

# Singular Value Decomposition: Reduced Density Matrix

We will introduce the singular value decomposition of a matrix in the context of the reduced density matrix of a quantum system connected to an environment.

## REDUCED DENSITY MATRIX<sup>1</sup>

Let us consider a quantum system (block)  $B$ , which is spanned by the  $N$ -dimensional orthonormal basis set  $\{|i\rangle \mid i = 1, \dots, N\}$ , surrounded by an environment  $E$ , which is spanned by the  $M$ -dimensional orthonormal basis set  $\{|j\rangle \mid j = 1, \dots, M\}$  (see the figure below).



The ground state of the total (= block + environment) system can be represented as

$$|\psi\rangle = \sum_{i=1}^N \sum_{j=1}^M \psi_{ij} |i\rangle |j\rangle. \quad (1)$$

Now consider the expectation value of an arbitrary operator,  $A$ , which acts only within the block:

$$\begin{aligned} \langle A \rangle &= \sum_i \sum_j \psi_{ij}^* \langle i|A \sum_{i'} \sum_{j'} \psi_{i'j'} |i'\rangle |j'\rangle \\ &= \sum_i \sum_j \sum_{i'} \sum_{j'} \psi_{i'j'} \psi_{ij}^* \langle i|A|i'\rangle \langle j|j'\rangle \\ &= \sum_i \sum_{i'} \sum_j \psi_{i'j} \psi_{ij}^* \langle i|A|i'\rangle \\ &\equiv \sum_i \sum_{i'} \rho_{i'i} A_{ii'} = \text{tr}_B(\rho A) \end{aligned} \quad (2)$$

where the reduced density matrix is defined as

$$\rho_{i'i} \equiv \sum_j \psi_{i'j} \psi_{ij}^* \quad (3)$$

and the matrix element of the operator is  $A_{ii'} \equiv \langle i|A|i'\rangle$ .

## SINGULAR VALUE DECOMPOSITION (SVD)

**Problem:** What is the optimal reduced density matrix  $\rho$  of rank- $m$  ( $\ll N$ )?

**Solution:** Singular value decomposition (SVD) of  $\psi \in \mathbf{R}^N \times \mathbf{R}^M$ .

(Theorem) An  $N \times M$  matrix  $\psi$  (assume  $N \geq M$ ) can be decomposed as (see [Appendix A](#) for proof of SVD and associated polar decomposition)

$$\begin{bmatrix} \psi \end{bmatrix} = \begin{bmatrix} U \end{bmatrix} \begin{bmatrix} d_1 & & \\ & \ddots & \\ & & d_M \end{bmatrix} \begin{bmatrix} V^T \end{bmatrix}, \quad (4)$$

or

$$\psi = U D V^T, \quad (5)$$

where  $U = [U_{iv} = u_i^{(v)}] \in \mathbf{R}^N \times \mathbf{R}^M$  is column orthogonal, i.e.,

$$\sum_{i=1}^N u_i^{(v)} u_i^{(v')} = \delta_{vv'}, \quad (6)$$

or

<sup>1</sup> R. P. Feynman, *Statistical Mechanics* (Benjamin/Cummings, Reading, MA, 1972) Chap. 2.

$$U^T U = I_M, \quad (7)$$

and  $V = [V_{iv} = v_i^{(v)}] \in \mathbf{R}^{M \times \mathbf{R}^M}$  is column orthogonal, *i.e.*,

$$\sum_{i=1}^M v_i^{(v)} v_i^{(v')} = \delta_{vv'}, \quad (8)$$

or

$$V^T V = I_M. \quad (9)$$

The columns of  $U$ , whose same-numbered elements  $d_v$  are *nonzero*, are an orthonormal set of basis vectors that span the range (see [Appendix B](#) for the range); the columns of  $V$ , whose same-numbered elements  $d_v$  are *zero*, are an orthonormal basis for the nullspace that is mapped to zero, *i.e.*, the subspace of  $x \in \mathbf{R}^M$ , where  $\psi x = 0$ . The program, singular.c, in the source code directory of the class home page demonstrates this automatic construction of orthonormal bases for the range and the nullspace.

### TRUNCATED SVD AS OPTIMAL APPROXIMATION

(Theorem) Let  $\psi = UDV^T$  be the SVD of  $\psi$  with the diagonal elements in descending order  $d_1 \geq d_2 \geq \dots \geq d_M$  and let

$$\psi^{(m)} \equiv \sum_{v=1}^m u^{(v)} d_v v^{(v)T}, \quad (10)$$

be the rank- $m$  truncation of the SVD. Then

$$\min_{\text{rank}(A)=m} \|A - \psi\|_2 = \|\psi^{(m)} - \psi\|_2 = d_{m+1}, \quad (11)$$

where the matrix 2-norm is defined in terms of the vector 2-norm as  $\|A\|_2 = \min_{\|x\|_2=1} \|Ax\|_2$ .

Therefore,  $\psi^{(m)}$  is the optimal rank- $m$  approximation to  $\psi$ .

Equation (10) shows that SVD is a representation of a matrix as a sum of outer products of two vectors, just as a density matrix is.

### LOW-RANK APPROXIMATION TO THE REDUCED DENSITY MATRIX

Substituting the rank- $m$  approximation (10) in the definition of the reduced density matrix, Eq. (3),

$$\begin{aligned} \rho &= \psi \psi^T \\ &= \sum_{v=1}^m \sum_{v'=1}^m u^{(v)} d_v (v^{(v)T} v^{(v')}) d_{v'} u^{(v')T} \\ &= \sum_{v=1}^m \sum_{v'=1}^m u^{(v)} d_v (d_{vv'}) d_{v'} u^{(v')T} \\ &= \sum_{v=1}^m u^{(v)} d_v^2 u^{(v)T} \end{aligned} \quad (12)$$

(Summary) The rank- $m$  truncation of the SVD of the global (= block + environment) ground state wave function,

$$\psi^{(m)} = \sum_{v=1}^m u^{(v)} d_v v^{(v)T}, \quad (13)$$

or

$$\psi_{ij}^{(m)} = \sum_{v=1}^m u_i^{(v)} d_v v_j^{(v)}, \quad (14)$$

produces the rank- $m$  approximation to the reduced density matrix,

$$\rho^{(m)} = \sum_{v=1}^m u^{(v)} w_v u^{(v)T}, \quad (15)$$

or

$$\rho_{ii'}^{(m)} = \sum_{v=1}^m u_{i(v)} w_v u_{i'(v)}, \quad (16)$$

where  $w_v = d_v^2$ . The rank- $m$  approximation  $\rho^{(m)}$  is optimal in the least square sense.

## DENSITY MATRIX RENORMALIZATION GROUP

The density matrix renormalization group (DMRG) algorithm by Steven White<sup>2</sup> is a systematic procedure to accurately obtain a quantum ground state with a modest computational cost. The DMRG incrementally add environments to the block, solve the global (= block + environment) ground state, and construct a low-rank block density matrix to represent the block with reduced degrees of freedom.

## Appendix A — Polar and Singular-Value Decompositions

### A.1 POLAR DECOMPOSITION

(Theorem) Let  $\mathbf{A}$  be a real  $N \times M$  matrix, where  $N \geq M$  (i.e., mapping from an  $M$ -dimensional source vector space to a larger  $N$ -dimensional target vector space). Then, there exists a column-wise orthogonal matrix  $\mathbf{S}$  ( $\in \mathfrak{R}^{N \times M}$  and) such that

$$\mathbf{A} = \mathbf{S}\mathbf{J}, \quad (A1)$$

$$\mathbf{S}^T \mathbf{S} = \mathbf{I}^{M \times M}, \quad (A2)$$

where  $\mathbf{I}^{M \times M}$  is the identity matrix and the unique nonnegative matrix  $\mathbf{J}$  is

$$\mathbf{J} = \sqrt{\mathbf{A}^T \mathbf{A}} \in \mathfrak{R}^{M \times M}. \quad (A3)$$

(Proof) Consider a spectral (or eigen) decomposition of  $\mathbf{J}$ :

$$\mathbf{J} = \sum_{i=1}^M \lambda_i |i\rangle \langle i|, \quad (A4)$$

where  $\lambda_i (\geq 0)$  is the  $i$ -th eigenvalue and  $\{|i\rangle \mid i = 1, \dots, M\}$  is an orthonormal set of eigenvectors. Define

$$|\psi_i\rangle = \mathbf{A}|i\rangle (\in \mathfrak{R}^N), \quad (A5)$$

then

$$\langle \psi_i | \psi_i \rangle = \langle i | \mathbf{A}^T \mathbf{A} | i \rangle = \langle i | \mathbf{J}^2 | i \rangle = \lambda_i^2. \quad (A6)$$

For those eigenvectors with  $\lambda_i \neq 0$ , define

$$|e_i\rangle = |\psi_i\rangle / \lambda_i (\in \mathfrak{R}^N), \quad (A7)$$

so that these vectors are orthonormal. For those eigenvectors with  $\lambda_i = 0$ , we use the Gram-Schmidt procedure to construct an orthonormal basis set and append it to the above basis set. Define a column-wise orthogonal matrix,

$$\mathbf{U} = \sum_{i=1}^M |e_i\rangle \langle i| \in \mathfrak{R}^{N \times M}. \quad (A8)$$

When  $\lambda_i \neq 0$ , we have

$$\mathbf{U}\mathbf{J}|i\rangle = \sum_{j=1}^M |e_j\rangle \lambda_i \underbrace{\langle j|i\rangle}_{\delta_{ji}} = \lambda_i |e_i\rangle = |\psi_i\rangle = \mathbf{A}|i\rangle. \quad (A9)$$

When  $\lambda_i = 0$ ,

$$\mathbf{U}\mathbf{J}|i\rangle = \sum_{j=1}^M |e_j\rangle \underbrace{\lambda_i}_{0} \underbrace{\langle j|i\rangle}_{\delta_{ji}} = 0 |e_i\rangle = 0 = |\psi_i\rangle = \mathbf{A}|i\rangle. \quad (A10)$$

Namely,  $\mathbf{U}\mathbf{J}$  is identical to  $\mathbf{A}$  as a mapping for the entire  $M$ -dimensional source vector space. //

<sup>2</sup> S. R. White, “Density-matrix algorithms for quantum renormalization groups,” *Physical Review B* **48**, 10345 (1993).

## A.2 SINGULAR VALUE DECOMPOSITION

(Theorem) Let  $\mathbf{A}$  be a real  $N \times M$  matrix, where  $N \geq M$  as above. Then, there exists column-wise orthogonal matrices  $\mathbf{U}$  ( $\in \mathbb{R}^{N \times M}$ ) and  $\mathbf{V}$  ( $\in \mathbb{R}^{M \times M}$ ), such that

$$\mathbf{A} = \mathbf{U}\mathbf{D}\mathbf{V}^T, \quad (\text{A11})$$

$$\mathbf{U}^T\mathbf{U} = \mathbf{V}^T\mathbf{V} = \mathbf{I}^{M \times M}, \quad (\text{A12})$$

where  $\mathbf{D}$  ( $\in \mathbb{R}^{M \times M}$ ) is a nonnegative diagonal matrix.

(Proof) Consider the polar decomposition,  $\mathbf{A} = \mathbf{S}\mathbf{J}$ , in Eq. (A1). We perform the eigen-decomposition of  $\mathbf{J}$  as

$$\mathbf{J} = \mathbf{V}\mathbf{D}\mathbf{V}^T, \quad (\text{A13})$$

where  $\mathbf{D}$  is the diagonal matrix such that its matrix elements are

$$D_{ij} = \lambda_i \delta_{ij}, \quad (\text{A14})$$

and  $\mathbf{V}$  ( $\in \mathbb{R}^{M \times M}$ ) is an orthogonal matrix, *i.e.*,  $\mathbf{V}^T\mathbf{V} = \mathbf{I}^{M \times M}$ . Substituting Eq. (A13) in Eq. (A1), we have

$$\mathbf{A} = \mathbf{S}\mathbf{V}\mathbf{D}\mathbf{V}^T \equiv \mathbf{U}\mathbf{D}\mathbf{V}^T, \quad (\text{A15})$$

Note that  $\mathbf{U} = \mathbf{S}\mathbf{V}$  ( $\in \mathbb{R}^{N \times M}$ ) is a column-wise orthogonal, since

$$\mathbf{U}^T\mathbf{U} = \mathbf{V}^T\mathbf{S}^T\mathbf{S}\mathbf{V} = \mathbf{V}^T \underbrace{\mathbf{S}^T\mathbf{S}}_{\mathbf{I} \in \mathbb{R}^{M \times M}} \mathbf{V} = \mathbf{V}^T\mathbf{V} = \mathbf{I}^{M \times M}. //$$

## Appendix B — Rank and Range of a Matrix

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For an  $N \times M$  matrix  $A$ , consider the mapping,

$$x(\in R^M) \xrightarrow{A} b = Ax(\in R^N). \quad (\text{B1})$$

The *range* of matrix  $A$  is the vector space spanned by all linearly independent vectors  $\{b\}$ , which are mapped from some  $x$ . The *rank* of matrix  $A$  is the size (*i.e.*, the number of linearly independent vectors) of its range.