

The Singular Value Decomposition

We are interested in more than just sym+def matrices. But the eigenvalue decompositions discussed in the last section of notes will play a major role in solving general systems of equations

$$\mathbf{y} = \mathbf{A}\mathbf{x}, \quad \mathbf{y} \in \mathbb{R}^M, \quad \mathbf{A} \text{ is } M \times N, \quad \mathbf{x} \in \mathbb{R}^N.$$

We have seen that a symmetric positive definite matrix can be decomposed as $\mathbf{A} = \mathbf{V}\mathbf{\Lambda}\mathbf{V}^T$, where \mathbf{V} is an orthogonal matrix ($\mathbf{V}^T\mathbf{V} = \mathbf{V}\mathbf{V}^T = \mathbf{I}$) whose columns are the eigenvectors of \mathbf{A} , and $\mathbf{\Lambda}$ is a diagonal matrix containing the eigenvalues of \mathbf{A} . Because both orthogonal and diagonal matrices are trivial to invert, this eigenvalue decomposition makes it very easy to solve systems of equations $\mathbf{y} = \mathbf{A}\mathbf{x}$ and analyze the stability of these solutions.

The **singular value decomposition** (SVD) takes apart an arbitrary $M \times N$ matrix \mathbf{A} in a similar manner. The SVD of a $M \times N$ matrix \mathbf{A} with rank¹ R is

$$\mathbf{A} = \mathbf{U}\mathbf{\Sigma}\mathbf{V}^T$$

where

1. \mathbf{U} is a $M \times R$ matrix

$$\mathbf{U} = [\mathbf{u}_1 \mid \mathbf{u}_2 \mid \cdots \mid \mathbf{u}_R],$$

whose columns $\mathbf{u}_m \in \mathbb{R}^M$ are orthogonal. Note that while $\mathbf{U}^T\mathbf{U} = \mathbf{I}$, in general $\mathbf{U}\mathbf{U}^T \neq \mathbf{I}$ when $R < M$. The columns of \mathbf{U} are an orthobasis for the range space of \mathbf{A} .

¹Recall that the rank of a matrix is the number of linearly independent columns of a matrix (which is always equal to the number of linearly independent rows).

2. \mathbf{V} is a $N \times R$ matrix

$$\mathbf{V} = [\mathbf{v}_1 \mid \mathbf{v}_2 \mid \cdots \mid \mathbf{v}_R],$$

whose columns $\mathbf{v}_n \in \mathbb{R}^N$ are orthonormal. Again, while $\mathbf{V}^T \mathbf{V} = \mathbf{I}$, in general $\mathbf{V} \mathbf{V}^T \neq \mathbf{I}$ when $R < N$. The columns of \mathbf{V} are an orthobasis for the range space of \mathbf{A}^T (recall that $\text{Range}(\mathbf{A}^T)$ consists of everything which is orthogonal to the nullspace of \mathbf{A}).

3. $\mathbf{\Sigma}$ is a $R \times R$ diagonal matrix with positive entries:

$$\mathbf{\Sigma} = \begin{bmatrix} \sigma_1 & 0 & 0 & \cdots \\ 0 & \sigma_2 & 0 & \cdots \\ \vdots & & \ddots & \\ 0 & \cdots & \cdots & \sigma_R \end{bmatrix}.$$

We call the σ_r the **singular values** of \mathbf{A} . By convention, we will order them such that $\sigma_1 \geq \sigma_2 \geq \cdots \geq \sigma_R$.

4. The $\mathbf{v}_1, \dots, \mathbf{v}_R$ are eigenvectors of the positive semi-definite matrix $\mathbf{A}^T \mathbf{A}$. Note that

$$\mathbf{A}^T \mathbf{A} = \mathbf{V} \mathbf{\Sigma} \mathbf{U}^T \mathbf{U} \mathbf{\Sigma} \mathbf{V}^T = \mathbf{V} \mathbf{\Sigma}^2 \mathbf{V}^T,$$

and so the singular values $\sigma_1, \dots, \sigma_R$ are the square roots of the non-zero eigenvalues of $\mathbf{A}^T \mathbf{A}$.

5. Similarly,

$$\mathbf{A} \mathbf{A}^T = \mathbf{U} \mathbf{\Sigma}^2 \mathbf{U}^T,$$

and so the $\mathbf{u}_1, \dots, \mathbf{u}_R$ are eigenvectors of the positive semi-definite matrix $\mathbf{A} \mathbf{A}^T$. Since the non-zero eigenvalues of $\mathbf{A}^T \mathbf{A}$

and $\mathbf{A}\mathbf{A}^T$ are the same, the σ_r are also square roots of the eigenvalues of $\mathbf{A}\mathbf{A}^T$.

The rank R is the dimension of the space spanned by the columns of \mathbf{A} , this is the same as the dimension of the space spanned by the rows. Thus $R \leq \min(M, N)$. We say \mathbf{A} is **full rank** if $R = \min(M, N)$.

As before, we will often times find it useful to write the SVD as the sum of R rank-1 matrices:

$$\mathbf{A} = \mathbf{U}\mathbf{\Sigma}\mathbf{V}^T = \sum_{r=1}^R \sigma_r \mathbf{u}_r \mathbf{v}_r^T.$$

When \mathbf{A} is **overdetermined** ($M > N$), the decomposition looks like this

$$\begin{bmatrix} \mathbf{A} \end{bmatrix} = \begin{bmatrix} \mathbf{U} \end{bmatrix} \begin{bmatrix} \sigma_1 & & \\ & \cdots & \\ & & \sigma_R \end{bmatrix} \begin{bmatrix} \mathbf{V}^T \end{bmatrix}$$

When \mathbf{A} is **underdetermined** ($M < N$), the SVD looks like this

$$\begin{bmatrix} \mathbf{A} \end{bmatrix} = \begin{bmatrix} \mathbf{U} \end{bmatrix} \begin{bmatrix} \sigma_1 & & \\ & \cdots & \\ & & \sigma_R \end{bmatrix} \begin{bmatrix} \mathbf{V}^T \end{bmatrix}$$

When \mathbf{A} is **square** and full rank ($M = N = R$), the SVD looks like

$$\begin{bmatrix} \mathbf{A} \end{bmatrix} = \begin{bmatrix} \mathbf{U} \end{bmatrix} \begin{bmatrix} \sigma_1 & & \\ & \cdots & \\ & & \sigma_N \end{bmatrix} \begin{bmatrix} \mathbf{V}^T \end{bmatrix}$$

Technical Details: Existence of the SVD

In this section we will prove that any $M \times N$ matrix \mathbf{A} with $\text{rank}(\mathbf{A}) = R$ can be written as

$$\mathbf{A} = \mathbf{U}\mathbf{\Sigma}\mathbf{V}^T$$

where \mathbf{U} , $\mathbf{\Sigma}$, \mathbf{V} have the five properties listed at the beginning of the last section.

Since $\mathbf{A}^T\mathbf{A}$ is symmetric positive semi-definite, we can write:

$$\mathbf{A}^T\mathbf{A} = \sum_{n=1}^N \lambda_n \mathbf{v}_n \mathbf{v}_n^T,$$

where the \mathbf{v}_n are orthonormal and the λ_n are real and non-negative. Since $\text{rank}(\mathbf{A}) = R$, we also have $\text{rank}(\mathbf{A}^T\mathbf{A}) = R$, and so $\lambda_1, \dots, \lambda_R$ are all strictly positive above, and $\lambda_{R+1} = \dots = \lambda_N = 0$.

Set

$$\mathbf{u}_m = \frac{1}{\sqrt{\lambda_m}} \mathbf{A}\mathbf{v}_m, \quad \text{for } m = 1, \dots, R, \quad \mathbf{U} = [\mathbf{u}_1 \ \dots \ \mathbf{u}_R].$$

Notice that these \mathbf{u}_m are orthonormal, as

$$\langle \mathbf{u}_m, \mathbf{u}_\ell \rangle = \frac{1}{\sqrt{\lambda_m \lambda_\ell}} \mathbf{v}_\ell^T \mathbf{A}^T \mathbf{A} \mathbf{v}_m = \sqrt{\frac{\lambda_m}{\lambda_\ell}} \mathbf{v}_\ell^T \mathbf{v}_m = \begin{cases} 1, & m = \ell, \\ 0, & m \neq \ell. \end{cases}$$

These \mathbf{u}_m also happen to be eigenvectors of $\mathbf{A}\mathbf{A}^T$, as

$$\mathbf{A}\mathbf{A}^T \mathbf{u}_m = \frac{1}{\sqrt{\lambda_m}} \mathbf{A}\mathbf{A}^T \mathbf{A}\mathbf{v}_m = \sqrt{\lambda_m} \mathbf{A}\mathbf{v}_m = \lambda_m \mathbf{u}_m.$$

Now let $\mathbf{u}_{R+1}, \dots, \mathbf{u}_M$ be an orthobasis for the null space of \mathbf{U}^T — concatenating these two sets into $\mathbf{u}_1, \dots, \mathbf{u}_M$ forms an orthobasis for all of \mathbb{R}^M .

Let

$$\mathbf{V} = [\mathbf{v}_1 \ \mathbf{v}_2 \ \cdots \ \mathbf{v}_R], \quad \mathbf{V}_0 = [\mathbf{v}_{R+1} \ \mathbf{v}_{R+2} \ \cdots \ \mathbf{v}_N], \quad \mathbf{V}_{\text{full}} = [\mathbf{V} \ \mathbf{V}_0]$$

and

$$\mathbf{U}_0 = [\mathbf{u}_{R+1} \ \mathbf{u}_{R+2} \ \cdots \ \mathbf{u}_M], \quad \mathbf{U}_{\text{full}} = [\mathbf{U} \ \mathbf{U}_0].$$

It should be clear that \mathbf{V}_{full} is an $N \times N$ orthonormal matrix and \mathbf{U}_{full} is a $M \times M$ orthonormal matrix. Consider the $M \times N$ matrix $\mathbf{U}_{\text{full}}^T \mathbf{A} \mathbf{V}_{\text{full}}$ — the entry in the m th rows and n th column of this matrix is

$$\begin{aligned} (\mathbf{U}_{\text{full}}^T \mathbf{A} \mathbf{V}_{\text{full}})[m, n] &= \mathbf{u}_m^T \mathbf{A} \mathbf{v}_n = \begin{cases} \sqrt{\lambda_n} \mathbf{u}_m^T \mathbf{u}_n & n = 1, \dots, R \\ 0, & n = R + 1, \dots, N. \end{cases} \\ &= \begin{cases} \sqrt{\lambda_n}, & m = n = 1, \dots, R \\ 0, & \text{otherwise.} \end{cases} \end{aligned}$$

Thus

$$\mathbf{U}_{\text{full}}^T \mathbf{A} \mathbf{V}_{\text{full}} = \mathbf{\Sigma}_{\text{full}}$$

where

$$\Sigma_{\text{full}}[m, n] = \begin{cases} \sqrt{\lambda_n}, & m = n = 1, \dots, R \\ 0, & \text{otherwise.} \end{cases}$$

Since $\mathbf{U}_{\text{full}} \mathbf{U}_{\text{full}}^T = \mathbf{I}$ and $\mathbf{V}_{\text{full}} \mathbf{V}_{\text{full}}^T = \mathbf{I}$, we have

$$\mathbf{A} = \mathbf{U}_{\text{full}} \mathbf{\Sigma}_{\text{full}} \mathbf{V}_{\text{full}}^T.$$

Since $\mathbf{\Sigma}_{\text{full}}$ is non-zero only in the first R locations along its main diagonal, the above reduces to

$$\mathbf{A} = \mathbf{U} \mathbf{\Sigma} \mathbf{V}^T, \quad \mathbf{\Sigma} = \begin{bmatrix} \sqrt{\lambda_1} & & & & \\ & \sqrt{\lambda_2} & & & \\ & & \ddots & & \\ & & & \ddots & \\ & & & & \sqrt{\lambda_R} \end{bmatrix}.$$

The Least-Squares Problem

We can use the SVD to “solve” the general system of linear equations

$$\mathbf{y} = \mathbf{A}\mathbf{x}$$

where $\mathbf{y} \in \mathbb{R}^M$, $\mathbf{x} \in \mathbb{R}^N$, and \mathbf{A} is an $M \times N$ matrix.

Given \mathbf{y} , we want to find \mathbf{x} in such a way that

1. when there is a unique solution, we return it;
2. when there is no solution, we return something reasonable;
3. when there are an infinite number of solutions, we choose one to return in a “smart” way.

The **least-squares** framework revolves around finding an \mathbf{x} that minimizes the length of the residual

$$\mathbf{r} = \mathbf{y} - \mathbf{A}\mathbf{x}.$$

That is, we want to solve the optimization problem

$$\underset{\mathbf{x} \in \mathbb{R}^N}{\text{minimize}} \quad \|\mathbf{y} - \mathbf{A}\mathbf{x}\|_2^2, \quad (1)$$

where $\|\cdot\|_2$ is the standard Euclidean norm. We will see that the SVD of \mathbf{A} :

$$\mathbf{A} = \mathbf{U}\mathbf{\Sigma}\mathbf{V}^T, \quad (2)$$

plays a pivotal role in solving this problem.

To start, note that we can write any $\mathbf{x} \in \mathbb{R}^N$ as

$$\mathbf{x} = \mathbf{V}\mathbf{\alpha} + \mathbf{V}_0\mathbf{\alpha}_0. \quad (3)$$

Here, \mathbf{V} is the $N \times R$ matrix appearing in the SVD decomposition (2), and \mathbf{V}_0 is a $N \times (N - R)$ matrix whose columns are orthogonal to one another and to the columns in \mathbf{V} . We have the relations

$$\mathbf{V}^T \mathbf{V} = \mathbf{I}, \quad \mathbf{V}_0^T \mathbf{V}_0 = \mathbf{I}, \quad \mathbf{V}^T \mathbf{V}_0 = \mathbf{0}.$$

You can think of \mathbf{V}_0 as an orthobasis for the null space of \mathbf{A} . Of course, \mathbf{V}_0 is not unique, as there are many orthobases for $\text{Null}(\mathbf{A})$, but any such set of vectors will serve our purposes here. The decomposition (3) is possible since $\text{Range}(\mathbf{A}^T)$ and $\text{Null}(\mathbf{A})$ partition \mathbb{R}^N for any $M \times N$ matrix \mathbf{A} . Taking

$$\boldsymbol{\alpha} = \mathbf{V}^T \mathbf{x}, \quad \boldsymbol{\alpha}_0 = \mathbf{V}_0^T \mathbf{x},$$

we see that (3) holds since

$$\mathbf{x} = \mathbf{V} \mathbf{V}^T \mathbf{x} + \mathbf{V}_0 \mathbf{V}_0^T \mathbf{x} = (\mathbf{V} \mathbf{V}^T + \mathbf{V}_0 \mathbf{V}_0^T) \mathbf{x} = \mathbf{x},$$

where we have made use of the fact that $\mathbf{V} \mathbf{V}^T + \mathbf{V}_0 \mathbf{V}_0^T = \mathbf{I}$, because $\mathbf{V} \mathbf{V}^T$ and $\mathbf{V}_0 \mathbf{V}_0^T$ are ortho-projectors onto complementary subspaces² of \mathbb{R}^N . So we can solve for $\mathbf{x} \in \mathbb{R}^N$ by solving for the pair $\boldsymbol{\alpha} \in \mathbb{R}^R$, $\boldsymbol{\alpha}_0 \in \mathbb{R}^{N-R}$.

Similarly, we can decompose \mathbf{y} as

$$\mathbf{y} = \mathbf{U} \boldsymbol{\beta} + \mathbf{U}_0 \boldsymbol{\beta}_0, \tag{4}$$

where \mathbf{U} is the $M \times R$ matrix from the SVD decomposition, and \mathbf{U}_0 is a $M \times (M - R)$ complementary orthogonal basis. Again,

$$\mathbf{U}^T \mathbf{U} = \mathbf{I}, \quad \mathbf{U}_0^T \mathbf{U}_0 = \mathbf{I}, \quad \mathbf{U}^T \mathbf{U}_0 = \mathbf{0},$$

²Subspaces \mathcal{S}_1 and \mathcal{S}_2 are **complementary** in \mathbb{R}^N if $\mathcal{S}_1 \perp \mathcal{S}_2$ (everything in \mathcal{S}_1 is orthogonal to everything in \mathcal{S}_2) and $\mathcal{S}_1 \oplus \mathcal{S}_2 = \mathbb{R}^N$. You can think of $\mathcal{S}_1, \mathcal{S}_2$ as a partition of \mathbb{R}^N into two orthogonal subspaces.

and we can think of \mathbf{U}_0 as an orthogonal basis for everything in \mathbb{R}^M that is not in the range of \mathbf{A} . As before, we can calculate the decomposition above using

$$\boldsymbol{\beta} = \mathbf{U}^T \mathbf{y}, \quad \boldsymbol{\beta}_0 = \mathbf{U}_0^T \mathbf{y}.$$

Using the decompositions (2), (3), and (4) for \mathbf{A} , \mathbf{x} , and \mathbf{y} , we can write the residual $\mathbf{r} = \mathbf{y} - \mathbf{A}\mathbf{x}$ as

$$\begin{aligned} \mathbf{r} &= \mathbf{U}\boldsymbol{\beta} + \mathbf{U}_0\boldsymbol{\beta}_0 - \mathbf{U}\boldsymbol{\Sigma}\mathbf{V}^T(\mathbf{V}\boldsymbol{\alpha} + \mathbf{V}_0\boldsymbol{\alpha}_0) \\ &= \mathbf{U}\boldsymbol{\beta} + \mathbf{U}_0\boldsymbol{\beta}_0 - \mathbf{U}\boldsymbol{\Sigma}\boldsymbol{\alpha} \quad (\text{since } \mathbf{V}^T\mathbf{V} = \mathbf{I} \text{ and } \mathbf{V}^T\mathbf{V}_0 = \mathbf{0}) \\ &= \mathbf{U}_0\boldsymbol{\beta}_0 + \mathbf{U}(\boldsymbol{\beta} - \boldsymbol{\Sigma}\boldsymbol{\alpha}). \end{aligned}$$

We want to choose $\boldsymbol{\alpha}$ that minimizes the energy of \mathbf{r} :

$$\begin{aligned} \|\mathbf{r}\|_2^2 &= \langle \mathbf{U}_0\boldsymbol{\beta}_0 + \mathbf{U}(\boldsymbol{\beta} - \boldsymbol{\Sigma}\boldsymbol{\alpha}), \mathbf{U}_0\boldsymbol{\beta}_0 + \mathbf{U}(\boldsymbol{\beta} - \boldsymbol{\Sigma}\boldsymbol{\alpha}) \rangle \\ &= \langle \mathbf{U}_0\boldsymbol{\beta}_0, \mathbf{U}_0\boldsymbol{\beta}_0 \rangle + 2\langle \mathbf{U}_0\boldsymbol{\beta}_0, \mathbf{U}(\boldsymbol{\beta} - \boldsymbol{\Sigma}\boldsymbol{\alpha}) \rangle \\ &\quad + \langle \mathbf{U}(\boldsymbol{\beta} - \boldsymbol{\Sigma}\boldsymbol{\alpha}), \mathbf{U}(\boldsymbol{\beta} - \boldsymbol{\Sigma}\boldsymbol{\alpha}) \rangle \\ &= \|\boldsymbol{\beta}_0\|_2^2 + \|\boldsymbol{\beta} - \boldsymbol{\Sigma}\boldsymbol{\alpha}\|_2^2 \end{aligned}$$

where the last equality comes from the facts that $\mathbf{U}_0^T\mathbf{U}_0 = \mathbf{I}$, $\mathbf{U}^T\mathbf{U} = \mathbf{I}$, and $\mathbf{U}^T\mathbf{U}_0 = \mathbf{0}$. We have no control over $\|\boldsymbol{\beta}_0\|_2^2$, since it determined entirely by our observations \mathbf{y} . Therefore, our problem has been reduced to finding $\boldsymbol{\alpha}$ that minimizes the second term $\|\boldsymbol{\beta} - \boldsymbol{\Sigma}\boldsymbol{\alpha}\|_2^2$ above, which is non-negative. We can make it zero (i.e. as small as possible) by taking

$$\hat{\boldsymbol{\alpha}} = \boldsymbol{\Sigma}^{-1}\boldsymbol{\beta}.$$

Finally, the \mathbf{x} which minimizes the residual (solves (1)) is

$$\hat{\mathbf{x}} = \mathbf{V}\hat{\boldsymbol{\alpha}} = \mathbf{V}\boldsymbol{\Sigma}^{-1}\boldsymbol{\beta} = \mathbf{V}\boldsymbol{\Sigma}^{-1}\mathbf{U}^T\mathbf{y}. \quad (5)$$

Thus we can calculate the solution to (1) simply by applying the linear operator $\mathbf{V}\Sigma^{-1}\mathbf{U}^T$ to the input data \mathbf{y} . There are two interesting facts about the solution $\hat{\mathbf{x}}$ in (5):

1. When $\mathbf{y} \in \text{span}(\{\mathbf{u}_1, \dots, \mathbf{u}_M\})$, we have $\beta_0 = \mathbf{U}_0^T \mathbf{y} = \mathbf{0}$, and so the residual $\mathbf{r} = \mathbf{0}$. In this case, there is at least one exact solution, and the one we choose satisfies $\mathbf{A}\hat{\mathbf{x}} = \mathbf{y}$.
2. Note that if $R < N$, then the solution is not unique. In this case, \mathbf{V}_0 has at least one column, and any part of a vector \mathbf{x} in the range of \mathbf{V}_0 is not seen by \mathbf{A} , since

$$\mathbf{A}\mathbf{V}_0\boldsymbol{\alpha}_0 = \mathbf{U}\Sigma\mathbf{V}^T\mathbf{V}_0\boldsymbol{\alpha}_0 = \mathbf{0} \quad (\text{since } \mathbf{V}^T\mathbf{V}_0 = \mathbf{0}).$$

As such,

$$\mathbf{x}' = \hat{\mathbf{x}} + \mathbf{V}_0\boldsymbol{\alpha}_0$$

for *any* $\boldsymbol{\alpha}_0 \in \mathbb{R}^{N-R}$ will have exactly the same residual, since $\mathbf{A}\mathbf{x}' = \mathbf{A}\hat{\mathbf{x}}$. In this case, our solution $\hat{\mathbf{x}}$ is the solution with smallest norm, since

$$\begin{aligned} \|\mathbf{x}'\|_2^2 &= \langle \hat{\mathbf{x}} + \mathbf{V}_0\boldsymbol{\alpha}_0, \hat{\mathbf{x}} + \mathbf{V}_0\boldsymbol{\alpha}_0 \rangle \\ &= \langle \hat{\mathbf{x}}, \hat{\mathbf{x}} \rangle + 2\langle \hat{\mathbf{x}}, \mathbf{V}_0\boldsymbol{\alpha}_0 \rangle + \langle \mathbf{V}_0\boldsymbol{\alpha}_0, \mathbf{V}_0\boldsymbol{\alpha}_0 \rangle \\ &= \|\hat{\mathbf{x}}\|_2^2 + 2\langle \mathbf{V}\Sigma^{-1}\mathbf{U}^T\mathbf{y}, \mathbf{V}_0\boldsymbol{\alpha}_0 \rangle + \|\boldsymbol{\alpha}_0\|_2^2 \quad (\text{since } \mathbf{V}_0^T\mathbf{V}_0 = \mathbf{I}) \\ &= \|\hat{\mathbf{x}}\|_2^2 + \|\boldsymbol{\alpha}_0\|_2^2 \quad (\text{since } \mathbf{V}^T\mathbf{V}_0 = \mathbf{0}) \end{aligned}$$

which is minimized by taking $\boldsymbol{\alpha}_0 = \mathbf{0}$.

To summarize, $\hat{\mathbf{x}} = \mathbf{V}\Sigma^{-1}\mathbf{U}^T\mathbf{y}$ has the desired properties stated at the beginning of this module, since

1. when $\mathbf{y} = \mathbf{A}\mathbf{x}$ has a unique exact solution, it must be $\hat{\mathbf{x}}$,
2. when an exact solution is not available, $\hat{\mathbf{x}}$ is the solution to (1),

3. when there are an infinite number of minimizers to (1), $\hat{\mathbf{x}}$ is the one with smallest norm.

Because the matrix $\mathbf{V}\mathbf{\Sigma}^{-1}\mathbf{U}^T$ gives us such an elegant solution to this problem, we give it a special name: the **pseudo-inverse**.

The Pseudo-Inverse

The **pseudo-inverse** of a matrix \mathbf{A} with singular value decomposition (SVD) $\mathbf{A} = \mathbf{U}\mathbf{\Sigma}\mathbf{V}^T$ is

$$\mathbf{A}^\dagger = \mathbf{V}\mathbf{\Sigma}^{-1}\mathbf{U}^T. \quad (6)$$

Other names for \mathbf{A}^\dagger include **natural inverse**, **Lanczos inverse**, and **Moore-Penrose inverse**.

Given an observation \mathbf{y} , taking $\hat{\mathbf{x}} = \mathbf{A}^\dagger\mathbf{y}$ gives us the **least squares** solution to $\mathbf{y} = \mathbf{A}\mathbf{x}$. The pseudo-inverse \mathbf{A}^\dagger always exists, since every matrix (with rank R) has an SVD decomposition $\mathbf{A} = \mathbf{U}\mathbf{\Sigma}\mathbf{V}^T$ with $\mathbf{\Sigma}$ as an $R \times R$ diagonal matrix with $\Sigma[r, r] > 0$.

When \mathbf{A} is full rank ($R = \min(M, N)$), then we can calculate the pseudo-inverse without using the SVD. There are three cases:

- When \mathbf{A} is square and invertible ($R = M = N$), then

$$\mathbf{A}^\dagger = \mathbf{A}^{-1}.$$

This is easy to check, as here

$$\mathbf{A} = \mathbf{U}\mathbf{\Sigma}\mathbf{V}^T \quad \text{where both } \mathbf{U}, \mathbf{V} \text{ are } N \times N,$$

and since in this case $\mathbf{V}\mathbf{V}^T = \mathbf{V}^T\mathbf{V} = \mathbf{I}$ and $\mathbf{U}\mathbf{U}^T = \mathbf{U}^T\mathbf{U} = \mathbf{I}$,

$$\begin{aligned}\mathbf{A}^\dagger\mathbf{A} &= \mathbf{V}\boldsymbol{\Sigma}^{-1}\mathbf{U}^T\mathbf{U}\boldsymbol{\Sigma}\mathbf{V}^T \\ &= \mathbf{V}\boldsymbol{\Sigma}^{-1}\boldsymbol{\Sigma}\mathbf{V}^T \\ &= \mathbf{V}\mathbf{V}^T \\ &= \mathbf{I}.\end{aligned}$$

Similarly, $\mathbf{A}\mathbf{A}^\dagger = \mathbf{I}$, and so \mathbf{A}^\dagger is both a left and right inverse of \mathbf{A} , and thus $\mathbf{A}^\dagger = \mathbf{A}^{-1}$.

- When \mathbf{A} more rows than columns and has full column rank ($R = N \leq M$), then $\mathbf{A}^T\mathbf{A}$ is invertible, and

$$\mathbf{A}^\dagger = (\mathbf{A}^T\mathbf{A})^{-1}\mathbf{A}^T. \quad (7)$$

This type of \mathbf{A} is “tall and skinny”

$$\begin{bmatrix} \mathbf{A} \end{bmatrix},$$

and its columns are linearly independent. To verify equation (7), recall that

$$\mathbf{A}^T\mathbf{A} = \mathbf{V}\boldsymbol{\Sigma}\mathbf{U}^T\mathbf{U}\boldsymbol{\Sigma}\mathbf{V}^T = \mathbf{V}\boldsymbol{\Sigma}^2\mathbf{V}^T,$$

and so

$$(\mathbf{A}^T\mathbf{A})^{-1}\mathbf{A}^T = \mathbf{V}\boldsymbol{\Sigma}^{-2}\mathbf{V}^T\mathbf{V}\boldsymbol{\Sigma}\mathbf{U}^T = \mathbf{V}\boldsymbol{\Sigma}^{-1}\mathbf{U}^T,$$

which is exactly the content of (6).

- When \mathbf{A} has more columns than rows and has full row rank ($R = M \leq N$), then $\mathbf{A}\mathbf{A}^T$ is invertible, and

$$\mathbf{A}^\dagger = \mathbf{A}^T(\mathbf{A}\mathbf{A}^T)^{-1}. \quad (8)$$

This occurs when \mathbf{A} is “short and fat”

$$\begin{bmatrix} & & \\ & \mathbf{A} & \\ & & \end{bmatrix},$$

and its rows are linearly independent. To verify equation (8), recall that

$$\mathbf{A}\mathbf{A}^T = \mathbf{U}\mathbf{\Sigma}\mathbf{V}^T\mathbf{V}\mathbf{\Sigma}\mathbf{U}^T = \mathbf{U}\mathbf{\Sigma}^2\mathbf{U}^T,$$

and so

$$\mathbf{A}^T(\mathbf{A}\mathbf{A}^T)^{-1} = \mathbf{V}\mathbf{\Sigma}\mathbf{U}^T\mathbf{U}\mathbf{\Sigma}^{-2}\mathbf{U}^T = \mathbf{V}\mathbf{\Sigma}^{-1}\mathbf{U}^T,$$

which again is exactly (6).

\mathbf{A}^\dagger is as close to an inverse of \mathbf{A} as possible

As discussed in the last section, when \mathbf{A} is square and invertible, \mathbf{A}^\dagger is exactly the inverse of \mathbf{A} . When \mathbf{A} is not square, we can ask if there is a better right or left inverse. We will argue that there is not.

Left inverse Given $\mathbf{y} = \mathbf{A}\mathbf{x}$, we would like $\mathbf{A}^\dagger\mathbf{y} = \mathbf{A}^\dagger\mathbf{A}\mathbf{x} = \mathbf{x}$ for any \mathbf{x} . That is, we would like \mathbf{A}^\dagger to be a *left inverse* of \mathbf{A} : $\mathbf{A}^\dagger\mathbf{A} = \mathbf{I}$. Of course, this is not always possible, especially when \mathbf{A} has more columns than rows, $M < N$. But we can ask if any other matrix \mathbf{H} comes closer to being a left inverse

than \mathbf{A}^\dagger . To find the “best” left-inverse, we look for the matrix which minimizes

$$\min_{\mathbf{H} \in \mathbb{R}^{N \times M}} \|\mathbf{H}\mathbf{A} - \mathbf{I}\|_F^2. \quad (9)$$

Here, $\|\cdot\|_F$ is the *Frobenius norm*, defined for an $N \times M$ matrix \mathbf{Q} as the sum of the squares of the entries:

$$\|\mathbf{Q}\|_F^2 = \sum_{n=1}^M \sum_{m=1}^N |Q[m, n]|^2$$

(It is also true, and you can and should prove this at home, that $\|\mathbf{Q}\|_F^2$ is the sum of the squares of the singular values of \mathbf{Q} : $\|\mathbf{Q}\|_F^2 = \lambda_1^2 + \dots + \lambda_p^2$.) With (9), we are finding \mathbf{H} such that $\mathbf{H}\mathbf{A}$ is as close to the identity as possible in the least-squares sense.

The pseudo-inverse \mathbf{A}^\dagger minimizes (9). To see this, recognize (see the exercise below) that the solution $\hat{\mathbf{H}}$ to (9) must obey

$$\mathbf{A}\mathbf{A}^\top \hat{\mathbf{H}}^\top = \mathbf{A}. \quad (10)$$

We can see that this is indeed true for $\hat{\mathbf{H}} = \mathbf{A}^\dagger$:

$$\mathbf{A}\mathbf{A}^\top \mathbf{A}^{\dagger\top} = \mathbf{U}\Sigma\mathbf{V}^\top \mathbf{V}\Sigma\mathbf{U}^\top \mathbf{U}\Sigma^{-1}\mathbf{V}^\top = \mathbf{U}\Sigma\mathbf{V}^\top = \mathbf{A}.$$

So there is no $N \times M$ matrix that is closer to being a left inverse than \mathbf{A}^\dagger .

Right inverse If we re-apply \mathbf{A} to our solution $\hat{\mathbf{x}} = \mathbf{A}^\dagger \mathbf{y}$, we would like it to be as close as possible to our observations \mathbf{y} . That is,

we would like $\mathbf{A}\mathbf{A}^\dagger$ to be as close to the identity as possible. Again, achieving this goal exactly is not always possible, especially if \mathbf{A} has more rows than columns. But we can attempt to find the “best” right inverse, in the least-squares sense, by solving

$$\underset{\mathbf{H} \in \mathbb{R}^{N \times M}}{\text{minimize}} \quad \|\mathbf{A}\mathbf{H} - \mathbf{I}\|_F^2. \quad (11)$$

The solution $\hat{\mathbf{H}}$ to (11) (see the exercise below) must obey

$$\mathbf{A}^\top \mathbf{A} \hat{\mathbf{H}} = \mathbf{A}^\top. \quad (12)$$

Again, we show that \mathbf{A}^\dagger satisfies (12), and hence is a minimizer to (11):

$$\mathbf{A}^\top \mathbf{A} \mathbf{A}^\dagger = \mathbf{V} \boldsymbol{\Sigma}^2 \mathbf{V}^\top \mathbf{V} \boldsymbol{\Sigma}^{-1} \mathbf{U}^\top = \mathbf{V} \boldsymbol{\Sigma} \mathbf{U}^\top = \mathbf{A}^\top.$$

Moral:

$\mathbf{A}^\dagger = \mathbf{V} \boldsymbol{\Sigma}^{-1} \mathbf{U}^\top$ is as close (in the least-squares sense) to an inverse of \mathbf{A} as you could possibly have.

Exercise:

1. Show that the minimizer $\hat{\mathbf{H}}$ to (9) must obey (10). Do this by using the fact that the derivative of the functional $\|\mathbf{H}\mathbf{A} - \mathbf{I}\|_F^2$ with respect to an entry $H[k, \ell]$ in \mathbf{H} must obey

$$\frac{\partial \|\mathbf{H}\mathbf{A} - \mathbf{I}\|_F^2}{\partial H[k, \ell]} = 0, \quad \text{for all } 1 \leq k \leq N, 1 \leq \ell \leq M,$$

to be a solution to (9). Do the same for (11) and (12).

Stability Analysis of the Pseudo-Inverse

We have seen that if we make indirect observations $\mathbf{y} \in \mathbb{R}^M$ of an unknown vector $\mathbf{x}_0 \in \mathbb{R}^N$ through a $M \times N$ matrix \mathbf{A} , $\mathbf{y} = \mathbf{A}\mathbf{x}_0$, then applying the pseudo-inverse of \mathbf{A} gives us the least squares estimate of \mathbf{x}_0 :

$$\hat{\mathbf{x}}_{\text{ls}} = \mathbf{A}^\dagger \mathbf{y} = \mathbf{V}\mathbf{\Sigma}^{-1}\mathbf{U}^\text{T}\mathbf{y},$$

where $\mathbf{A} = \mathbf{U}\mathbf{\Sigma}\mathbf{V}^\text{T}$ is the singular value decomposition (SVD) of \mathbf{A} .

We will now discuss what happens if our measurements contain *noise* — the analysis here will be very similar to when we looked at the stability of solving square sym+def systems, and in fact this is one of the main reasons we introduced the SVD.

Suppose we observe

$$\mathbf{y} = \mathbf{A}\mathbf{x}_0 + \mathbf{e},$$

where $\mathbf{e} \in \mathbb{R}^M$ is an unknown perturbation. Say that we again apply the pseudo-inverse to \mathbf{y} in an attempt to recover \mathbf{x} :

$$\hat{\mathbf{x}}_{\text{ls}} = \mathbf{A}^\dagger \mathbf{y} = \mathbf{A}^\dagger \mathbf{A}\mathbf{x}_0 + \mathbf{A}^\dagger \mathbf{e}$$

What effect does the presence of the noise vector \mathbf{e} had on our estimate of \mathbf{x}_0 ? We answer this question by comparing $\hat{\mathbf{x}}_{\text{ls}}$ to the reconstruction we would obtain if we used standard least-squares on perfectly noise-free observations $\mathbf{y}_{\text{clean}} = \mathbf{A}\mathbf{x}_0$. This noise-free recon-

struction can be written as

$$\begin{aligned}
 \mathbf{x}_{\text{pinv}} &= \mathbf{A}^\dagger \mathbf{y}_{\text{clean}} = \mathbf{A}^\dagger \mathbf{A} \mathbf{x}_0 \\
 &= \mathbf{V} \boldsymbol{\Sigma}^{-1} \mathbf{U}^T \mathbf{U} \boldsymbol{\Sigma} \mathbf{V}^T \mathbf{x}_0 \\
 &= \mathbf{V} \mathbf{V}^T \mathbf{x}_0 \\
 &= \sum_{r=1}^R \langle \mathbf{x}_0, \mathbf{v}_r \rangle \mathbf{v}_r.
 \end{aligned}$$

The vector \mathbf{x}_{pinv} is the orthogonal projection of \mathbf{x}_0 onto the row space (everything orthogonal to the null space) of \mathbf{A} . If \mathbf{A} has full column rank ($R = N$), then $\mathbf{x}_{\text{pinv}} = \mathbf{x}_0$. If not, then the application of \mathbf{A} destroys the part of \mathbf{x}_0 that is not in \mathbf{x}_{pinv} , and so we only attempt to recover the “visible” components. In some sense, \mathbf{x}_{pinv} contains all of the components of \mathbf{x}_0 that \mathbf{A} does not completely remove, and has them preserved perfectly.

The reconstruction error (relative to \mathbf{x}_{pinv} is)

$$\|\hat{\mathbf{x}}_{\text{ls}} - \mathbf{x}_{\text{pinv}}\|_2^2 = \|\mathbf{A}^\dagger \mathbf{e}\|_2^2 = \|\mathbf{V} \boldsymbol{\Sigma}^{-1} \mathbf{U}^T \mathbf{e}\|_2^2. \quad (13)$$

Now suppose for a moment that the error has unit norm, $\|\mathbf{e}\|_2^2 = 1$. Then the worst case for (13) is given by

$$\underset{\mathbf{e} \in \mathbb{R}^M}{\text{maximize}} \quad \|\mathbf{V} \boldsymbol{\Sigma}^{-1} \mathbf{U}^T \mathbf{e}\|_2^2 \quad \text{subject to} \quad \|\mathbf{e}\|_2 = 1.$$

Since the columns of \mathbf{U} are orthonormal, $\|\mathbf{U}^T \mathbf{e}\|_2^2 \leq \|\mathbf{e}\|_2^2$, and the above is equivalent to

$$\max_{\boldsymbol{\beta} \in \mathbb{R}^R: \|\boldsymbol{\beta}\|_2=1} \|\mathbf{V} \boldsymbol{\Sigma}^{-1} \boldsymbol{\beta}\|_2^2. \quad (14)$$

Also, for any vector $\mathbf{z} \in \mathbb{R}^R$, we have

$$\|\mathbf{V}\mathbf{z}\|_2^2 = \langle \mathbf{V}\mathbf{z}, \mathbf{V}\mathbf{z} \rangle = \langle \mathbf{z}, \mathbf{V}^T \mathbf{V}\mathbf{z} \rangle = \langle \mathbf{z}, \mathbf{z} \rangle = \|\mathbf{z}\|_2^2,$$

since the columns of \mathbf{V} are orthonormal. So we can simplify (14) to

$$\underset{\boldsymbol{\beta} \in \mathbb{R}^R}{\text{maximize}} \|\boldsymbol{\Sigma}^{-1}\boldsymbol{\beta}\|_2^2 \quad \text{subject to} \quad \|\boldsymbol{\beta}\|_2 = 1.$$

The worst case $\boldsymbol{\beta}$ (you should verify this at home) will have a 1 in the entry corresponding to the largest entry in $\boldsymbol{\Sigma}^{-1}$, and will be zero everywhere else. Thus

$$\max_{\boldsymbol{\beta} \in \mathbb{R}^R: \|\boldsymbol{\beta}\|_2=1} \|\boldsymbol{\Sigma}^{-1}\boldsymbol{\beta}\|_2^2 = \max_{r=1, \dots, R} \sigma_r^{-2} = \frac{1}{\sigma_R^2}.$$

(Recall that by convention, we order the singular values so that $\sigma_1 \geq \sigma_2 \geq \dots \geq \sigma_R$.)

Returning to the reconstruction error (13), we now see that

$$\|\hat{\mathbf{x}}_{\text{ls}} - \mathbf{x}_{\text{pinv}}\|_2^2 = \|\mathbf{V}\boldsymbol{\Sigma}^{-1}\mathbf{U}^T\mathbf{e}\|_2^2 \leq \frac{1}{\sigma_R^2} \|\mathbf{e}\|_2^2.$$

Since \mathbf{U} is an $M \times R$ matrix, it is possible when $R < M$ that the reconstruction error is zero. This happens when \mathbf{e} is orthogonal to every column of \mathbf{U} , i.e. $\mathbf{U}^T\mathbf{e} = \mathbf{0}$. Putting this together with the work above means

$$0 \leq \frac{1}{\sigma_1^2} \|\mathbf{U}^T\mathbf{e}\|_2^2 \leq \|\hat{\mathbf{x}}_{\text{ls}} - \mathbf{x}_{\text{pinv}}\|_2^2 \leq \frac{1}{\sigma_R^2} \|\mathbf{U}^T\mathbf{e}\|_2^2 \leq \frac{1}{\sigma_R^2} \|\mathbf{e}\|_2^2.$$

Notice that if σ_R is small, the worst case reconstruction error can be **very bad**.

We can also relate the “average case” error to the singular values. Say that \mathbf{e} is additive Gaussian white noise, that is each entry $e[m]$ is a random variable independent of all the other entries, and distributed

$$e[m] \sim \text{Normal}(0, \nu^2).$$

Then, as we have argued before, the average measurement error is

$$\mathbb{E}[\|\mathbf{e}\|_2^2] = M\nu^2,$$

and the average reconstruction error³ is

$$\begin{aligned} \mathbb{E} \left[\|\mathbf{A}^\dagger \mathbf{e}\|_2^2 \right] &= \nu^2 \cdot \text{trace}(\mathbf{A}^{\dagger T} \mathbf{A}^\dagger) = \nu^2 \cdot \left(\frac{1}{\sigma_1^2} + \frac{1}{\sigma_2^2} + \cdots + \frac{1}{\sigma_R^2} \right) \\ &= \frac{1}{M} \left(\frac{1}{\sigma_1^2} + \frac{1}{\sigma_2^2} + \cdots + \frac{1}{\sigma_R^2} \right) \cdot \mathbb{E}[\|\mathbf{e}\|_2^2]. \end{aligned}$$

Again, if σ_R is tiny, $1/\sigma_R^2$ will dominate the sum above, and the average reconstruction error will be quite large.

Exercise: Let \mathbf{D} be a diagonal $R \times R$ matrix whose diagonal elements are positive. Show that the maximizer $\hat{\boldsymbol{\beta}}$ to

$$\underset{\boldsymbol{\beta} \in \mathbb{R}^R}{\text{maximize}} \quad \|\mathbf{D}\boldsymbol{\beta}\|_2^2 \quad \text{subject to} \quad \|\boldsymbol{\beta}\|_2 = 1$$

has a 1 in the entry corresponding to the largest diagonal element of \mathbf{D} , and is 0 elsewhere.

³We are using the fact that if \mathbf{e} is vector of iid Gaussian random variables, $\mathbf{e} \sim \text{Normal}(\mathbf{0}, \nu^2 \mathbf{I})$, then for any matrix \mathbf{M} , $\mathbb{E}[\|\mathbf{M}\mathbf{e}\|_2^2] = \nu^2 \text{trace}(\mathbf{M}^T \mathbf{M})$. We will argue this carefully as part of the next homework.