## **Singular Value Decomposition**

An application to Big Data

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# Singular Value Decomposition

#### Theorem

Given a matrix  $A \in \mathbb{R}^{m \times n}$ , it can always be found a decomposition such that

$$A = U\Sigma V^T \tag{1}$$

where  $U \in \mathbb{R}^{m \times m}$ ,  $V \in \mathbb{R}^{n \times n}$  and  $\Sigma \in \mathbb{R}^{m \times n}$ .

U and V are two orthogonal matrices and  $\Sigma$  is a diagonal matrix, namely:

$$(\Sigma)_{ij} = \begin{cases} \sigma_i, & i = j \\ 0, & i \neq j \end{cases}$$

where  $\sigma_1 \geq \sigma_2 \geq \ldots \geq \sigma_p \geq 0$ ,  $p = \min\{m, n\}$ .

The non-zero entries of  $\Sigma$ , denoted by  $\sigma_i$ , are called *singular values*.

They are arrenged in a nonincreasing order by convention.

The column vectors  $u_i$  of U are called *left singular vectors* and those  $v_i$  of V are called *right singular vectors*.

Since in general  $m \neq n$ , we have:

$$A = \sum_{i=1}^{p} \boldsymbol{u_i} \sigma_i \boldsymbol{v_i}^T$$

#### **Theorem**

*If for some* r *such that*  $1 \le r < p$  *we have* 

$$\sigma_1 \ge \ldots \ge \sigma_r > \sigma_{r+1} = \ldots = \sigma_p = 0$$

then

- rank(A) = r
- $A = \sum_{i=1}^{r} \boldsymbol{u_i} \sigma_i \boldsymbol{v_i}^T$

It means that all other p-r dimensions of matrix A are linear combinations of the first r.

#### Lower rank approximation

Let  $A \in \mathbb{R}^{m \times n}$  be a matrix whose rank is  $\operatorname{rank}(A) = r$ .

If for a fixed integer value k < r we define

$$A_k = \sum_{i=1}^k \sigma_i \boldsymbol{u_i} \boldsymbol{v_i}^T \tag{2}$$

and

$$\mathcal{B} = \left\{ B \in \mathbb{R}^{m \times n} : \operatorname{rank}(B) = k \right\}$$

then

$$\min_{B \in \mathcal{B}} \|A - B\|_2 = \|A - A_k\|_2 = \sigma_{k+1}$$

This result tell us that  $A_k$  represents the best approximation (considering the *spectral norm*) of rank k of matrix A.

### Singular values computation

To compute the singular values, consider the transponse of  $\cal A$  given its decomposition:

$$A^T = (U\Sigma V^T)^T = V\Sigma^T U^T$$

The symmetric matrix  $A^TA$  is equal to:

$$A^T A = (V \Sigma^T U^T)(U \Sigma V^T) = V \Sigma^T \Sigma V^T$$

Furthermore, this equation can be written as:

$$A^TAV = V\Sigma^T\Sigma$$

It means that the diagonal entries of the square matrix  $\Sigma^T\Sigma$ , which are the square of the singular values, are the eigenvalues of matrix  $A^TA$  and V is the matrix of eigenvectors.

### Singular values computation

Similarly, consider the product of  $AA^T$ . It is equal to:

$$AA^T = (U\Sigma V^T)(V\Sigma^T U^T) = U\Sigma \Sigma^T U^T$$

Which means that:

$$AA^TU = U\Sigma\Sigma^T$$

Hence U is the matrix of eigenvectors of  $AA^T$ .

Since  $\operatorname{rank}(A)=r,$  only the first r eigenvalues of  $AA^T$  and  $A^TA$  are non-zero.

# Finding eigenvalues and

eigenvectors

### **QR** Method

A possible method to find eigenvalues and eigenvectors of a matrix is based on  ${\it QR}$  decompositions and this theorem:

#### Theorem

Suppose  $A \in \mathbb{R}^{n \times n}$  is a matrix having eigenvalues  $\lambda_1, \lambda_2, \dots \lambda_n$  satisfying

$$|\lambda_1| > |\lambda_2| > \dots > |\lambda_n| \tag{3}$$

then the following sequence for  $A_1 = A$  and k = 1, 2, ...

$$\begin{cases}
A_k = Q_k R_k \\
A_{k+1} = R_k Q_k
\end{cases}$$
(4)

converges to an upper triangular matrix where  $(A_k)_{ii} = \lambda_i$ ,  $i = 1, 2, \ldots, n$ . In case (3) is not satisfied, this sequence converges to a triangular matrix with square blocks of order at most 2 along the diagonal. If A is symmetric, then the sequence converges to a diagonal matrix.

### **QR** Method

So a basic implementation would be like this:

```
while err > tol1
[Q, R] = qr(A);
A = R * Q;

err = max( max( abs( tril(A, -1) ) ) );
end
```

This method can be sped up by using a technique called *shifting*:

```
n = length( A );
while err > toll
% A(n, n) is a usual choice, it can be any real number
T = A(n, n) * eye(n);
[Q, R] = qr( A - T );
A = R * Q + T;

err = max( max( abs( tril(A, -1) ) ) );
end
```

### **QR** Method

In our case, this method must be applied to  $AA^T$  and  $A^TA$ , so if (4) converges then we have a diagonal matrix.

Now, consider the diagonalization of B, where  $B = AA^T$  or  $B = A^TA$ :

$$B = P\Lambda P^{-1} = P\Lambda P^{T}$$

As B can be factored using (4), the matrix containing the eigenvectors must be equal to:

$$P = \prod_{i} Q_i = Q_1 Q_2 Q_3 \cdots$$

Hence, for every iteration  $B_k = Q_k R_k$  and  $B_{k+1} = R_k Q_k$ , requiring each step to have a computational cost equal to  $O(\frac{2n^3}{3})$ .

### **Hessemberg Reduction**

To achieve a lower computational cost, one solution consists in transforming B in a similar tridiagonal matrix using Householder matrices. A triangular matrix can be obtained because B is symmetric.

In general, the process of trasforming a matrix in a similar matrix using Householder matrices allows us to obtain a matrix in which  $B_{ij}=0$ , for all i>j. Such a matrix is called a *Hessemberg matrix*.

Consequently, after B has been transformed, each step of the previous code can be realized using  $\it Givens \ rotation \ matrices.$ 

#### **Givens rotation matrices**

A Givens rotation can be represented with the following matrix:

$$G_{ij} = \begin{pmatrix} 1 & \cdots & 0 & \cdots & 0 & \cdots & 0 \\ \vdots & \ddots & \vdots & & \vdots & & \vdots \\ 0 & \cdots & c & \cdots & -s & \cdots & 0 \\ \vdots & & \vdots & \ddots & \vdots & & \vdots \\ 0 & \cdots & s & \cdots & c & \cdots & 0 \\ \vdots & & \vdots & & \vdots & \ddots & \vdots \\ 0 & \cdots & 0 & \cdots & 0 & \cdots & 1 \end{pmatrix} \leftarrow \mathbf{j}$$

where  $c = \cos \theta$  and  $s = \sin \theta$ , for a particular value of  $\theta \in [0, 2\pi]$ . A rotation occurs in the plane spanned by the two coordinates axes i and j.

#### **Givens rotation matrices**

Using the following values:

$$\cos \theta = \frac{|b_{ii}|}{\sqrt{b_{ii}^2 + b_{ji}^2}}$$

$$\sin \theta = \operatorname{sign}\left(\frac{b_{ji}}{b_{ii}}\right) \frac{|b_{ji}|}{\sqrt{b_{ii}^2 + b_{ji}^2}}$$

when B gets left multiplied by  $G_{ij}$ , the final matrix will have the element in position  $(j, i)^1$  equal to 0.

Therefore, instead of factoring B at each step by the QR method, if B is tridiagonal, we can simply iterate n-1 times by constructing  $G_{k\,k+1}$  and eliminating the elements in position  $(k+1,\,k)$  and  $(k,\,k+1)$ .

<sup>&</sup>lt;sup>1</sup>Note that the indices are reversed

### Pseudocode for finding eigenvalues and eigenvectors

A pseudocode of what to do should clarify what we have done so far (excluding shifting for brevity):

```
\begin{split} B, P \leftarrow \mathsf{hessemberg}(B) \\ \mathbf{while} \ \mathsf{error} > \mathsf{toll} \ \mathbf{do} \\ \mathbf{for} \ k = 1, 2, \dots, n \ \mathbf{do} \\ & \mathsf{construct} \ G_{k \ k+1} \\ B \leftarrow G_{k \ k+1} B \\ \mathbf{end} \ \mathbf{for} \\ \mathbf{for} \ k = 1, 2, \dots, n \ \mathbf{do} \\ B \leftarrow BG_{k \ k+1}^T \\ P \leftarrow PG_{k \ k+1}^T \\ \mathbf{end} \ \mathbf{for} \\ \mathsf{update} \ \mathsf{error} \\ \mathbf{end} \ \mathbf{while} \end{split}
```

The function implementing the Hessemberg reduction must also return the matrix used for the transformation, meaning that  $B = PHP^T$ . Note that P is updated in this way because  $Q_i = G_{12}^{(i)}G_{23}^{(i)}\cdots G_{n-1}^{(i)}$ ,

where i denotes the generic iteration for error minimization.

# The Algorithm

### Calculation of left singular vectors

Suppose we start computing V by the method explained in the previous slides.

We apply the method to  $A^TA$  and get P=V and  $B=\Sigma^2$ .

Once we have take the square root of the singular values, by explicitly writing  $\Sigma$ , the matrix containing the left singular vectors can be calculated as:

$$A = U\Sigma V^T \Rightarrow U = AV\Sigma^{-1}$$

Unlike (1), U is a rectangular matrix, and  $\Sigma$  and V are square matrices. The algorithm is now complete. The next slides will illustrate the code.

#### **Function to calculate Givens rotation matrices**

```
function [c, s] = givens(A, i, j)

%GIVENS Function to compute cos and sin of Gij

% Given a matrix A, this function returns a vector containing the values

% of givens rotation matrix Gij such that Gij * A is equal to A except

% for the element (j, i), which will be zero.

c = abs( A(i, i) ) / sqrt( A(i, i)^2 + A(j, i)^2 );

s = - sign( A(j, i) / A(i, i) ) * abs( A(j, i) ) / sqrt( A(i, i)^2 + A(j, i)^2 );

end
```

### Function to calculate the Hessemberg reduction i

```
function [A, P] = hessemberg(A)
2 %HESSEMBERG Fuction to compute the Hessember reduction of a matrix
3 %
      Given a matrix A, this function trasform A in a similar hessemberg
      matrix. It will be produced as output the hessember matrix and the
4 %
      matrix P used for the transformation.
5 %
      n = size(A, 2);
      P = eve(n):
      for k=1:n-2
8
          sigma = sign(A(k+1, k)) * norm(A(k+1:n, k));
9
          v = [sigma + A(k+1, k); A(k+2:n, k)];
10
          beta = 1 / (sigma * (sigma + A(k+1, k)));
          % transform lower triangular matrix
          for i=k:n
              tau = beta * (v' * A(k+1:n, j));
14
              A(k+1:n, j) = A(k+1:n, j) - tau * v;
          end
16
          % transform upper triangular matrix
          for j=1:n
18
```

### Function to calculate the Hessemberg reduction ii

```
tau = beta * (A(j, k+1:n) * v);

A(j, k+1:n) = A(j, k+1:n) - tau * v';

modify P in a similar way

tau = beta * (P(j, k+1:n) * v);

P(j, k+1:n) = P(j, k+1:n) - tau * v';

end

end

end

end
```

### Function to calculate singular values and singular vectors i

```
function [M, H] = singular vectors(A, toll, left)
2 %SINGULAR VECTORS Function to compute the left or right singular vectors
3 %
      Given a matrix A and a tollerance for the stopping criterion, this
      function computes the left or right singular vectors of A and its
4 %
5 %
      corrisponding singular values. If left is true then U is calculated, V
      otherwise.
6 %
      if left
          [H, P] = hessemberg(A * A.');
8
      else
9
          [H, P] = hessemberg(A.' * A);
10
      end
      n = length(H);
14
      % G will hold c and s for each step
      G = zeros(n - 1, 2);
16
      % G aux will be equal to [c, -s; s, c] for each step
      G aux = zeros(2);
18
```

### Function to calculate singular values and singular vectors ii

```
19
       % M will contain the eigenvectors
20
      M = P;
21
       err = toll + 1:
22
       while err > toll
24
           H1 = H;
25
           T = H(n, n) * eve(n);
26
           % shifting
28
           H = H - T;
29
30
           for k = 1:n-1
31
                [G(k, 1), G(k, 2)] = givens(H, k, k+1);
32
                G \text{ aux}(1, 1) = G(k, 1);
33
                G \text{ aux}(1, 2) = -G(k, 2);
34
                G aux(2, 1) = G(k, 2);
35
                G_{aux}(2, 2) = G(k, 1);
36
37
```

### Function to calculate singular values and singular vectors iii

```
% nullify the elements below the diagonal
38
                H(k:k+1, k:n) = G \text{ aux } * H(k:k+1, k:n):
39
           end
40
41
           for k = 1:n-1
42
                G \text{ aux}(1, 1) = G(k, 1):
43
                G aux(1, 2) = G(k, 2);
44
                G \text{ aux}(2, 1) = -G(k, 2);
45
                G aux(2, 2) = G(k, 1);
46
47
                % reconstruct the Hessemberg form
48
                H(1:k+1, k:k+1) = H(1:k+1, k:k+1) * G aux;
49
                % update M as M * O = M * G12' * G23' * ...
50
                M(1:n, k:k+1) = M(1:n, k:k+1) * G aux;
           end
52
           % shifting
54
           H = H + T;
55
56
```

### Function to calculate singular values and singular vectors iv

```
err = norm(diag(H - H1), 1);
      end
58
59
      % calculate square root for singular values
60
      H = sqrt( diag( diag( H ) ) );
61
62
      % discard imaginary parts if too small
63
      discard imag = 10^-5:
64
65
       if all( imag( diag(H) ) < discard_imag )</pre>
66
           H = real(H);
67
       end
68
       if all( imag( diag(M) ) < discard_imag )</pre>
69
           M = real(M);
70
       end
72 end
```

#### Function to calculate SVD i

```
function [U, S, V] = custom svd(A, toll)
2 %CUSTOM SVD Function to compute the SVD decomposition of a matrix
3 %
      Given a matrix A as input and a tollerance for the stopping criterion,
      this function computes the SVD decomposition of A.
4 %
      [m, n] = size(A);
      [V, S1] = singular vectors(A, toll, false);
      % sort singular values and right singular vectors if they are not
8
      % already
9
      if ~ issorted( diag(S1) )
10
           [New Diag, s order] = sort(diag(S1), 'descend');
          S = diag(New Diag);
14
          new order = 1:n;
          for h = 1:length(s order)
16
              if h ~= s order(h)
                  new order(h) = s order(h);
18
```

#### Function to calculate SVD ii

```
end
19
           end
20
21
           V = V(:, new order);
22
       end
24
      % take only non-zero diagonal entries
25
       arr = New Diag;
26
       arr(arr > 0) = NaN;
       [zero, index] = max(arr, [], 'omitnan');
28
       if index ~= 1
29
           arr = New Diag(1 : index - 1);
30
       else
31
           arr = New_Diag;
32
       end
33
34
      % modify S1 = inv(S) to fit dimensions
35
      S1 = diag( 1./ arr );
36
      S1(n, 1) = 0;
37
```

#### Function to calculate SVD iii

```
38 S1(1, n) = 0;
39
40 % compute U
41 U = A*V*S1;
42 end
```

# Application to Big Data

### **Dimensionality reduction**

The SVD decomposition can be applied to Big Data in order to reduce the dimensionality of data sets.

As an example, consider the data set [5] containing 33 different attributes for 145 students, which includes student ID, personal information and data, and higher education performance ratings<sup>2</sup>.

Dimensionality can be reduced as follows:

- 1. truncate the SVD retaining top-k singular values,  $A\approx U_k\Sigma_k V_k^T;$
- 2. compute the modified data set as  $D' = DV_k = U_k \Sigma_k$ .

 $<sup>^2</sup> https://archive.ics.uci.edu/ml/datasets/Higher+Education+Students+Performance+Evaluation+Dataset$ 

### **Dimensionality reduction**

 $D^\prime$  represents a new data set that corresponds to the projection of the data set D on a k-dimensional basis system of right singular vectors.

#### Number of singular values

As indicated in [3], a usual choice of k is dictated by maintaining enough singular values to make up 90% of the energy in  $\Sigma$ .

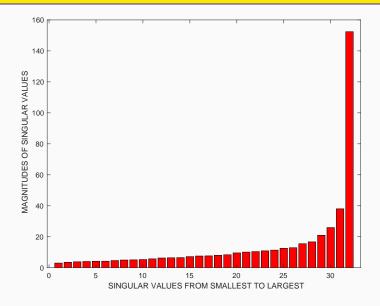
The total energy of a data set is:

$$\mathrm{E}(\Sigma) = \sum_{i=1}^r \sigma_i^2$$

where r = rank(D).

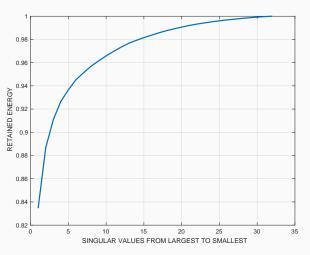
By plotting the singular values we can get an insight of how they are distributed and how many of them we can retain.

### Magnitudes of singular values



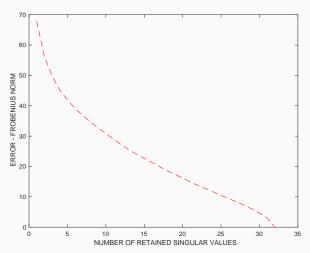
### **Retained energy**

From what we can see, with 3 singular values we can get about 25,346 of energy, which is enough to exceed 90% of  $E(\Sigma)=27,830$ .



### Decreasing error as singular values increase

By increasing the number of retained singular values, the following graph is obtained:



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