

International Master in Computer Vision

Fundamentals of machine learning for computer vision

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Machine learning (I)

- **Machine learning**: synthesize a function from a set of selected values (training examples).
- **Simple case**: linear regression (it concerns two-dimensional sample points with one independent variable and one dependent variable, $(x_1, y_1), \dots, (x_N, y_N)$, and finds a linear function that, as accurately as possible (minimizing the mean squared error), predicts the dependent variable values as a function of the independent variable.

Machine learning (II)

- The coefficients of the line or polynomial can be calculated minimizing the mean squared error. They can be used to predict a function for $x \neq x_i$.
- The polynomial coefficients are the learned or trained parameters. For higher polynomial degrees, more trained parameters.
- Other possibilities would be to do interpolation (lineal, spline, etc.) instead of adjustment.

Machine learning (III)

- So, we always need to learn a set of parameters.
- In some cases, the learning of parameters leads to analytical expressions (mathematical formulas like in linear regression).
- If the problem complexity or number of trained parameters increase, more sophisticated and efficient methods may be needed instead of analytical methods.

Classification or regression

- How are **the values of the function** to predict or approximate?
 - **Continuous**: then **regression**
 - **Discrete** (categorical): (integer values with or without order) then **classification**

Supervised and unsupervised learning

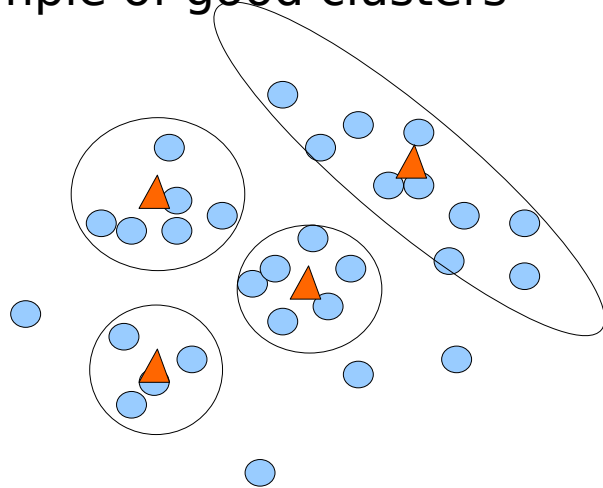
- **Supervised learning**: the value of the function to be predicted is already unknown, like in the linear regression, interpolation or classification.
- **Unsupervised learning**: the objective is to extract information from the data using criteria like similarity or distances, as in the case of **clustering** or **dimensionality reduction**.

Clustering

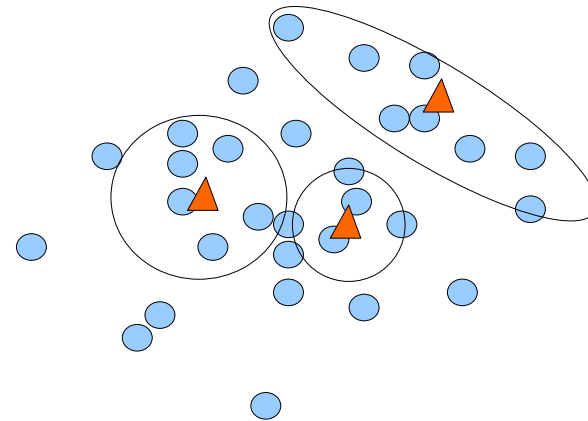
- Objective: to group a set of patterns into clusters (groups of patterns) using a similarity or distance measurement.
- To learn the data structure (the pattern groups).
- Similar patterns (low distances among them) are assigned to the same cluster and used to create a cluster prototype.
- Different patterns (high distances) are assigned to different clusters.
- The threshold to split the distances between similar and different patterns depends of the number of clusters.
- Many clusters: each cluster will have similar patterns.
- Few clusters: each cluster can have very different patterns.

Clustering

Example of good clusters



Example of difficult clustering



- The pattern distribution determines the clustering difficulty.
- The quality of the learned clustering: to what extent do cluster prototypes represent patterns?
- Are there pattern-free (empty) regions between clusters?
- In high dimensions, distances among patterns are very similar.

Clustering

- The results of the clustering are the cluster prototypes. The whole prototype set constitutes a compressed data representation.
- The prototypes should be placed on the regions in the input space with higher pattern density.
- Non-supervised learning: there is no pattern label (discrete nor continuous).
- There is only a distance measurement, which indicates how far two patterns are from each other or from a prototype.

Clustering: k-means

- The number K of clusters that you want to create must be known.
- Set randomly K prototypes $\{\mathbf{p}_k\}_{k=1}^K$ in K patterns.
- For each pattern $\{\mathbf{x}_i\}_{i=1}^N$:
 - Calculate the distance between \mathbf{x}_i and $\{\mathbf{p}_k\}_{k=1}^K$
 - Select the closest prototype.
- For each prototype $\{\mathbf{p}_k\}_{k=1}^K$, update the prototype using only the closest patterns and return the calculation of the prototype.
- Repeat the process until the prototypes do not change significantly or a specified number of times.

K-means algorithm

Data: $K > 1$, $\{\mathbf{x}_i\}_{i=1}^N$, $\varepsilon > 0$

Output: $\{\mathbf{p}_j, E_j\}_{j=1}^K$: prototypes and pattern list for each cluster

$\{\mathbf{p}_k = \mathbf{x}_{s(k)}\}_{k=1}^K$ # $s(k)$: random number in $\{1..N\}$

repeat

$\mathbf{P} = \{\mathbf{p}_k\}_{k=1}^K$; $\{N_k\}_{k=1}^K = 0$; $\{E_k = \emptyset\}_{k=1}^K$

for $i=1:N$

$m = \underset{k=1..K}{\operatorname{argmin}} \{ \|\mathbf{x}_i - \mathbf{p}_k\| \}$

$N_k = N_k + 1$; $E_k = E_k \cup \{i\}$

endfor

for $k=1:K$

$\mathbf{p}_k' = \frac{1}{N_k} \sum_{i \in E_k} \mathbf{x}_i$

endfor

$\mathbf{P}' = \{\mathbf{p}_k'\}_{k=1}^K$; $\{\mathbf{p}_k = \mathbf{p}_k'\}_{k=1}^K$

until $|\mathbf{P}' - \mathbf{P}| < \varepsilon$

Distance measurements between patterns and prototypes

- Euclidean distance: $d(\mathbf{x}_i, \mathbf{x}_j) = \sqrt{\sum_{k=1}^N (x_{ik} - x_{jk})^2}$

- Mahalanobis distance: $d(\mathbf{x}_i, \mathbf{x}_j) = \sqrt{(\mathbf{x}_i - \mathbf{x}_j)^T \Sigma^{-1} (\mathbf{x}_i - \mathbf{x}_j)}$

Σ is the data covariance matrix

- Similarity: $s(\mathbf{x}_i, \mathbf{x}_j) = \frac{\mathbf{x}_i^T \mathbf{x}_j}{\|\mathbf{x}_i\| \|\mathbf{x}_j\|}$

- Tanimoto similarity: $s(\mathbf{x}_i, \mathbf{x}_j) = \frac{\mathbf{x}_i^T \mathbf{x}_j}{\|\mathbf{x}_i\|^2 + \|\mathbf{x}_j\|^2 - \mathbf{x}_i^T \mathbf{x}_j}$

Measures of error and clustering quality (I)

- To minimize the quadratic error between patterns and prototypes:

$$J = \sum_{k=1}^K \sum_{x \in E_k} |\mathbf{x} - \mathbf{p}_k|^2$$

- To maximize the trace of the between-cluster covariance matrix \mathbf{S}_B :

$$\text{tr}(\mathbf{S}_B) = \sum_{k=1}^K N_k |\mathbf{p}_k - \mathbf{p}|^2 \quad \mathbf{p} = \frac{1}{N} \sum_{i=1}^N \mathbf{x}_i$$

- To minimize the determinant of within-cluster cov. matrix \mathbf{S}_W :

$$\det(\mathbf{S}_W) = \det\left(\sum_{k=1}^K \mathbf{S}_k\right) \quad \mathbf{S}_k = \sum_{x \in E_k} (\mathbf{x} - \mathbf{m}_k)(\mathbf{x} - \mathbf{m}_k)^T$$

\mathbf{S}_W : within-cluster covariance
 \mathbf{S}_B : between-cluster covariance
 $\mathbf{S}_T = \mathbf{S}_W + \mathbf{S}_B$: total covariance

Measures of error and cluster quality (II)

- To maximize the trace of $\mathbf{S}_W^{-1}\mathbf{S}_B$:

$$tr(\mathbf{S}_W^{-1}\mathbf{S}_B) = \sum_{i=1}^n \lambda_i$$

λ_i : i -th eigenvalue
of $\mathbf{S}_W^{-1}\mathbf{S}_B$

- To maximize the trace of $\mathbf{S}_T^{-1}\mathbf{S}_B$:

$$tr(\mathbf{S}_T^{-1}\mathbf{S}_B) = \sum_{i=1}^n \frac{1}{1+\lambda_i}$$

- To maximize the determinant ratio between \mathbf{S}_W and \mathbf{S}_T :

$$\frac{|\mathbf{S}_W|}{|\mathbf{S}_T|} = \prod_{i=1}^n \frac{1}{1+\lambda_i}$$

- Post-processing: the clusters of lower quality (for example, high quadratic error) can be splitted to reduce the error.

The number K of clusters

- It is a hyper-parameter which must be set before clustering.
- It can be tuned if there is some quality measure of the clustering.
- Sometimes, it is pre-defined by the problem.
- PCA (Principal Component Analysis) can be used to visualize the data in 2D and to select an appropriated number of clusters.
- Instead of setting K , we can set a maximum distance D :
 - 1) If $d < D$, the pattern is assigned to the cluster and update its prototype.
 - 2) If $d > D$ for all clusters, a new cluster is created. The final number K of clusters depends on data.

Example: image segmentation

- To divide a RGB image in regions with similar appearance.
- The K-means method shows the images with K colors.

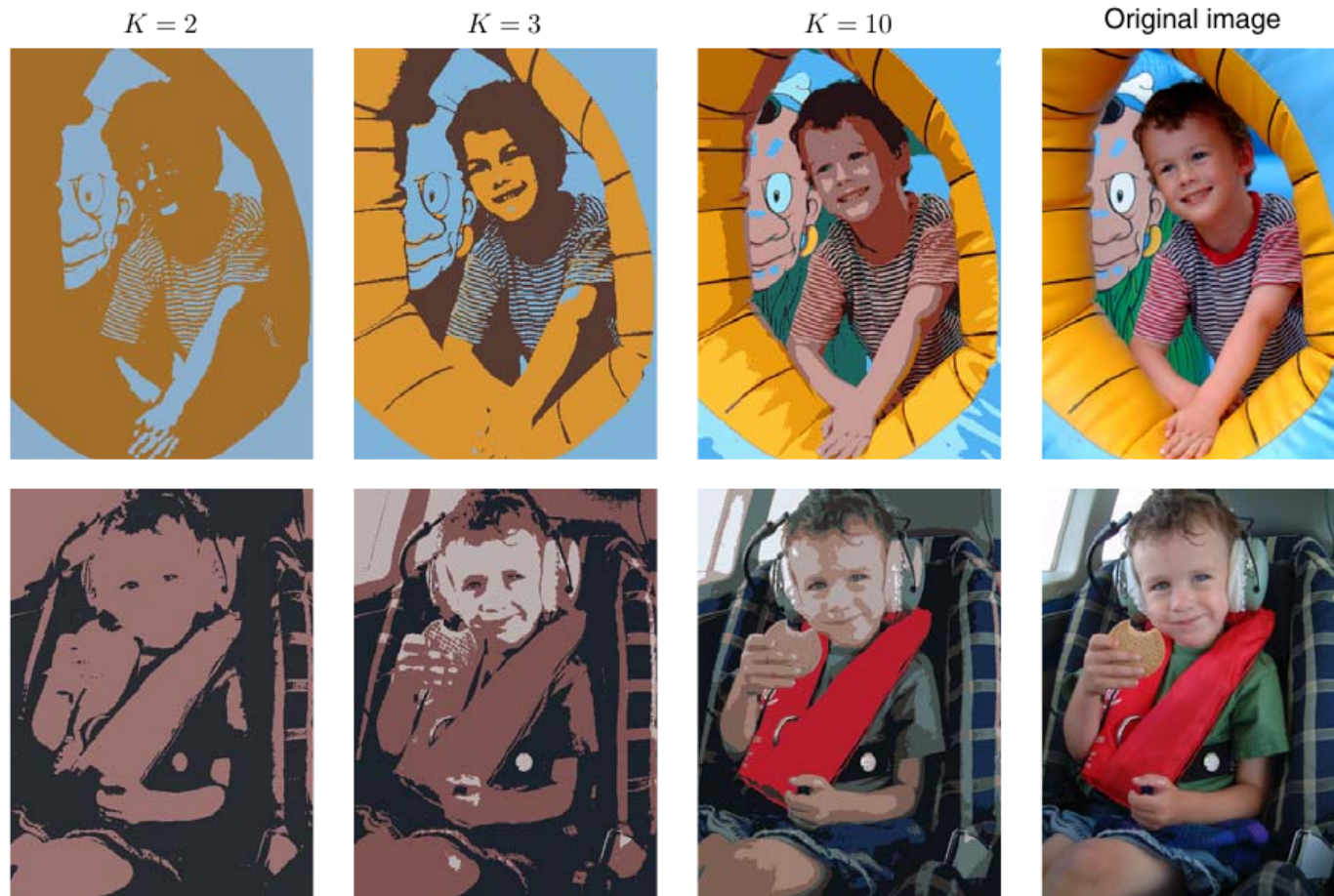


Image segmentation: performance

- **Pixel level evaluation:**
 - Assume that each pixel is a pattern
 - The image segmentation process is a two-class classification problem
 - Class 0 and 1 for background and foreground pixels
 - Use the quality measures for classification:
 - Kappa and accuracy
 - Precision, recall and F1
- The K-means method shows the images with K colors.

Image segmentation: performance

- **Region level evaluation:**
 - Both examples are rather bad

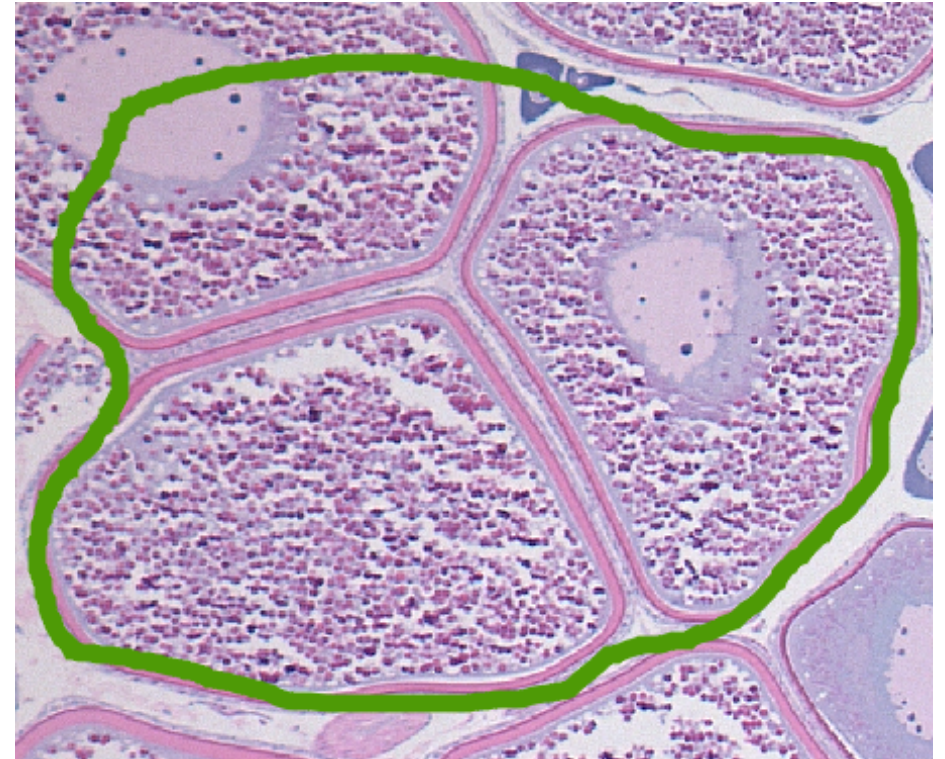


Image segmentation: performance

- **Region level evaluation:**

- Quantify the agreement of expert and algorithm at region level.
- Measure the overlap between the expert and computer to segment objects.
- D^i and A^i the number of segmented and true objects in image I^i .
- P_d^i , number of pixels of recognized object R_d^i , $d=1..D^i$
- P_a^i , number of pixels of true object R_a^i , $a=1..A^i$

Image segmentation: performance

- **Region level evaluation:**

- $O_{da}^i = P_d^i \cap P_a^i$ number of overlapped pixels between R_d^i and R_a^i
- $O_{da}^i = \emptyset$ both regions are not overlapped.
- $O_{da}^i = R_d^i = R_a^i$, the overlapping is complete.

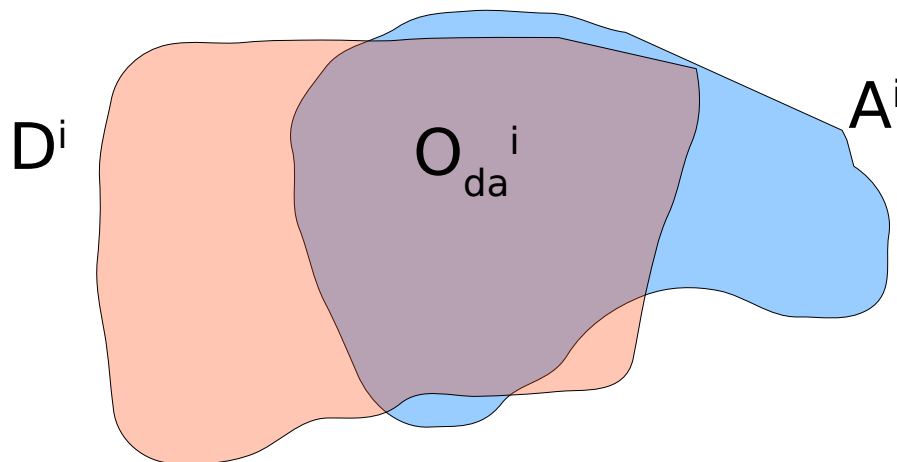
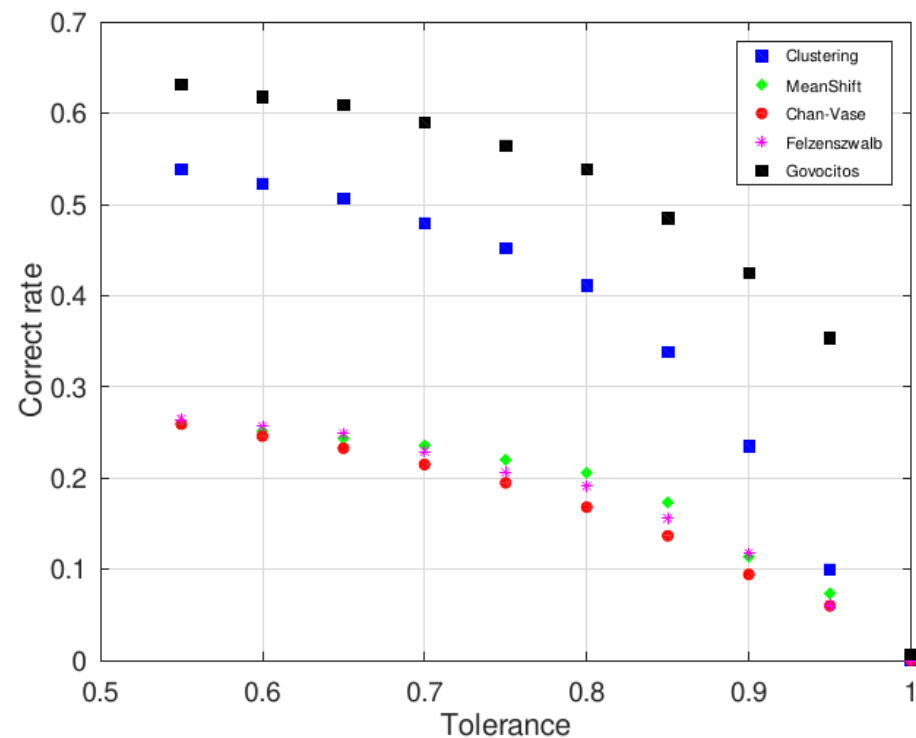
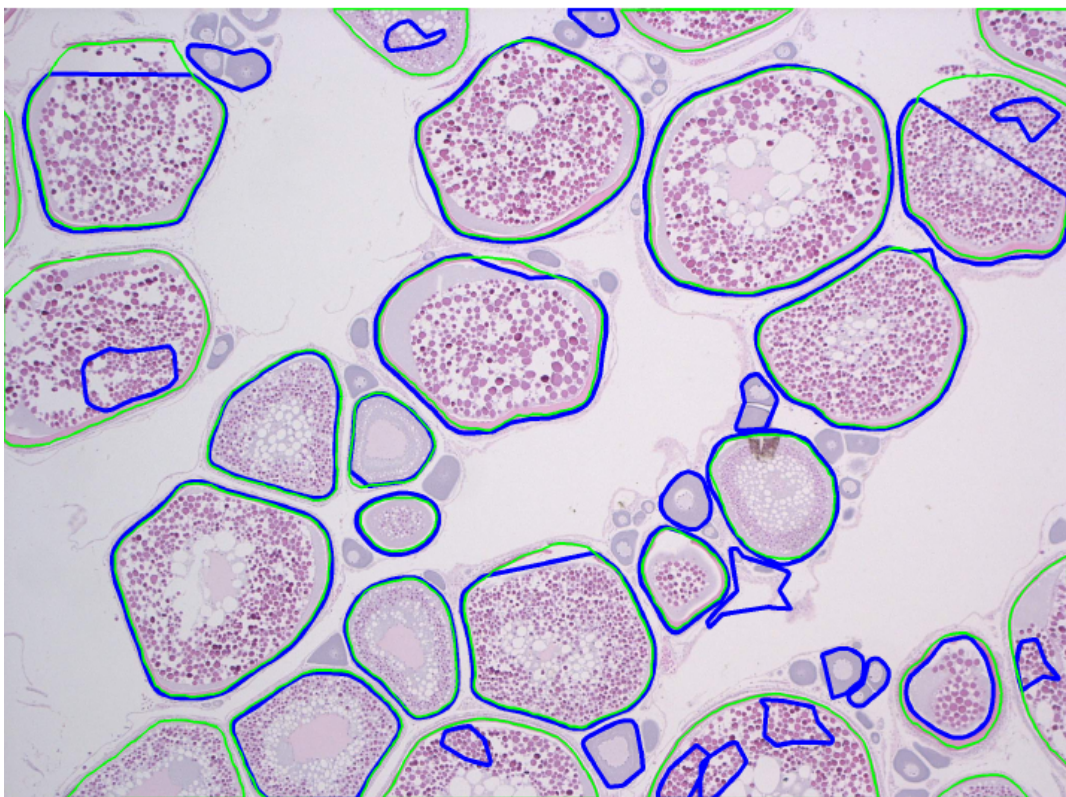


Image segmentation: performance

- **A region is classified in the following types:**
 - Let $0 \leq T \leq 1$ is a strictness criterion
 - **Correct:** a region R_a^i is correctly detected if
$$O_{da}^i \geq P_d^i T$$
 - **Missed:** a region R_a^i that it is not correctly detected is missed.
 - **Noise:** a region R_d^i that does not participate in any instance of correct detection is classified as noise.

Image segmentation: performance

- **Example:** “MSCF: Multi-Scale Canny Filter to Recognize Cells in Microscopic Images”, 2023.
- <https://doi.org/10.3390/su151813693>



Example: image segmentation

- Matlab: function **imsegkmeans()**

K-means clustering based image segmentation

[collapse all in page](#)

Syntax

```
L = imsegkmeans(I,k)
[L,centers] = imsegkmeans(I,k)
L = imsegkmeans(I,k,Name,Value)
```

Description

`L = imsegkmeans(I,k)` segments image `I` into `k` clusters by performing k-means clustering and returns the segmented labeled output in `L`.

[example](#)

`[L,centers] = imsegkmeans(I,k)` also returns the cluster centroid locations, `centers`.

[example](#)

`L = imsegkmeans(I,k,Name,Value)` uses name-value arguments to control aspects of the k-means clustering algorithm.

Example: K-means clustering

- Matlab/Octave: function **kmeans()**

kmeans

K-means clustering

Syntax

```
IDX = kmeans(X,k)
[IDX,C] = kmeans(X,k)
[IDX,C,sumd] = kmeans(X,k)
[IDX,C,sumd,D] = kmeans(X,k)
[...] = kmeans(...,param1,val1,param2,val2,...)
```

Example: image segmentation

- OpenCV: function `kmeans()` (see OpenCV help)

Functions

```
double cv::kmeans (InputArray data, int K, InputOutputArray bestLabels, TermCriteria criteria, int attempts, int flags, OutputArray centers=noArray())  
    Finds centers of clusters and groups input samples around the clusters. More...
```

```
template<typename _Tp, class _EqPredicate >
```

```
int partition (const std::vector< _Tp > &_vec, std::vector< int > &labels, _EqPredicate predicate=_EqPredicate())  
    Splits an element set into equivalency classes. More...
```

Detailed Description

Function Documentation

◆ `kmeans()`

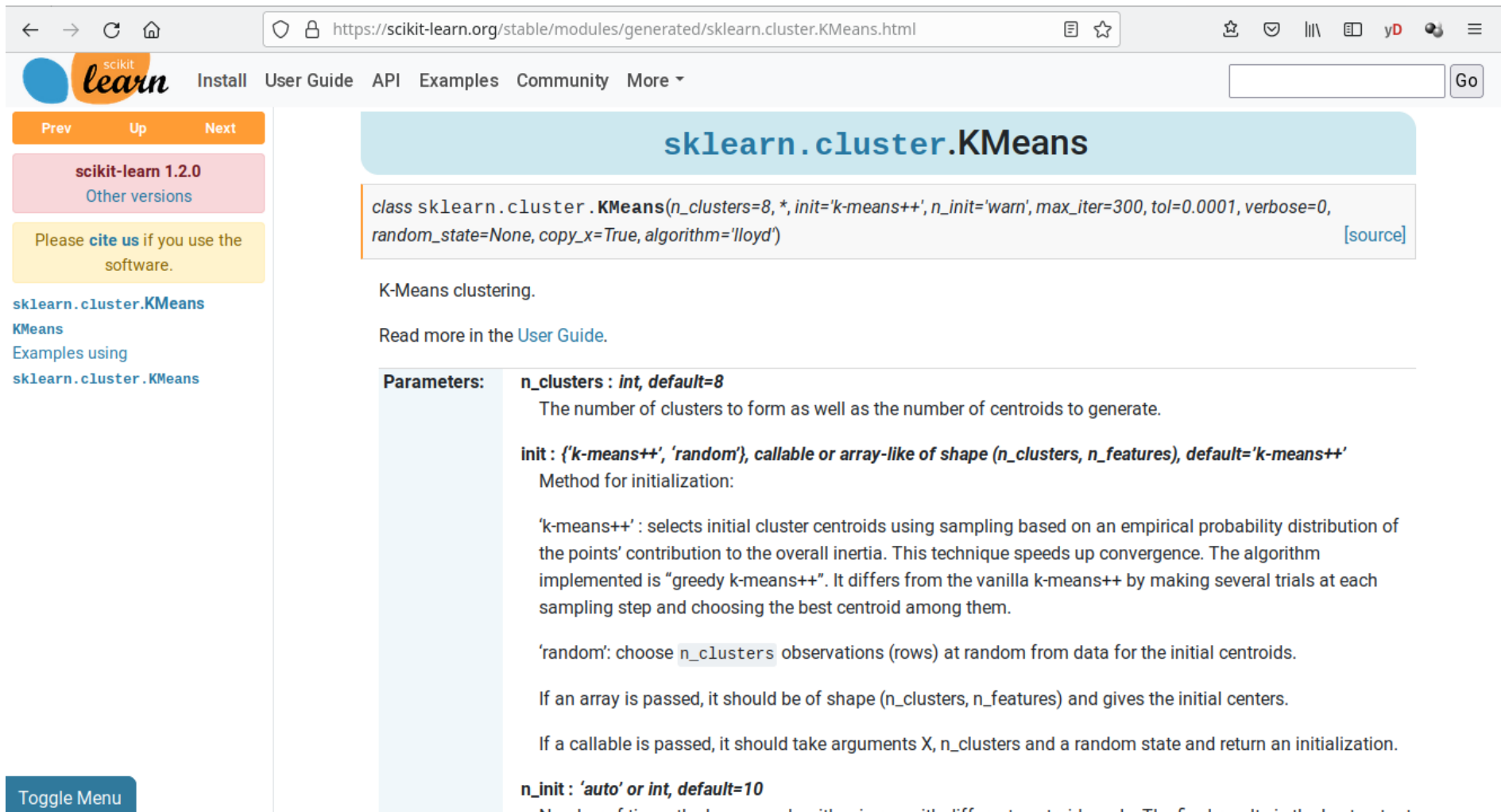
```
double cv::kmeans ( InputArray      data,  
                   int             K,  
                   InputOutputArray bestLabels,  
                   TermCriteria    criteria,  
                   int             attempts,  
                   int             flags,  
                   OutputArray     centers = noArray()  
                   )
```

Python:

```
retval, bestLabels, centers = cv.kmeans( data, K, bestLabels, criteria, attempts, flags[, centers] )
```

Example: image segmentation

- Sklearn: module cluster (python)



The screenshot shows the sklearn.org website documentation for `sklearn.cluster.KMeans`. The page title is `sklearn.cluster.KMeans`. The main content area displays the class signature: `class sklearn.cluster.KMeans(n_clusters=8, *, init='k-means++', n_init='warn', max_iter=300, tol=0.0001, verbose=0, random_state=None, copy_x=True, algorithm='lloyd')` with a [\[source\]](#) link. Below the signature, it states "K-Means clustering." and "Read more in the [User Guide](#)." The "Parameters:" section lists `n_clusters` (int, default=8), `init` (tuple of ('k-means++', 'random'), callable or array-like of shape (n_clusters, n_features), default='k-means++'), and `n_init` (auto or int, default=10). The `init` parameter description includes details about the 'k-means++' and 'random' methods. A "Toggle Menu" button is visible in the bottom left corner.

`sklearn.cluster.KMeans`

```
class sklearn.cluster.KMeans(n_clusters=8, *, init='k-means++', n_init='warn', max_iter=300, tol=0.0001, verbose=0, random_state=None, copy_x=True, algorithm='lloyd')
```

K-Means clustering.

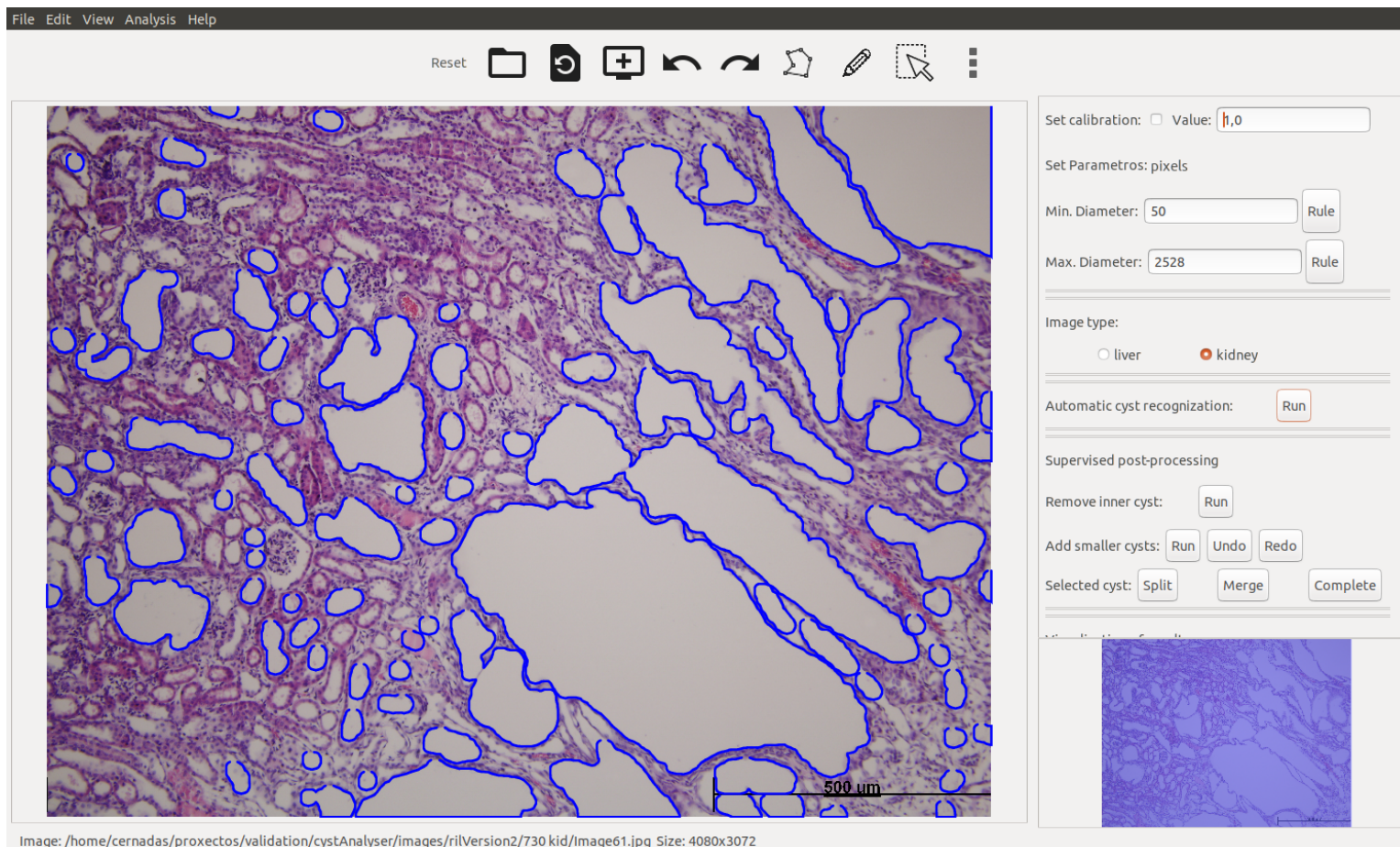
Read more in the [User Guide](#).

Parameters:

- n_clusters** : *int, default=8*
The number of clusters to form as well as the number of centroids to generate.
- init** : *(('k-means++', 'random'), callable or array-like of shape (n_clusters, n_features), default='k-means++'*
Method for initialization:
 - 'k-means++': selects initial cluster centroids using sampling based on an empirical probability distribution of the points' contribution to the overall inertia. This technique speeds up convergence. The algorithm implemented is "greedy k-means++". It differs from the vanilla k-means++ by making several trials at each sampling step and choosing the best centroid among them.
 - 'random': choose `n_clusters` observations (rows) at random from data for the initial centroids.If an array is passed, it should be of shape (n_clusters, n_features) and gives the initial centers.
If a callable is passed, it should take arguments X, n_clusters and a random state and return an initialization.
- n_init** : *'auto' or int, default=10*
Number of times the k-means algorithm is run with different centroid seeds. The final results is the best output

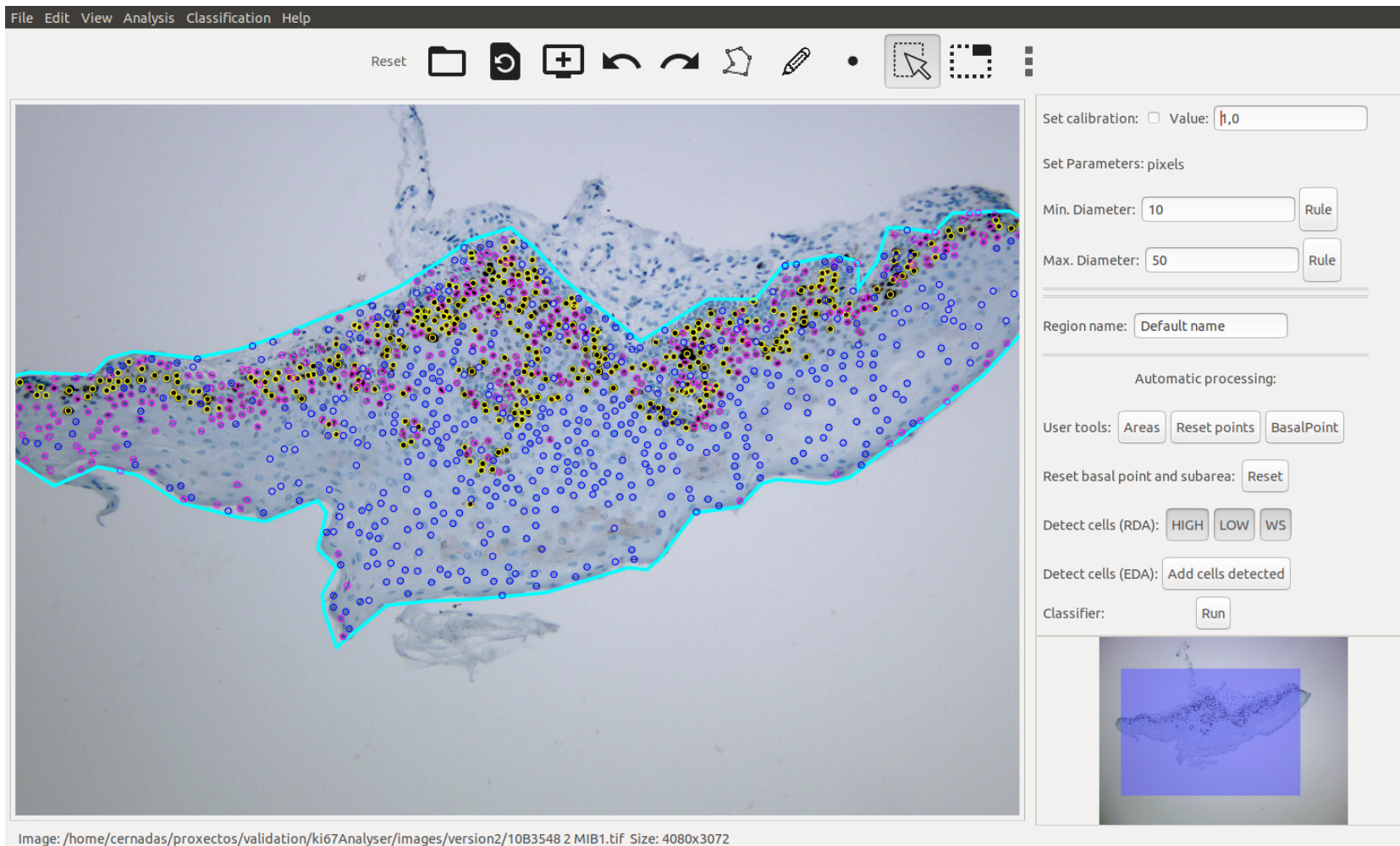
Real examples of clustering

- **CystAnalyser:** [10.1371/journal.pcbi.1008337](https://doi.org/10.1371/journal.pcbi.1008337) or <https://citius.usc.es/transferecia/software/cystanalyser>



Real examples of clustering

- **OralImmunoAnalyser:**



The screenshot displays the OralImmunoAnalyser software interface. The main window shows a histological image of tissue with a cyan boundary and numerous colored dots (yellow, purple, blue) representing detected cells. The interface includes a menu bar (File, Edit, View, Analysis, Classification, Help) and a toolbar with icons for Reset, file operations, zoom, and navigation. On the right, a control panel contains the following settings:

- Set calibration: Value: 1,0
- Set Parameters: pixels
- Min. Diameter: 10
- Max. Diameter: 50
- Region name: Default name
- Automatic processing:
- User tools:
- Reset basal point and subarea:
- Detect cells (RDA):
- Detect cells (EDA):
- Classifier:

At the bottom left, the image path is shown: `Image: /home/cernadas/proxectos/validation/ki67Analyser/images/version2/10B3548 2 MIB1.tif Size: 4080x3072`. A small thumbnail of the processed image is visible in the bottom right corner of the control panel.

Example: image compression

- Image N pixels, 1 pixel=3 colors x 8 bits/color=24 bits: size= $24N$ bits.
- Using K ($\ll N$) prototypes, an input pattern is a pixel $\mathbf{x}_i=(R_i, G_i, B_i)$.
- Clustering: K prototypes $\{\mathbf{p}_k\}_{k=1}^K$, with $\mathbf{p}_k=(R_k, G_k, B_k)$, stored only once.
- Compression: pixel \mathbf{x}_i (8x3 bits) \rightarrow prototype index \mathbf{p}_k of size $\log_2 K$
- Size of compressed image: number N of pixels multiplied by the number of bits required to code the number K of prototypes ($N \cdot \log_2 K$ bits) and the K prototypes (24K bits)
- Compression factor: $\alpha = \frac{T_0}{T_1} = \frac{24 N}{24 K + N \log_2 K} \approx \frac{24}{\log_2 K}$ for large N
- With $K=10$, $\alpha=7.2$