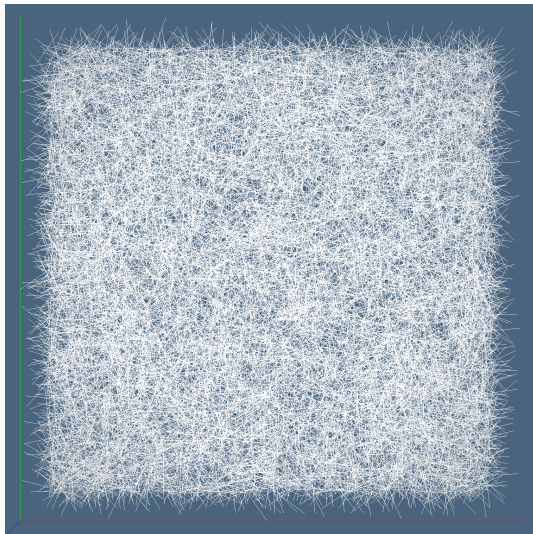


Computational topology: Lecture 8

Alberto Paoluzzi

March 28, 2019

Problem: compute the \mathbb{E}^2 partition



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- 1 Reduction of arrangements to segment intersection
- 2 Segment intersection
- 3 Planar graph by congruence
- 4 Maximal biconnected components

Reduction of arrangements to segment intersection

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Definition

An input collection \mathcal{S} of geometric complexes that mutually intersect, will partition \mathbb{E}^d into a cellular complex $X = \bigcup X_p$ ($0 \leq p \leq d$), called the arrangement $\mathcal{A}(\mathcal{S})$ induced by \mathcal{S}

Computational pipeline

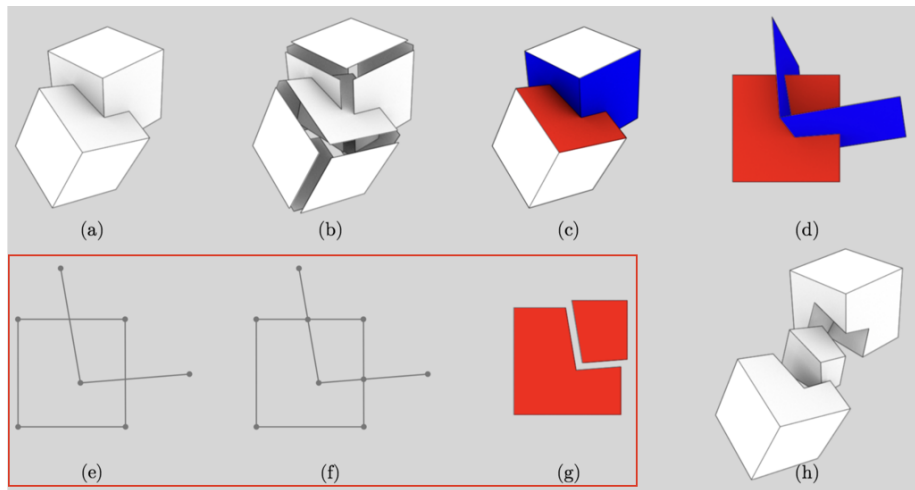


Figure 2: Pipeline segment in 2D

From segment intersection to \mathbb{E}^2 arrangement

Software construction workflow

testing

segment intersection

2D arrangement

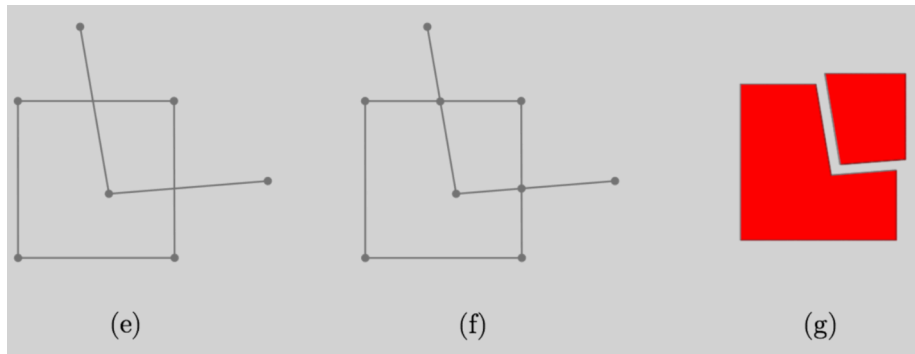


Figure 3: Iterate over 2-cells

Correctness and robustness

- The computation is **correct** and **robust** because the boundaries of adjacent decomposed 2-cells are **compatible as cellular complexes** by **construction**

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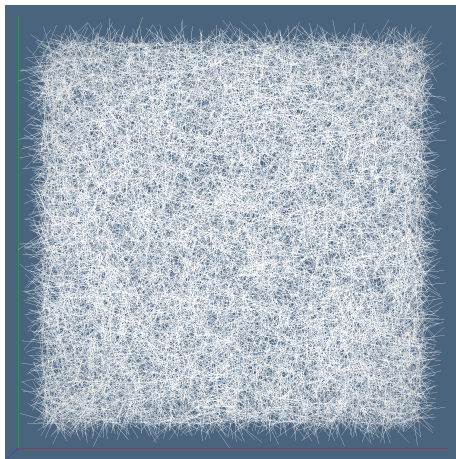
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- This fact is induced by **continuity of topological spaces**, mathematically modeled here by **chains of cells** of a complex
- A requirement of the **standard definition** of a cellular complex demands **boundary compatibility** to hold
- This fact is guaranteed here, since **abutting subsets of 1-cells** have **non-empty intersection**, so they generate **congruent 0-, 1-cells**

Segment intersection

Test-driven development

Always start by parametrically generating some test data

In our case start looking to a **script file** in `/repo/examples`



PARAMETRIC !! generation of random test data

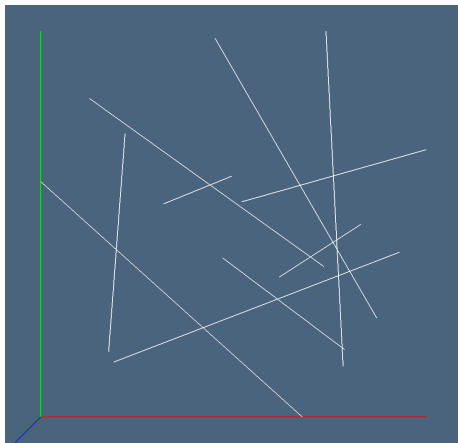


Figure 5: Some examples

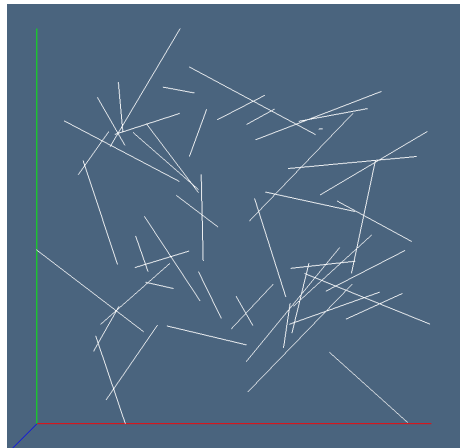


Figure 6: Some examples

Code: examples/randomlines.jl

```
using LinearAlgebraicRepresentation
Lar = LinearAlgebraicRepresentation
using Plasm
```

```
n = 50 #1000 #1000 #20000
```

```
t = 0.5 #0.15 #0.4 #0.15
```

```
V = zeros(Float64,2,2*n)
```

```
EV = [zeros(Int64,2) for k=1:n]
```

```
for k=1:n
```

```
    v1 = rand(Float64,2)
```

```
    v2 = rand(Float64,2)
```

```
    vm = (v1+v2)/2
```

```
    transl = rand(Float64,2)
```

```
    V[:,k] = (v1-vm)*t + transl
```

```
    V[:,n+k] = (v2-vm)*t + transl
```

```
    EV[k] = [k,n+k]
```

```
end
```

```
V = Plasm.normalize(V)
```

```
model = (V,EV)
```

```
Plasm.view(model)
```

<https://github.com/cvdlab/LinearAlgebraicRepresentation.jl/1.0/src/refactoring.jl>

```
function fragmentlines(model)
    V,EV = model

    # acceleration via spatial index computation
    # actual parametric intersection of each line with the cl
    # initialization of local data structures
    # generation of intersection points
    # normalization of output

    return V,EV
end
```

Acceleration via spatial index computation

$\text{Sigma} = \text{Lar.spaceindex}(\text{model})$

```
function spaceindex(model::Lar.LAR)::Lar.Cells
    V,CV = model[1:2]
    dim = size(V,1)

    cellpoints = [ V[:,CV[k]]::Lar.Points
        for k=1:length(CV) ]
    bboxes = [hcat(Lar.boundingBox(cell)...)
        for cell in cellpoints]
    xboxdict = Lar.coordintervals(1,bboxes)
    yboxdict = Lar.coordintervals(2,bboxes)
    # xs,ys are IntervalTree type
    xs = IntervalTrees.IntervalMap{Float64, Array}()
    for (key,boxset) in xboxdict
        xs[tuple(key...)] = boxset
    end
    ys = IntervalTrees.IntervalMap{Float64, Array}()
    for (key,boxset) in yboxdict
        ys[tuple(key...)] = boxset
    end
    xcovers = Lar.boxcovering(bboxes, 1, xs)
    ycovers = Lar.boxcovering(bboxes, 2, ys)
    covers = [intersect(pair...)
        for pair in zip(xcovers,ycovers)]

    if dim == 3
        zboxdict = Lar.coordintervals(3,bboxes)
        zs = IntervalTrees.IntervalMap{Float64, Array}()
        for (key,boxset) in zboxdict
            zs[tuple(key...)] = boxset
        end
        zcovers = Lar.boxcovering(bboxes, 3, ys)
        covers = [intersect(pair...) for pair in zip(zcovers, covers)]
    end
    # remove each cell from its cover
    for k=1:length(covers)
        covers[k] = setdiff(covers[k],[k])
    end
    return covers
end
```

Parametric intersection of lines w the closest ones

```
lineparams = linefragments(V,EV,Sigma)
```

```
julia> Sigma = Lar.spaceindex(model)
20-element Array{Array{Int64,1},1}:
 [19]
 [19]
 [7]
 []
 []
 [19, 13]
 [3]
 [17]
 [12]
 []
 []
 [9]
 [19, 6]
 []
 []
 []
 [8]
 []
 [13, 6, 2, 1]
 []
```

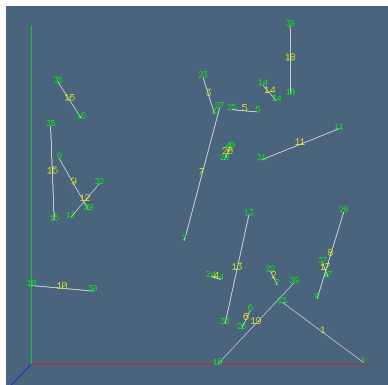


Figure 7: Spatial index computation

Accelerated line intersection

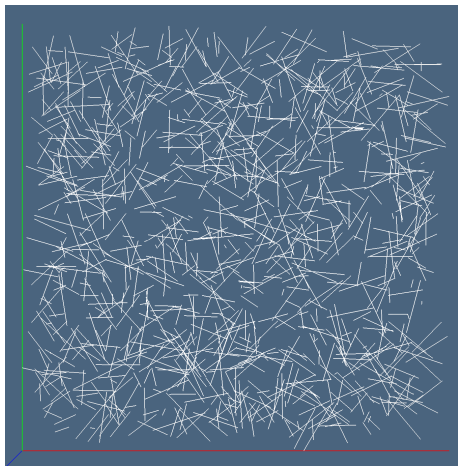


Figure 8: Spatial index computation

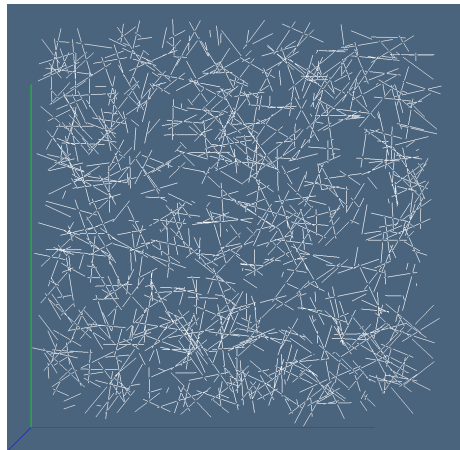


Figure 9: Line segment intersection

Planar graph by congruence

From independently generated line segments to their graph

Homology relation

Two $(d - 1)$ -spaces (curves, surfaces, etc.) embedded in \mathbb{E}^d are topologically *homologous* when their boundaries can be glued, enclosing a portion of the ambient space, and subdivide \mathbb{E}^d in two parts, inner and outer.

From independently generated line segments to their graph

Homology relation

Two $(d - 1)$ -spaces (curves, surfaces, etc.) embedded in \mathbb{E}^d are topologically **homologous** when their boundaries can be glued, enclosing a portion of the ambient space, and subdivide \mathbb{E}^d in two parts, inner and outer.

Congruence relation

Two geometric figures are geometrically **congruent** iff one can be transformed into the other by an isometry.

Congruences are equivalence relations

Congruences R_p between p -cells of geometric complexes are equivalence relations, so we may compute the chain complex of quotient chain spaces:

$$C_2(U_2/R_2) \xrightarrow{\partial_2} C_1(U_1/R_1) \xrightarrow{\partial_1} C_0(U_0/R_0),$$

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- $*C_p(U_p/R_p)$ stands for the chain space generated by $X_p = U_p/R_p$.
- in this stage we compute, for each $\sigma \in \mathcal{S}_2$, the quotient sets and the maps ∂_p in-between, for $p = 0, 1, 2$.

PARAMETRIC !! generation of random test data

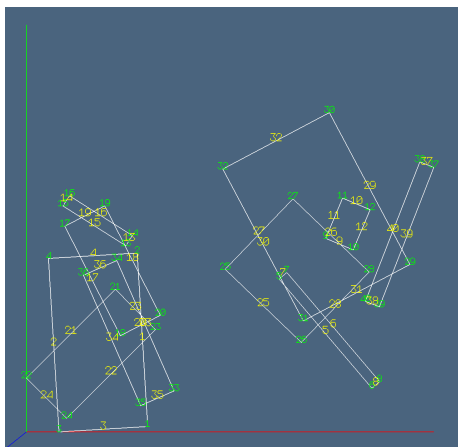


Figure 10: Some examples

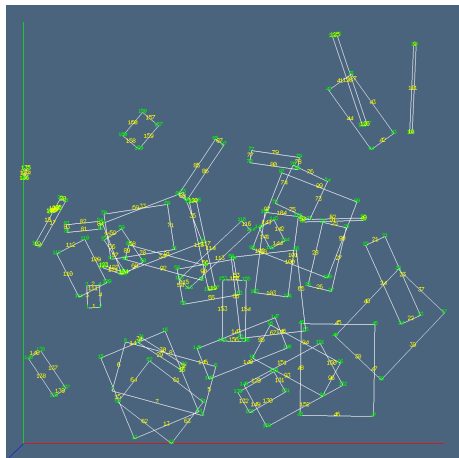


Figure 11: Some examples

Random 2D cuboids in `examples/randomshapes.jl`

using LinearAlgebraicRepresentation Lar = LinearAlgebraicRepresentation
using Plasm

```
function cuboids(n,scale=1.)
    assembly = []
    for k=1:n
        corner = rand(Float64, 2)
        sizes = rand(Float64, 2)
        V,(_,EV,_) = Lar.cuboid(corner,true,corner+sizes)
        center = (corner + corner+sizes)/2
        angle = rand(Float64)*2*pi
        obj = Lar.Struct([ Lar.t(center...), Lar.r(angle),
                           Lar.s(scale,scale), Lar.t(-center...), (V,EV)
                           ])
        push!(assembly, obj)
    end
    Lar.struct2lar(Lar.Struct(assembly))
end
```

end

Generation of parametric test data

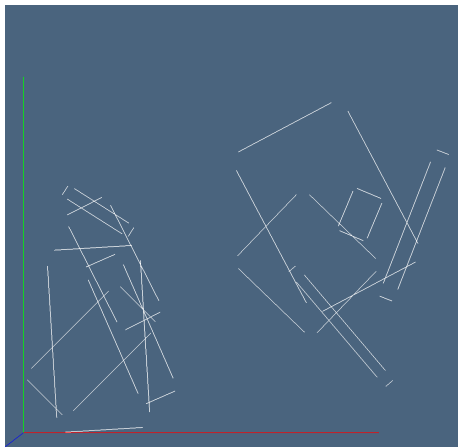


Figure 12: Some example

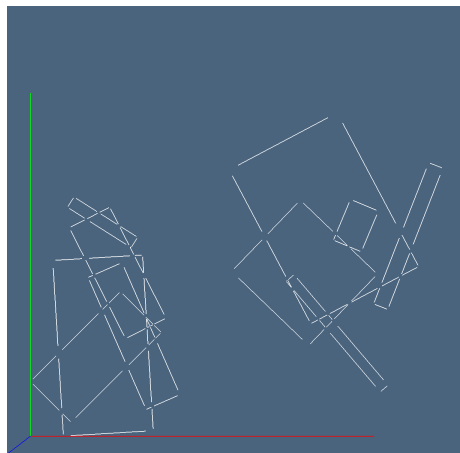
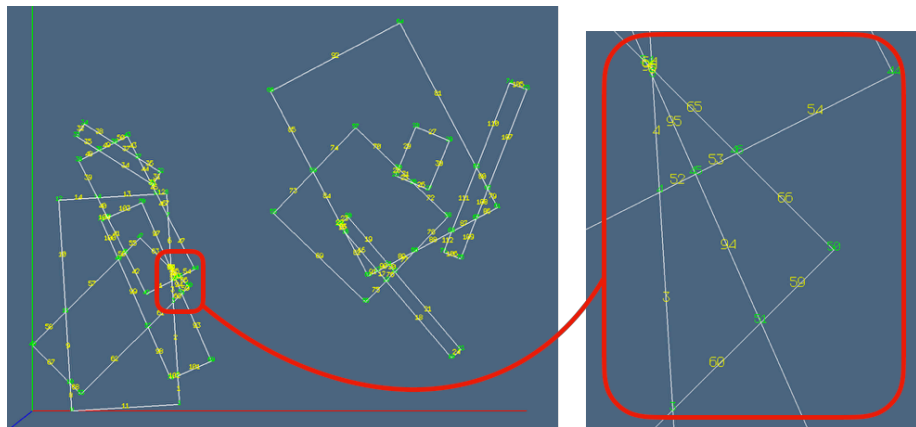


Figure 13: Some example

Visual correctness test

A **formal** (statistical) **proof** will be only possible using the **Euler formula**
 $V - E + F$ after the **construction of the 2-complex** is completed ...



Merge results via search for local neighbor

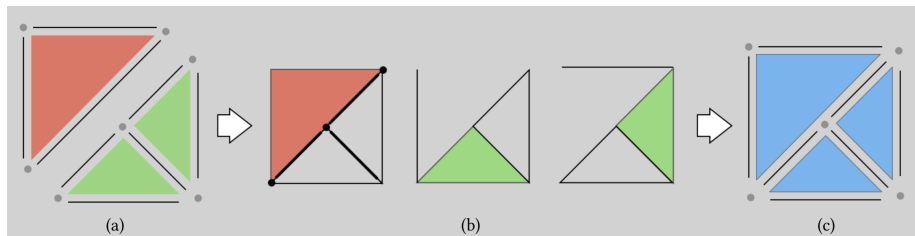


Figure 15: two 2D complexes with incompatible boundaries

Maximal biconnected components

Checks of correctness via graph algorithms

The **planar processing** of each 2-cell continues by pairwise executing the line segment intersection algorithm, and producing a **correct linear graph**

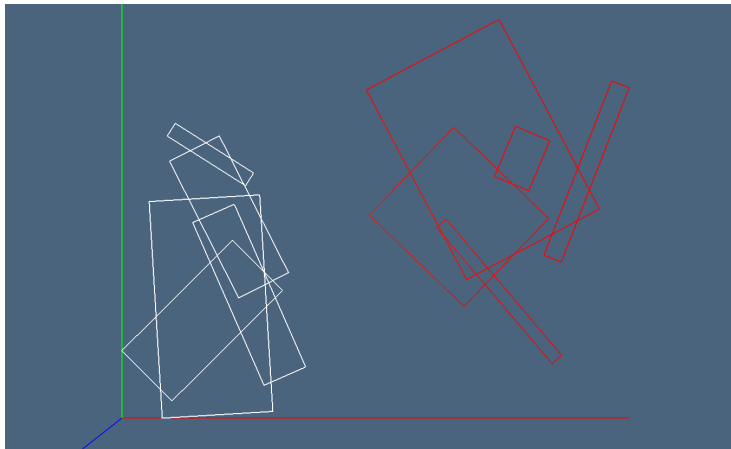


Figure 16: **connected components**

Removing dangling subcomplexes

- In a d -complex, **dangling cells** are p -cells, $p < d$, that are **not contained** in some **boundary cycle** of a d -cell

$$Aop^*B = \overline{(A\overset{\circ}{o}pB)}$$

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- Dangling edges are removed using the **Hopcroft's and Tarjan's algorithm [1974]** for computing the **maximal 2-vertex-connected subgraphs**

bbbbbb

A connected graph G is **2-vertex-connected** if it has at least three vertices and no articulation points

A vertex is an **articulation point** if its removal increases the number of connected components of G

bbbbbb

bbbbbb

bbbbbb