

# Parallelizing zigzag persistence

Mikael Vejdemo-Johansson

Royal Institute of Technology  
Stockholm, Sweden



KTH Computer Science  
and Communication

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# Outline

- 1 Zigzag persistence
- 2 Pullbacks and Parallelization
- 3 Implemented algorithm
- 4 Getting rid of linearity

# Persistent homology

## Classical persistence

A diagram of spaces

$$X_0 \hookrightarrow X_1 \hookrightarrow \dots \hookrightarrow X_n$$

produces a diagram of homology groups

$$H_* X_0 \rightarrow H_* X_1 \rightarrow \dots \rightarrow H_* X_n$$

Decomposes (over a field) into *intervals* of 1-dimensional spaces with identity maps.

# Zigzag persistence

But what if spaces do not include cleanly? What if simplices appear and disappear?

## Zigzag homology

Diagram of Dynkin type  $A_n$ ; arrows can take arbitrary directions.

$$X_0 \hookrightarrow X_1 \hookleftarrow \dots \hookrightarrow X_n$$

produces

$$H_* X_0 \rightarrow H_* X_1 \leftarrow \dots \rightarrow H_* X_n$$

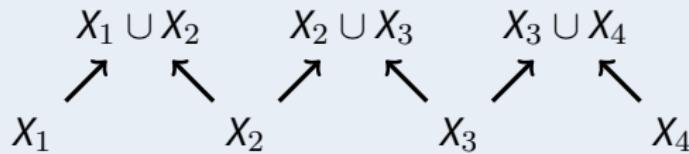
**Gabriel (1970)** proved that these, too, decompose into intervals of 1-dimensional spaces with identity maps.

## Union zigzag

Carlsson & de Silva introduced zigzag persistence, and described a *bootstrap* type technique:

### Union zigzag

Suppose  $X_1, \dots, X_n$  are approximations of a space  $X$ . The zigzag diagram:

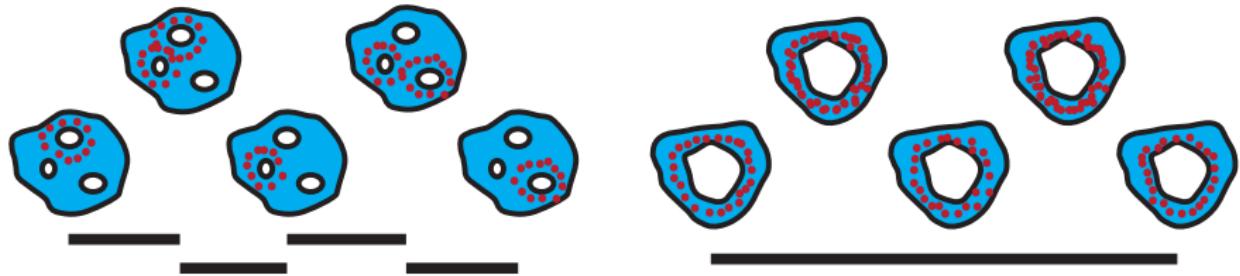


gives rise to a topological style bootstrap: topological features that are artifacts of the approximations live locally; features that are in  $X$  exist everywhere.

# Bootstrap example



# Bootstrap example



## Previous zigzag algorithm

Morozov already presented a parallelizable but different algorithm for computing zigzag persistent homology.

### Basic formulation

Examines the act of adding and removing a simplex, demonstrates how these primitive operations influence the results.

### Advanced formulation

This can be recast into essentially a matrix multiplication, on the order of # insertions and deletions of simplices.

For the bootstrap case, this is prohibitively expensive. Other use cases, this is close to optimal.

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# Pullbacks of linear maps

## Pullbacks in category theory

The pullback  $P$  of two maps  $A \xrightarrow{f} C \xleftarrow{g} B$  is a least specific object with maps to  $A$  and  $B$  such that the diagram

$$\begin{array}{ccc} P & \xrightarrow{\quad \text{---} \quad} & A \\ \downarrow & & \downarrow f \\ B & \xrightarrow{g} & C \end{array}$$

commutes.

## Pullbacks in vector spaces

The pullback has a particularly easy formulation in the category of vector spaces:

$P$  is the subspace of  $A \oplus B$  where the maps agree:

$$P = \{(a, b) \in A \oplus B : fa = gb\}$$

## Pullbacks can compute zigzags

The key recognition here is that the pullback computes the subspace where functions agree.

Consider the diagram, where all maps are induced from inclusions:

$$\begin{array}{ccc} & H_*(X_1 \cup X_2) & \\ \nearrow & & \nwarrow \\ H_*X_1 & & H_*X_2 \end{array}$$

## Pullbacks can compute zigzags

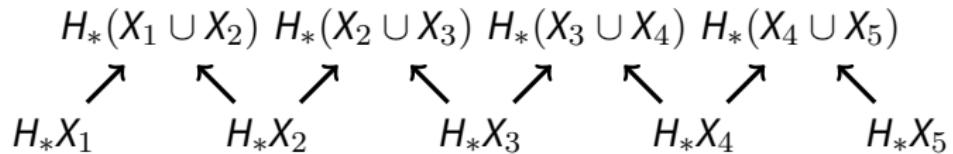
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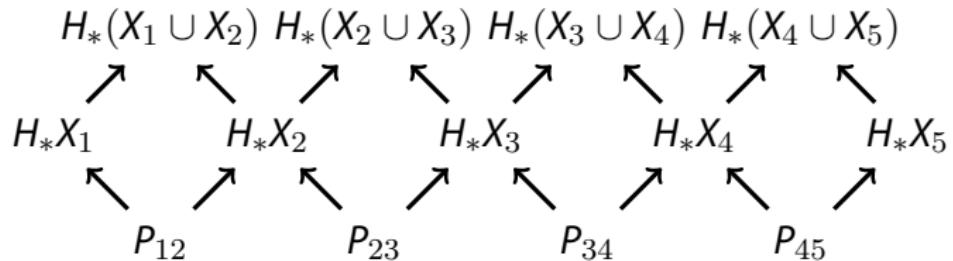
$$\begin{array}{ccc} & H_*(X_1 \cup X_2) & \\ & \swarrow & \nwarrow \\ H_*X_1 & & H_*X_2 \\ & \nwarrow & \swarrow \\ P & & \end{array}$$

Basis vectors of  $P$  encode classes in  $H_*X_1$  and  $H_*X_2$  that agree, up to homology, in  $X_1 \cup X_2$ .

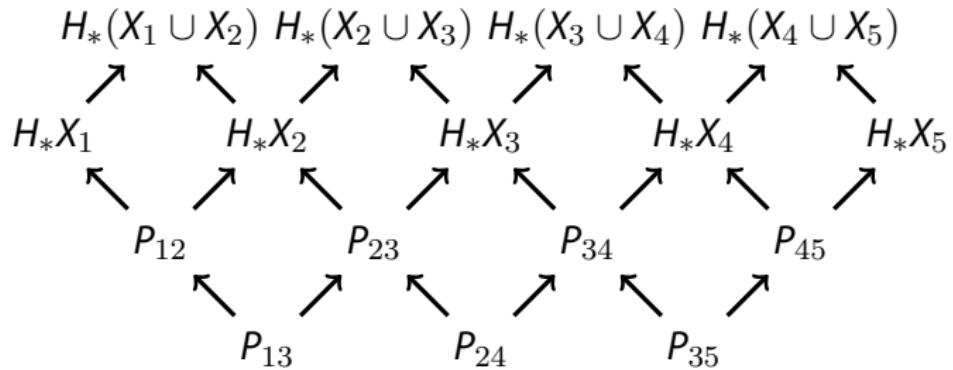
# Algorithm overview



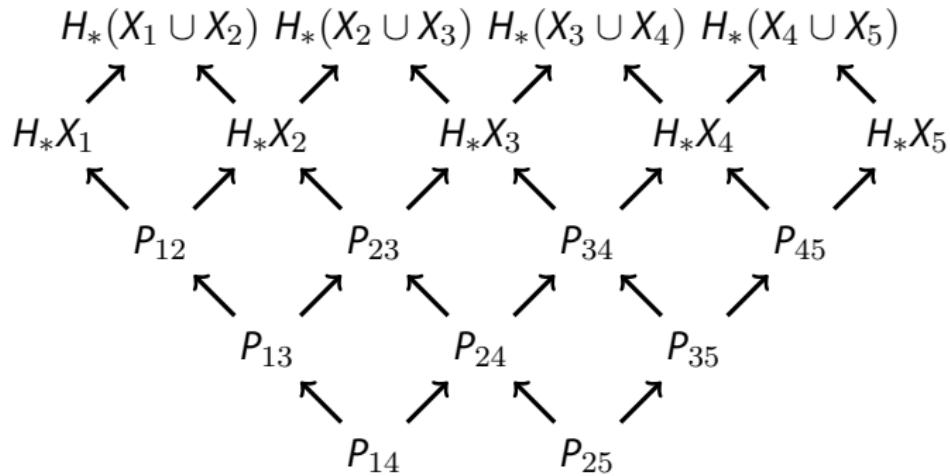
# Algorithm overview



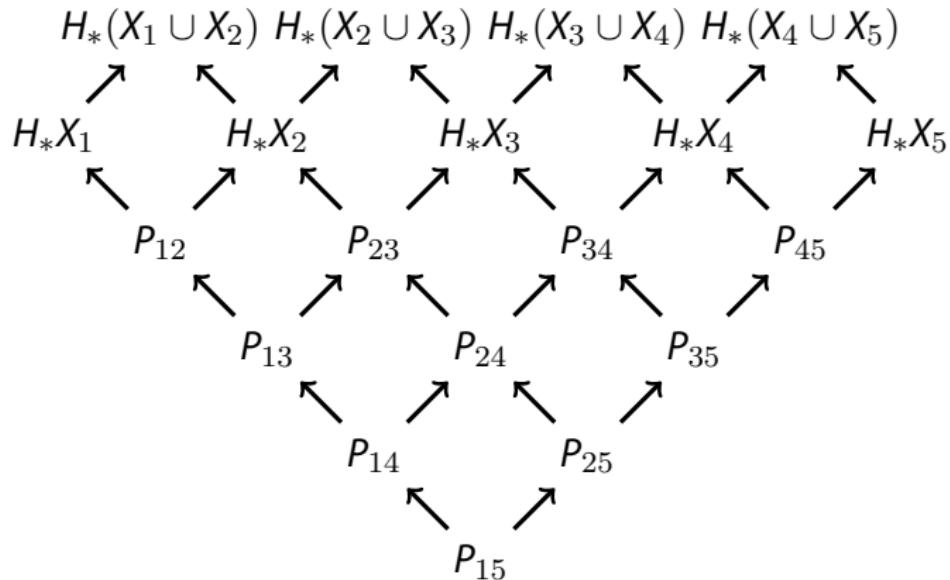
# Algorithm overview



# Algorithm overview



# Algorithm overview



# Algorithm overview

- ① Compute all homology groups.
- ② Compute all induced maps  $H_*X_i \rightarrow H_*(X_i \cup X_j)$ .
- ③ Compute pullbacks repeatedly until all layers processed.
- ④ Compute cokernels of the pullback maps.  
This picks out generators where they take effect.

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## Sample code

A serial implementation in GAP with comment annotations for parallelization with HPC-GAP exists.

Access with Mercurial: `hg://hg.gap-system.org/geometry`

Annotated code resides in `hpczz.g`

Shared but provably disjoint array write access. Shared array read access.

# Code: Assumptions

The functionality is parametrized on:

- A sequence of pointclouds pts.
- A globally chosen Vietoris-Rips radius eps.
- A globally chosen top interesting dimension d.
- Some field field.

## Homology setup

Computing homology is embarrassingly parallel; no dependencies between the spaces. ParList or dependency-free task launches. Needs to be done both for each single space, and each union of neighbouring spaces.

```
SpaceKernel := function(array, index, pts)
    local graph, complex;

    graph := VietorisRipsGraph(pts, eps);
    cpx := VietorisRipsIncremental(graph, d);
    array[index] := rec( cc := CreateChainComplex(cpx, field),
                         spxs := cpx,
                         graph := graph );
end;
```

## Inclusion maps setup

This step is specific to the union zigzag; assumptions on vertex numbering come from how unions are coded. Step `index` depends on the `SpaceKernel` for `index` and `index+1`.

```
UnionKernel := function(array, index, spaces, unions)
    local uu;

    uu := CreateUnion(spaces[index].spxs,
                      spaces[index].graph.nV,
                      spaces[index+1].spxs,
                      unions[index].spxs,
                      field);
    array[index].right := uu[1];
    array[index+1].left := uu[2];
end;
```

# Homology computation

Homology computation locally is still internally dependency-free; but relies on the corresponding SpaceKernel having finished.

```
HomologyKernel := function(array, index, spaces)
  local p, h, z;
  p := NaturalHomomorphismBySubspace(
    Kernel(spaces[index].cc),
    Image(spaces[index].cc));
  h := Image(x1p);
  z := List(Basis(x1h), v -> PreImagesRepresentative(x1p, v));
  array[index].space := h;
  array[index].cycles := z;
  array[index].proj := p;
end;
```

# Induced maps

```
CreateMap := function(from,to,unionmap,proj)
  return LeftModuleHomomorphismByImages(
    from.space, to.space, Basis(from.space), List(List(from.cycles,
      v -> Image(unionmap, v)), w -> Image(proj, w)));
end;

HomologyMapKernel := function(array, idx, state)
  local f, g, sl;
  sl := state.layers; f := fail; g := fail;
  if index < Length(array) then
    f := CreateMap(sl[2][idx], sl[1][idx],
                  state.unionmaps[idx][1], sl[1][idx].proj);
  fi;
  if index > 1 then
    g := CreateMap(sl[2][idx], sl[1][idx],
                  state.unionmaps[idx-1][2], sl[1][idx-1].proj);
  fi;
  array[index].left := g; array[index].right := f;
end;
```

## Pullbacks

The function here is almost trivial; but the dependency structure more intricate.

```
PullbackKernel := function(state, depth, index)
    state.layers[depth][index] := Pullback(
        state.layers[depth-1][index].right,
        state.layers[depth-1][index+1].left);
end;
```

The depth/index invocation of the `PullbackKernel` depends on the `PullbackKernel` invocation for `depth-1/index` and `index+1`.

The first layer depends on the correspondingly indexed induced maps having been computed.

# Cokernels

A final step, for a nice and complete presentation, is to compute cokernels, isolating the essential content of the computed barcode.

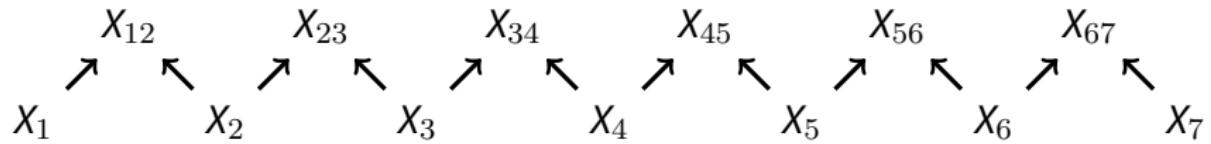
```
CokernelKernel := function(state, depth, index)
local quot, hf, hh;

quot := Subspace(state.layers[depth][index].space, []);
if index <> Length(state.layers[depth]) then
    quot := quot + ImagesSet(state.layers[depth+1][index].left,
                                state.layers[depth+1][index].space);
fi;
if index <> 1 then
    quot := quot + ImagesSet(state.layers[depth+1][index-1].right,
                                state.layers[depth+1][index-1].space);
fi;
hf := NaturalHomomorphismBySubspace(
    state.layers[depth][index].space,
    quot);
hh := Image(hf);
state.zz[depth][index].hf := hf;
state.zz[depth][index].hh := hh;
end;
```

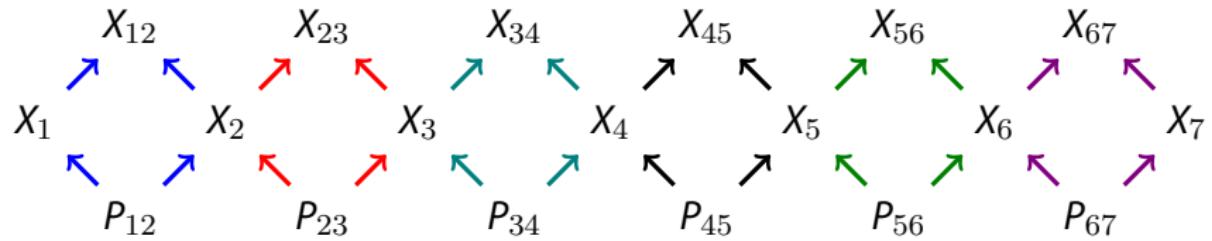
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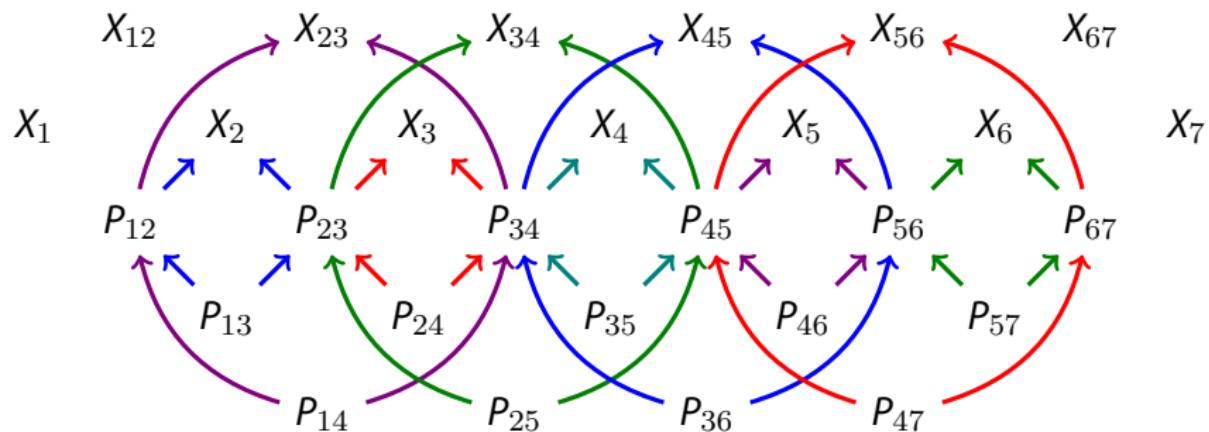
# Logarithmic time; assuming quadratic resources



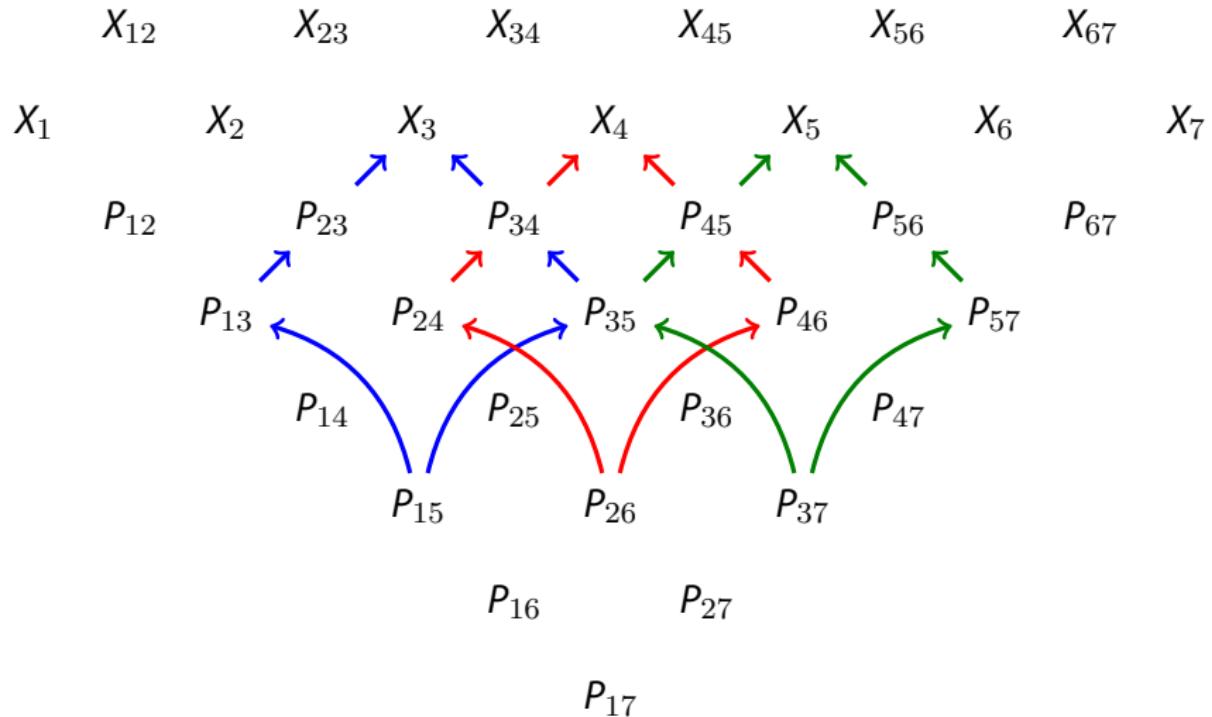
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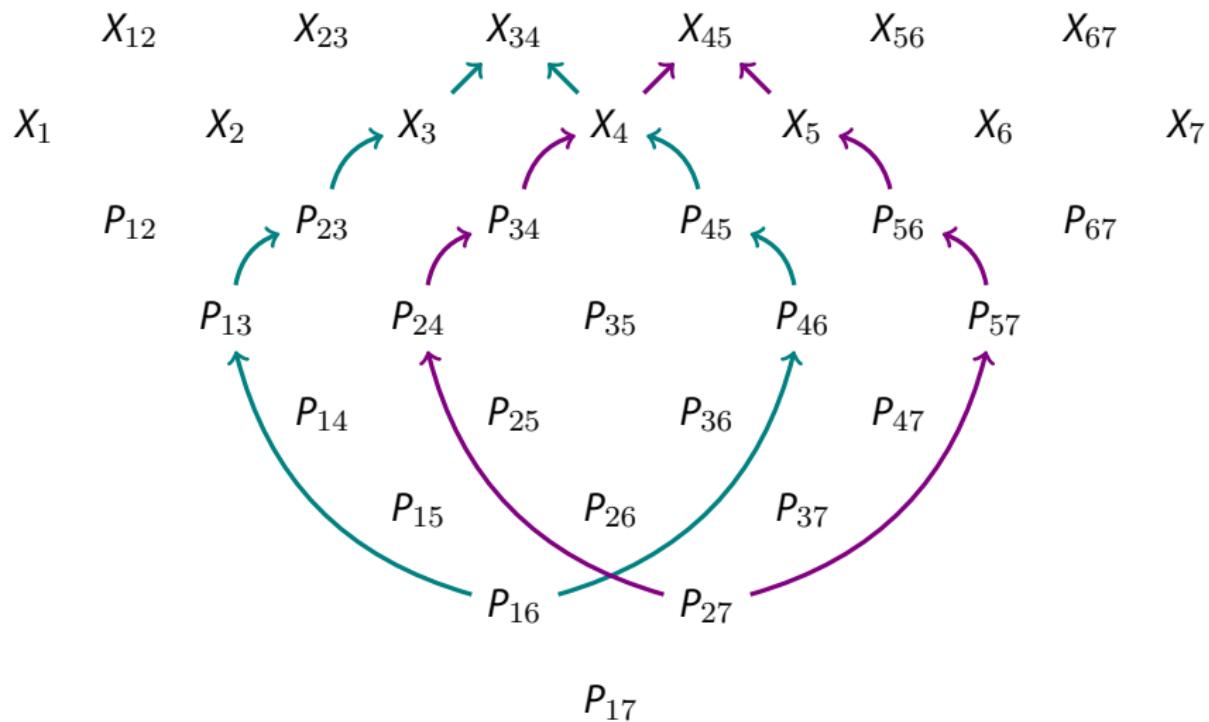
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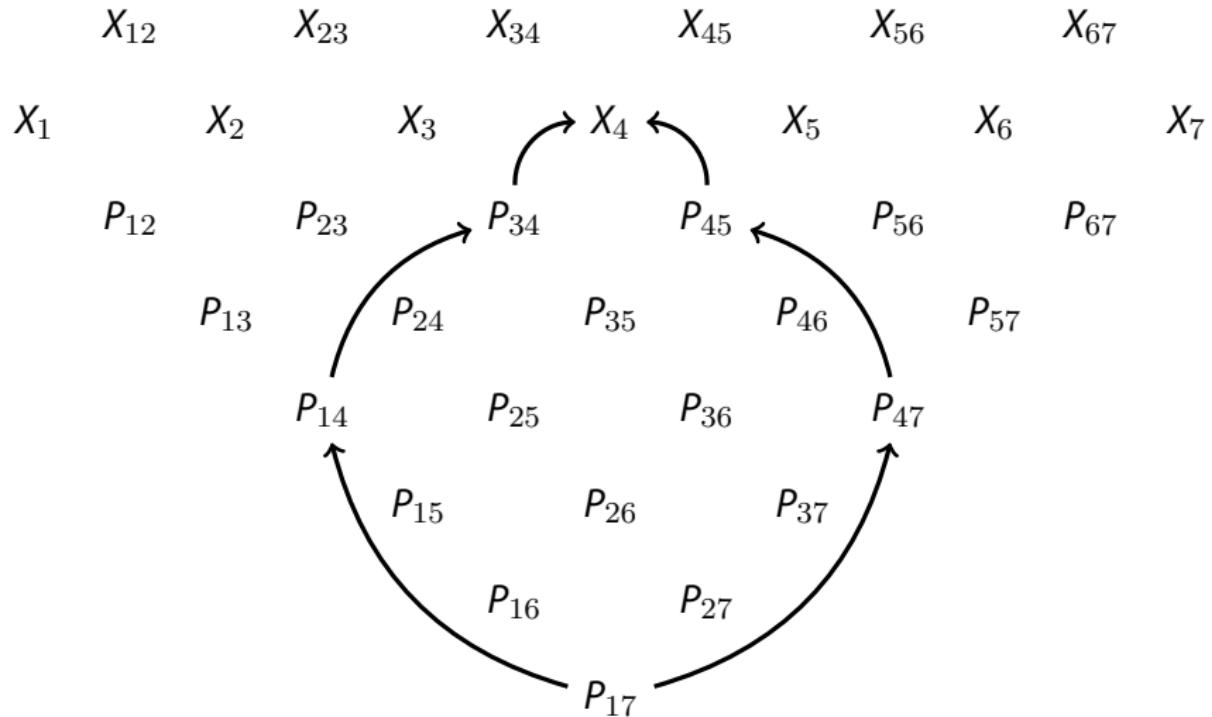
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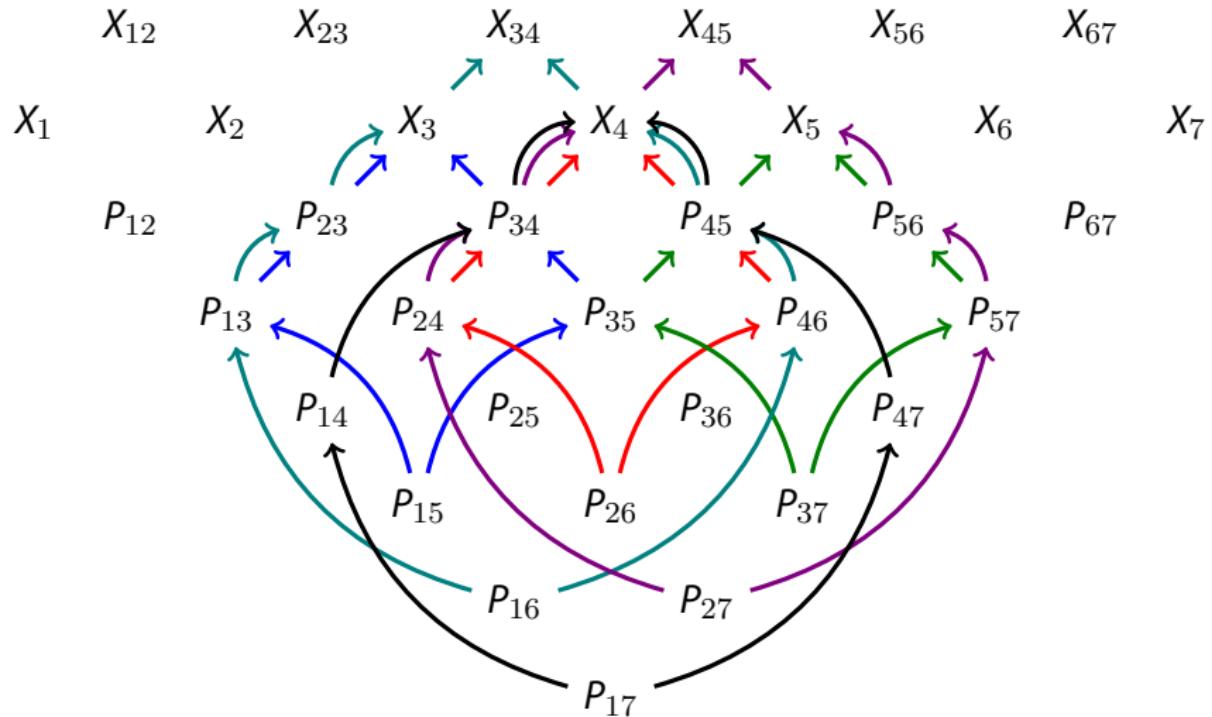
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# Complexity analysis

Assume  $n$  spaces;  $n - 1$  unions. Largest space has  $m$  simplices. Suppose that  $b_k$  is the upper bound on the number of classes persisting at least  $k$  steps.

Homology computation Time  $O(m^\omega)$ . Processors  $2n - 1$ .

Maps computation Time  $O(m^\omega)$ . Processors  $2n - 2$ .

Pullbacks, 1st round Time  $O(b_1^\omega)$ . Processors  $n - 1$ .

Pullbacks, 2nd round Time  $O(b_2^\omega + b_3^\omega)$ . Processors  $(n - 2) + (n - 3)$ .

Pullbacks,  $k$ th round Time  $O(\sum_{i=2^{k-1}}^{2^k-1} b_i^\omega)$ . Processors  $\sum_{i=2^{k-1}}^{2^k-1} n - i$ .

The computation ends after  $\lceil \log_2 n \rceil$  rounds.

$\omega$  is the exponent of matrix multiplication complexity.