

RAGDNet

A Region-Adjacency Graph for Semantic Segmentation of Mechanical Drawings Using Graph Neural Networks

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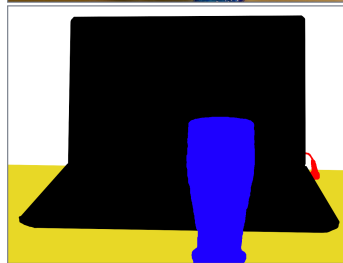
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What is semantic segmentation?

- Assign a class label to every pixel of an image.
- Standard recipe: an encoder–decoder over the pixel grid (*U-Net*) or a transformer over patches (*SegFormer*).

The question is always the same: *what* is it at this location?

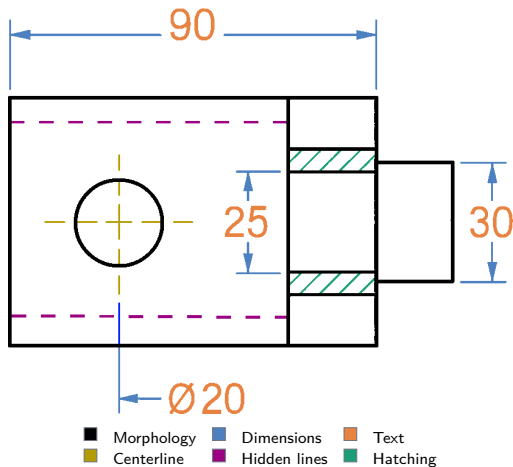


- Laptop
- Table
- Glass
- Cable
- Background

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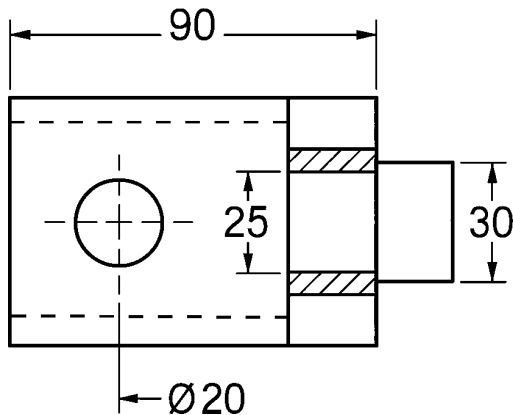
Why segment mechanical drawings?

- Standard medium of technical communication across many industrial domains.
- Semantics are lost in exchange formats: raster is a matrix of pixels (PNG, JPG), vector, a set of geometric primitives with no meaning attached (DXF, DWG).
- Downstream automation:
 - retrieving similar drawings in large databases;
 - automated cost estimation;
 - comparison of information between two versions.



Limitations of current approaches

- In our setting, drawings are **binary and extremely sparse**: in the drawing on the right, ink covers **7.3%** of the pixels.
- Yet in a CNN as in a ViT, **every pixel becomes a variable** in the computation.



829 × 665 px, 40 380 inked, 7.3%

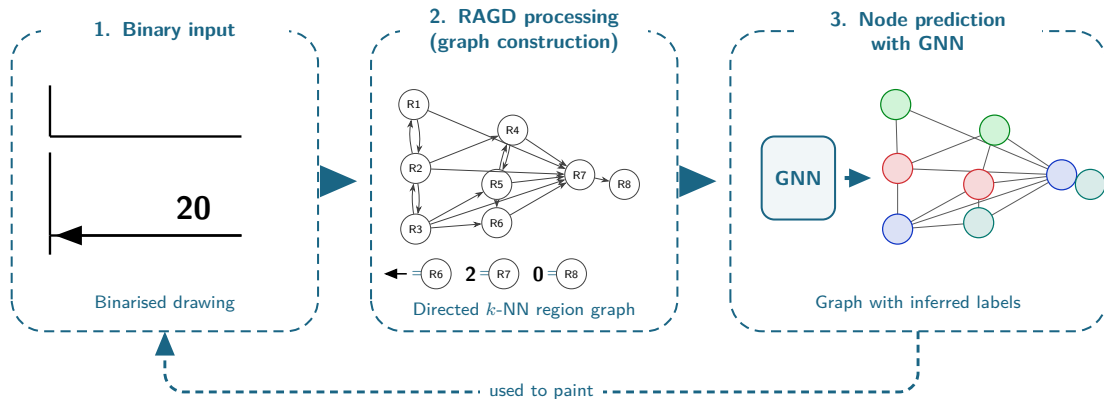
Contributions

- RAGDNet splits into two parts:
 - RAGD, the algorithm that turns a binarised drawing into a region-adjacency graph: regions become nodes, adjacency becomes edges;
 - the network that predicts one class per node of that graph, using a graph neural network.
- A labeled dataset of 2D CAD mechanical drawings, with four foreground classes.

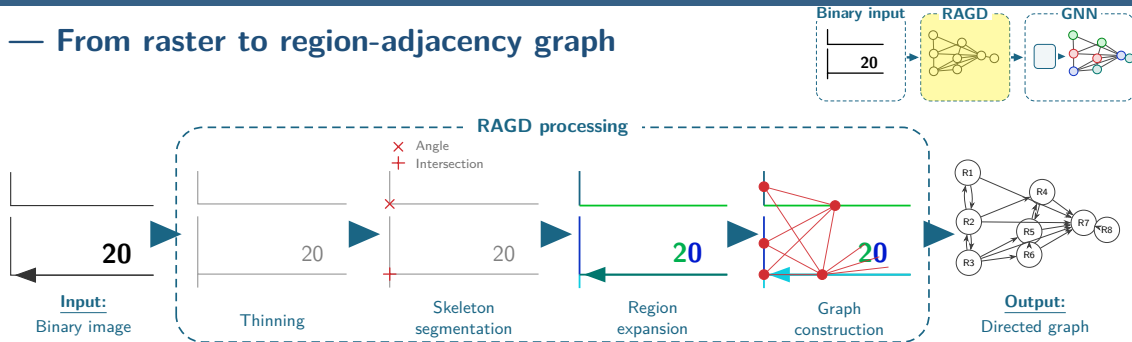
Outline

- 1 Method (RAGDNet)
- 2 Dataset
- 3 Experiments

Pipeline overview (key idea)



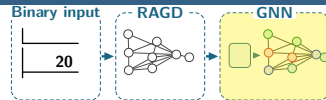
1 — From raster to region-adjacency graph



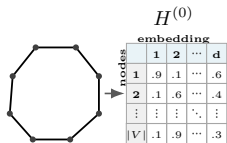
A directed k -NN region graph $G = (V, E)$, one node per stroke region.

- **Parameters:** corner angle range $[\tau_{\min}, \tau_{\max}]$ and k , the neighbours kept per node.
- **Node features:** the region polygon P_i , its simplified outer contour.
- **Edge features:** normalised distance $\hat{d}_{ij}$ and unit direction $\hat{u}_{ij}$ to each of the k nearest regions.

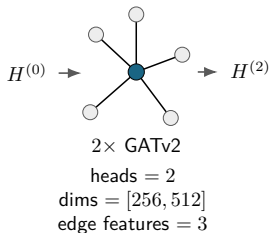
2 — Node prediction with GNN



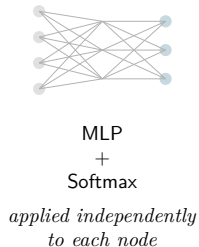
1. Polygon embedding



2. Message passing



3. Node prediction



4. Output

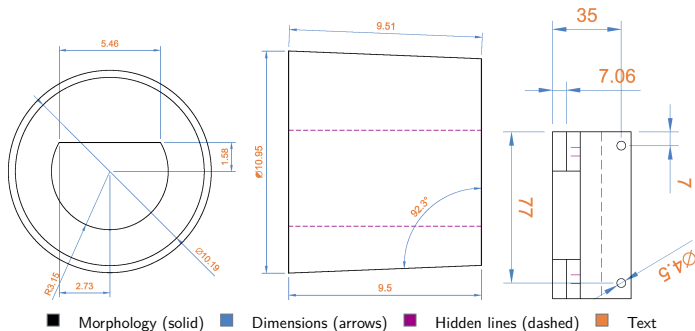
classes		1	2	3	4
nodes	1	0.59	0.13	0.20	0.08
	2	0.12	0.60	0.18	0.10
	3	0.15	0.13	0.53	0.19
	⋮	⋮	⋮	⋮	⋮
	V	0.08	0.16	0.61	0.15

node-wise class scores

C classes

A CAD-derived annotated dataset

- Our dataset consists of 422 drawings.
- They come from a 3D dataset of mechanical parts: Autodesk Fusion flattens each part to 2D and generates the annotations.



Cross-validation results

Method	Morphology	Dimensions	Hidden lines	Text	mIoU
RAGDNet-GAT (ours)	0.90 ± 0.02	0.87 ± 0.01	0.74 ± 0.04	0.96 ± 0.01	0.869 ± 0.019
RAGDNet-MLP	0.87 ± 0.01	0.78 ± 0.03	0.51 ± 0.04	0.88 ± 0.01	0.761 ± 0.007
U-Net	0.94 ± 0.02	0.88 ± 0.03	0.56 ± 0.04	0.93 ± 0.04	0.826 ± 0.030
SegFormer-B0	0.92 ± 0.02	0.83 ± 0.03	0.32 ± 0.15	0.92 ± 0.04	0.747 ± 0.060
SegFormer-B2	0.95 ± 0.02	0.88 ± 0.07	0.56 ± 0.19	0.96 ± 0.03	0.838 ± 0.075

- 5-fold experiments: 80% training, 20% validation.
- Best mIoU: 0.869 vs 0.838.
- Where each wins: a thin stroke gives a pixel model almost nothing to see, but it is still one node on the graph; on the thick classes the pixel models keep the edge, 0.95 vs 0.90.

Model size & inference speed

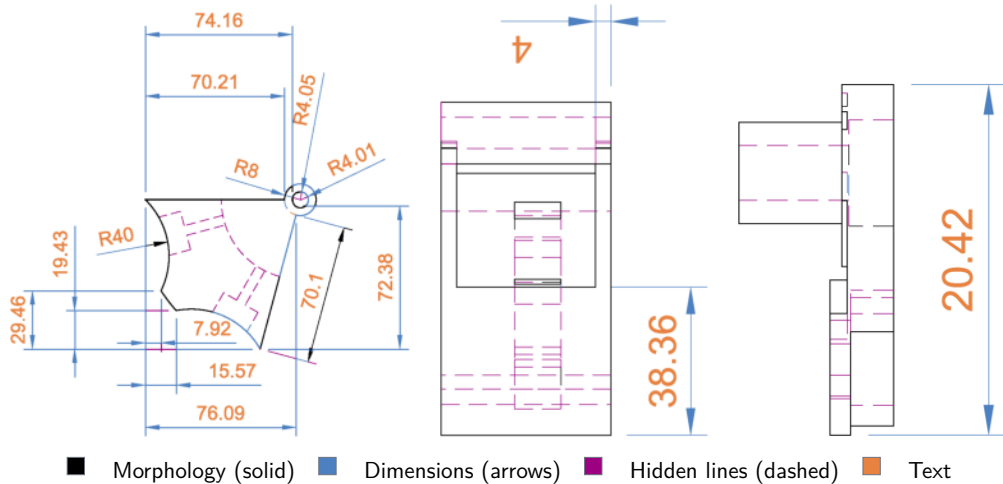
Table: Model size and CPU timing on an Intel Core i7-10610U CPU over 422 drawings.

Method	Params	Pre* (s)	Inf (s)	Total (s)
RAGDNet-GAT	0.881M	11.86±10.90	0.03 ±0.03	11.89±10.90
U-Net	31.0M	–	12.26±5.24	12.26±5.24
SegFormer-B0	3.75M	–	7.29±6.17	7.29 ±6.17
SegFormer-B2	27.3M	–	18.08±13.26	18.08±13.26

* A recent optimisation of the graph construction cuts this preprocessing time by a factor of ≈ 49 , down to 0.26 s, a new total of ≈ 0.29 s.

- Inference with RAGDNet is much faster
- The model is also lightweight

Qualitative results



Conclusion

- We exploit the **sparse, structured nature** of mechanical drawings: RAGDNet shifts the prediction from the pixel grid to a **Region Adjacency Graph**. A dataset comes with it.
- This topological representation achieves **pixel-based accuracy** with a **far more compact** architecture and **faster inference**.
- Three limitations:
 - the dataset covers a **limited set of classes and drawing conventions**, short of industrial diversity;
 - the **multi-stage image-to-graph** construction lets early errors propagate into the final region graph;
 - instance segmentation is not supported yet.

Future work

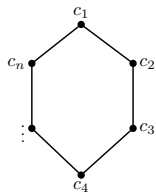
- **Broader coverage**: an industrial partner dataset, re-annotated to our label definitions, extending the taxonomy beyond the four foreground classes.
- **Learnable graph extraction** instead of fixed preprocessing heuristics: a differentiable construction, trainable **end-to-end from raw pixels**.
- **Panoptic segmentation**, to distinguish individual objects within each class.
- Longer term: **legacy scanned drawings**, with image restoration as a pre-processing step, ideally without extensive retraining.

Thank you

Dataset & code: github.com/Alewep/RAGDNet · alexandre.monnier@univ-grenoble-alpes.fr

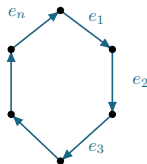
Polygon encoder

1 Polygon



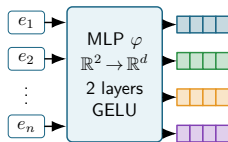
$$p_i = (c_1, \dots, c_n)$$

2 Contour vectors



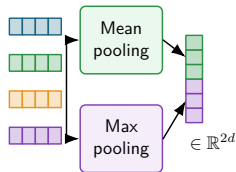
$$e_j = c_{j+1} - c_j$$

3 Shared MLP



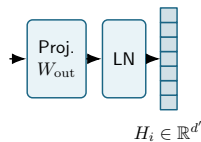
n descriptors
in $\mathbb{R}^d$

4 Aggregation



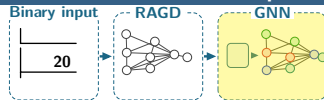
no longer
depends on n

5 Output



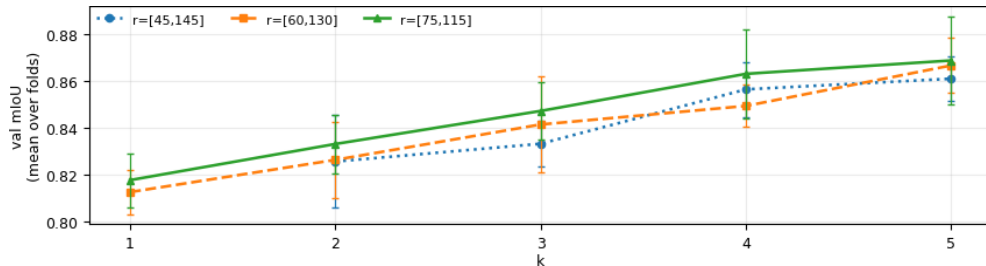
fixed size d'

- Contour vectors $e_j = c_{j+1} - c_j$ pass through a shared MLP.
- Pooling yields one feature vector per region.



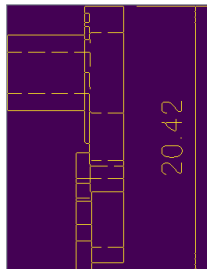
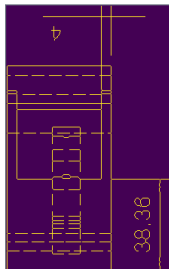
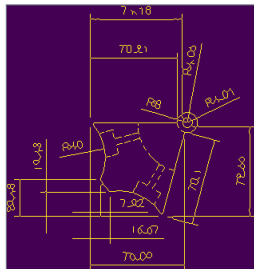
RAGD parameters

Parameter	Tested	Retained
$[\tau_{\min}, \tau_{\max}]$, corner angle range	(75, 115), (60, 130), (45, 145)	$[75^\circ, 115^\circ]$
k , directed k -NN size, $\text{indegree}(v) = k$	$1, \dots, 5$	5



- Mean validation mIoU across folds, one curve per angle range.
- Increasing k helps, with diminishing returns.
- $[75^\circ, 115^\circ]$ is best for small k ; the ranges converge as k grows.

Thinning



- We use [Lee's thinning algorithm](#) (Lee, Kashyap & Chu, 1994).
- It erodes the foreground iteratively, deleting a point only when it is *simple*: its removal changes neither connectivity nor topology.
- The result is a one-pixel-wide, 8-connected skeleton centred in the stroke. The same three drawings as in the qualitative results.

Junction detection

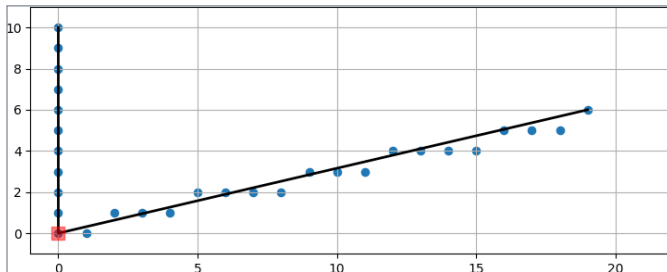
Convolve the skeleton with the 8-neighbour kernel: $D = \text{Skel} * K$ counts the foreground neighbours of each pixel.

$$K = \begin{pmatrix} 1 & 1 & 1 \\ 1 & 0 & 1 \\ 1 & 1 & 1 \end{pmatrix}$$

0 0 0	0 0 0	1 0 0	1 0 1
0 1 1	1 1 1	0 1 1	0 1 0
0 0 0	0 0 0	1 0 0	1 0 1
Endpoint, $D = 1$	Internal, $D = 2$	T-junction, $D = 3$	Cross, $D = 4$

- On a skeleton pixel, D is the **degree** of the branch point.
- $D = 1$ endpoint, $D = 2$ internal pixel, $D \geq 3$ **junction**.
- Cutting at every $D \geq 3$ splits the skeleton into independent branches.

Branch simplification & corners



- A branch is an ordered run of 8-connected skeleton pixels between two junctions, or a junction and an endpoint.
- Pixel quantisation makes raw branches zigzag, so each one is simplified by [Douglas–Peucker \(1973\)](#) with tolerance $\varepsilon \approx 2.12$ px.
- An interior point is a [corner](#) when its turning angle θ_i falls in $[\tau_{\min}, \tau_{\max}]$. The branch is split there.

Where the tolerance ε comes from

ε is not tuned: it is the largest displacement the **pixel grid alone** can produce, so Douglas–Peucker absorbs quantisation noise and nothing else.

- **One axis**: mapping a continuous point to the nearest pixel centre costs at most half a pixel, in x as in y .
- **One point**: the worst case is a diagonal displacement,

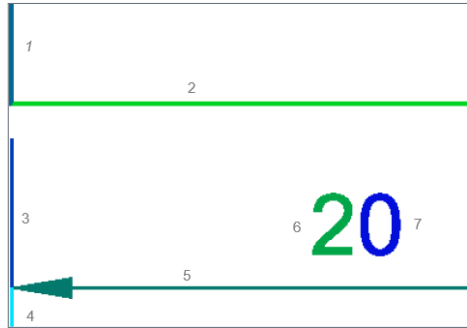
$$\eta = \sqrt{\left(\frac{1}{2}\right)^2 + \left(\frac{1}{2}\right)^2} = \frac{\sqrt{2}}{2} \approx 0.71 \text{ px.}$$

- **Three points**: the criterion measures the distance from a branch point to the segment carried by two others. All three are quantised, each off by up to η , so in the worst case the errors add up:

$$\varepsilon = 3\eta = \frac{3\sqrt{2}}{2} \approx 2.12 \text{ px.}$$

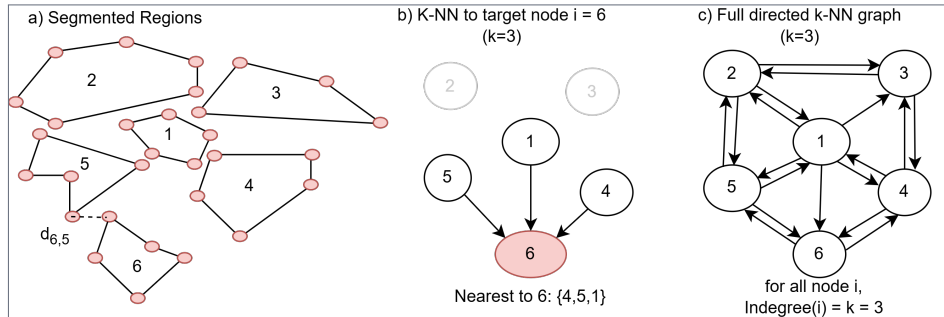
A conservative bound: it kills grid oscillation without touching a real direction change.

Watershed



- **Relief**: the negated distance transform, deepest at the centre of a stroke (left, colour bar $0 \rightarrow -8$).
- The skeleton branches are the **seeds**; basins flood outward until they meet.
- One region per branch (right), and those regions become the nodes of the graph.

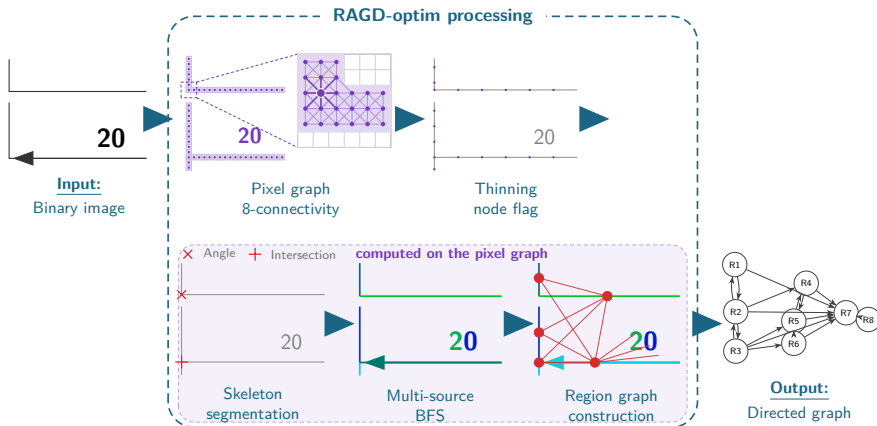
Graph construction



- One **node** per region; its feature is the simplified outer polygon (a).
- For each node i , the k nearest regions each send an edge *into* i (b), so $\text{indegree}(i) = k$ for every node (c).
- **Edge features:** normalised distance $\hat{d}_{ij}$ and unit direction $\hat{u}_{ij}$.

RAGD-optim pipeline

Key idea: compute the graph earlier, to avoid processing the blank pixels that carry no information.



Graph construction cost

Table: Comparison of the mean execution time and standard deviation between RAGD and RAGD Optim.

Step	RAGD (s)	RAGD Optim (s)
Del Peak	3.534 ± 3.742	0.057 ± 0.025
Watershed	6.254 ± 6.033	0.013 ± 0.007
Region Graph	2.945 ± 2.850	0.086 ± 0.072
Skeletonize	0.068 ± 0.050	0.065 ± 0.049
Del Junction	0.027 ± 0.012	0.003 ± 0.001
Pixel Graph	n/a	0.039 ± 0.015
Total	12.827 ± 11.606	0.264 ± 0.117