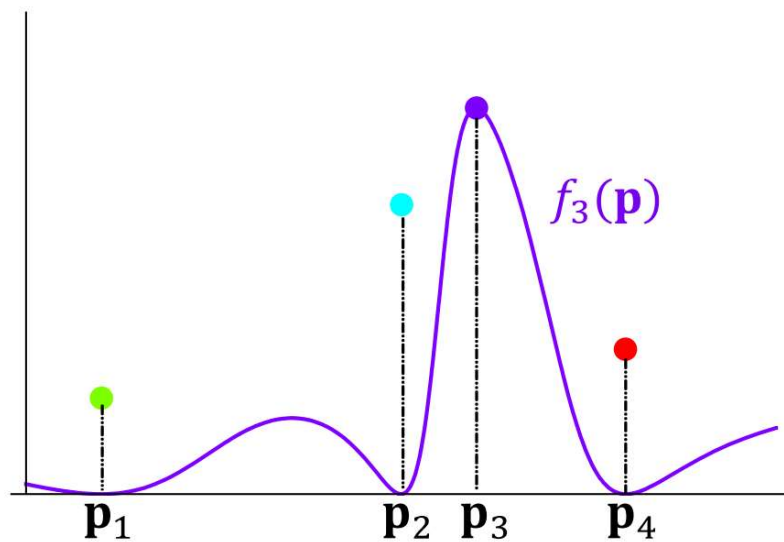
$$f_i(\mathbf{p}) = \frac{(\|\mathbf{p} - \mathbf{p}_i\|^{-\alpha})}{\sum_{j=1}^n (\|\mathbf{p} - \mathbf{p}_j\|^{-\alpha})} v_i$$

$$f(\mathbf{p}) = \sum_{i=1}^n f_i(\mathbf{p})$$

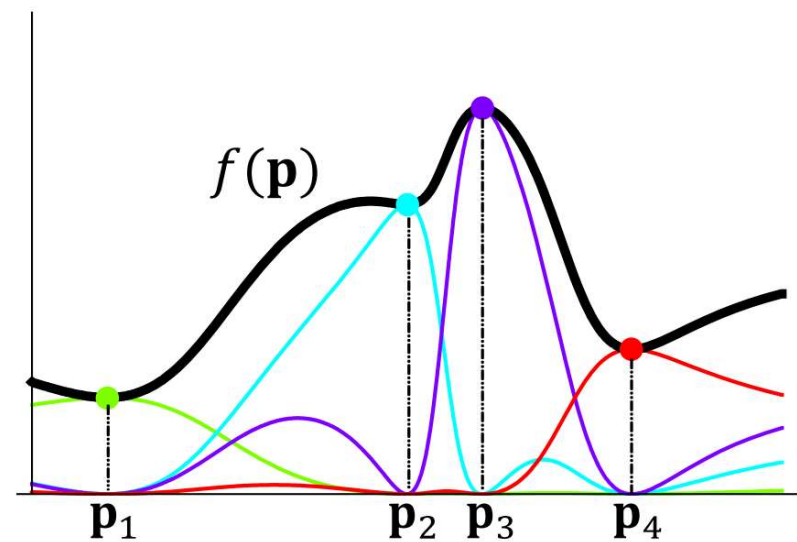
- When approaching  $\mathbf{p}_i$ , the bump function  $f_i(\mathbf{p})$  has both numerator and denominator approach infinity
  - When approaching  $\mathbf{p}_j$  then  $f_i(\mathbf{p})$  approaches 0
- Not confined to convex hull of points...can extrapolate
- Not ideal...flattens at  $\mathbf{p}_i$  and has more waviness than seems natural

# Shepard's Interpolation: Example

Using  $\alpha=2$



(a)



(b)



# Radial Basis Functions (RBFs)

- Any function dependent on distance from a center is *radial*
- We can compute an approximate function as a weighted sum..

$$\phi(x, p) = \phi(\|x - p\|)$$

$$f(x) \approx \sum_{i=1}^N w_i \phi(x, p_i)$$

- Some popular RBFs include

$$\phi(r) = e^{-\lambda r^2} \quad \text{Gaussian}$$

$$\phi(r) = \frac{1}{1 + r^2} \quad \text{Inverse distance}$$

$\lambda$  is a parameter you choose. What behavior does it control?

$r$  is a distance

# Radial Basis Functions (RBFs)

- We have the freedom to choose the weights  $w_i$
- But we have to interpolate the data
- We can find the weights that allow us to do that by solving a system of equations

$$f(p_j) = \sum_{i=1}^N w_i \phi(p_j, p_i)$$

$$Aw = p$$

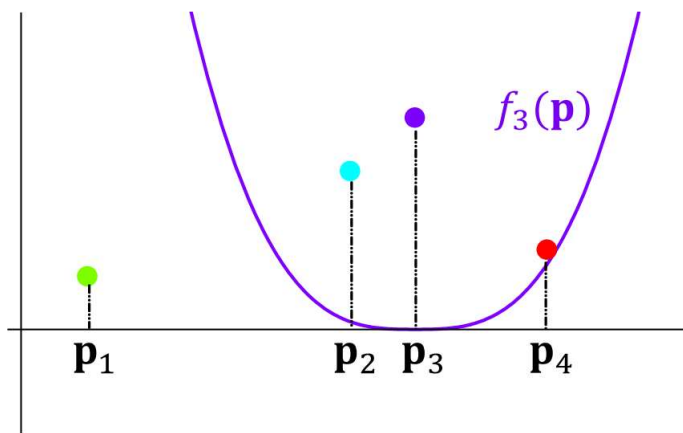
$$A = \begin{bmatrix} \phi(p_1, p_1) & \dots & \phi(p_1, p_N) \\ \dots & \dots & \dots \\ \phi(p_N, p_1) & \dots & \phi(p_N, p_N) \end{bmatrix}$$

$$w = \begin{bmatrix} w_1 \\ \dots \\ w_N \end{bmatrix}$$

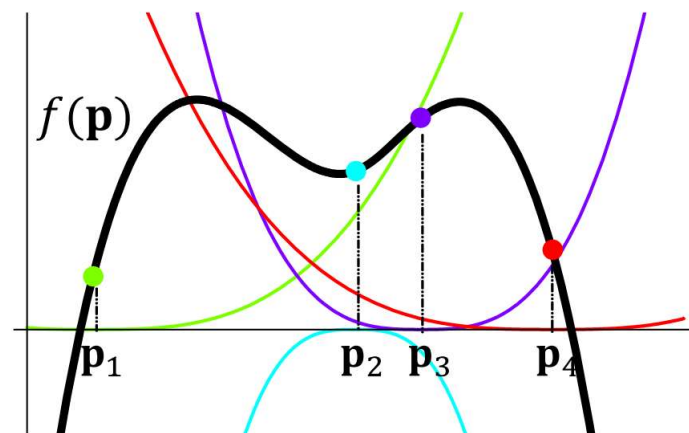
$$p = \begin{bmatrix} f(p_1) \\ \dots \\ f(p_N) \end{bmatrix}$$

# Example: RBF

- Using triharmonic radial functions



(a)



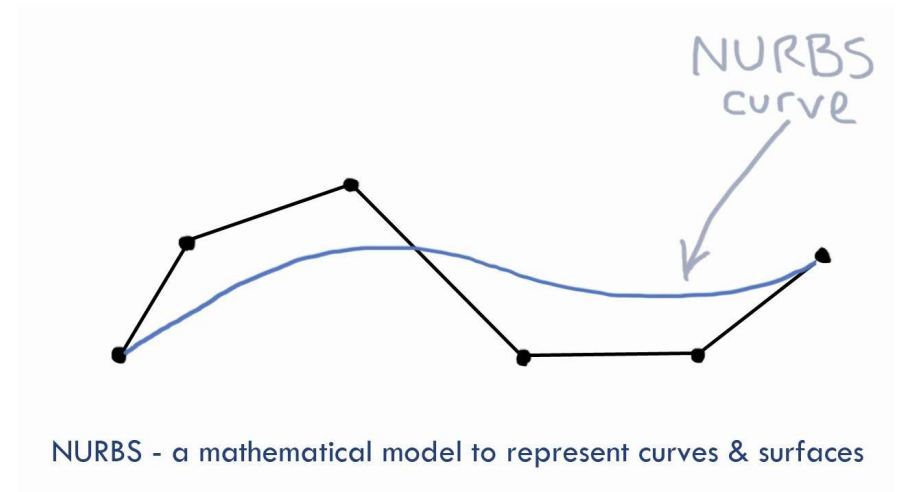
(b)

# RBF Issues

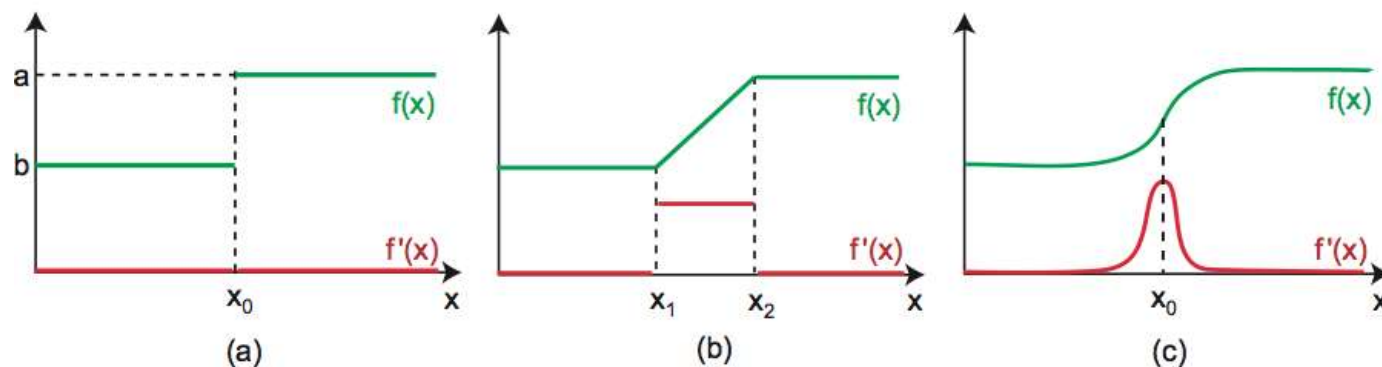
- ❑ Doesn't scale well
- ❑ For large number of points  $A$  becomes ill-conditioned
  - ❑ What does that mean?
- ❑ Also solving system takes time
  - ❑ How much for general Gaussian Elimination on an  $n \times n$ ?
- ❑ Just evaluating the interpolant takes time
- ❑ For large data sets, need to use RBFs with local support
  - ❑ Creates a sparse system more amenable to fast solvers

# Filling in Missing Data

- ▣ There's lots of other options
- ▣ Especially polynomial interpolants
  - ▣ Especially piecewise polynomial interpolants
  - ▣ Splines (need extra control points)
- ▣ Could also use approximation methods
  - ▣ Least squares best fit....



# Reviewing Assigned Reading: Continuous Data



**Figure 3.1.** Function continuity. (a) Discontinuous function. (b) First-order  $C^0$  continuous function. (c) High-order  $C^k$  continuous function.

## Cauchy definition of continuity

A function  $f$  is continuous iff

$$\forall \epsilon > 0, \exists \delta > 0 \text{ such that if } \|x - p\| < \delta, x \in C \text{ then } \|f(x) - f(p)\| < \epsilon.$$



- $C^0$  graph of **derivative** of the function has “holes”
- $C^1$  graph of **derivative** of function has “kinks”
- $C^k$  first  $k$  derivatives of the function are continuous

# Interpolation

Interpolation: Fundamental tool for signal reconstruction

## 1. **Reconstruction** formula

$$\tilde{f} = \sum_{i=1}^N f_i \phi_i \quad \phi_i : D \rightarrow \mathbb{C} \text{ are basis (or interpolation) functions}$$

## 2. **Interpolation**: reconstruction passes through (interpolates) the sampled values

$$\sum_{i=1}^N f_i \phi_i(p_j) = f_j, \forall j. \quad \text{because } \tilde{f}(p_i) = f(p_i) = f_i$$

## 3. **Orthogonality** of basis functions

$$\phi_i(p_j) = \begin{cases} 1, & i = j, \\ 0, & i \neq j. \end{cases} \quad \text{why? Just apply (2) to } f = \begin{cases} 1, & p = p_j \\ 0, & p \neq p_j \end{cases}$$

## 4. **Normality** of basis functions

$$\sum_{i=1}^N \phi_i(x) = 1, \forall x \in D \quad \text{why? } \sum_{i=1}^N \phi_i(p_j) = 1, \forall p_j \text{ (sum (3) over } i = 1..N)$$

and apply above to all  $p_i \in D$

# Piecewise Interpolation

Recall the interpolation formula

$$\tilde{f} = \sum_{i=1}^N f_i \phi_i$$

This becomes **very inefficient** if

- $N$  is very large and we have to evaluate  $\phi_i$  at all these  $N$  points
- $\phi_i$  have complicated expressions

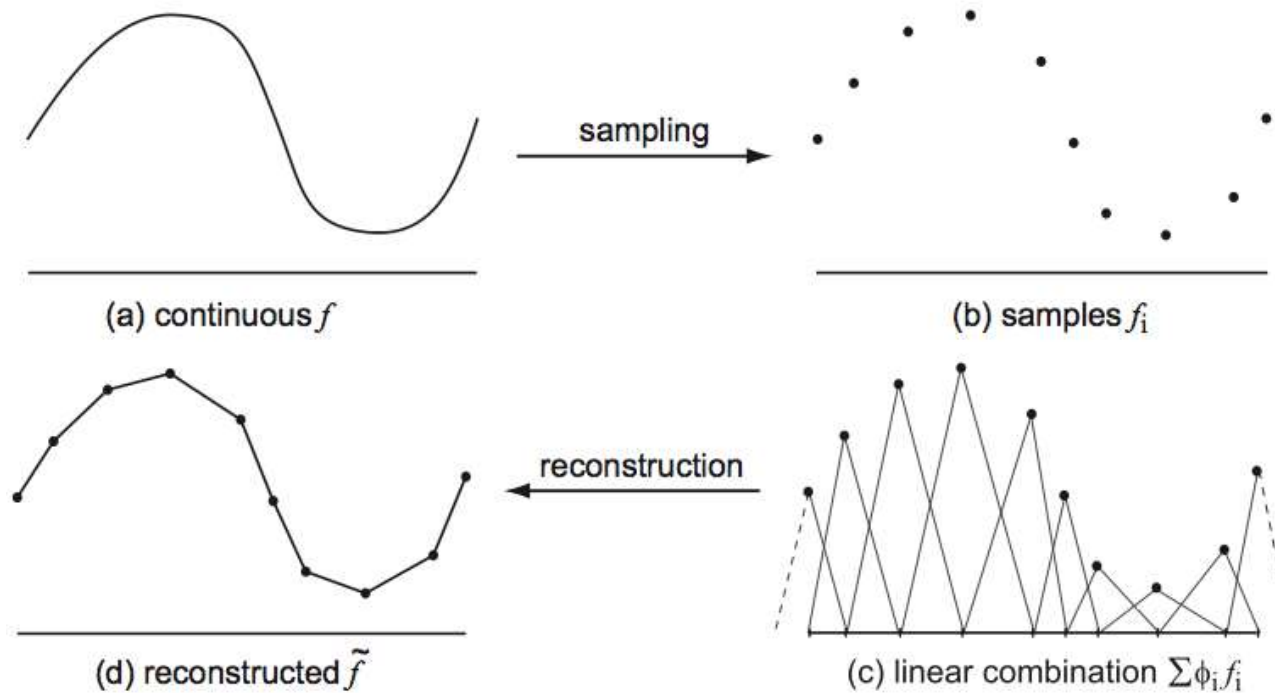
## Practical basis functions

- are non-zero over small spatial ‘pieces’ of  $D$  only (limited support)
- have the same simple formula at all sample points  $p_i$

➡ Additional motivation for discretizing our spatial domain  $D$  into **cells**



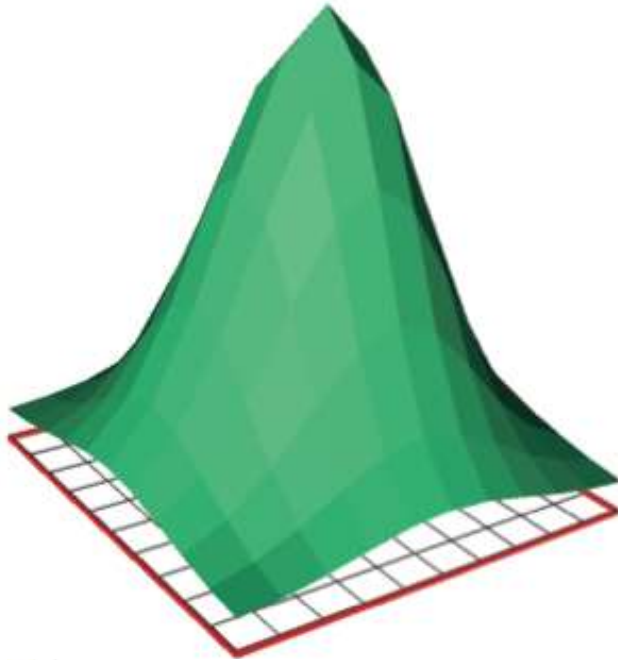
# 1D Example



## Remarks

- interpolation & reconstruction goes cell-by-cell
- only need sample points at a cell vertices to interpolate over that cell

## Bilinear interpolation



$$\Phi_1^1(r, s) = (1 - r)(1 - s),$$

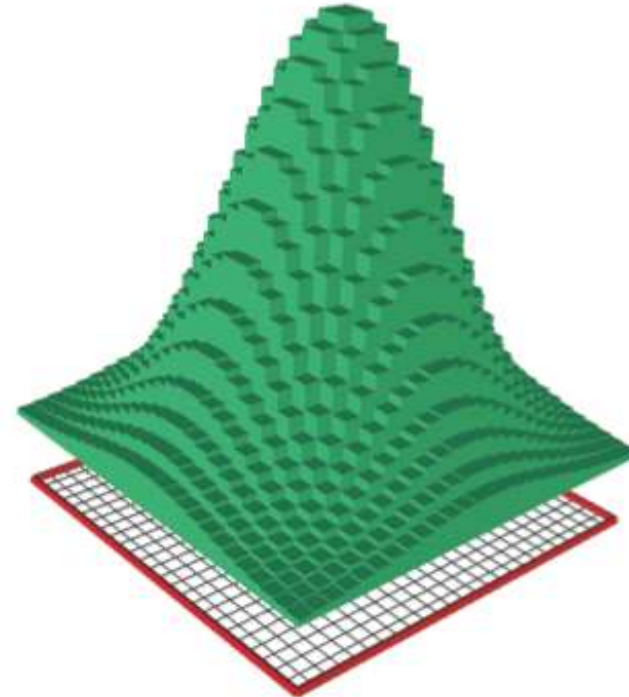
$$\Phi_2^1(r, s) = r(1 - s),$$

$$\Phi_3^1(r, s) = rs,$$

$$\Phi_4^1(r, s) = (1 - r)s;$$

- 4 functions, one **per vertex**
- result:  $C^0$  but never  $C^1$  (why?)
- good for **vertex-based** samples

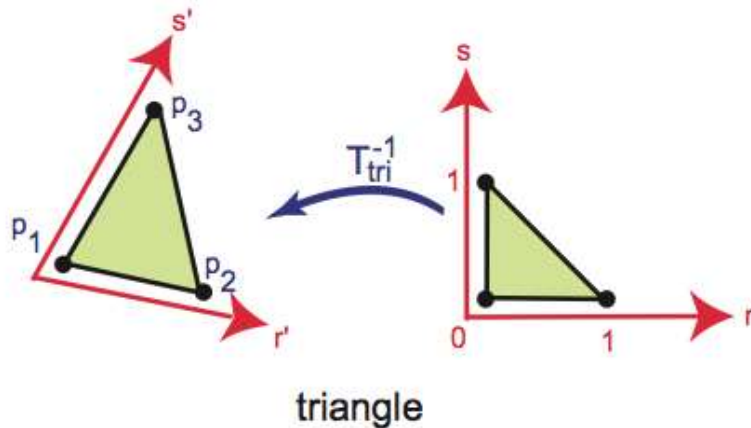
## Constant interpolation



$$\phi_i^0(x) = \begin{cases} 1, & x \in c_i, \\ 0, & x \notin c_i. \end{cases}$$

- 1 functions per **whole cell**
- result: not even  $C^0$
- good for **cell-based** samples

# 2D Cells: Triangles



$$\Phi_1^1(r, s) = 1 - r - s,$$

$$\Phi_2^1(r, s) = r,$$

$$\Phi_3^1(r, s) = s.$$

## Remarks

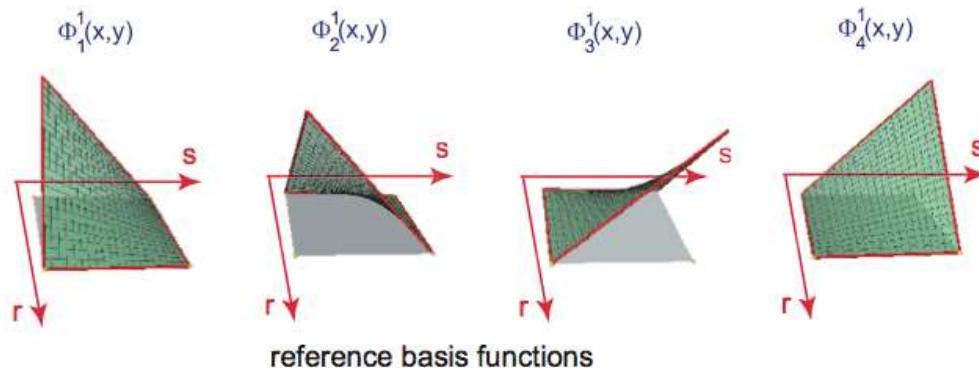
- in graphics/visualization, triangles used more often than quads
  - easier to cover complex shapes with triangles than quads
  - same computational complexity

# From the Book: 2D Cells (Quads)

Same as in 1D case, but

- we have to decide on different cells; say we take quads
- quads  $\rightarrow$  4 vertices, 4 basis functions
- particular case: square cells = pixels

Bilinear basis functions



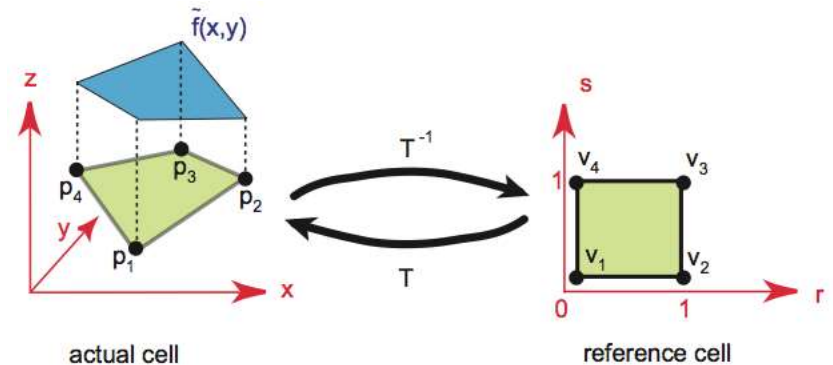
$$\Phi_1^1(r, s) = (1 - r)(1 - s),$$

$$\Phi_2^1(r, s) = r(1 - s),$$

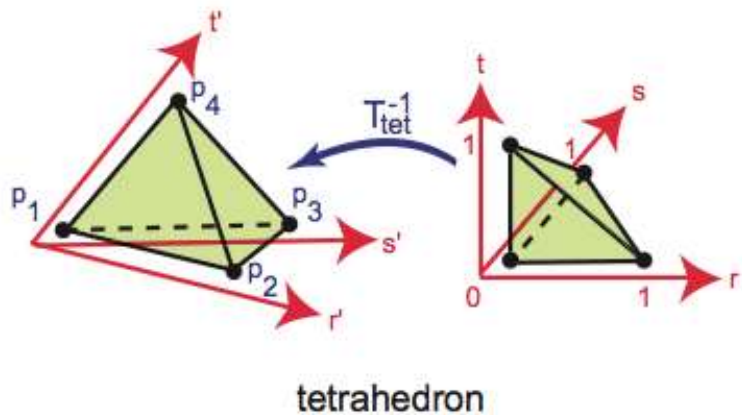
$$\Phi_3^1(r, s) = rs,$$

$$\Phi_4^1(r, s) = (1 - r)s;$$

Bilinear transforms



# 3D Cells: Tetrahedra



$$\Phi_1^1(r, s, t) = 1 - r - s - t,$$

$$\Phi_2^1(r, s, t) = r,$$

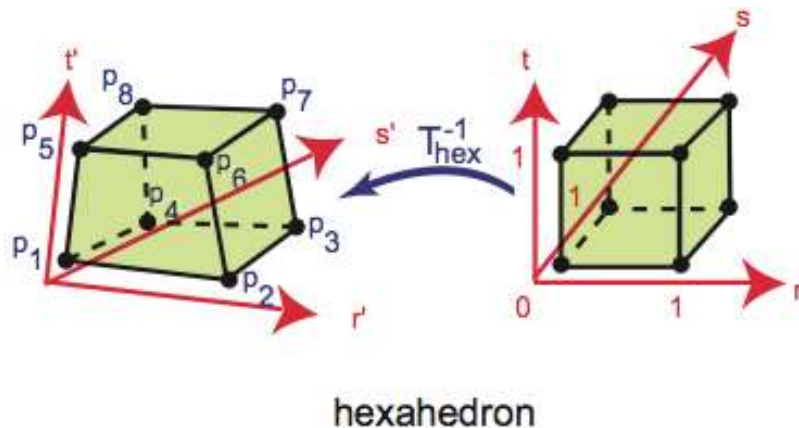
$$\Phi_3^1(r, s, t) = s,$$

$$\Phi_4^1(r, s, t) = t.$$

## Remarks

- counterparts of triangles in 3D
- interpolate **volumetric** functions  $f: \mathbf{R}^3 \rightarrow \mathbf{R}$
- three parametric coordinates  $r, s, t$
- **trilinear** interpolation

# 3D Cells: Hexahedra



$$\begin{aligned}\Phi_1^1(r, s, t) &= (1 - r)(1 - s)(1 - t), \\ \Phi_2^1(r, s, t) &= r(1 - s)(1 - t), \\ \Phi_3^1(r, s, t) &= rs(1 - t), \\ \Phi_4^1(r, s, t) &= (1 - r)s(1 - t), \\ \Phi_5^1(r, s, t) &= (1 - r)(1 - s)t, \\ \Phi_6^1(r, s, t) &= r(1 - s)t, \\ \Phi_7^1(r, s, t) &= rst, \\ \Phi_8^1(r, s, t) &= (1 - r)st.\end{aligned}$$

## Remarks

- counterparts of quads in 3D
- interpolate **volumetric** functions  $f: \mathbf{R}^3 \rightarrow \mathbf{R}$
- **trilinear** interpolation
- particular case: cubic cells or voxels

# Common Cell Types

## 0D

- point

## 1D

- line

## 2D

- triangle, quad, rectangle

## 3D

- tetrahedron, parallelepiped, box, pyramid, prism, ...

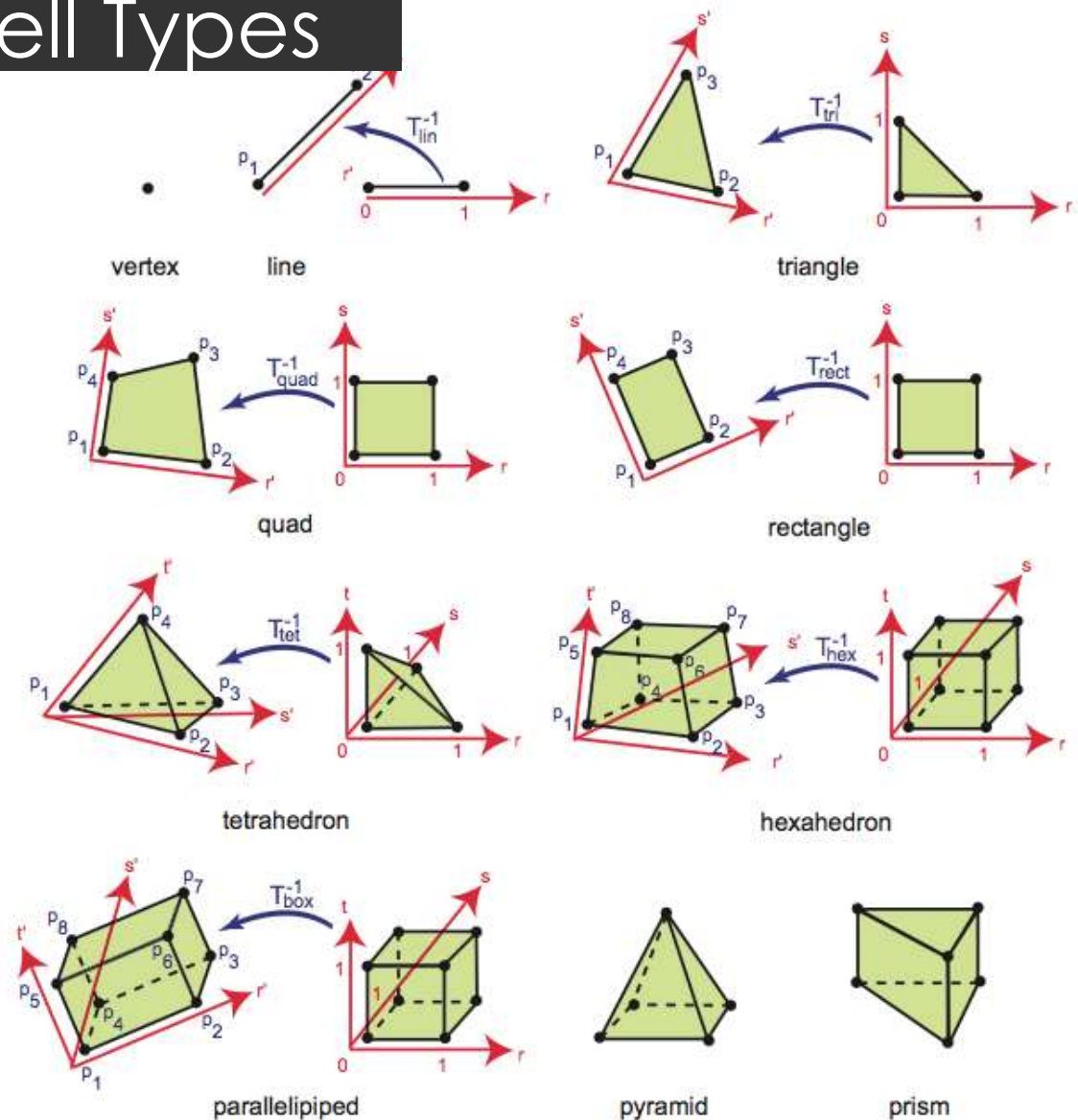
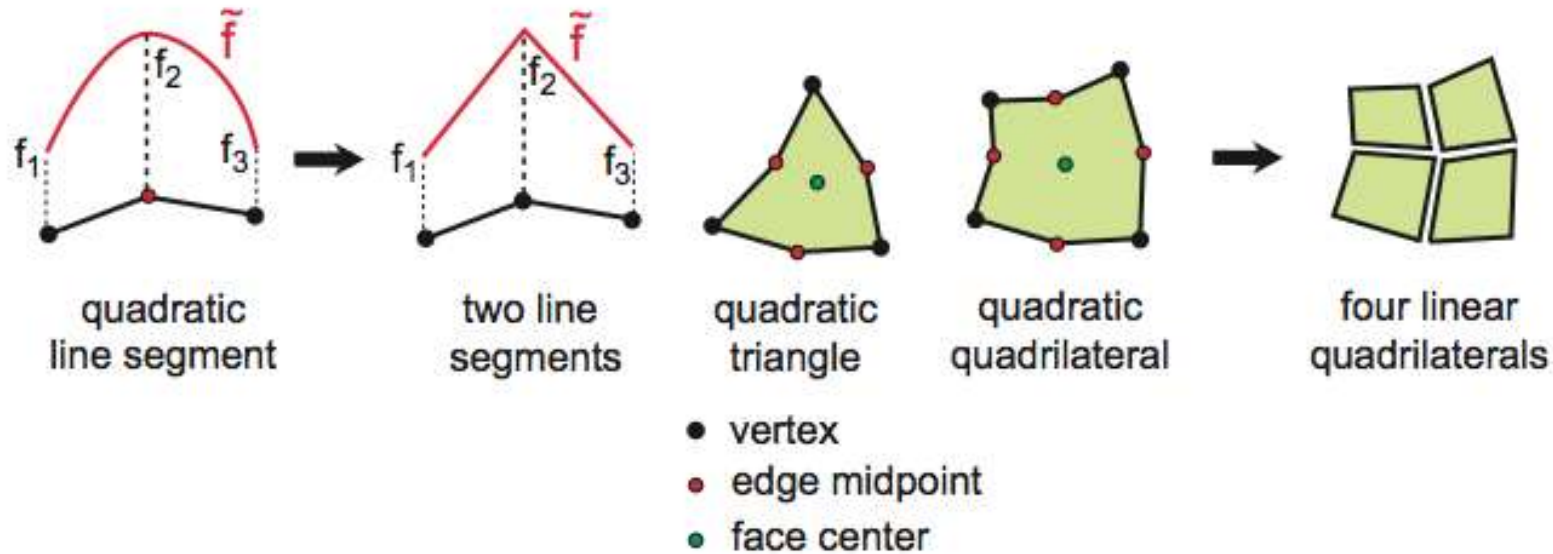


Figure 3.5. Cell types in world and reference coordinate systems.



# Quadratic Cells



**Figure 3.6.** Converting quadratic cells to linear cells.

- allow defining **quadratic** basis functions
- higher **precision** for interpolation
- however, we need data samples at extra midpoints, not just vertices
- used in more complex numerical simulations (e.g. finite elements)
- split into linear cells for visualization purposes