

Assignment 1 — Math Fundamentals for Robotics 16-811, Fall 2024

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1. Implement the $PA = LDU$ decomposition algorithm discussed in class. Do so yourself (in other words, do not merely use predefined Gaussian elimination code in MatLab or Python). Simplifications: (i) You may assume that the matrix A is square and invertible. (ii) Do not worry about column interchanges, just row interchanges. Demonstrate (in your pdf) that your implementation works properly, on some examples.

Solution:

```
def LDU(A: np.ndarray):
    """PA = LDU decomposition algorithm

    Args:
        A (np.ndarray): input array of shape (m, n) where m >= n and A is of rank n

    Returns:
        P (np.ndarray): permutation matrix
        L (np.ndarray): lower triangular matrix
        D (np.ndarray): diagonal matrix
        U (np.ndarray): upper triangular matrix
    """

    # const A
    A = A.copy().astype(float)
    print("input matrix A:\n", A, end="\n\n")

    # verify the input matrix A is valid for LDU decomposition
    assert len(A.shape) == 2, "Input matrix A must be a 2D array"

    n_row, n_col = A.shape
    print("Input matrix A is a {}x{} matrix\n".format(n_row, n_col))

    assert n_row >= n_col, "Input matrix A must have more rows than columns"
    assert np.linalg.matrix_rank(A) == n_col, "Input matrix A must be of full rank"

    # initialize the permutation matrix P, lower triangular matrix L,
    # diagonal matrix D, and upper triangular matrix U
    P = np.eye(n_row, dtype=float)
    L = np.eye(n_row, dtype=float)
    D = np.eye(n_row, dtype=float)
    U = np.zeros((n_row, n_col), dtype=float)

    # solve for PA = LA'
    print("\nSolve for PA = LA':\n")
    A_prime = A.copy()
    for col in range(n_col):
        if A_prime[col, col] == 0:
            if np.max(A_prime[col:, col]) == 0:
                # Should not happen since A is of full rank
                continue
            # find the row with the largest absolute value in the current column
            max_row = np.argmax(np.abs(A_prime[col:, col])) + col
            print("switching row", col + 1, "with row", max_row + 1)
            # swap the current row with the row of the largest value
            # in the current column
            P[[col, max_row]] = P[[max_row, col]]
            A_prime[[col, max_row]] = A_prime[[max_row, col]]
        for row in range(col + 1, n_row):
            L[row, col] = A_prime[row, col] / A_prime[col, col]
            A_prime[row] -= L[row, col] * A_prime[col]
```

```

    print(
        "After eliminating column",
        col + 1,
        ":\nP:\n",
        P,
        "\nL:\n",
        L,
        "\nA':\n",
        A_prime,
        end="\n\n",
    )

print("After solving PA = LA'\nP:\n", P, "\nL:\n", L, "\nA':\n", A_prime)
print("PA =\n", P @ A, "\nLA' =\n", L @ A_prime, end="\n\n")

# Solve for A' = DU
print("\nSolve for A' = DU:\n")
for row in range(n_col):
    D[row, row] = A_prime[row, row]
    U[row] = A_prime[row] / D[row, row]
    print(
        "After eliminating row",
        row + 1,
        ":\nD:\n",
        D,
        "\nU:\n",
        U,
        end="\n\n",
    )

print("After solving A' = DU\nD:\n", D, "\nU:\n", U)
print("A'=\n", A_prime, "\nDU =\n", D @ U)

# Check if the decomposition is correct
assert np.array_equal(P @ A, L @ D @ U), "Decomposition is incorrect"

# Print the final result
print("\nFinal result:", "\nP:\n", P, "\nL:\n", L, "\nD:\n", D, "\nU:\n", U)

return P, L, D, U

```

Refer to my jupyter notebook for the full implementation and demonstration of the LDU decomposition algorithm other than the example given next page.

<p>P1, L1, D1, U1 = LDU(A1)</p> <hr/> <p>input matrix A:</p> <pre>[[10. -10. 0.] [0. -4. 2.] [2. 0. -5.]]</pre> <p>Input matrix A is a 3x3 matrix</p> <p>Solve for PA = LA':</p> <p>After eliminating column 1 :</p> <p>P:</p> <pre>[[1. 0. 0.] [0. 1. 0.] [0. 0. 1.]]</pre> <p>L:</p> <pre>[[1. 0. 0.] [0. 1. 0.] [0.2 0. 1.]]</pre> <p>A':</p> <pre>[[10. -10. 0.] [0. -4. 2.] [0. 2. -5.]]</pre> <p>After eliminating column 2 :</p> <p>P:</p> <pre>[[1. 0. 0.] [0. 1. 0.] [0. 0. 1.]]</pre> <p>L:</p> <pre>[[1. 0. 0.] [0. 1. 0.] [0.2 -0.5 1.]]</pre> <p>A':</p> <pre>[[10. -10. 0.] [0. -4. 2.] [0. 0. -4.]]</pre> <p>After eliminating column 3 :</p> <p>P:</p> <pre>[[1. 0. 0.] [0. 1. 0.] [0. 0. 1.]]</pre> <p>L:</p> <pre>[[1. 0. 0.] [0. 1. 0.] [0.2 -0.5 1.]]</pre> <p>A':</p> <pre>[[10. -10. 0.] [0. -4. 2.] [0. 0. -4.]]</pre>	<p>After solving PA = LA'</p> <p>P:</p> <pre>[[1. 0. 0.] [0. 1. 0.] [0. 0. 1.]]</pre> <p>L:</p> <pre>[[1. 0. 0.] [0. 1. 0.] [0.2 -0.5 1.]]</pre> <p>A':</p> <pre>[[10. -10. 0.] [0. -4. 2.] [0. 0. -4.]]</pre> <p>PA =</p> <pre>[[10. -10. 0.] [0. -4. 2.] [2. 0. -5.]]</pre> <p>LA' =</p> <pre>[[10. -10. 0.] [0. -4. 2.] [2. 0. -5.]]</pre> <p>Solve for A' = DU:</p> <p>After eliminating row 1 :</p> <p>D:</p> <pre>[[10. 0. 0.] [0. 1. 0.] [0. 0. 1.]]</pre> <p>U:</p> <pre>[[1. -1. 0.] [0. 0. 0.] [0. 0. 0.]]</pre> <p>After eliminating row 2 :</p> <p>D:</p> <pre>[[10. 0. 0.] [0. -4. 0.] [0. 0. 1.]]</pre> <p>U:</p> <pre>[[1. -1. 0.] [-0. 1. -0.5] [0. 0. 0.]]</pre> <p>After eliminating row 3 :</p> <p>D:</p> <pre>[[10. 0. 0.] [0. -4. 0.] [0. 0. -4.]]</pre> <p>U:</p> <pre>[[1. -1. 0.] [-0. 1. -0.5] [-0. -0. 1.]]</pre>	<p>After solving A' = DU</p> <p>D:</p> <pre>[[10. 0. 0.] [0. -4. 0.] [0. 0. -4.]]</pre> <p>U:</p> <pre>[[1. -1. 0.] [-0. 1. -0.5] [-0. -0. 1.]]</pre> <p>A' =</p> <pre>[[10. -10. 0.] [0. -4. 2.] [0. 0. -4.]]</pre> <p>DU =</p> <pre>[[10. -10. 0.] [0. -4. 2.] [0. 0. -4.]]</pre> <p>Final result:</p> <p>P:</p> <pre>[[1. 0. 0.] [0. 1. 0.] [0. 0. 1.]]</pre> <p>L:</p> <pre>[[1. 0. 0.] [0. 1. 0.] [0.2 -0.5 1.]]</pre> <p>D:</p> <pre>[[10. 0. 0.] [0. -4. 0.] [0. 0. -4.]]</pre> <p>U:</p> <pre>[[1. -1. 0.] [-0. 1. -0.5] [-0. -0. 1.]]</pre>
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Fig. 1: Function LDU run on A_1 of question 2

2. Compute the $PA = LDU$ decomposition and the SVD decomposition for each of the following matrices: (In fact, here it is enough to let P be an identity matrix, so $A = LDU$.)

$$A_1 = \begin{pmatrix} 10 & -10 & 0 \\ 0 & -4 & 2 \\ 2 & 0 & -5 \end{pmatrix} \quad A_2 = \begin{pmatrix} 5 & -5 & 0 & 0 \\ 5 & 5 & 5 & 0 \\ 0 & -1 & 4 & 1 \\ 0 & 4 & -1 & 2 \\ 0 & 0 & 2 & 1 \end{pmatrix} \quad A_3 = \begin{pmatrix} 1 & 1 & 1 \\ 10 & 2 & 9 \\ 8 & 0 & 7 \end{pmatrix}$$

You may use any method or tools you wish to solve this problem, including pre-defined routines from MatLab or Python, your code, and/or hand calculations. (Hand calculations may be easiest for computing the $PA = LDU$ decompositions of these examples. We recommend using pre-defined code to compute the SVD decompositions.)

Show how you obtained your solutions, that is, show your work, including intermediate steps.

Solution:

All the intermediate steps of LDU decomposition can be found in the jupyter notebook. Steps of LDU of A_1 is shown in Fig. 1.

a) **LDU Decomposition:**

i) A_1 :

$$L_1 = \begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0.2 & -0.5 & 1 \end{pmatrix} \quad D_1 = \begin{pmatrix} 10 & 0 & 0 \\ 0 & -4 & 0 \\ 0 & 0 & -4 \end{pmatrix} \quad U_1 = \begin{pmatrix} 1 & -1 & 0 \\ 0 & 1 & -0.5 \\ 0 & 0 & 1 \end{pmatrix}$$

ii) A_2 :

$$L_2 = \begin{pmatrix} 1 & 0 & 0 & 0 & 0 \\ 1 & 1 & 0 & 0 & 0 \\ 0 & -0.1 & 1 & 0 & 0 \\ 0 & 0.4 & -0.667 & 1 & 0 \\ 0 & 0 & 0.444 & 0.2083 & 1 \end{pmatrix} \quad D_2 = \begin{pmatrix} 5 & 0 & 0 & 0 & 0 \\ 0 & 10 & 0 & 0 & 0 \\ 0 & 0 & 4.5 & 0 & 0 \\ 0 & 0 & 0 & 2.667 & 0 \\ 0 & 0 & 0 & 0 & 1 \end{pmatrix} \quad U_2 = \begin{pmatrix} 1 & -1 & 0 & 0 & 0 \\ 0 & 1 & 0.5 & 0 & 0 \\ 0 & 0 & 1 & 0.222 & 0 \\ 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 1 \end{pmatrix}$$

iii) A_3 :

$$L_3 = \begin{pmatrix} 1 & 0 & 0 \\ 10 & 1 & 0 \\ 8 & 1 & 1 \end{pmatrix} \quad D_3 = \begin{pmatrix} 1 & 0 & 0 \\ 0 & -8 & 0 \\ 0 & 0 & 0 \end{pmatrix} \quad U_3 = \begin{pmatrix} 1 & 1 & 1 \\ 0 & 1 & 0.125 \\ 0 & 0 & 0 \end{pmatrix}$$

b) **SVD Decomposition:**

i) A_1 :

$$U_1 = \begin{pmatrix} -0.97502551 & 0.01928246 & -0.22125424 \\ -0.19998937 & -0.50949125 & 0.83691273 \\ -0.09658937 & 0.86025976 & 0.50062326 \end{pmatrix}$$

$$\Sigma_1 = \begin{pmatrix} 14.49778417 & 0 & 0 \\ 0 & 5.94733738 & 0 \\ 0 & 0 & 1.85564877 \end{pmatrix}$$

$$V_1^T = \begin{pmatrix} -0.68585887 & 0.72771207 & 0.00572281 \\ 0.3217144 & 0.31024647 & -0.89456524 \\ -0.65276141 & -0.61170439 & -0.44690075 \end{pmatrix}$$

ii) A_2 :

$$U_2 = \begin{pmatrix} 0.1126266 & 0.86708016 & 0.37480574 & -0.30750061 & -0.0212435 \\ -0.93215705 & 0.15225202 & 0.1626033 & 0.2846251 & 0.0212435 \\ -0.20196646 & 0.2257811 & -0.74973268 & -0.31691184 & -0.4956816 \\ -0.2398716 & -0.41225956 & 0.38334628 & -0.77085017 & -0.17702914 \\ -0.14166742 & 0.06368894 & -0.35217519 & -0.36026217 & 0.84973988 \end{pmatrix}$$

$$\Sigma_2 = \begin{pmatrix} 9.14492811 & 0 & 0 & 0 \\ 0 & 7.79814769 & 0 & 0 \\ 0 & 0 & 4.42070712 & 0 \\ 0 & 0 & 0 & 2.23976139 \\ 0 & 0 & 0 & 0 \end{pmatrix}$$

$$V_2^T = \begin{pmatrix} -0.44807922 & -0.65407164 & -0.60275097 & -0.09003647 \\ 0.65357327 & -0.69875055 & 0.28263403 & -0.06861233 \\ 0.60783153 & 0.27645026 & -0.74051747 & -0.07582844 \\ -0.05106685 & 0.08667875 & 0.09188657 & -0.99067443 \end{pmatrix}$$

iii) A_3 :

$$U_3 = \begin{pmatrix} -0.08686637 & -0.57077804 & -0.81649658 \\ -0.78592384 & -0.4643889 & 0.40824829 \\ -0.6121911 & 0.67716718 & -0.40824829 \end{pmatrix}$$

$$\Sigma_3 = \begin{pmatrix} 17.2832333 & 0 & 0 \\ 0 & 1.51322343 & 0 \\ 0 & 0 & 0 \end{pmatrix}$$

$$V_3^T = \begin{pmatrix} -0.74312678 & -0.09597244 & -0.66223249 \\ 0.1339329 & -0.99096789 & -0.0066797 \\ 0.65561007 & 0.09365858 & -0.74926865 \end{pmatrix}$$

Here Σ_3 's last singular value is smaller than 10^{-15} , so it's regarded as zero.

3. Solve the systems of equations $Ax = b$ for the values of A and b given below. For each system, specify whether the system has zero, one, or many exact solutions. If a system has zero exact solutions, give “the SVD solution” (as defined in class) and explain what this solution means. If a system has a unique exact solution, compute that solution. If a system has more than one exact solution, specify both “the SVD solution” and all solutions, using properties of the SVD decomposition of the matrix A , as discussed in class.

Show your work, including verifying that your answers are correct.

$$\begin{aligned} \text{a) } A &= \begin{pmatrix} 10 & -10 & 0 \\ 0 & -4 & 2 \\ 2 & 0 & -5 \end{pmatrix} \quad b = \begin{pmatrix} 10 \\ 2 \\ 13 \end{pmatrix} \\ \text{b) } A &= \begin{pmatrix} 1 & 1 & 1 \\ 10 & 2 & 9 \\ 8 & 0 & 7 \end{pmatrix} \quad b = \begin{pmatrix} 1 \\ 3 \\ 1 \end{pmatrix} \\ \text{c) } A &= \begin{pmatrix} 1 & 1 & 1 \\ 10 & 2 & 9 \\ 8 & 0 & 7 \end{pmatrix} \quad b = \begin{pmatrix} 3 \\ 2 \\ 2 \end{pmatrix} \end{aligned}$$

Parts (b) and (c) have the same A matrix but different b vectors. You will see that “the SVD solution” is the same for both parts. Explain why that makes sense. Your explanation should be based on the meaning of the columns of U that arose in A ’s SVD decomposition.

Solution:

```
def solve_systems_of_equations(A: np.ndarray, b: np.ndarray) -> np.ndarray:
    """Solve a system of equations of the form  $Ax = b$  for  $x$ .

    Args:
        A (np.ndarray): Input matrix A
        b (np.ndarray): Target vector b

    Returns:
        np.ndarray: Solution vector x
    """

    # copy A and b to avoid modifying the original matrices
    A = A.copy()
    b = b.copy()

    # reshape b into a column vector
    b = b.reshape(-1, 1)

    # verify that b has the same number of rows as A
    assert A.shape[0] == b.shape[0], "A and b have different numbers of rows"

    # perform SVD on A
    U, Sigma, Vt = np.linalg.svd(A)

    # factor out extremely small singular values of Sigma
    Sigma = Sigma[Sigma > 1e-10]

    # construct the pseudo-inverse of A
    M_sigma_inverse = np.zeros((U.shape[1], Vt.shape[0]))
    for i in range(len(Sigma)):
        M_sigma_inverse[i][i] = 1 / Sigma[i]
    M_pseudo_inverse = Vt.T @ M_sigma_inverse @ U.T
    print("pseudo inverse of matrix A:\n", M_pseudo_inverse)

    # compute the "solution" to the system of equations
    x = M_pseudo_inverse @ b
    return x
```

- a) $A = \begin{pmatrix} 10 & -10 & 0 \\ 0 & -4 & 2 \\ 2 & 0 & -5 \end{pmatrix}$ ’s a full rank matrix, so its column space spans \mathbb{R}^3 , and thus its systems of equations all have a unique exact solution.

$$A^\dagger = \begin{pmatrix} 0.125 & -0.3125 & -0.125 \\ 0.025 & -0.3125 & -0.125 \\ 0.05 & -0.125 & -0.25 \end{pmatrix} x = \begin{pmatrix} -1 \\ -2 \\ -3 \end{pmatrix}$$

b) SVD decomposition of A :

$$U = \begin{pmatrix} -0.08686637 & -0.57077804 & -0.81649658 \\ -0.78592384 & -0.4643889 & 0.40824829 \\ -0.6121911 & 0.67716718 & -0.40824829 \end{pmatrix}$$

$$\Sigma = \begin{pmatrix} 17.2832333 & 0 & 0 \\ 0 & 1.51322343 & 0 \\ 0 & 0 & 0 \end{pmatrix}$$

$$V^T = \begin{pmatrix} -0.74312678 & -0.09597244 & -0.66223249 \\ 0.1339329 & -0.99096789 & -0.0066797 \\ 0.65561007 & 0.09365858 & -0.74926865 \end{pmatrix}$$

Thus the pseudo-inverse of A is:

$$A^\dagger = \begin{pmatrix} -0.04678363 & -0.00730994 & 0.08625731 \\ 0.37426901 & 0.30847953 & -0.44005848 \\ 0.00584795 & 0.03216374 & 0.02046784 \end{pmatrix}$$

Σ of A has 2 non-zero diagonal entries and 1 zero diagonal entry, so the first 2 columns of U span the column space of A , and the last column of U spans the left null space of A . The last column of V spans the null space of A .

The SVD solution is:

$$A^\dagger = \begin{pmatrix} -0.04678363 & -0.00730994 & 0.08625731 \\ 0.37426901 & 0.30847953 & -0.44005848 \\ 0.00584795 & 0.03216374 & 0.02046784 \end{pmatrix} \bar{x} = \begin{pmatrix} 0.01754386 \\ 0.85964912 \\ 0.12280702 \end{pmatrix}$$

\bar{x} is exact because $A\bar{x} = b$ by verification, and there are infinitely many solutions to the system of equations, because we can add any vector in the null space of A to \bar{x} to get another solution.

c) The SVD solution is:

$$A^\dagger = \begin{pmatrix} -0.04678363 & -0.00730994 & 0.08625731 \\ 0.37426901 & 0.30847953 & -0.44005848 \\ 0.00584795 & 0.03216374 & 0.02046784 \end{pmatrix} \bar{x} = \begin{pmatrix} 0.01754386 \\ 0.85964912 \\ 0.12280702 \end{pmatrix}$$

\bar{x} is not an exact solution because $A\bar{x} = \begin{pmatrix} 1 \\ 3 \\ 1 \end{pmatrix} \neq b$, which means that b_c is not in the column space of A .

The value of $A\bar{x}$ indicates that b_c 's projection onto the column space of A is exactly b_b , in this case the \bar{x} is a least square solution, so b) and c) have the same SVD solution.

4. Suppose that u is an n -dimensional column vector of unit length in \mathbb{R}^n , and let u^T be its transpose. Then uu^T is a matrix. Consider the $n \times n$ matrix $A = I - uu^T$.
- Describe the action of the matrix A geometrically.
 - Give the eigenvalues of A .
 - Describe the null space of A .
 - What is A^2 ?

(As always, show your work.)

Solution:

- a) Ax represents the operation of projecting vector x onto the hyperplane orthogonal to u .

$$Ax = (I - uu^T)x = x - u(u^T x) = x - (u^T x)u$$

The amount of projection of x onto u is $u^T x$. Subtracting this projection from x gives the projection of x onto the hyperplane orthogonal to u .

- b) The eigenvalues of A are 1 and 0. Definition of eigenvalues:

$$\begin{aligned} \det(A - \lambda I) &= \det(I - uu^T - \lambda I) \\ &= \det((1 - \lambda)I - uu^T) = 0 \end{aligned}$$

- i) For any vector x orthogonal to u :

$$Ax = (I - uu^T)x = x - (u^T x)u = x$$

- ii) For any vector x parallel to u :

$$Ax = (I - uu^T)x = x - (u^T x)u = x - x = 0$$

- c) The null space of A is the subspace spanned by u , because any vector x parallel to u will be projected to $\vec{0}$.

$$Au = (I - uu^T)u = u - uu^T u = u - u = 0$$

- d) $A^2 = (I - uu^T)^2 = I - 2uu^T + uu^T uu^T = I - uu^T = A$.

5. The following problem arises in a large number of robotics and vision problems: Suppose p_1, \dots, p_n are the 3D coordinates of n points located on a rigid body in three-space. Suppose further that q_1, \dots, q_n are the 3D coordinates of these same points after the body has been translated and rotated by some unknown amount. Derive an algorithm in which SVD plays a central role for inferring the body's translation and rotation. (You may assume that the coordinate values are precise not noisy, but see comment and caution below.)

Show (in your pdf) that your algorithm works correctly by running it on some examples.

Comment: This problem requires some thought. There are different approaches. Although you can find a solution on the web or in a vision text book, try to solve the problem yourself before looking at any such sources. Spend some time on the problem. It is good practice to develop your analytic skills. Feel free to discuss among yourselves. (As always, cite any sources, including discussions with others.)

Requirement: Your algorithm should make use of all the information available. True, in principle you only need three pairs of points – but if you use more points your solution will be more robust, something that might come in handy some day when you need to do this for real with noisy data.

Caution: A common mistake is to derive an algorithm that finds the best affine transformation, rather than the best rigid body transformation. Even though you may assume precise coordinate values, imagine how your algorithm would behave with noise. Your algorithm should still produce a rigid body transformation.

Hint: Suppose for a moment that both sets of points have the origin as centroid. Assemble all the points p_i into a matrix P and all the points q_i into another matrix Q . Now think about the relationship between P and Q . You may wish to find a rigid body transformation that minimizes the sum of squared distances between the points q_i and the result of applying the rigid body transformation to the points p_i .

You may find the following facts useful (assuming the dimensions are sensible):

$$\|x\|^2 = x^T x, x^T R^T y = \text{Tr}(Rxy^T).$$

[Here x and y are column vectors (e.g., 3D vectors) and R is a matrix (e.g., a 3×3 rotation matrix). The superscript T means transpose, so $x^T x$ is a number and xy^T is a matrix. Also, Tr is the trace operator that adds up the diagonal elements of its square matrix argument.]

You will have more complicated expressions for x and y , involving the points p_i and q_i .

Solution: The realistic version of the problem's solution is the point-to-point *iterative closest point (ICP) algorithm* for point cloud matching/alignment, where the index of the points in the 2 point clouds are not matched, and even if they match, the points are not perfectly aligned due to imperfect perception. In more challenging scenario, we could even have missing chunks of points in one of the point clouds, where we need to use point-to-plane ICP algorithm.

Regardless, the objective is to find:

$$\arg \min_{T \in SE(3)} \sum_{i=1}^n \|q_i - T p_i\|^2$$

Where q_i and p_i are the homogeneous coordinates of q_i and p_i respectively.

This is equivalent to find:

$$\arg \min_{R \in SO(3), t \in \mathbb{R}^3} \sum_{i=1}^n \|q_i - (R p_i + t)\|^2$$

Now let \bar{q} and \bar{p} be the centroids of q_i and p_i respectively, i.e., $\bar{q} = \frac{1}{n} \sum_{i=1}^n q_i$, $\bar{p} = \frac{1}{n} \sum_{i=1}^n p_i$.

And let $\hat{q}_i = q_i - \bar{q}$, $\hat{p}_i = p_i - \bar{p}$.

Then the objective becomes:

$$\begin{aligned} & \arg \min_{R \in SO(3), t \in \mathbb{R}^3} \sum_{i=1}^n \|q_i - (R p_i + t)\|^2 \\ &= \arg \min_{R \in SO(3), t \in \mathbb{R}^3} \sum_{i=1}^n \|(\hat{q}_i + \bar{q}) - [R(\hat{p}_i + \bar{p}) + t]\|^2 \\ &= \arg \min_{R \in SO(3), t \in \mathbb{R}^3} \sum_{i=1}^n \|\hat{q}_i - R \hat{p}_i + (\bar{q} - R \bar{p} - t)\|^2 \\ &= \arg \min_{R \in SO(3), t \in \mathbb{R}^3} \left\{ \sum_{i=1}^n \|\hat{q}_i - R \hat{p}_i\|^2 + n \|\bar{q} - R \bar{p} - t\|^2 + 2 \sum_{i=1}^n (\hat{q}_i - R \hat{p}_i)^T (\bar{q} - R \bar{p} - t) \right\} \end{aligned}$$

It's easy to see that t is optimal when $t = \bar{q} - R\bar{p}$, i.e. when t is the difference of (rotated) centroids, as the second and the third term of the above equation become 0.

Now we have only R left to optimize:

$$\arg \min_{R \in SO(3)} \sum_{i=1}^n \|\hat{q}_i - R\hat{p}_i\|^2$$

Rewriting this into matrix form:

Let \hat{Q} be a $3 \times n$ matrix whose columns are \hat{q}_i , and \hat{P} be a $3 \times n$ matrix whose columns are \hat{p}_i .

$$\hat{Q} = \begin{pmatrix} \hat{q}_1 & \hat{q}_2 & \cdots & \hat{q}_n \end{pmatrix}, \hat{P} = \begin{pmatrix} \hat{p}_1 & \hat{p}_2 & \cdots & \hat{p}_n \end{pmatrix}$$

Then the objective becomes:

$$\begin{aligned} & \arg \min_{R \in SO(3)} \sum_{i=1}^n \|\hat{q}_i - R\hat{p}_i\|^2 \\ &= \arg \min_{R \in SO(3)} \|\hat{Q} - R\hat{P}\|_2^2 \\ &= \arg \min_{R \in SO(3)} \text{Tr}\{(\hat{Q} - R\hat{P})^T(\hat{Q} - R\hat{P})\} \\ &= \arg \min_{R \in SO(3)} \text{Tr}(\hat{Q}^T \hat{Q} - \hat{Q}^T R\hat{P} - \hat{P}^T R^T \hat{Q} + \hat{P}^T R^T R \hat{P}) \\ &= \arg \min_{R \in SO(3)} \text{Tr}(\hat{Q}^T \hat{Q}) - 2\text{Tr}(\hat{Q}^T R\hat{P}) + \text{Tr}(\hat{P}^T \hat{P}) \\ &= \arg \max_{R \in SO(3)} \text{Tr}(\hat{Q}^T R\hat{P}) \\ &= \arg \max_{R \in SO(3)} \text{Tr}(R\hat{P}\hat{Q}^T) \end{aligned}$$

Now let $SVD(\hat{P}\hat{Q}^T) = U\Sigma V^T$, then:

$$\begin{aligned} & \arg \max_{R \in SO(3)} \text{Tr}(R\hat{P}\hat{Q}^T) \\ &= \arg \max_{R \in SO(3)} \text{Tr}(RU\Sigma V^T) \\ &= \arg \max_{R \in SO(3)} \text{Tr}(V^T RU\Sigma) \end{aligned}$$

Since Σ is a diagonal matrix, $V^T RU$ is an orthogonal matrix, the above expression reaches maximum only when the diagonal term of $V^T RU$ reaches maximum, i.e. when $V^T RU = I$. Therefore, $R = VU^T$. However, VU^T may not be a rotation matrix, so we need to find the nearest rotation matrix to VU^T . That is, when $\det(VU^T) = -1$, we need to flip

the sign of the smallest singular value of VU^T , by $VU^T = V \begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & -1 \end{pmatrix} U^T$.

Examples of the algorithm can be found in the jupyter notebook.

```
def generate_random_points(n: int) -> np.ndarray:
    """Generate random points in 3D space.

    Args:
        n (int): number of points to generate

    Returns:
        np.ndarray: 3 x n array of random points, where each column is a point
    """
    return np.random.rand(3, n)
```

```
def generate_random_rigid_body_transformation():
    """Generate a random rigid body transformation.
```

```

Returns:
    R: 3 x 3 array representing the rotation matrix
    t: 3 x 1 array representing the translation vector
"""
# Generate a random rotation matrix
# Step 1: Generate a random axis (a unit vector)
axis = np.random.randn(3)
axis = axis / np.linalg.norm(axis) # Normalize the vector to make it a unit vector

# Step 2: Generate a random angle between 0 and 2pi
angle = np.random.uniform(0, 2 * np.pi)

# Step 3: Use Rodrigues' rotation formula to construct the rotation matrix
K = np.array([[0, -axis[2], axis[1]],
               [axis[2], 0, -axis[0]],
               [-axis[1], axis[0], 0]]) # Cross-product matrix of the axis

I = np.eye(3)
R = I + np.sin(angle) * K + (1 - np.cos(angle)) * (K @ K)

assert np.allclose(np.linalg.det(R), 1), "The generated matrix is not a rotation matrix"

# Generate a random translation vector
t = np.random.rand(3, 1)

return R, t

```

```

def ICP(P: np.ndarray, Q: np.ndarray):
    """point-to-point iterative closest point algorithm with aligned point indexes

    Args:
        P (np.ndarray): source point cloud
        Q (np.ndarray): target point cloud
    """
    P = P.copy()
    Q = Q.copy()

    assert P.shape[0] == 3 and Q.shape[0] == 3, "Input point clouds must be 3D"
    assert (
        P.shape[1] == Q.shape[1]
    ), "Input point clouds must have the same number of points"

    # Center the point clouds
    P_centered = P - np.mean(P, axis=1, keepdims=True)
    Q_centered = Q - np.mean(Q, axis=1, keepdims=True)

    U, _, Vt = np.linalg.svd(P_centered @ Q_centered.T)

    if np.linalg.det(U @ Vt) < 0:
        Vt[-1] *= -1

    R = Vt.T @ U.T
    t = np.mean(Q, axis=1, keepdims=True) - R @ np.mean(P, axis=1, keepdims=True)

    return R, t

```

```

P = generate_random_points(100)
R, t = generate_random_rigid_body_transformation()
print("R =\n", R, "\nt =\n", t)
Q = R @ P + t

```

```

R_ICP, t_ICP = ICP(P, Q)
print("R_ICP =\n", R_ICP, "\nt_ICP =\n", t_ICP)

```

```
assert np.allclose(R, R_ICP, atol=1e-3), "The rotation matrix is not correct"
assert np.allclose(t, t_ICP, atol=1e-3), "The translation vector is not correct"
```

```
P = generate_random_points(100)
R, t = generate_random_rigid_body_transformation()
print("R =\n", R, "\nt =\n", t)
Q = R @ P + t
```

[71] ✓ 0.0s

```
... R =
[[ 0.78760427  0.53794562  0.30048965]
 [-0.53690088  0.83842385 -0.09371707]
 [-0.30235237 -0.0875212   0.94916968]]
t =
[[0.16885856]
 [0.39316203]
 [0.61804859]]
```

```
R_ICP, t_ICP = ICP(P, Q)
print("R_ICP =\n", R_ICP, "\nt_ICP =\n", t_ICP)
```

[72] ✓ 0.0s

```
... R_ICP =
[[ 0.78760427  0.53794562  0.30048965]
 [-0.53690088  0.83842385 -0.09371707]
 [-0.30235237 -0.0875212   0.94916968]]
t_ICP =
[[0.16885856]
 [0.39316203]
 [0.61804859]]
```

```
assert np.allclose(R, R_ICP, atol=1e-3), "The rotation matrix is not correct"
assert np.allclose(t, t_ICP, atol=1e-3), "The translation vector is not correct"
```

[73] ✓ 0.0s

Fig. 2: ICP algorithm run on random points and transformation