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Missing Values and Outliers; Removing Outliers using k-means Algorithm

Introduction

The aim of this tutorial is to emphasize the importance of data preprocessing. You will learn basic methods how to cope with missing values and outliers.

1 Missing values

The occurrence of missing values has various causes, be it a broken sensor or people who forgot to fill in some questionnaire. In some cases some attributes make no sense for some types of objects (for example: pregnancy for males).

It is very important to identify type of the missing value, because otherwise the analysis could lead to meaningless conclusions. If there are missing values, we often suppose that they are *Missing Completely At Random*. Such an error can be imagined as a random ink blot on the paper with the printed data. Contrarily, *Missing At Random* model works with the assumption, that probability of a missing value depends on some other variable. Thermometer may measure incorrect values with probability $p > 0$ just in case of rain, otherwise it may work well. *Nonignorable Missing Values* are the worst. An example could be a broken thermometer which may not measure temperatures under 0°C .

Imputation

Missing values cannot be always excluded. The simplest way is to replace a missing value by the mean value in case of numeric attributes, or by modus in case of categorical attributes. Another option is to replace the missing value by the respective value of its nearest neighbour.

2 Normalisation

Normalisation is a procedure used to eliminate unequal contribution of attributes. Typical approaches are *min-max normalisation* and *z-score standardisation*.

- **min-max normalisation:** For a numeric attribute x the min-max normalisation is defined as:

$$x \mapsto \frac{x - \min_x}{\max_x - \min_x},$$

where \min_x and \max_x are minimum and maximum values of the attribute

- **Z-score standardisation:** For a numeric attribute x the z-score standardisation is defined as:

$$x \mapsto \frac{x - \hat{\mu}_x}{\hat{\sigma}_x},$$

where $\hat{\mu}_x$ is mean and $\hat{\sigma}_x$ is standard deviation of the attribute

3 Outliers

Outliers usually result from erroneous measurements. If it is not possible to correct such data, it is necessary to remove it from the dataset. It can happen that some outliers are not really outliers, their values are significantly different just by chance. Despite this fact, it can be suitable to remove them and analyse them separately.

Outliers depend on the distribution function. When detecting outliers for categorical data, we assume that outliers occur with extremely low frequency. The situation is more difficult when dealing with numerical attributes. The problem is that in case of large datasets there will be always some points marked as outliers.

Univariate outliers

Let \bar{x} be average value and s be standard deviation of a normal distribution. We say that an observation is an outlier if lies outside of the interval

$$(\bar{x} - ks, \bar{x} + ks),$$

where k is usually 2 or 3.

Multivariate outliers

When dealing with multi-dimensional data, we usually do not use any special assumptions about the distribution function. It is possible to exploit dimensionality reduction methods and try to identify the outliers by visual inspection of a low-dimensional plot of these data. Alternatively, we can use clustering techniques, where we mark the data as outliers when they cannot be meaningfully assigned to any cluster. An example of such an approach is the algorithm COR [?].

Algorithm COR

IN		OUT
X	data	X
m	number of clusters	data without outliers
th	threshold	
R	number of iterations	

```

i ← 0
repeat
  i ← i + 1
  len ← numOfSamples(X)
   $P_{new}, C_{new} \leftarrow k\_means(X, m)$ 
  {BEGIN: outlier removal}
   $X_{new} \leftarrow X$ 
  for cluster  $c \subseteq X_{new}$  do
    if numOfSamples(c) > 1 then
       $s_{max} \leftarrow maxDistantSampleFromCentroid(c)$ 
       $s_{min} \leftarrow minDistantSampleFromCentroid(c)$ 
       $distortion \leftarrow distanceFromCentroid(s_{min}, c) / distanceFromCentroid(s_{max}, c)$ 
      if  $distortion < th$  then
         $P_{new}, C_{new}, X_{new} \leftarrow removeSample(s_{max}, X_{new})$ 
      end if

```

```

    else
       $P_{new}, C_{new}, X_{new} \leftarrow removeCluster(c, X_{new})$ 
       $m \leftarrow m - 1$ 
    end if
  end for
  {END: outlier removal}
   $len_{new} \leftarrow numOfSamples(X_{new})$ 
   $X \leftarrow X_{new}$ 
until  $i > R \wedge len_{new} == len$ 
return( $X$ )

```

Work assignment

1. Replace missing values using nearest neighbours method.
2. Normalize data from the previous point using z-score standardisation. Discuss influence of standardisation on PCA reduction according to the two-dimensional plots showing data projected to the first two principal components.
3. Find and remove outliers in `outlier_data.mat` using algorithm COR.

Expected results

1. Replacing missing values using nearest neighbours method should yield results which look more consistently. A disadvantage of this approach might be undesirable influence of the location of the data centroids.
2. Some of the algorithms for data analysis are sensitive to relative scale of the input variables. If we have two variables with equal variance and they are positively correlated, PCA finds transformation, which rotates the original space by 45° . If we multiply all values of the first variable by 100, then the transformation will respect mostly the first variable. Thus the rotation will be minimal. As a result of this task we can expect visible elimination of the unequal contribution of the variables.
3. The data contain five artificially added outliers. It is possible to detect three of them by univariate methods. The remaining two objects are

visibly outliers, but it is not possible to remove them by univariate approaches.



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