

# K-means Clustering

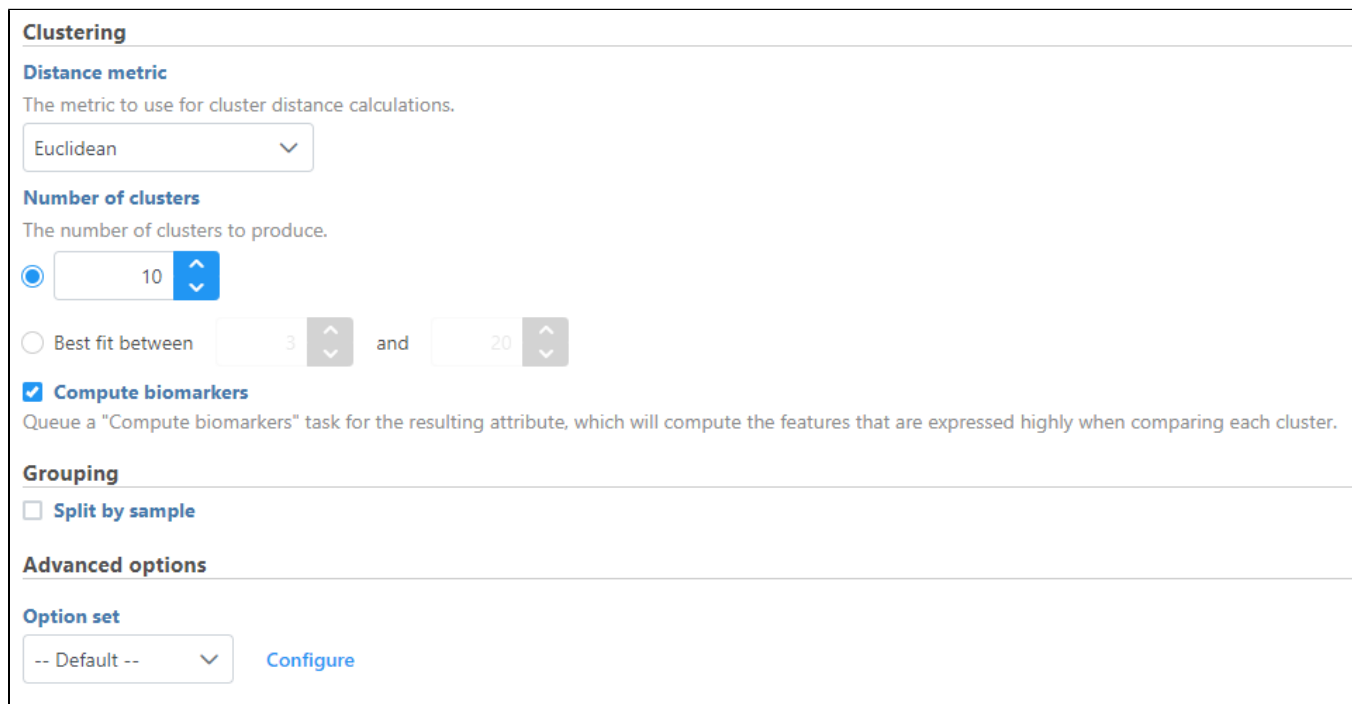
## What is K-means clustering?

K-means clustering is a method for identifying groups of similar observations, i.e. cells or samples. K-means clustering aims to group observations into a pre-determined number of clusters (k) so that each observation belongs to the cluster with the nearest mean. An important aspect of K-means clustering is that it expects clusters to be of similar size (equal variance) and shape (distribution of variance is spherical). The [Compare Clusters task](#) can also be used to help determine the optimal number of K-means clusters.

## Running K-means clustering

We recommend normalizing your data prior to running *K-means clustering*, but the task will run on any counts data node.

- Click the counts data node
- Click the **Exploratory analysis** section of the toolbox
- Click **K-means clustering**
- Configure the parameters
- Click **Finish** to run (Figure 1)



The image shows a configuration dialog for K-means clustering. It has several sections: 'Distance metric' with a dropdown set to 'Euclidean'; 'Number of clusters' with a radio button selected for a fixed number of 10 clusters, and an option for 'Best fit between' 3 and 20 clusters which is unselected; a checked checkbox for 'Compute biomarkers' with a description; a 'Grouping' section with an unchecked 'Split by sample' checkbox; and an 'Advanced options' section with an 'Option set' dropdown set to '-- Default --' and a 'Configure' button.

**Clustering**

**Distance metric**  
The metric to use for cluster distance calculations.  
Euclidean

**Number of clusters**  
The number of clusters to produce.

☒ 10

☐ Best fit between 3 and 20

☒ **Compute biomarkers**  
Queue a "Compute biomarkers" task for the resulting attribute, which will compute the features that are expressed highly when comparing each cluster.

**Grouping**

☐ **Split by sample**

**Advanced options**

**Option set**  
-- Default -- Configure

Figure 1. K-means clustering configuration dialog

K-means clustering produces a *K-means Clusters* result data node; double-click to open the task report which lists the cluster statistics (Figure 2). If *Compute biomarkers* was enabled, top markers will be available by double-clicking the *Biomarkers* result data node. If clustering was run with *Split by sample* enabled on a single cell counts data node, the cluster results table displays the number of clusters found for each sample and clicking the sample name opens the sample-level report.

Cluster statistics		
Total number of clusters 10		
Cluster ↑⇅	Size ↑↓	Size % ↑↓
1	10528	18.22%
2	10190	17.64%
3	8525	14.76%
4	8162	14.13%
5	7065	12.23%
6	4240	7.34%
7	3658	6.33%
8	2798	4.84%
9	2416	4.18%
10	188	0.33%
Task details		

Figure 2. *K-means* clustering task report

## Cluster statistics

The total number of clusters is listed along with the number and percentage of cells in each cluster.

The *K-means Clustering* result data node includes the input values for each gene and adds cluster assignment as a new attribute, *K-means*, for each observation. If the *K-means clusters* data node is visualized by *Scatter plot*, *PCA*, *t-SNE*, or *UMAP*, the plot can be colored by the *K-means* attribute (Figure 3).

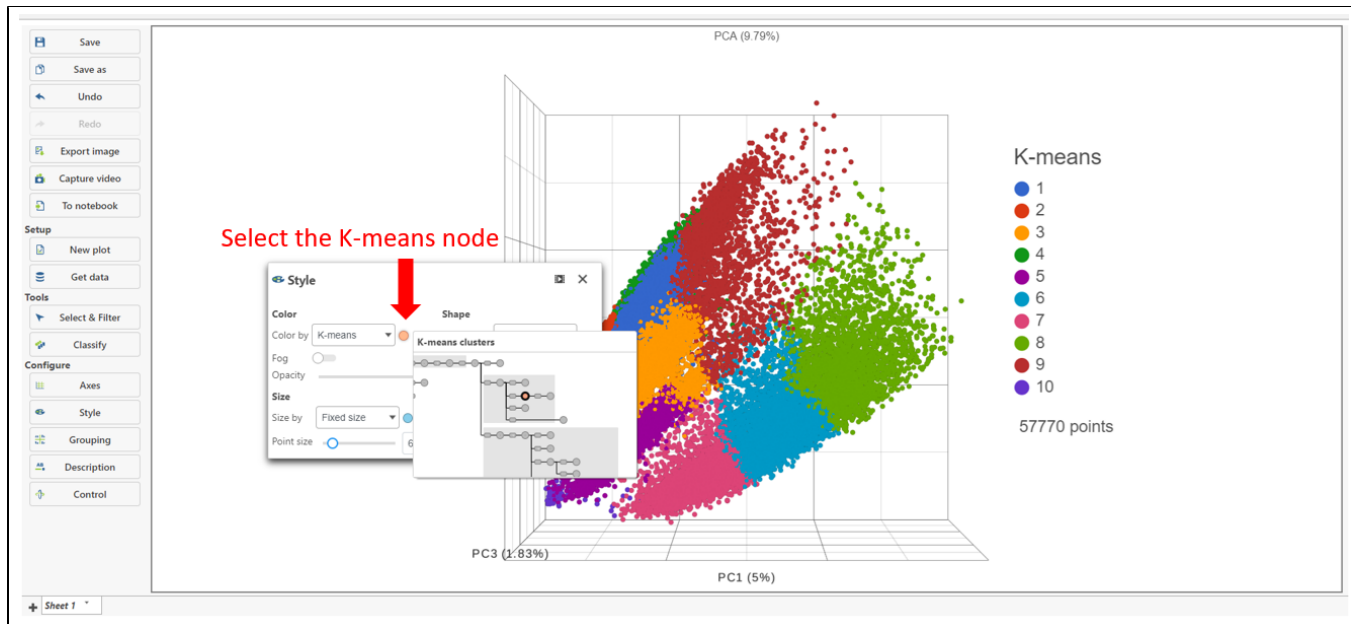


Figure 3. Visualizing K-means cluster results

## Basic K-means clustering parameters

### Distance metric

Choose which distance metric to use for cluster distance calculations. Options include *Euclidean*, *Absolute Value*, *Euclidean Squared*, *Kendall Correlation*, *Max Value*, *Min Value*, *Pearson Correlation*, *Rank Correlation*, *Average Euclidean*, *Shape*, *Cosine*, *Canberra*, *Bray Curtis*, *Tanimoto*, *Pearson Correlation Absolute*, *Rank Correlation Absolute*, and *Kendall Correlation Absolute*. The default is *Euclidean*.

### Number of clusters

Choose between specifying a set number of clusters or a range to test for the best fit number of clusters. The best fit is determined by the number of clusters with the lowest Davies–Bouldin index. The default is set to 10 for a fixed number of clusters. The initial values for the range option are 3 to 20 clusters.

### Compute biomarkers

Choose whether to run the ANOVA test comparing each cluster to all other observations to identify features that have higher values in that cluster. Default is *Enabled*.

### Split cells by sample

This option is present in single cell data. If enabled, K-means clustering will be run separately for each sample. If disabled, K-means clustering will be run on all cells from the input data. Default is set by the *Split single cell by sample* option in the user preference page.

## Advanced K-means clustering parameters

### Random cluster initialization

If enabled, the initial cluster centroids will be selected randomly from among the data points. If disabled, the initial cluster centroids will be selected to optimize distance between clusters. Default is *Disabled*.

### Random seed

This sets the random seed used if *Random cluster initialization* is enabled. Use the same random seed to reproduce results.

### Batch centroid computations

If enabled, all cluster centroids will be recomputed at the end of each iteration. If disabled, each cluster centroid will be recomputed as the members of the cluster change. Default is *Enabled*.

### Max iterations

The maximum number of iterations to perform before setting on a set of clusters. Default is 1000.

# Additional Assistance

If you need additional assistance, please visit [our support page](#) to submit a help ticket or find phone numbers for regional support.



Your Rating: Results: 22 rates