Data Mining and Matrix Factorization

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Outline

- Introduction
- Basic Linear Model
 - $\diamond\,$ Factor Analysis
 - \diamond Latent Semantic Indexing
 - $\diamond\,$ Cluster Analysis
 - $\diamond\,$ Receptor Analysis
- Factors versus Clusters
- Centroid Method
 - $\diamond\,$ Modified Centroid
 - $\diamond\,$ Centroid Decomposition
 - $\diamond\,$ Integer Programming on Hypercubes
- Truncated Data
 - $\diamond\,$ Statistical Meaning of SVD
 - $\diamond\,$ Truncation and Lower Rank Approximation
 - $\diamond\,$ Centroid versus SVD

Introduction

- Data mining is about extracting interesting information from raw data.
- What constitutes "information"?
 - $\diamond\,$ Patterns of appearance.
 - $\diamond\,$ Association rules between sets of items.
 - $\diamond\,$ Clustering of the data points.
 - $\diamond\,$ Concepts or categories.
 - $\diamond\,$ Principal components or factors.
 - ♦ ...
- What should be counted as "interesting"?
 - $\diamond\,$ Confidence and support.
 - $\diamond\,$ Information content.
 - \diamond Unexpectedness.
 - ♦ Actionability The ability to suggest concrete and profitable decision-making.

 $\diamond \ldots$

- For different information retrievals, different techniques should be used.
 - ♦ Factors Rank reduction or lower dimension approximation.
 - $\diamond~$ Clusters Centroids or k-means.

♦ ...

From Complexity to Simplicity

• Data analysis:

- ♦ An indispensable task in almost every discipline of science.
- \diamond Search for relationships between a set of externally caused and internal variables.
- \diamond Especially important in this era of information and digital technologies when massive amounts of data are generated at almost all levels of applications.
- Data observed from complex phenomena:
 - \diamond Often represent the integrated result of several interrelated variables acting together.
 - $\diamond\,$ These variables sometimes are less precisely defined.
- What to distinguish which variable is related to which and how the variables are related.

Two Classical Approaches

• Factor analysis:

- \diamond A class of procedures that can help identify and test what *constructs*, or *factors*, might be used to explain the interrelationships among the variables.
- ♦ Each construct itself is a complex image, idea, or theory formed from a number of simpler elements.
- Cluster analysis:
 - ♦ A procedure used to organize information about cases so that relatively *homogenous groups*, or *clusters*, can be formed.
 - ▷ Group members should be highly internally homogenous (members are "similar" to one another in their characteristics) and highly externally heterogenous (members are not "like" members of other clusters).
 - $\triangleright\,$ Need a measurement of similarity or dissimilarity.
 - $\triangleright\,$ Need a decision on how many clusters to keep.
- Either analysis is meant to bring forth the effect of reducing the size of the data table.

Basic Model

- Making observation, data gathering and processing:
 - \diamond Assume ℓ entities and n variable.
 - \diamond Record raw scores that entity *j* received from all variables.
 - ♦ Normalize raw scores to have mean 0 and standard deviation 1 (standardized score).
 - ♦ Let $Y = [y_{ij}] \in \mathbb{R}^{n \times \ell}$ denote the matrix of observed data.
 - $\triangleright y_{ij} = standard \ score \ of \ entity \ j \ on \ variable \ i.$
- Correlation matrix of all *n* variables:

$$R := \frac{1}{\ell} Y Y^T. \tag{1}$$

Linear Relationship (Only an Assumption)

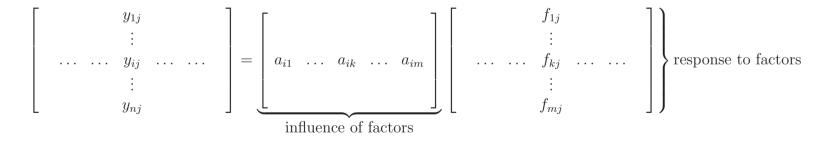
• Assume that y_{ij} is a linearly weighted score of entity j on several factors.

$$Y = AF.$$

• $A = [a_{ik}] \in \mathbb{R}^{n \times m}$ is the factor loading matrix.

 $\diamond a_{ik}$ = the loading of variable *i* on factor *k*, or the influence of factor *k* on variable *i*.

- $F = [f_{kj}] \in \mathbb{R}^{m \times \ell}$ is the factor scoring matrix.
 - $\diamond f_{kj}$ = the score of factor k on entity j, or the response of entity j to factor k.



Example 1: What Factors Affect Students' Academic Performance?

- Y represents the transcripts of ℓ college students (the entities) at the end of freshman year. Each column denote one student's grades on n fixed subjects (the variables), e.g., Calculus, English, Chemistry, and so on.
- A college freshman's academic performance depends on a number of factors including, for instance, family social status, finance, high school GPA, cultural background, and so on.
- Upon entering the college, each student could be asked to fill out a questionnaire inquiring these factors of his/her background. In turn, individual responses to those factors are translated into scores and placed in the corresponding column of the scoring matrix F.
- What is not clear to the educators/administrators is how to choose the factors to compose the questionnaire or how each of the chosen factors would be weighted (the loadings) to reflect the effect on each particular subject.
- In practice, we usually do not have a priori knowledge about the number and character of underlying factors in A. Sometimes we do not even know the factor scores in F.
- Only the data matrix Y is observable.
- Explaining the complex phenomena observed in Y with the help of a minimal number of factors extracted from the data matrix is the primary and most important goal of factor analysis.

Factor Analysis and Matrix Decomposition

- Two additional assumptions:
 - ♦ All sets of factors being considered are uncorrelated with each other.
 - \diamond Similar to Y, the scores in F for each factor are normalized.

$$\frac{1}{\ell}FF^T = I_m. \tag{3}$$

• The correlation matrix R can be expressed directly in terms of the loading matrix A, i.e.,

$$R = AA^T.$$
(4)

- \diamond Factor extraction now becomes a problem of decomposing the correlation matrix R into the product AA^{T} .
- \diamond Would like to use as few factors as possible.

Interpretation of the Loading Matrix A

- a_{i*} = how the data variable *i* is weighted across/influenced by the list of current factors.
 - $\diamond ||a_{i*}||_2^2 = \text{the communality of variable } i.$
 - \triangleright If $||a_{i*}||_2$ is small, variable *i* is of little consequence to the current factors.
- a_{*k} = correlations of the data variables with that particular kth factor.
 - $\diamond ||a_{*k}|| = \text{the significance of factor } k.$
 - ▷ Variables with high factor loadings are more "like" the factor in some sense.
 - \triangleright Variables with zero or near-zero loadings are treated as being unlike the factor.

Tasks to Do in Factor Analysis

- Want to rewrite the loadings of variables over some *newly selected* factors.
 - $\diamond\,$ Fewer factors.
 - $\diamond\,$ Manifest more clearly the correlation between variables and factors.
- Represent the loading of each variable (each row of A) as a single point in the factor space \mathbb{R}^m .
 - $\diamond\,$ What does it mean if these points cluster around a certain direction?
 - $\diamond\,$ How to find the clustering direction?

What Is Going On?

• Suppose the newly selected factors are expressed in terms of columns of the orthogonal matrix

$$V := [\mathbf{v}_1, \dots, \mathbf{v}_m] \in \mathbb{R}^{m \times m}.$$
(5)

- \diamond Determine some appropriate new basis for V.
- \diamond The rewriting of factor loadings with respect to V is mathematically equivalent to a change of basis, i.e., A is now written as B := AV.
- \diamond Because $Y = AF = (AV)(V^TF) = BG$,
 - $\triangleright B = AV$ denotes new factor loadings.
 - $\triangleright G = V^T F$ denotes new factor scores.
- ♦ The correlation matrix $\mathbf{R} = AA^T = BB^T \in \mathbb{R}^{n \times n}$ is independent of factors selected.
 - \triangleright Would like that the significance of factors concentrates on "fewer" columns of B.
 - \triangleright Lower rank approximation of A.
- In the process of defining new factors it is often desirable to retrieve information ...
 - \diamond Directly from the correlation matrix R rather than from any particular loading matrix A, if A is not readily available; or
 - \diamond Approximate A, if A is too large or too expensive.

Example 2: Latent Semantic Indexing

- Indexing matrix $H = [h_{ik}] \in \mathbb{R}^{n \times m}$:
 - $\diamond\,$ Each document is represented by one row in H.
 - $\diamond h_{ik}$ = the *weight* of one particular *term* k in document i.
 - \triangleright Each term could be just one single word or a string of phrases.
 - \triangleright The weight h_{ik} could simply be the number of occurrence of term k in document i.
 - ▷ More elaborate weighting schemes are available and yield better performance.

 $h_{ik} = t_{ik}g_kn_i,$

 \triangleright Each row should be normalized to unit length.

$$documents \rightarrow \begin{bmatrix} h_{1k} & & \\ & \vdots & \\ h_{i1} & \dots & h_{ik} & \dots & h_{im} \\ & & \vdots & \\ & & h_{nk} & \end{bmatrix}$$

Search Similarities

- Queries $\mathbf{q}_j = [q_{1j}, \dots, q_{mj}]^T \in \mathbb{R}^m$:
 - $\diamond q_{kj}$ = the weight of term k in the query j.
- Would like to find documents relevant to given queries.
 - \diamond To measure how the query \mathbf{q}_i matches the documents,
 - \triangleright Calculate the dots product

$$\mathbf{s}_j = H\mathbf{q}_j. \tag{6}$$

 \triangleright Rank the relevance of documents to \mathbf{q}_j according to the *scores* in \mathbf{s}_j .

Comparison of LSI with Linear Model

• Analogies:

| indexing matrix H | \longleftrightarrow | loading matrix A |
|--------------------------|-----------------------|---|
| document i | \longleftrightarrow | variable i |
| term k | \longleftrightarrow | factor k |
| weight h_{ik} | \longleftrightarrow | loading of factor k on variable i |
| one query \mathbf{q}_j | \longleftrightarrow | one column in scoring matrix ${\cal F}$ |
| weights q_{kj} | \longleftrightarrow | scores of entity j on factor k |
| scores in \mathbf{s}_i | \longleftrightarrow | scores in j column of data matrix Y |

• Differences:

- ♦ In LSI, terms/factors are predetermined.
 - \triangleright How are the terms/factors predetermined?
 - \triangleright What is the notion of "orthogonal words"?
 - \triangleright What is the notion of "term/factor reduction"?
- ♦ LSI is not trying to compute factors based on the scores in \mathbf{s}_j , $j = 1, ..., \ell$.
 - ▷ Though, this information may be used as a learning process for selecting terms/factors.
- \diamond LSI emphasizes effective vector-matrix multiplication (6).
 - ▷ Want to represent the indexing matrix and the queries in a more *compact form* so as to facilitate the computation of the scores.

Example 3: Applications of Cluster Analysis

- Split a shuffled deck of cards into two parts (k = 2).
- Classify living organisms into groups based on the shared possession of characteristics (taxonomy and dendrogram).
- AMS 1991 Mathematics Subject Classification
- Switch manufacturing in the telecommunication industry.
 - $\diamond\,$ A cabinet consists of m slots.
 - $\diamond\,$ Each slot may be filled with a selection from r types of boards.
 - \diamond History of past *n* customer orders have been recorded into a matrix $A \in \mathbb{R}^{n \times m}$.
 - \diamond Would like to preassemble q semi-finished cabinet models.
 - \diamond Determine the model configurations and the corresponding customer-to-model assignment of semi-finished cabinets based on A so as to minimize the total number of insertions required to manufacture the entire order.

Example 4: Receptor Model

- An observational technique within the air pollution research community.
- Make use of the ambient data and source profile data to apportion sources or source categories.
- Fundamental principle:
 - $\diamond\,$ Mass conservation can be assumed.
 - ♦ Mass balance analysis can identify and apportion sources of airborne particulate matter in the atmosphere.

- Obtain a large number of chemical constituents such as elemental concentrations in a number of samples.
- Mass balance equation describes the relationships between p sources which contribute m chemical species to n samples.

$$y_{ij} = \sum_{k=1}^{p} a_{ik} f_{kj},\tag{7}$$

- $\diamond y_{ij}$ = the elemental concentration of the *i*th chemical measured in the *j*th sample.
- $\diamond a_{ik}$ = the gravimetric concentration of the *i*th chemical in the *k*th source.
- $\diamond f_{kj}$ = the airborne mass concentration that the kth source has contributed to the *j*th sample.
- Application:
 - \diamond Typically, only values of y_{ij} are observable.
 - \diamond Neither the sources are known nor the compositions of the local particulate emissions are measured.
 - \diamond A critical question is to estimate the number p, the compositions a_{ik} , and the contributions f_{ki} of the sources.
- The source compositions a_{ik} and the source contributions f_{kj} must all be nonnegative. The identification and apportionment, therefore, becomes a nonnegative matrix factorization problem of Y.

Comparison of Cluster Analysis with Factor Analysis

- Consider a document-term matrix $A \in \mathbb{R}^{n \times m}$ with n documents and m terms.
- Analogies:

 $\begin{array}{rcl} (\text{variable } i) &\longleftrightarrow & \text{document } i/\text{case } i \\ (\text{factor } k) &\longleftrightarrow & \text{term } k/\text{characteristic } k \\ (\text{row } h_{i*}) &\longleftrightarrow & \text{samples of document/case } i \text{ by terms/characteristics} \\ (\text{column } h_{*k}) &\longleftrightarrow & \text{samples of term/characteristic } k \text{ by document/cases} \\ && (\text{These samples are not centered nor normalized.}) \end{array}$

• Differences:

♦ Factor analysis seeks to uncover the underlying structure of the set of factors/terms/characteristic.

- ▷ Which original factors are highly correlated to the principal component factors?
- $\triangleright\,$ Seek to reproduce the intercorrelation among the factors.
- ▷ Reduce columns/terms/characteristics.
- \triangleright Generally known as the *R*-mode factor analysis or simply factor analysis.
- ♦ Cluster analysis seeks to uncover the underlying structure of the set of variables/documents/cases.
 - \triangleright Seek the intercorrelation among the variables.
 - \triangleright Reduce rows/documents/cases.
 - \triangleright Cluster analysis is also known as the *Q*-mode factor analysis or the inverse factor analysis.
- Same factor retrieval techniques can be applied to the transpose of the data table to retrieve clusters.

Centroid Method

- Temporarily assuming that a loading matrix $A_1 \in \mathbb{R}^{n \times m}$ is given.
 - \diamond Coordinate axes in \mathbb{R}^m represent a set of *m* abstractly defined factors.
 - ♦ Define a new coordinate system representing the *centroid factors*.
 - \diamond Loadings with respect to the centroid factors can be calculated without the knowledge of A_1 or even the centroid factors. No need to know A_1 a priori.
- Denote each row of A_1 as a point in the factor space \mathbb{R}^m .
 - ♦ The arithmetic mean of these points, the *centroid*, is used to indicate a collective trend of the variables.
 - $\diamond\,$ Generally variables that tend to vary together form clusters.
 - \triangleright Truly uncorrelated variables form no clusters at all.
 - \triangleright If all variables depend on the same factor, then a single cluster should be formed.
 - \triangleright Do not know a priori how many clusters are to be expected.

Compute the Centroid

• The centroid of these n variables:

$$\mathbf{c}_1 := \frac{A_1^T \mathbf{1}_n}{n} = \left[\frac{\sum_{i=1}^n a_{i1}}{n}, \dots, \frac{\sum_{i=1}^n a_{im}}{n}\right]^T.$$
(8)

♦ First *centroid factor*:

$$\mathbf{v}_1 := \frac{\mathbf{c}_1}{\|\mathbf{c}_1\|}.\tag{9}$$

 $\diamond\,$ New loadings of variables with respect to ${\bf v}_1 :$

$$\mathbf{b}_{1} = A_{1}\mathbf{v}_{1} = A_{1}\frac{A_{1}^{T}\mathbf{1}_{n}}{\|A_{1}^{T}\mathbf{1}_{n}\|} = \frac{R_{1}\mathbf{1}_{n}}{\sqrt{\mathbf{1}_{n}^{T}R_{1}\mathbf{1}_{n}}}.$$
(10)

 \triangleright Loading vector \mathbf{b}_1 is extracted directly from R_1 . No reference to A_1 or \mathbf{v}_1 is needed.

- Remove the influence \mathbf{v}_1 from the the loading matrix:
 - ♦ Orthogonally reduced loading matrix:

$$A_2 := A_1 - A_1 \mathbf{v}_1 \mathbf{v}_1^T. \tag{11}$$

 $\diamond A_2$ inherits most of the loading information of the original A_1 except for the loadings along the direction \mathbf{v}_1 .

Update the Loading Matrix

• The product moment of A_2 can be computed via

$$\begin{aligned} \mathbf{R}_{2} &= A_{2}A_{2}^{T} &= \left(A_{1} - A_{1}\mathbf{v}_{1}\mathbf{v}_{1}^{T}\right)\left(A_{1}^{T} - \mathbf{v}_{1}\mathbf{v}_{1}^{T}A_{1}^{T}\right) \\ &= A_{1}A_{1}^{T} - A_{1}\mathbf{v}_{1}\mathbf{v}_{1}^{T}A_{1}^{T} \\ &= R_{1} - \frac{R_{1}\mathbf{1}_{n}\mathbf{1}_{n}^{T}R_{1}}{\mathbf{1}_{n}^{T}R_{1}\mathbf{1}_{n}}, \end{aligned}$$

- \diamond No explicit reference to A_2 is needed.
- $\diamond R_2$ is exactly one rank less than R_1 (Wedderburn formula).
- Repeat the procedure to extract the next centroid factor for A_2 , to introduce the next reduced loading matrix, and so on.
 - \diamond No! It cannot be done.
 - \diamond The procedure cannot be repeated because $A_2^T \mathbf{1}_n = \mathbf{0}_m$.
 - \diamond The centroid of A_2 is residing squarely at the origin of \mathbb{R}^m . No factor is retrieved.
 - \diamond Really bad?

Modified Centroid Factor

• Heuristic reasons for modification:

- \diamond Signs of loadings indicates positive or negative linear correlation between the variable and the factor \mathbf{v}_1 . Either sign is fine.
- \diamond Asymmetrically distributed points in \mathbb{R}^m identifies the centroid factor more easily.
 - $\triangleright \|\mathbf{c}_1\|$ measures the *eccentricity* of the system of variables with respect to the origin.
 - \triangleright The farther \mathbf{c}_1 is away from the origin, the more variables are qualitatively scattered in a general area surrounding \mathbf{c}_1 .
 - \triangleright The larger $||\mathbf{c}_1||$ is, the better an essential factor \mathbf{v}_1 represents.
- Replacing one particular variable by its negative does not cause trouble in the identification of an essential factor.
 - \diamond Would change the sign of certain rows, if that helps to bring out the eccentricity.
- Observe the relationship:

$$\mathbf{l}_{n}^{T}R\mathbf{1}_{n} = \|A_{1}^{T}\mathbf{1}_{n}\|^{2} = n^{2}\|\mathbf{c}_{1}\|^{2}.$$
(12)

 $\diamond\,$ The problem is now changed to solving the integer programming problem

$$\max_{|\mathbf{z}|=1} \mathbf{z}^T R_1 \mathbf{z},\tag{13}$$

- $\triangleright |\mathbf{z}| = 1$ means components of the column vector \mathbf{z} are either 1 or -1.
- \triangleright There are only 2^n many sign vectors for a fixed n.

Ready to Go!

• Modified centroid and factor:

$$\mathbf{c}_1 := \frac{A_1^T \mathbf{z}_1}{n},\tag{14}$$

$$\mathbf{v}_1 := \frac{A_1^{\mathsf{T}} \mathbf{z}_1}{\|A_1^T \mathbf{z}_1\|},\tag{15}$$

- $\diamond \mathbf{z}_1$ is the optimizer of (13).
- Centroid value of A_1 :

$$\mu_1 := \frac{1}{n} \max_{|\mathbf{z}|=1} \, \mathbf{z}^T R_1 \mathbf{z}. \tag{16}$$

- Updating information:
 - \diamond New loading: $\mathbf{b}_1 = A_1 \mathbf{v}_1$ and can be computed by

$$\mathbf{b}_1 = A_1 \mathbf{v}_1 = \frac{R_1 \mathbf{z}_1}{\sqrt{\mathbf{z}_1^T R_1 \mathbf{z}_1}}.$$
(17)

 \diamond New product moment R_2 :

$$R_2 = R_1 - \frac{R_1 \mathbf{z}_1 \mathbf{z}_1^T R_1}{\mathbf{z}_1^T R_1 \mathbf{z}_1}.$$
(18)

 $\diamond\,$ Repeat the procedure.

Centroid Decomposition

- Each application of this centroid factor retrieval reduces the rank of the loading matrix by one.
 - \diamond The procedure therefore has to come to a stop in finitely many steps.
 - $\diamond\,$ With the recurrence

$$A_{i} = A_{i-1} - A_{i-1} \mathbf{v}_{i-1} \mathbf{v}_{i-1}^{T}, \quad i = 2, \dots, \gamma,$$
(19)

we may write

$$A = A_1 = \mathbf{b}_{\gamma} \mathbf{v}_{\gamma}^T + \ldots + \mathbf{b}_1 \mathbf{v}_1^T, \tag{20}$$

- \triangleright **v**_i = the modified centroid factor of A_i .
- $\triangleright \gamma = \operatorname{rank} \operatorname{of} A_1.$

$$\triangleright$$
 b_i = A_i**v**_i.

- \diamond This is called a *centroid decomposition* of A.
 - \triangleright Closely related to the singular value decomposition (SVD).
- The modified centroid vectors (and factors) are mutually orthogonal, even though they are not explicitly calculated.

$$\mathbf{c}_{1}^{T}\mathbf{c}_{2} = \frac{1}{n^{2}} \left(\mathbf{z}_{1}^{T}A_{1} \right) \left(A_{2}^{T}\mathbf{z}_{2} \right) = \frac{1}{n^{2}} \left(\mathbf{z}_{1}^{T}A_{1} \right) \left[A_{1}^{T} \left(\mathbf{z}_{2} - \frac{\mathbf{z}_{1}^{T}R_{1}\mathbf{z}_{2}}{\mathbf{z}_{1}^{T}R_{1}\mathbf{z}_{1}} \mathbf{z}_{1} \right) \right] = 0.$$

- When $\|\mathbf{b}_i\|$ is small, the factor \mathbf{v}_i is less significant.
 - $\diamond\,$ Less significant factors can be discarded.
 - \diamond Closely related to the truncated SVD (TSVD).

Topology of *n***-dimensional Hypercubes**

- To perform the centroid decomposition, a sequence of integer programming problems must be solved.
 - \diamond The feasible set consists of 2^n sign vectors.
 - $\diamond\,$ An exhaustive search would be expensive.
- Representing hypercubes:
 - \diamond Identifying -1 as 0 and keeping 1 as 1, a unique binary tag can be assigned to each sign vector.
 - \diamond Each binary tag translated into a unique integer between 0 and $2^n 1$ provides a natural ordering of the sign vectors.
 - Each sign vector as one node connected only to those sign vectors whose binary tags differ from its own by exactly one bit (Hamming metric 1).
 - \diamond The set of 2^n sign vectors can be identified as an *n*-dimensional hypercube.

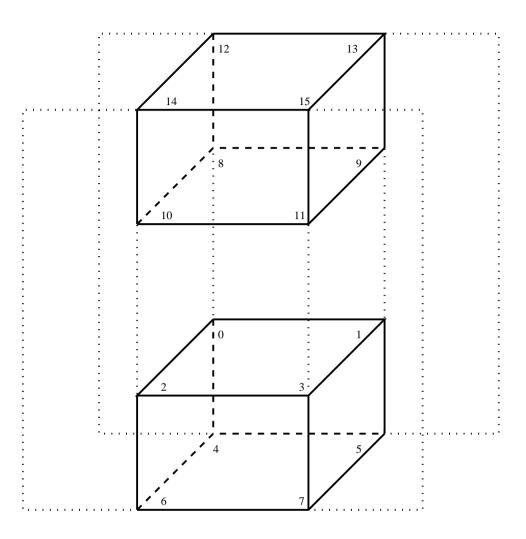


Figure 1: Topology of a 4-dimensional hypercube

- Each *n*-dimensional hypercube consists of two (n-1)-dimensional sub-hypercubes.
 - $\diamond\,$ One sub-hypercube is simply a $bit\ reversal$ of the other.
 - $\diamond\,$ The objective values z^TRz therefore always appear in pairs.

Integer Programming on Hypercubes

- Write $R = [r_{ij}] = P + diag(diag(R)).$
 - $\diamond \mathbf{z}^T R \mathbf{z} = \mathbf{z}^T P \mathbf{z} + \sum_{i=1}^n r_{ii}.$
 - $\diamond\,$ Suffice to maximize

$$f(\mathbf{z}) := \mathbf{z}^T P \mathbf{z}$$

with $|\mathbf{z}| = 1$.

- Classical centroid method:
 - $\diamond\,$ Given any sign vector $\mathbf{z}^{(0)}$ and machine zero threshold $\epsilon.$
 - $\diamond \text{ Define } \mathbf{w}^{(0)} := P \mathbf{z}^{(0)}.$
 - $\diamond\,$ Repeat the following steps for $i=0,1,\ldots$:
 - 1. If $\operatorname{sgn}(\mathbf{w}_k^{(i)}) = \operatorname{sgn}(\mathbf{z}_k^{(i)})$ for all $k = 1, \dots, n$, then stop; otherwise, choose k so that $|\mathbf{w}_k^{(i)}| > \epsilon$ and is the largest among all $|\mathbf{w}_j^{(i)}|$'s where $\operatorname{sgn}(\mathbf{w}_j^{(i)}) \neq \operatorname{sgn}(\mathbf{z}_j^{(i)})$.
 - 2. Define $\mathbf{z}^{(i+1)}$ by simply changing the sign of $\mathbf{z}_k^{(i)}$.
 - 3. Define $\mathbf{w}^{(i+1)} := \mathbf{w}^{(i)} + 2\text{sgn}(\mathbf{z}_k^{(i+1)})P(:,k).$

• Main ideas:

♦ At each given node on the hypercube, check through its neighboring nodes and move to the node with highest bit.

♦ At most one bit is changed in each cycle.

$$\mathbf{z}^{(i+1)} := \mathbf{z}^{(i)} - 2\mathrm{sgn}(\mathbf{z}_k^{(i)})\mathbf{e}_k$$

 \diamond The sequence $\{f(\mathbf{z}^{(i)})\}$ generated by the centroid method from any starting value $\mathbf{z}^{(0)}$ is finite and increasing.

$$f(\mathbf{z}^{(i+1)}) = \left(\mathbf{z}^{(i)} - 2\operatorname{sgn}(\mathbf{z}_k^{(i)})\mathbf{e}_k\right)^T P\left(\mathbf{z}^{(i)} - 2\operatorname{sgn}(\mathbf{z}_k^{(i)})\mathbf{e}_k\right)$$
$$= f(\mathbf{z}^{(i)}) - 4\operatorname{sgn}(\mathbf{z}_k^{(i)})(\mathbf{e}_k^T P \mathbf{z}^{(i)})$$
$$= f(\mathbf{z}^{(i)}) - 4\operatorname{sgn}(\mathbf{z}_k^{(i)})\mathbf{w}_k^{(i)}.$$

♦ The centroid method is a steepest ascent method along the nodes of the hypercube.

• Cost:

 $\diamond\,$ It takes at most n iterations to locate a maximum.

 \diamond Tje expected number of iterations required for convergence is n/2.

Statistical Meaning of TSVD

- Let $\mathcal{X} \in \mathbb{R}^n$ denote a random column vector.
 - $\diamond \ C := \mathcal{E}[(\mathcal{X} \mathcal{E}[\mathcal{X}])(\mathcal{X} \mathcal{E}[\mathcal{X}])^T] \in \mathbb{R}^{n \times n} \text{ is defined as the covariance matrix of } \mathcal{X}.$
 - ♦ In practice, often only random samples are available. A data matrix $X = [x_{ij}] \in \mathbb{R}^{n \times m}$ is collected where each column represents one sample of \mathcal{X} .
 - \diamond Sample mean and sample covariance matrix approximate the true mean and the true covariance matrix of the random variable \mathcal{X} , if m is large enough.
- Assume that C has a spectral decomposition

$$cov(\mathcal{X}) = \sum_{j=1}^{n} \lambda_j \mathbf{u}_j \mathbf{u}_j^T.$$

- $\diamond \mathbf{u}_1, \ldots, \mathbf{u}_p$ are deterministic and form an orthonormal basis for \mathbb{R}^n .
- $\diamond\,$ The random column vector ${\mathcal X}$ can be expressed as

$$\mathcal{X} = \sum_{j=1}^{n} (\mathbf{u}_{j}^{T} \mathcal{X}) \mathbf{u}_{j}$$

- \diamond Each coefficient $\alpha_i := \mathcal{X}^T u_i$ itself is a random variable.
- Properties of α :

$$\mathcal{E}[\alpha] = U^T \mathcal{E}[\mathcal{X}],$$

$$cov(\alpha) = \text{diag}\{\lambda_1, \dots, \lambda_n\}.$$

Meaning of Truncation

- Since \mathbf{u}_j , j = 1, ..., n, are deterministic, stochastic properties of \mathcal{X} are caused only by the stochastic properties of coefficients α_j , j = 1, ..., n.
 - \diamond The randomness of \mathcal{X} is due to the randomness of α .
 - $\diamond\,$ Variance measures the unpredictability of a random variable.
 - \diamond Random variables $\alpha_j, j = 1, \ldots, n$, are mutually stochastically independent.
- The larger the eigenvalue λ_j is, the larger the variance of α_j is and, hence, the more randomness it contributes.
 - \diamond Those coefficients with larger variances and the corresponding directions are the more important components in representing the stochastic nature of \mathcal{X} .
 - \diamond Rank the importance of corresponding eigenvectors \mathbf{u}_i as essential components for the variable \mathcal{X} according to the magnitude of λ_i .
 - ◊ If truncation is necessary, those eigenvectors corresponding to smaller variances should be thrown away first.

Lower Dimensional Minimum-Variance Approximation

• Given a random vector $\mathcal{X} \in \mathbb{R}^n$ with mean zero, let its covariance matrix be spectrally decomposed as

$$cov(\mathcal{X}) = \sum_{j=1}^{n} \lambda_j \mathbf{u}_j \mathbf{u}_j^T$$

Then among all unbiased variables restricted to any r-dimensional subspaces in \mathbb{R}^n , the random variable

$$\hat{\mathcal{X}} := \sum_{j=1}^{r} (\mathbf{u}_{j}^{T} \mathcal{X}) \mathbf{u}_{j}$$
(21)

is the best linear minimum-variance estimate of \mathcal{X} in the sense that $\mathcal{E}[\|\mathcal{X} - \hat{\mathcal{X}}\|^2]$ is minimized.

Truncation in Sample Space

- The distribution of a random variable is often simulated by a collection of ℓ random samples.
 - \diamond Samples are recorded in a $n \times \ell$ matrix X.
 - \diamond Each column of X represents one random sample of the underlying random (column vector) variable $\mathcal{X} \in \mathbb{R}^n$.
 - \diamond When ℓ is large enough, many of the stochastic properties of \mathcal{X} can be recouped from X.
- How to retrieve a sample data matrix from X to represent the minimum-variance approximation $\hat{\mathcal{X}}$ of \mathcal{X} ?
 - $\diamond\,$ Sample covariance:

$$R = \frac{XX^T}{\ell}.$$

♦ Spectral decomposition of sample variance:

$$R = \sum_{i=1}^{n} \mu_i \mathbf{u}_i \mathbf{u}_i^T.$$
 (22)

 \diamond Best low dimensional minimum-variance estimate $\hat{\mathcal{X}}$ to \mathcal{X} :

$$\hat{X} := \sum_{j=1}^{r} \mathbf{u}_j(\mathbf{u}_j^T X).$$
(23)

- The low dimension estimate $\hat{\mathcal{X}}$ to the (continuous) random variable \mathcal{X} is ntranslated into a low rank approximation \hat{X} to the (discrete) random sample matrix X.
- The singular value decomposition of X:

$$X = U\Sigma V^T = \sum_{i=1}^n \sigma_i \mathbf{u}_i \mathbf{v}_i^T$$
(24)

- \diamond Share the same eigenvectors of R as its left singular vectors, i.e., $U = [\mathbf{u}_1, \dots, \mathbf{u}_n]$.
- \diamond Singular values $\sigma_i = \sqrt{\ell \mu_i}$ are ranked in the same ordering as eigenvalues $\mu_i, i = 1, \dots n$.
- ♦ The notion of the truncated singular value decomposition of X is simply the partial sum $\sum_{i=1}^{r} \sigma_i \boldsymbol{u}_i \boldsymbol{v}_i^T$.
- The TSVD of a give data matrix X representing random samples of an unknown random variable \mathcal{X} has a statistical meaning.
 - \diamond The truncated rank-*r* SVD represents random samples of the best minimum-variance linear estimate $\hat{\mathcal{X}}$ to \mathcal{X} among all possible *r*-dimensional subspaces.

Centroid Decomposition versus SVD

• Observe that that

$$\lambda_{1}(R_{1}) = (\mathbf{u}_{1}(R_{1}))^{T} R_{1} \mathbf{u}_{1}(R_{1}) = \max_{\|\mathbf{u}\|=1} \mathbf{u}^{T} R_{1} \mathbf{u} = \max_{\|\mathbf{u}\|=1} \|A_{1}^{T} \mathbf{u}\|^{2}$$

$$\geq \mu_{1} = \frac{1}{n} \mathbf{z}_{1}^{T} R_{1} \mathbf{z}_{1} = \frac{1}{n} \max_{\|\mathbf{z}\|=1} \mathbf{z}^{T} R_{1} \mathbf{z} = \frac{1}{n} \max_{\|\mathbf{z}\|=1} \|A_{1}^{T} \mathbf{z}\|^{2}, \qquad (25)$$

where \mathbf{z}_1 is the sign vector defining the first modified centroid.

- \diamond The sign vector \mathbf{z}_1 and the centroid value μ_1 is *minicking* the left singular vector \mathbf{u}_1 and the square of the singular value λ_1 of A_1 , respectively.
- ♦ Geometric meaning of the variational formulation.

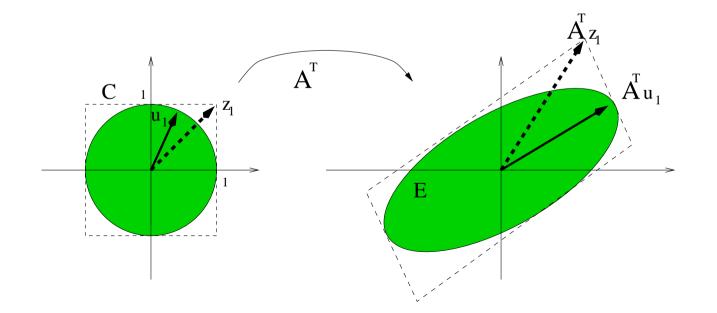


Figure 2: Comparison of geometric meanings of \mathbf{z}_1 and $\mathbf{u}_1(R_1)$ when n = 2.

| Centroid Decomposition | Singular value Decomposition |
|--|---|
| $\mu_1 = \frac{1}{n} \max_{ \mathbf{z} =1} \mathbf{z}^T R_1 \mathbf{z}$ (centroid value) | $\lambda_1 = \max_{\ \mathbf{x}\ =1} \mathbf{x}^T R_1 \mathbf{x}$ (eigenvalue) |
| $\mathbf{z}_1 = \arg \max_{ \mathbf{z} =1} \mathbf{z}^T R_1 \mathbf{z}$ (sign vector for modified centroid) | $\mathbf{u}_1 = \arg \max_{\ \mathbf{x}\ =1} \mathbf{x}^T R_1 \mathbf{x}$ (left singular vector) |
| easy to obtain \mathbf{z}_1 in $O(n)$ steps (tranverse hypercube) | not easy to obtain \mathbf{u}_1 via iterations (nonlinear iteration) |
| $\mathbf{v}_1 = \frac{A_1^T \mathbf{z}_1}{\sqrt{n\mu_1}}$ (centroid factor) | $\hat{\mathbf{v}}_1 = \frac{A_1^T \mathbf{u}_1}{\sqrt{\lambda_1}}$ (right singular vector) |
| $\gamma_1 = \ A_1 \mathbf{v}_1\ $ (significance) | $\sigma_1 = \sqrt{\lambda_1} = \ A_1 \hat{\mathbf{v}}_1\ $ (largest singular value) |
| $b_1 = A_1 \mathbf{v}_1$ (loading vector) | $\sigma_1 \mathbf{u}_1 = A_1 \hat{\mathbf{v}}_1$ (internal relation) |
| $A_1 = \sum b_i \mathbf{v}_i^T$ (centroid decomposition) | $A_1 = \sum \sigma_i \mathbf{u}_i \hat{\mathbf{v}}_i^T$ (singular value decomposition) |
| $R = \sum_{i} b_{i} b_{i}^{T} = \sum_{i} \gamma_{i}^{2} \frac{b_{i}}{\ b_{i}\ } \left(\frac{b_{i}}{\ b_{i}\ }\right)^{T}$ (factor decomposition) | $R = \sum \lambda_i \mathbf{u}_i \mathbf{u}_i^T = \sum \sigma_i^2 \mathbf{u}_i \mathbf{u}_i^T$ (spectral decomposition) |
| $R_{2} = R_{1} - \frac{R_{1}\mathbf{z}_{1}\mathbf{z}_{1}^{T}R_{1}}{\mathbf{z}_{1}^{T}R_{1}\mathbf{z}_{1}} = R_{1} - \gamma_{1}^{2}\frac{b_{1}}{\ b_{1}\ } \left(\frac{b_{1}}{\ b_{1}\ }\right)^{T}$ (rank reduction) | $\overline{R}_2 = R_1 - \frac{R_1 \mathbf{u}_1 \mathbf{u}_1^T R_1}{\mathbf{u}_1^T R_1 \mathbf{u}_1} = R_1 - \lambda_1 \mathbf{u}_1 \mathbf{u}_1^T$ (rank reduction) |

Table 1: Comparison of centroid decomposition and singular value decomposition.

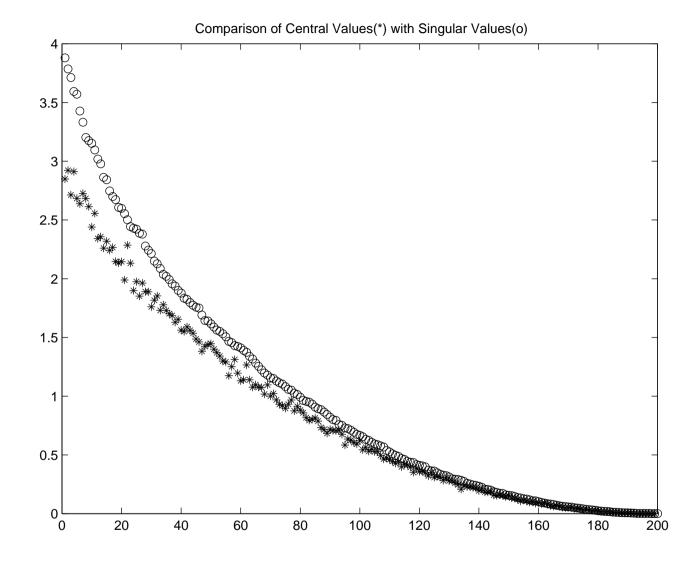


Figure 3: Comparison of centroid values and singular values for correlation matrix of n = 200.

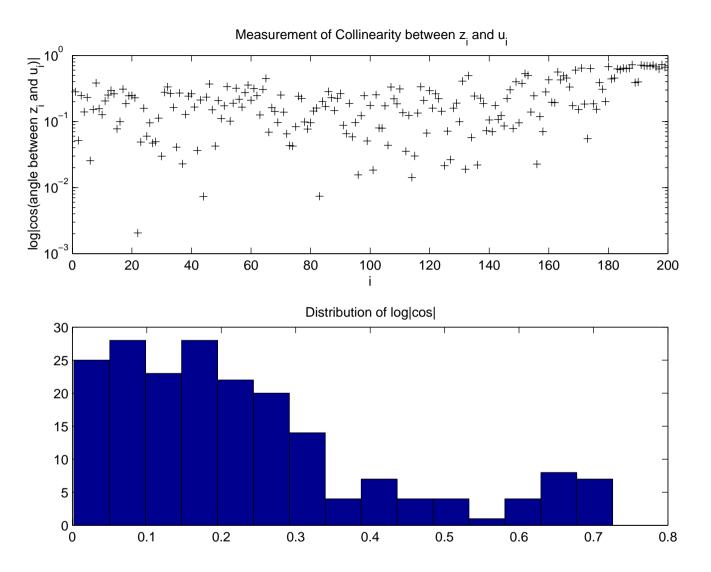


Figure 4: Degree of Alignment between \mathbf{z}_i and u_i .

Conclusion

- We try to clarify the notions of objectives in data mining.
- We compare similarities and differences between factor analysis and cluster analysis.
- The centroid method is cast as an $\mathcal{O}(n)$ -step optimization problem on a hypercube.
- Centroid decomposition is a cheaper simulator of the SVD.
- We offer the insight explaining why, how, and when a low rank approximation makes sensible approximation to the original matrix.
- We show empirically that the centroid decomposition provides a measurement of second order statistical information of the original data.
- The information of significance of a loading vector provides a decision-making on when principal factors have been found.