## Quantitative text analysis: Describing and Comparing Text

Blake Miller

MY 459: Quantitative Text Analysis

February 27, 2023

Course website: lse-my459.github.io

- 1. Overview and Fundamentals
- 2. Descriptive Statistical Methods for Text Analysis
- 3. Automated Dictionary Methods

8. Similarity and Clustering Methods

- 4. Machine Learning for Texts5. Supervised Scaling Models for Texts
- 6 Reading Week
- 6. Reading Week
- 7. Unsupervised Models for Scaling Texts
- 9. Topic models
- 10. Word embeddings
- 11. Working with Social Media

### Overview of text as data methods



### Outline

- Describing a single document
- ► Comparing documents
  - ▶ Similarity metrics: cosine, Euclidean, Jaccard, edit distance
  - Clustering methods: k-means clustering, hierarchical clustering

## Quantities for describing a document

Length in characters, words, lines, sentences, paragraphs, pages, sections, chapters, etc.

Word (relative) frequency counts or proportions of words

Lexical diversity (At its simplest) involves measuring a type-to-token ratio (TTR) where unique words are types and the total words are tokens

Readability statistics Use a combination of syllables and sentence length to indicate "readability" in terms of complexity

### Outline

- Describing a single document
- Comparing documents
  - ► Similarity metrics: cosine, Euclidean, Jaccard, edit distance
  - Clustering methods: k-means clustering, hierarchical clustering

## Comparing documents

- The idea is that (weighted) features form a vector for each document, and that these vectors can be judged using metrics of similarity
- A document's vector for us is simply (for us) the row of the document-feature matrix
- ► The question is: how do we measure distance or similarity between the vector representation of two (or more) different documents?

## Characteristics of similarity measures

Let A and B be any two documents in a set and d(A, B) be the distance between A and B.

- 1.  $d(x,y) \ge 0$  (the distance between any two points must be non-negative)
- 2. d(A, B) = 0 iff A = B (the distance between two documents must be zero if and only if the two objects are identical)
- 3. d(A, B) = d(B, A) (distance must be symmetric: A to B is the same distance as from B to A)
- 4.  $d(A, C) \le d(A, B) + d(B, C)$  (the measure must satisfy the triangle inequality)

### Euclidean distance

Between document A and B where j indexes their features, where  $y_{ij}$  is the value for feature j of document i

- Euclidean distance is based on the Pythagorean theorem
- Formula

$$\sqrt{\sum_{j=1}^{j} (y_{Aj} - y_{Bj})^2} \tag{1}$$

In vector notation:

$$\|\mathbf{y}_A - \mathbf{y}_B\| \tag{2}$$

► Can be performed for any number of features *J* (where *J* is the number of columns in of the dfm, same as the number of feature types in the corpus)

## Cosine similarity

- Cosine distance is based on the size of the angle between the vectors
- Formula

$$\frac{\mathbf{y}_A \cdot \mathbf{y}_B}{\|\mathbf{y}_A\| \|\mathbf{y}_B\|} \tag{3}$$

- ▶ The · operator is the dot product, or  $\sum_{i} y_{Aj} y_{Bj}$
- ▶ The  $\|\mathbf{y}_A\|$  is the vector norm of the (vector of) features vector  $\mathbf{y}$  for document A, such that  $\|\mathbf{y}_A\| = \sqrt{\sum_j y_{Aj}^2}$
- Nice property: independent of document length, because it deals only with the angle of the vectors
- Ranges from -1.0 to 1.0 for term frequencies, or 0 to 1.0 for normalized term frequencies (or tf-idf)

### Jaccard coefficient

- ► Similar to the Cosine similarity
- ► Formula

$$\frac{\mathbf{y}_A \cdot \mathbf{y}_B}{\|\mathbf{y}_A\| + \|\mathbf{y}_B\| - \mathbf{y}_A \cdot \mathbf{y}_B} \tag{4}$$

▶ Ranges from 0 to 1.0

### Edit distances

- ► Edit distance refers to the number of operations required to transform one string into another for strings of equal length
- ► Common edit distance: the Levenshtein distance
- Example: the Levenshtein distance between "kitten" and "sitting" is 3
  - ▶ kitten → sitten (substitution of "s" for "k")
  - sitten → sittin (substitution of "i" for "e")
  - Sittin → sitting (insertion of "g" at the end).
- Hamming distance: for two strings of equal length, the Hamming distance is the number of positions at which the corresponding characters are different
- Not common, as at a textual level this is hard to implement and possibly meaningless

### Other uses, extensions

- Used extensively in information retrieval
- ➤ Summary measures of how far apart two texts are but be careful exactly how you define "features"
- Some but not many applications in social sciences to measure substantive similarity — scaling models are generally preferred
- Can be used to generalize or represent features in machine learning, by computing similarities between textual (sub)sequences without extracting the features explicitly (as we will do in a second)

### Outline

- Describing a single document
- Comparing documents
  - ▶ Similarity metrics: cosine, Euclidean, Jaccard, edit distance
  - ► Clustering methods: *k*-means clustering, hierarchical clustering

### The idea of "clusters"

- Essentially: groups of items such that inside a cluster they are very similar to each other, but very different from those outside the cluster
- "unsupervised classification": cluster is not to relate features to classes or latent traits, but rather to estimate membership of distinct groups
- groups are given labels through post-estimation interpretation of their elements
- typically used when we do not and never will know the "true" class labels
- issues:
  - how many clusters?
  - which features to include?
  - how to compute distance is arbitrary

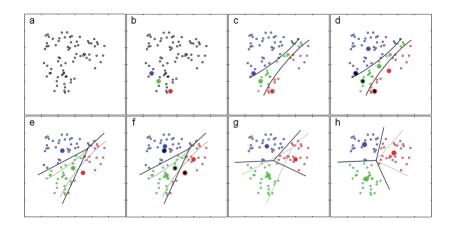
### k-means clustering

- Essence: assign each item to one of k clusters, where the goal is to minimised within-cluster difference and maximize between-cluster differences
- Uses random starting positions and iterates until stable
- k-means clustering treats feature values as coordinates in a multi-dimensional space
- Advantages
  - simplicity
  - highly flexible
  - efficient
- Disadvantages
  - no fixed rules for determining k
  - uses an element of randomness for starting values

### algorithm details

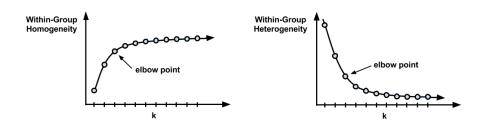
- 1. Choose starting values
  - ▶ assign random positions to k starting values that will serve as the "cluster centres", known as "centroids"; or,
  - assign each feature randomly to one of k classes
- 2. assign each item to the class of the centroid that is "closest"
  - Euclidean distance is most common
  - any others may also be used (Manhattan, Minkowski, Mahalanobis, etc.)
  - (assumes feature vectors are normalized within document)
- 3. update: recompute the cluster centroids as the mean value of the points assigned to that cluster
- 4. repeat reassignment of points and updating centroids
- 5. repeat 2-4 until some stopping condition is satisfied
  - e.g. when no items are reclassified following update of centroids

# k-means clustering illustrated



## choosing the appropriate number of clusters

- very often based on prior information about the number of categories sought
  - for example, you need to cluster people in a class into a fixed number of (like-minded) tutorial groups
- ▶ a (rough!) guideline: set  $k = \sqrt{N/2}$  where N is the number of items to be classified
  - ▶ usually too big: setting *k* to large values will improve within-cluster similarity, but risks *overfitting*
- "elbow plots": fit multiple clusters with different k values, and choose k beyond which are diminishing gains

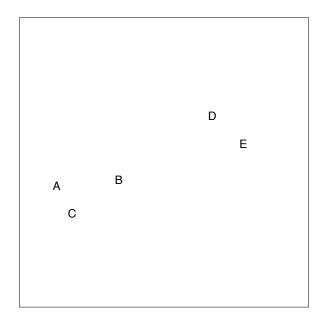


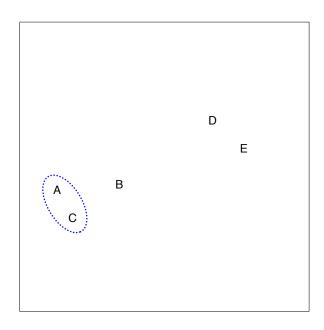
## Other clustering methods: hierarchical clustering

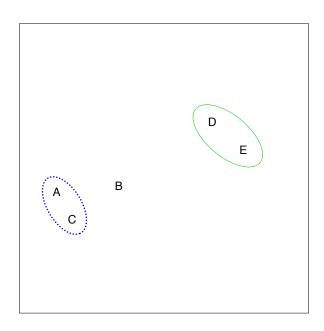
- agglomerative: works from the bottom up to create clusters
- ▶ like k-means, usually involves projection: reducing the features through either selection or projection to a lower-dimensional representation
  - 1. SVD methods, such PCA on a normalised feature matrix
  - usually simple threshold-based truncation is used (keep all but 100 highest frequency or tf-idf terms)
- Can be done at the document level, but also at the feature level (creating clusters of features)

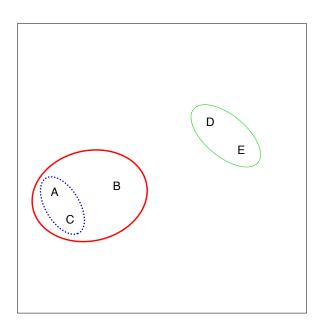
## Hierarchical Clustering

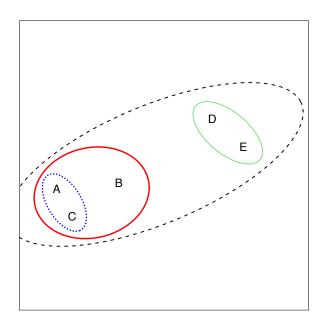
- ► K-means clustering requires us to pre-specify the number of clusters K. This can be a disadvantage (later we discuss strategies for choosing K)
- ► Hierarchical clustering is an alternative approach which does not require that we commit to a particular choice of K.
- ▶ In this section, we describe bottom-up or agglomerative clustering. This is the most common type of hierarchical clustering, and refers to the fact that a dendrogram is built starting from the leaves and combining clusters up to the trunk.







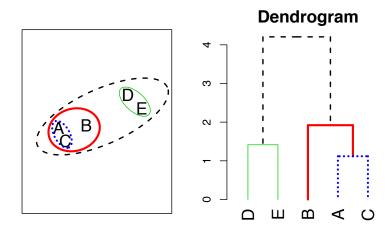




## Hierarchical Clustering Algorithm

### The approach in words:

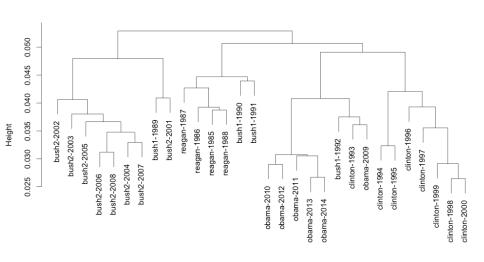
- Start with each point in its own cluster.
- ▶ Identify the closest two clusters and merge them. -Repeat.
- ► Ends when all points are in a single cluster.



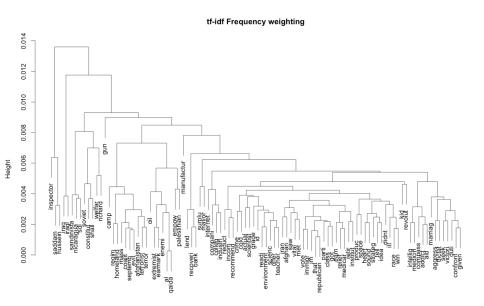
## Hierarchical Clustering Algorithm

- 1. start by considering each item as its own cluster, for n clusters
- 2. calculate the N(N-1)/2 pairwise distances between each of the n clusters, store in a matrix  $D_0$
- 3. find smallest (off-diagonal) distance in  $D_0$ , and merge the items corresponding to the i,j indexes in  $D_0$  into a new "cluster"
- 4. recalculate distance matrix  $D_1$  with new cluster(s). Options for determining the location of a cluster include:
  - centroids (mean)
  - most dissimilar objects
  - Ward's measure(s) based on minimising variance
- 5. repeat 3-4 until a stopping condition is reached
  - ▶ i.e. all items have been merged into a single cluster
- 6. to plot the *dendrograms*, need decisions on ordering, since there are  $2^{(N-1)}$  possible orderings

## Dendrogram: Presidential State of the Union addresses



# Dendrogram: Presidential State of the Union addresses



## pros and cons of hierarchical clustering

#### advantages

- deterministic, unlike k-means
- ▶ no need to decide on k in advance (although can specify as a stopping condition)
- allows hierarchical relations to be examined (usually through dendrograms)

#### disadvantages

- more complex to compute: quadratic in complexity:  $O(n^2)$ 
  - whereas k-means has complexity that is O(n)
- the decision about where to create branches and in what order can be somewhat arbitrary, determined by method of declaring the "distance" to already formed clusters
- for words, tends to identify collocations as base-level clusters (e.g. "saddam" and "hussein")

## Dendrogram: Presidential State of the Union addresses

