

CIS PA3

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1 Cartesian Math Package

For our program, we used native Python functions and various libraries—namely Numpy, Pandas, and time— to help perform vector/matrix mathematics, frame transformations, rotations, and ICP. For performing our ICP algorithms, we were able to use a combination of these libraries, the mathematical functions we wrote in PA 1, the mathematical functions from PA 3, and our newly written functions to compute the answers. For instance, we used our frame class and registration algorithm from the first programming assignment in writing our ICP algorithm. As always, we relied on the homework instructions and the ways we learned to approach this algorithm in class to complete the assignment.

2 Algorithmic and Mathematical Approach

In the following subsections, we will break down our approach to our complex computations first algorithmically, then it mathematically, and finally we will show how it is reflected in the code. The explanation for our ICP algorithm (iterative), Find Closest Point (on surface mesh), Linear ICP algorithm, Project On Segment computation, Efficient ICP (using Covariance trees) algorithm, Computing the covariance Frame, Registration Algorithm, Construct Subtrees and Split Sort Algorithm, Transformation computation, and Composition computation are all below. Our linear, or simple ICP, will cover the behavior for all of the functions in `icp.py`, except for `project_on_segment()`, which is covered in section 2.3 of this report.

2.1 Iterative Closest Point

Our ICP approach can be explained algorithmically by the following:

2.1.1 Algorithmic Approach

1. Find the closest points on the mesh using the previous closest points (using a tree search with the threshold as the previous match).
2. Compute the registration between `d_ks`, the tip in tracker coordinates, and `c_ks`, the closest points on A to B, below the threshold
3. Calculate the error statistics, σ_n , $\bar{\epsilon}_n$, and $(\epsilon_n)_{max}$ for the current frame.
4. Adjust the threshold. First set the threshold to $3\bar{\epsilon}_n$, and if the number of valid matched points drops significantly, increase the threshold.
5. Check termination conditions and terminate, or go back to the start. As we discussed in class, the conditions for termination are not necessarily concrete. We drew inspiration from the lecture slides, and stopped our iterations when σ_n , $\bar{\epsilon}_n$, and $(\epsilon_n)_{max}$ are less than the desired threshold (0.005, 0.001, and 0.001 respectively). We also check that $\gamma \leq \frac{\bar{\epsilon}_n}{\epsilon_{n-1}} \leq 1$ over the course of several iterations, where γ is set to 0.95. Of course, we also do not iterate past a maximum number of iterations.

2.1.2 Mathematical Approach

Mathematically speaking, the approach looked like this:

1. $c_ks = \text{FindClosestPoint}$
2. $F_{reg} = \text{registration}(A, B)$, explained in section 2.9.

3. Calculate the error by:

$$s_{ks} = F_{reg} \cdot d_{ks}$$

$$\sigma_n = \frac{\sum_k \vec{e}_k \cdot \vec{e}_k}{numElements(E)}, \text{ where } E = \{\dots, \vec{e}_k, \dots\}, \vec{e}_k = \vec{c}_{ks} - \vec{s}_{ks}$$

$$(\epsilon_n)_{max} = max_k(\sqrt{\vec{e}_k \cdot \vec{e}_k})$$

$$\bar{\epsilon}_n = \frac{\sum_k \sqrt{\vec{e}_k \cdot \vec{e}_k}}{numElements(E)}$$

4. The last two steps of the algorithm are conditionals discussed above in the algorithmic approach and will be displayed in the code in the programming approach, but cannot be displayed as easily mathematically.

2.1.3 Programming Approach

Reflected in our code, this looks like: Steps 1 and 2:

```
s_ks = F_reg.compose_transform(d_ks)
prev_A = len(A)
prev_B = len(B)
# find the closest points
new_c_ks, A, B = match_points(d_ks, c_ks, s_ks, cov_tree, thresh)
c_ks = new_c_ks
# compute registration between d_ks and c_ks below the threshold
F_reg = registration(A, B)
```

Step 3 is

```
eps_n_v, e_max_v, sig_n_v = compute_error_stats(F_reg, d_ks, c_ks)
```

where compute_error_stats is defined as:

```
s_ks = F_reg.compose_transform(d_ks)
# compute the error between the tip and the surface mesh
for i in range(len(s_ks)):
    E.append((c_ks[i] - s_ks[i]))
E = np.array(E)
dot_E = []
for i in range(len(E)):
    dot_E.append(np.dot(E[i], E[i]))
for i in range(len(E)):
    eps_n += dot_E[i] ** 0.5
    sig_n += dot_E[i]
# compute the error statistics
eps_n = eps_n / len(E)
e_max = np.max(dot_E) ** 0.5
sig_n = sig_n ** 0.5 / len(E)
return eps_n, e_max, sig_n
```

Steps 4 and 5 are:

```
thresh = 3 * eps_n[-1]
if (len(A) < 0.8 * prev_A) and (len(B) < 0.8 * prev_B):
    thresh = 15 * eps_n[-1]

if eps_n[-1] < 0.005 and e_max[-1] < 0.01 and sig_n[-1] < 0.001:
    break

if n_iter > 1:
    change = eps_n[-1]/eps_n[-2]
    if 0.95 < change < 1:
        term_stack.append(change)
    elif len(term_stack) > 1:
        term_stack.pop()
    if len(term_stack) > 10:
        break
```

2.2 Linear Iterative Closest Point

Our Linear ICP approach/finding the closest point can be explained algorithmically by the following:

2.2.1 Algorithmic Approach

1. Perform a registration to calculate the poses of $F_{A,k}$ and $F_{B,k}$ (the registration algorithm and mathematical approach are explained in the registration section). Using the newly computed poses, calculate \vec{d}_k to find the position of the pointer tip with respect to rigid body B .
2. After finding a \vec{d}_k for each frame, find sample points estimated to be on the surface mesh, \vec{s}_k , such that $\vec{s}_k = F_{reg} \cdot \vec{d}_k$
3. Find points \vec{c}_k on the surface mesh that are closest to \vec{s}_k . We essentially find the closest point to the triangle by performing a least squares operation (and projecting our point onto the plane of the triangle if it is not).
4. Our last step is to use this computed point to find the distance from the point on the triangle to the mesh. If this distance is less than our bound, it is the closest distance.

2.2.2 Mathematical Approach

Mathematically, this approach looks like :

1. Using $F_{A,k}$ and $F_{B,k}$ the rigid body poses of A and B respectively, we compute \vec{d}_k , such that:

$$\vec{d}_k = F_{B,k}^{-1} \cdot F_{A,k} \cdot \vec{A}_{tip}$$

2. Compute sample points \vec{s}_k such that:

$$\vec{s}_k = F_{reg} \cdot \vec{d}_k$$

, where $F_{reg} = I$

3. Find the points \vec{c}_k on the surface mesh that are closest to \vec{s}_k , where $\vec{c}_k = F_{reg} \cdot \vec{d}_k$. This will look something like: Perform a least squares operation to solve for μ and λ in the equation

$$\vec{a} - \vec{p} = \lambda(\vec{q} - \vec{p}) + \mu(\vec{r} - \vec{p})$$

, where \vec{p} , \vec{q} , and \vec{r} are the vertices of the triangle and \vec{a} contains the sample points. Compute \vec{c} such that:

$$\vec{c} = \vec{p} + \lambda(\vec{q} - \vec{p}) + \mu(\vec{r} - \vec{p})$$

The results of λ and μ will determine where \vec{c} lies in the triangle, so if $\mu \geq 0$, $\lambda \geq 0$, and $\mu + \lambda \leq 1$, \vec{c} is in the triangle. Otherwise, find a point on the border (the approach to this is explained in the Project onto Segment section).

4. The last step to finding \vec{c}_k is to check if $\|\vec{c}_k - \vec{a}\| \leq \text{bound}$. If it is, then $\text{bound} = \|\vec{c}_k - \vec{a}\|$, and \vec{c}_k is the closest point.

2.2.3 Programming Approach

These same steps are reflected in the code by:

1. `d_ks = find_rigid_body_pose(a_read, b_read, a_tip, a_leds, b_leds)`

, where the function `find_rigid_body_pose` is :

```
Nf = len(a_frames)

# initialize the array to store the tip coordinates
d_k_cloud = np.zeros((Nf, 3))
# loop through the frames
for i in range(Nf):
    # find the rigid body poses
    F_ak = registration(a_leds, a_frames[i])
    F_bk = registration(b_leds, b_frames[i])
```

```

        # calculate the pointer tip location
        d_k = Frame.compose_transform(F_bk.invert(),
                                      Frame.compose_transform(F_ak, a_tip))
        d_k_cloud[i] = d_k
    return d_k_cloud

```

```

2.     for i in range(len(a_read)):
        F_reg = Frame(np.identity(3), np.zeros(3))
        s_k = find_sample_points(F_reg, d_ks[i])

```

where `find_sample_points` is defined as:

```

    sample_points = Frame.compose_transform(F_reg, d_k)
    return sample_points

```

```

3.         c_ks[i] = (find_closest_point(vertices, indices, s_k).reshape(1, 3))

```

where `find_closest_point` does

```

    for i in range(len(indices)):
        cur_c_k = find_closest_point_triangle(mesh_vertices[indices[i]], s_k)
        cur_d_k = find_euclidian_distance(cur_c_k, s_k)
        if cur_d_k < d_min:
            d_min = cur_d_k
            c_min = cur_c_k
    return c_min

```

and where `find_closest_point_triangle` computes:

```

    p, q, r = vertices
    A_minus_p = s_k - vertices[0]
    B = np.vstack(((q - p), (r - p))).T
    lam, mu = np.linalg.lstsq(B, A_minus_p.T, rcond=None)[0]
    c = p + lam * (q - p) + mu * (r - p)

    if lam < 0:
        c = project_on_segment(c, r, p)
    elif mu < 0:
        c = project_on_segment(c, p, q)
    elif lam + mu > 1:
        c = project_on_segment(c, q, r)
    return c

```

and `project_on_segment` does:

```

    if np.linalg.norm(p - q) == 0:
        return p

    t = np.dot(c - p, q - p) / np.dot(q - p, q - p)
    t = np.clip(t, 0, 1)
    return p + t * (q - p)

```

2.3 Project Onto Segment

Our Project Onto Segment computation is sometimes performed after finding the closest point on the triangle. If the closest point is out of bounds (not in the plane of the triangle), we must find a point on the border of the triangle instead. This section of the report covers the behavior in the function `project_on_segment` from `icp.py`.

2.3.1 Algorithmic Approach

Algorithmically, the approach looks like:

1. Perform an orthogonal projection where we project \vec{c} onto the plane of the triangle by projecting \vec{c} onto the bounding edge of the triangle between two of the vertices.
2. After computing the projection, λ , confine to the bounds between $[0,1]$ (essentially, if $\lambda \leq 0$, set it to 0, if $\lambda \geq 1$, set it to 1).
3. To finally get our correct closest point on the plane of the triangle, add λ , our corrected distance, to the vertex \vec{p} , and multiply by the bounding edge of the triangle, $\vec{q} - \vec{p}$.

2.3.2 Mathematical Approach

Mathematically speaking, the following steps are taken:

1.

$$\lambda = \frac{(\vec{c} - \vec{p}) \cdot (\vec{q} - \vec{p})}{(\vec{q} - \vec{p}) \cdot (\vec{q} - \vec{p})}$$

2.

$$\lambda^* = \text{Max}(0, \text{Min}(\lambda, 1))$$

3. Finally, we compute our \vec{c} ,

$$\vec{c} = \vec{p} + \lambda^* \times (\vec{q} - \vec{p})$$

2.3.3 Programming Approach

This is reflected in the code as:

```
def project_on_segment(c: np.ndarray, p: np.ndarray, q: np.ndarray):

    if np.linalg.norm(p - q) == 0:
        return p

    t = np.dot(c - p, q - p) / np.dot(q - p, q - p)
    t = np.clip(t, 0, 1)
    return p + t * (q - p)
```

2.4 Compute Covariance Frame

For our computation of the Covariance Frame in the Covariance Tree, we followed Dr. Taylor's "Finding Point-Pairs" slide to construct a Covariance Tree of Thing objects and used a modification of his algorithm to find the Frame.

2.4.1 Algorithmic Approach

The algorithmic approach is derived from the pseudocode on the "Finding Point-Pairs" slides. We compute the covariance frame using an approach virtually identical to the registration algorithm. We chose not to use the eigenvector/eigenvalue approach given on the slides because this approach is mathematically the same and allows for correction of the rotation if the determinant is negative. The approach is shown below:

1. Compute the mean of the corners of the triangles in the Covariance Tree
2. Calculate a matrix A by summing the outer products of the centered value of each corner
3. Find the SVD of A
4. Use the results of the SVD, specifically the V matrix to calculate the rotation matrix
5. Calculate the determinant of rotation: it should be 1, or correct the matrix if the determinant is negative one
6. Set the translation component of the frame to the mean of the corners (centroid)
- 7.

2.4.2 Mathematical Approach

Mathematically, the algorithmic approach can be represented as:

1. $centroid = corners - \bar{corners}$
2. $points = corners - centroid$

$$A = \sum_{i=1}^N \langle points, points \rangle$$
3. $A = U \Sigma V^T$
4. Steps 4 and 5 of the algorithm can be combined into:

$$R = VV$$
 where R is our Rotation matrix, assuming the determinant is 1
5. $p = centroid$
 where p is our translation component.

2.4.3 Programming Approach

Computing the covariance frame is reflected in our code in the function `compute_cov_frame`

```

1.    def compute_cov_frame(self, ts: np.ndarray):

        points = np.array([ts[i].sort_point() for i in range(len(ts))])
        n_p = len(points)
        centroid = np.mean(points, axis=0)

2.    for i in range(n_p):
        A += np.outer(points[i] - centroid, points[i] - centroid)

3.    u, s, vt = np.linalg.svd(A)
        u = u.T
        vt = vt.T

4.    comp_size = vt.shape[0]
        reflection_comp = np.eye(comp_size)
        reflection_comp[comp_size - 1][comp_size - 1] = np.linalg.det(np.dot(vt, u))

        R = np.dot(vt, np.dot(reflection_comp, vt))
        p = centroid

    return Frame(R, p)

```

2.5 Construct Subtrees and Split Sort

For our construction and sorting of the subtrees in the Covariance Tree, we followed Dr. Taylor's "Finding Point-Pairs" slide to construct a Covariance Tree of Thing objects.

2.5.1 Algorithmic Approach

The algorithmic approach to solving the subtree problem is summarized below:

1. If the number of triangles or the size of the bounds is less than our accepted limits, we do not construct further subtrees
2. We split the points in the Node into subtrees by putting each corner from the triangle into the frame of the node. If the x value is below zero, we add it to the left tree and otherwise, we add to the right tree.
3. We construct new subtrees from these subtrees in a recursive loop, ending the loop if the size of either the left or right tree is the same size as the number of points in the node (indicating that the subtrees are no longer splitting the node)

2.5.2 Mathematical Approach

Our mathematical approach for the construction and sorting of subtrees is as follows considering we have not reached the exit conditions described above:

1. Convert each corner of a triangle into the frame of the node.

$$v_x = (F_{node}^{-1} * [x, y, z]_{corner}).x$$

2. Add to the left or right subtree and return the subtrees

2.5.3 Programming Approach

Subtree construction and sorting is reflected in our code in the file `cov_tree.py`. Specifically, in the functions `construct_subtrees` and `split_sort`

1.

```
if n_t <= min_count or np.linalg.norm(self.UB - self.LB) <= min_diag:
    return None, None, False
```
2.

```
left_tree, right_tree = self.split_sort(n_t)
```

where `split_sort` is defined by:

- ```
ts = self.things
for i in range(n_t):
 # transform triangle corners into the frame of the node
 if self.F.invert().compose_transform(ts[i].sort_point().reshape(1, 3))[0][0] < 0:
 left_tree.append(ts[i])
 else:
 right_tree.append(ts[i])
return np.array(left_tree), np.array(right_tree)
```
3. 

```
if len(left_tree) == n_t or len(right_tree) == n_t:
 return None, None, False
left, right = CovTreeNode(left_tree, len(left_tree)),
 CovTreeNode(right_tree, len(right_tree))
return left, right, True
```

## 2.6 Efficient ICP

For our efficient ICP implementation, we followed Dr. Taylor's "Finding Point-Pairs" slide to construct a Covariance Tree of Thing objects to find the closest point.

### 2.6.1 Algorithmic/Mathematical Approach

I am combining our Algorithmic and Mathematical approach given that much of the algorithm involves if/else statements. I will give the mathematical representation of the statements along with their usage.

1. First, we find the  $d_k$ s as described in the Simple ICP Approach
2. We create an array of Thang (Thing) objects, representing the triangles of the surface mesh of the given object
3. We then create a CovTreeNode object using the computation of covariance frame and subtree construction algorithms described earlier. We initialize the previous closest value for this iterative approach to the first vertex in our surface mesh.
4. For the closest point computation, we iterate through each point in our  $d_k$  cloud, finding the norm between each point and our previous closest:  $\|s - d_k\|$
5. We now find the closest point:

- (a) Find the local frame of the point:

$$v_{local} = F_{node}^{-1} * v$$

- (b) Ensure  $v_{local}$  is in the bounds defined by the covariance tree

$$LB < v_{local} < UB$$

- (c) If the node has subtrees and  $v_{local}.xis < -bound$ , search the left subtree recursive for a closest point. Else, search the right.
- (d) If the node does not have subtrees, complete a linear search along the nodes points, updating the new closest point

### 2.6.2 Programming approach

Our code has the efficient ICP algorithm in `pa.three.py` in our `efficient_icp`, which calls the method `find_closest_point` in `cov_tree.py`. These two methods combined perform the efficient ICP algorithm. This looks like:

```

1. d_ks = icp.find_rigid_body_pose(a_read, b_read, a_tip, a_leds, b_leds)

2. ts = np.array([thang.Thang(vertices[indices[i]]) for i in
 range(len(indices))])

3. root = ct.CovTreeNode(ts, len(ts))
 previous_closest = ts[0].corners[0]

4. for _, s in enumerate(d_ks):
 bound = np.linalg.norm(s - previous_closest)
 closest.append(root.find_closest_point(s, bound, previous_closest))

5. previous_closest = closest[-1]
 c_ks = np.array(closest)

 mag_dif = icp.find_euclidian_distance(c_ks, d_ks)

```

where `find_closest_point` is defined as:

```

v_local = self.F.invert().compose_transform(v.reshape(1, 3))
for i in range(3):
 if v_local[0, i] > self.UB[0, i] + bound or v_local[0, i]
 < self.LB[0, i] - bound:
 return
if self.have_subtrees:
 if v_local[0, 0] < -bound:
 return self.left.find_closest_point(v, bound, closest)
 elif v_local[0, 0] > bound:
 return self.right.find_closest_point(v, bound, closest)
 else:
 left = self.left.find_closest_point(v, bound, closest)
 right = self.right.find_closest_point(v, bound, closest)
 if left is not None and right is None:
 return left
 elif left is None and right is not None:
 return right
 elif left is None and right is None:
 return closest
 else:
 return min(left, right, key=lambda x: np.linalg.norm(x - v))
else:
 for i in range(self.n_things):
 bound, closest = self.update_closest(self.things[i], v, bound, closest)
 return closest

```



## 2.7 Registration algorithm

For our registration algorithm, we used an approach developed by Arun et al., in which the authors explored a non-iterative algorithm which employs Singular Value Decomposition. Additionally, we compensated for rotation matrices with a determinant of negative one using an approach described by Sorkine-Hornung et al. This explanation will cover the behavior of the registration.py file.

### 2.7.1 Algorithmic Approach

The algorithmic approach using can be broken down into roughly five or so steps:

1. Center the point sets
2. Calculate a matrix  $H$  by multiplying one matrix by the transpose of the other
3. Find the SVD of  $H$
4. Use the results of the SVD to calculate a matrix  $X$
5. Calculate the determinant of  $X$ : it should be 1, or correct the matrix if the determinant is negative one

### 2.7.2 Mathematical Approach

Mathematically speaking, the function would look like

1.  $a_i = a - \bar{a}$   
 $b_i = b - \bar{b}$
2.  $H = \sum_{i=1}^N a_i b_i^T$
3.  $H = U \Sigma V^T$
4. Steps 4 and 5 of the algorithm can be combined into:  
 $X = V U^T$   
where  $X$  is our Rotation matrix, assuming the determinant is 1

### 2.7.3 Programming Approach

Reflected in the code, this looked like:

1. 

```
a_mean = np.mean(A.points, axis=1, keepdims=True)
b_mean = np.mean(B.points, axis=1, keepdims=True)
centered_a = A.points - a_mean
centered_b = B.points - b_mean
```
2. 

```
H = np.dot(centered_a, centered_b.transpose())
```
3. 

```
u, s, vt = np.linalg.svd(H)
```
4. 

```
u = u.transpose()
vt = vt.transpose()
```
5. 

```
comp_size = vt.shape[0]
reflection_comp = np.eye(comp_size)
reflection_comp[comp_size - 1][comp_size - 1] =
np.linalg.det(np.dot(vt, u))
```
6. And with the newly found rotation matrix, we returned a Frame:  

```
R = np.dot(vt, np.dot(reflection_comp, u))
p = b_mean - np.dot(R, a_mean)
return Frame(R, p)
```

## 2.8 Transformation

Our frame transformation function, `compose_transform`, in `frame.py` looked something like:

### 2.8.1 Algorithmic Approach

1. Iterate through each row in the rotation matrix of the frame,  $R$ , and compute the dot product of that vector with the points to transform and add to it the translation vector,  $p$ .

### 2.8.2 Mathematical Approach

Mathematically, this looks like:

1.  $v = F \cdot b$   
 $v = [R, p] \cdot b$   
 $v = R \cdot b + p$

### 2.8.3 Programming Approach

Reflected in the code, this looks like:

1. 

```
for i in range(frame_size):
 t_points[i] = np.dot(self.R, points[i]) + self.p
```

## 2.9 Composition

Our frame composition function, `compose_frame`, in `frame.py` looks something like:

### 2.9.1 Algorithmic Approach

1. Compute the dot product between Frame 1's rotation Matrix and Frame 2's rotation matrix
2. Compute the dot product between Frame 1's rotation Matrix and Frame 2's translation vector. Add to it Frame 1's translation vector

### 2.9.2 Mathematical Approach

Mathematically speaking, this looks like:

1.  $F_1 \cdot F_2 = [R_1, p_1] \cdot [R_2, p_2]$   
 $F_1 \cdot F_2 = [R_1 \cdot R_2, R_1 p_2 + p_1]$

### 2.9.3 Programming Approach

This was represented in the code with the lines:

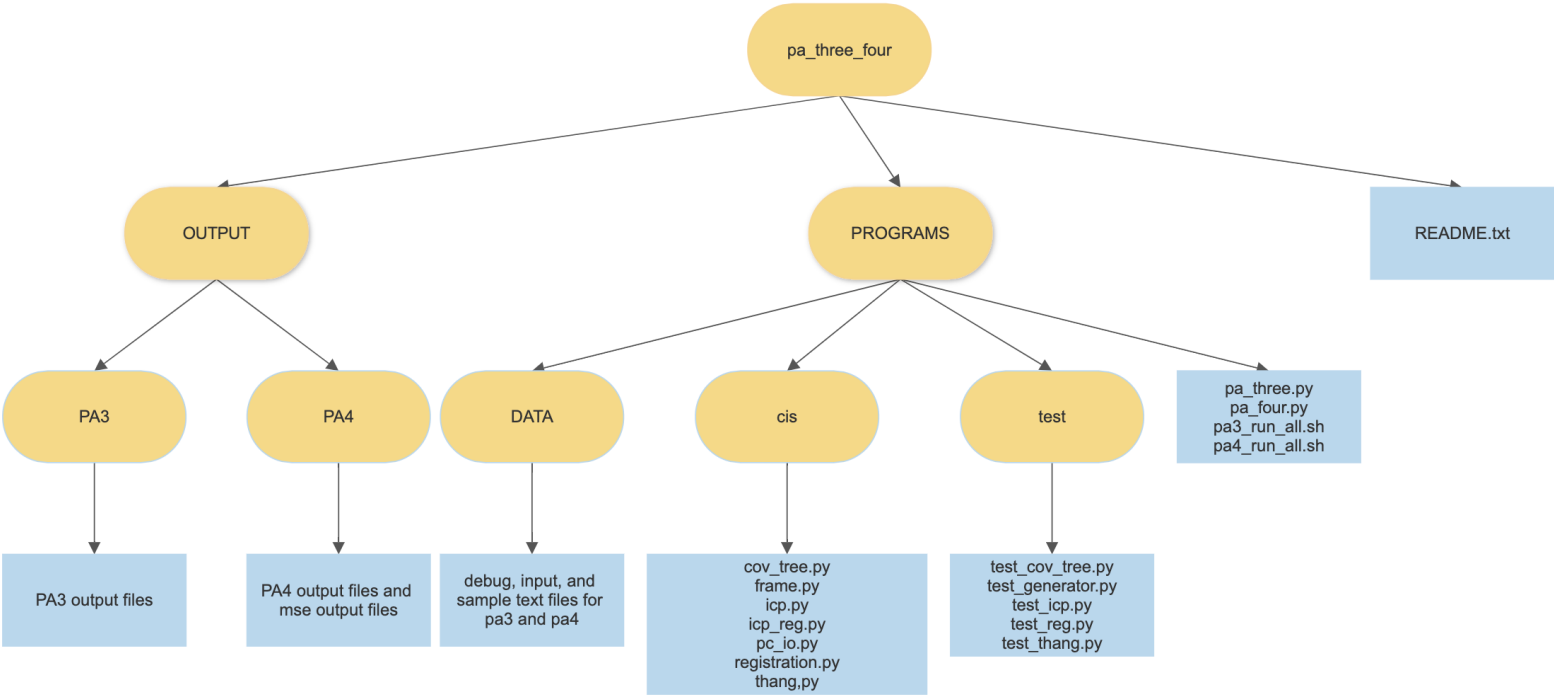
1. 

```
mat = np.dot(self.R, other_frame.R)
```
2. 

```
vec = np.dot(self.R, other_frame.p) + self.p
```

## 3 Program Structure

Our program structure is broken down into two main sections: the programs and input files, located in the `PROGRAMS` folder, and outputs, located in the `OUTPUTS` folder. Inside `PROGRAMS`, is our `README.txt`, main method for programming assignment 3, `pa_three.py`, and our main method for programming assignment 4, `pa_four.py`. We also have a folder containing all of the input files, called `DATA`, and we have a folder containing all of our test files, called `test`. Lastly, in `PROGRAMS` there is a folder titled `cis`; this is where the bulk of our program lies. `cov_tree.py`, `frame.py`, `icp.py`, `icp_reg.py`, `pc_io.py`, `registration.py`, and `thang.py`. The overall structure of our program can be seen in the diagram below, and is further elaborated following the diagram.



### 3.1 cov\_tree.py

This file contains the CovTreeNode class which is the class we used to represent Covariance Tree Functions. This class has the following fields:

1. things: a 3x3 matrix used to represent the triangles on the mesh
2. n\_things: an integer used to represent the number of things
3. F : the frame of the covariance tree node
4. UB: The upper bound of the covariance tree node
5. LB: The lower bound of the covariance tree node
6. left: The left subtree of the covariance tree node
7. right: The right subtree of the covariance tree node
8. have\_subtrees: A boolean representing whether or not the covariance tree node has subtrees

In this file there is the constructor and methods compute\_cov\_bounds, compute\_cov\_frame, construct\_subtrees, split\_sort, find\_closest\_point, and update\_closest.

1. The constructor takes in parameters ts, representing the triangles of the covariance tree node, and n\_t, the number of triangles in the covariance tree node. It initializes the fields above, and does with help from the methods compute\_cov\_frame and construct\_subtrees.
2. compute\_cov\_bounds is the method for computing the covariance bounds of a node. It takes in parameters ts, representing the triangles of the covariance tree node, and n\_t, the number of triangles in the covariance tree node, and returns UB, the upper bound of the covariance tree node, and LB, the lower bound of the covariance tree node. To compute the minimum and maximum bounds of the triangle points, we iterate through the triangles and call the method enlarge\_bounds.
3. compute\_cov\_frame is a method for computing the covariance frame of a node. It takes in the parameter ts, the triangles of the covariance tree node. To find the frame, we compute a covariance matrix  $A$  by centering the triangles. We then perform SVD resulting in vectors  $\vec{u}$  and  $\vec{v}_t$ , and validate the result by checking the determinant to account for any potential issues. Lastly, we return our newly computed frame.
4. construct\_subtrees is a method for constructing the subtrees of a node. It takes in parameters, n\_t, the number of triangles in the covariance tree node, min\_count, the minimum number of triangles in a subtree, and min\_diag, the minimum diagonal of a subtree. This function then returns the left and right subtrees of the covariance tree node.
5. split\_sort is a method for dividing the triangles of a node into two subtrees. Then transforming the triangle corners into the frame of the node, we return the left and right subtrees.
6. find\_closest\_point is the method for finding the closest point to a given point where the parameters are v, the point to find the closest point to, bound, the current closest distance, and closest, the current closest point.
7. update\_closest is the method for updating the closest point to a given point.

### 3.2 frame.py

this file was created in programming assignment 1 and contains the frame class. The functions of the class are `compose_frame`, `compose_transform`, and `invert`.

1. `compose_frame` is called on a frame and takes in another frame as the additional argument. It then computes a frame composition and returns the final frame.
2. `compose_transform` is also called on a frame and takes in a point set as the additional argument. The function computes a frame transform and returns the corresponding points.
3. `invert` is called on a frame and performs the frame inversion calculations. The final resulting frame is returned.

### 3.3 icp.py

this file contains the method (and corresponding helper methods) to our ICP algorithm. We have `find_rigid_body_pose`, `find_sample_points`, `find_closest_point`, `find_closest_point_triangle`, `project_on_segment`, `find_euclidian_distance`, and, of course, `ICP_linear`.

1. `find_rigid_body_pose` is, as the name suggests, the method for finding the rigid body pose given arguments `a_frames`, the xyz coordinates of A body LED markers in tracker coordinates, `b_frames`, the xyz coordinates of B body LED markers in tracker coordinates, `a_tip`, the xyz coordinates of the tip in tracker coordinates, `a_leds`, the xyz coordinates of A body LED markers in body coordinates, and `b_leds`, the xyz coordinates of B body LED markers in body coordinates. After iterating through the frames and finding the pose, we return `d_k_cloud`, the xyz coordinates of the pointer tip with respect to rigid body B. This method calls the registration method and the `compose_transform` method.
2. `find_sample_points` is a method for finding sample points to match to the surface mesh given `F_reg`, the frame transformation of the surface mesh from the pointer tip, and `d_k`, the xyz coordinates of the tip with respect to rigid body B. This method returns `sample_points`, which are the xyz coordinates of sample points estimated to be on the surface mesh. This method calls the `compose_transform` method.
3. `find_closest_point` is a method for finding the closest point on the surface mesh. It takes in parameters `mesh_vertices`, the xyz coordinates of the vertices of the surface mesh, `indices`, the indices of the vertices of the surface mesh, and `s_k`, the xyz coordinates of the sample points. After calculating the closest point, this method returns `c_min`, the xyz coordinates of the closest point on the surface mesh. This method also calls the methods `find_closest_point_triangle` and `find_euclidian_distance`.
4. `find_closest_point_triangle` is the method for finding the closest point on a triangle given arguments `vertices`, the xyz coordinates of the vertices of the triangle, and `s_k`, the xyz coordinates of the sample points. It returns `c` the xyz coordinates of the closest point on the triangle. This method also calls the helper method `project_on_segment`.
5. `project_on_segment` is the method for projecting a point on a segment given arguments `c`, the xyz coordinates of the point to be projected, `p`, the xyz coordinates of the first point on the segment, and `q`, the xyz coordinates of the second point on the segment. This then returns the xyz coordinates of the projected point.
6. `find_euclidian_distance` is the method for finding the euclidian distance between the sample points and the surface mesh given `c_k`, the xyz coordinates on the surface mesh found from `F_reg * d_k`, and `d_k`, the xyz coordinates of the tip with respect to rigid body B. This then returns `mag_dif`, the euclidian distance between the sample points and the surface mesh.

7. `ICP_linear` is our method for finding the rigid body pose using ICP. It takes in `mesh_vertices`, the xyz coordinates of the vertices of the surface mesh, `indices`, the indices of the vertices of the surface mesh, `a_frames`, the xyz coordinates of A body LED markers in tracker coordinates, `b_frames`, the xyz coordinates of B body LED markers in tracker coordinates, `a_tip`, the xyz coordinates of the tip in tracker coordinates, `a_leds`, the xyz coordinates of A body LED markers in body coordinates, `b_leds`, the xyz coordinates of B body LED markers in body coordinates. This then returns the xyz coordinates of the pointer tip with respect to rigid body B. This method makes calls to helper methods `find_rigid_body_pose`, `find_sample_points`, `find_closest_point`, and `find_euclidian_distance`.

### 3.4 `icp_reg.py`

This file contains all of the methods to perform an iterative ICP algorithm for PA4: `icp`, `match_points`, and `compute_error_stats`.

1. `icp` is the method for finding the rigid body pose using ICP (iteratively). It takes in the parameters `a_read`, the xyz coordinates of A body LED markers in tracker coordinates, `b_read`, the xyz coordinates of B body LED markers in tracker coordinates, `a_tip`, the xyz coordinates of the tip in tracker coordinates, `a_leds`, the xyz coordinates of A body LED markers in body coordinates, `b_leds`, the xyz coordinates of B body LED markers in body coordinates, `vertices`, the xyz coordinates of the vertices of the surface mesh, `indices`, the indices of the vertices of the surface mesh, and `max_iter`, the maximum number of iterations. It eventually returns the xyz coordinates of the pointer tip with respect to rigid body B. This function calls helper functions `find_rigid_body_pose` (from `icp.py`), `compose_transform`, `match_points`, `registration`, and `compute_error_stats`.
2. `match_points` is the method for finding the closest points on the mesh using the previous closest points. It takes in parameters `d_ks`, the xyz coordinates of the tip in tracker coordinates, `c_ks`, the estimated xyz coordinates of the vertices of the surface mesh, `s_ks`, the xyz coordinates of the tip in body coordinates, `root`, the root node of the covariance tree, and `threshold`, the threshold for the distance between the closest points. It then returns the updated closest points. This method calls the function `find_closest_points` (from `cov_tree.py`).
3. `compute_error_stats` is the method for computing the error statistics. It takes in the arguments `F_reg`, the rigid body transformation between the tip and the surface mesh, `d_ks`, the xyz coordinates of the tip in tracker coordinates, and `c_ks`, the estimated xyz coordinates of the vertices of the surface mesh. It returns the error statistics  $\sigma_n$ ,  $\bar{\epsilon}_n$ , and  $(\epsilon_n)_{max}$ . This method calls the `compose_transform` helper function.

### 3.5 `pc_io.py`

This file contains all of the methods to read from the input files and to write to the output files. We have the methods `import_rigid_body`, `import_surface_mesh`, `import_sample_readings`, `output_pa34`, `read_answer_pa3`, and `read_answer_pa4`.

1. `import_rigid_body` is the method for importing the rigid body design data. It takes in the parameter `fName`, which is the name of the data file, and, after reading the file, returns the point clouds representing the xyz coordinates of the marker LEDs in body coordinates, and the xyz coordinate of the tip in body coordinates.
2. `import_surface_mesh` is our method for importing body surface definition data. It also takes in the argument `fName`, which is the name of the data file, and after reading the file, returns the point clouds representing the xyz coordinates of vertices in CT coordinates and the xyz coordinates of the triangle indices.

3. `import_sample_readings` is our method for importing sample readings. It takes in the arguments `fName`, the name of the data file, `Na`, the number of A markers, and `Nb`, the number of B markers. After reading the file, it returns the point clouds representing frames of xyz coordinates of A body LED markers, B body LED markers, and D (unneeded) body LED markers.
4. `output_pa34` is our method for outputting PA34 data, and takes in parameters: `output_dir`, the directory to output the data, `name`, the name of the data output file, `cs`, the xyz coordinates on the surface mesh found from `F_reg * d_k`, `ds`, the xyz coordinates of the tip with respect to rigid body B, and `mag_dif`, the magnitude of the difference between the tip in CT coordinates and the tip in DCS coordinates.
5. `read_answer_pa3` is our method for reading the answer file for PA3. It takes in `fname`, the path of the answer file, and returns point clouds representing the xyz coordinates of the tip in CT coordinates.
6. `read_answer_pa4` is our method for reading the answer file for PA4. It takes in `fname`, the path of the answer file, and returns point clouds representing the xyz coordinates of the tip in CT coordinates.

### 3.6 registration.py

This file contains the registration function for our registration algorithm from assignment 1, `registration()`.

1. `registration` is our method performing a non-Iterative registration, employing Arun, Huang, and Blostein's algorithm. It takes in arguments A and B, the point cloud and the point cloud to be mapped to, respectively, and returns `Frame`, the point cloud transformation for the two point cloud inputs.

### 3.7 thang.py

This file contains the `thang` class, which is the class used to represent triangles in 3D space. This class has the field `corners`, a 3x3 matrix representation of the corners of the triangle. In addition to the constructor, it has the methods, `sort_point`, `closest_point_to`, `enlarge_bounds`, `bounding_box`, and `may_be_in_bounds`.

1. `sort_point` is the method for sorting the points of the triangle. Taking no parameters, this function returns one corner of the triangle.
2. `closest_point_to` is the method for finding the closest point on the triangle to a given point. It takes in the parameter `point`, which is the point to find the closest point on the triangle to. It calls the function `find_closest_point_triangle` from `icp.py`, and returns the resulting closest point on the triangle to the given point.
3. `enlarge_bounds` is a method for finding the bounding box of the triangle. It takes the parameters `frame`, the frame to be composed with, `LB`, the lower bound of the bounding box, and `UB`, the upper bound of the bounding box. This function then returns `LB`, the lower bound of the bounding box, and `UB`, the upper bound of the bounding box. It also calls the functions `invert` and `compose_transform`.
4. `bounding_box` is the method for finding the bounding box of the triangle. It takes in `frame`, the frame to be composed with, as an argument. It calls the function `enlarge_bounds` and then returns `LB`, the lower bound of the bounding box, and `UB`, the upper bound of the bounding box.

5. `may_be_in_bounds` is the method for checking if the triangle is in the bounding box. It takes in parameters `frame`, the frame to be composed with, `LB`, the lower bound of the bounding box, and `UB`, the upper bound of the bounding box. It calls the functions `invert` and `compose_transform` and eventually returns a boolean representing whether the triangle is in the bounding box.

### 3.8 `pa_three.py`

This file is our file that essentially runs the program. It contains our main method, the method `simple_ICP`, and the method `efficient_ICP`.

1. `main` is our main method for PA3 that runs our simple and efficient ICP algorithms. It takes in the parameters `data_dir`, the directory of the data files, `sample_readings_type`, the name of the sample readings file, `output_dir`, the directory to output the data, and `name`, the name of the data output file. This main method calls our other methods from `pc.io.py`, `import_surface_mesh`, `import_sample_readings`, and `output_pa34`, and also calls methods `simple_ICP`, and `efficient_ICP`.
2. `simple_ICP` is our method for performing a simple ICP. It takes in parameters `a_read`, the readings from the first rigid body, `b_read`, the readings from the second rigid body, `a_tip`, the tip of the first rigid body, `a_leds`, the LEDs of the first rigid body, `b_leds`, the LEDs of the second rigid body, `vertices`, the vertices of the surface mesh, and `indices`, the indices of the surface mesh. This function calls the `ICP_linear` function from `icp.py`, and returns the resulting `d_ks`, the points on the surface mesh, `c_ks`, the points on the rigid body, and `mag_dif`, the magnitude of the difference between the points.
3. `efficient_ICP` is our method for performing an efficient ICP. It takes in parameters `a_read`, the readings from the first rigid body, `b_read`, the readings from the second rigid body, `a_tip`, the tip of the first rigid body, `a_leds`, the LEDs of the first rigid body, `b_leds`, the LEDs of the second rigid body, `vertices`, the vertices of the surface mesh, and `indices`, the indices of the surface mesh. This function calls `find_rigid_body_pose` from `icp.py`, the `thang` constructor, the `CovTreeNode` constructor, `find_closest_point` from `cov_tree.py`, and `find_euclidian_distance` from `icp.py`. It returns the resulting `d_ks`, the points on the surface mesh, `c_ks`, the points on the rigid body, and `mag_dif`, the magnitude of the difference between the points.

### 3.9 `pa_four.py`

This file contains our main method for PA4, `main`, and the method to compute the mean squared error (`mse`), `mse`.

1. `main` is our main method for PA4. It takes in arguments `data_dir`, the directory of the DATA files, `sample_readings_type`, the name of the sample readings file, `output_dir`, the directory to output the data, and `name`, the name of the data output file. It calls methods `import_rigid_body`, `import_surface_mesh`, `import_sample_readings`, and `output_pa34` from `pc.io.py`. It also calls `compose_transform` from `frame.py`, `icp`, from `icp_reg.py`, `find_euclidean_distance` from `icp.py`, and `mse`.
2. `mse` is our method to compute the mean squared error of our output compared to the given output. It takes the parameters `name`, the name of the data output file, `sample_readings_type`, the type of sample readings, `output_dir`, the directory to output the data, `c_ks`, the computed `c_ks` from ICP, and `icp_time`, the time it took to run ICP. It calls on the helper method `read_answer_pa4` from `pc.io.py`.



## 4 Verification

We compared our output to the debug output to verify our program. Our results were nearly the same as the debugging output, with occasional micrometer differences that are negligible. In our main method, we added a print statement that computed and logged the MSE between the simple and efficient ICP output coordinates, and the MSE was always 0. Moreover in our output folder we have an MSE output text file for most of the results which show the MSE, which was 0, as well as the length of time it took for our efficient ICP algorithm to compute. In creating our algorithm, we originally had a perfect output every time, however the program took hours to run, and this was because there was a logical bug that caused it to build new subtrees on every iteration. So, even though our output was perfect, the program was inefficient and therefore not usable. We had to compromise a bit on accuracy in order to have a much quicker algorithm. By logging the time it took for our ICP to run, we ensured that our algorithm always finished in less than 2 minutes (closer to a minute and a half). We were able to verify not only accuracy in our computations, but speed as well, which is just as important. We also created unit tests to verify our algorithms.

## 5 Results

### 5.1 PA4-A-output.txt

| $s_x, s_y, s_z$      | $c_x, c_y, c_z$      | $ \vec{s}_k - \vec{c}_k $ |
|----------------------|----------------------|---------------------------|
| -4.77 20.39 13.36    | -4.76 20.39 13.36    | 0.002                     |
| -6.13 17.37 12.30    | -6.13 17.37 12.29    | 0.002                     |
| -0.25 4.92 -21.88    | -0.25 4.91 -21.88    | 0.003                     |
| -30.48 -21.33 -44.17 | -30.48 -21.34 -44.18 | 0.007                     |
| $\vdots$             | $\vdots$             | $\vdots$                  |

### 5.2 PA4-B-output.txt

| $s_x, s_y, s_z$      | $c_x, c_y, c_z$      | $ \vec{s}_k - \vec{c}_k $ |
|----------------------|----------------------|---------------------------|
| 6.68 -4.80 63.42     | 6.68 -4.80 63.42     | 0.001                     |
| -37.19 -17.17 -41.07 | -37.19 -17.17 -41.07 | 0.001                     |
| 28.12 19.27 12.82    | 28.12 19.27 12.82    | 0.001                     |
| -32.92 -28.12 -19.42 | -32.92 -28.12 -19.43 | 0.001                     |
| $\vdots$             | $\vdots$             | $\vdots$                  |

### 5.3 PA4-C-output.txt

| $s_x, s_y, s_z$      | $c_x, c_y, c_z$      | $ \vec{s}_k - \vec{c}_k $ |
|----------------------|----------------------|---------------------------|
| 3.14 -7.01 52.60     | 3.14 -7.02 52.60     | 0.005                     |
| 30.05 12.94 16.46    | 30.05 12.94 16.46    | 0.001                     |
| -33.87 -23.65 -13.48 | -33.87 -23.65 -13.48 | 0.001                     |
| 2.23 -12.95 1.39     | 2.23 -12.96 1.39     | 0.004                     |
| $\vdots$             | $\vdots$             | $\vdots$                  |

5.4 PA4-D-output.txt

| $s_x, s_y, s_z$      | $c_x, c_y, c_z$      | $ \vec{s}_k - \vec{c}_k $ |
|----------------------|----------------------|---------------------------|
| 4.64 10.65 -11.34    | 4.64 10.65 -11.33    | 0.002                     |
| -33.66 -26.58 -17.16 | -33.66 -26.58 -17.16 | 0.002                     |
| 3.95 19.38 0.63      | 3.96 19.38 0.63      | 0.002                     |
| -14.56 6.26 -40.59   | -14.56 6.26 -40.59   | 0.002                     |
| $\vdots$             | $\vdots$             | $\vdots$                  |

5.5 PA4-E-output.txt

| $s_x, s_y, s_z$    | $c_x, c_y, c_z$    | $ \vec{s}_k - \vec{c}_k $ |
|--------------------|--------------------|---------------------------|
| -1.54 23.96 21.66  | -1.53 24.11 21.65  | 0.154                     |
| -16.94 6.21 -13.17 | -16.93 6.17 -13.22 | 0.065                     |
| 6.56 16.61 -5.96   | 6.52 16.64 -5.99   | 0.055                     |
| 11.10 -6.58 56.98  | 11.11 -6.58 56.97  | 0.006                     |
| $\vdots$           | $\vdots$           | $\vdots$                  |

5.6 PA4-F-output.txt

| $s_x, s_y, s_z$     | $c_x, c_y, c_z$     | $ \vec{s}_k - \vec{c}_k $ |
|---------------------|---------------------|---------------------------|
| -3.42 4.74 -23.90   | -3.43 4.72 -23.90   | 0.021                     |
| -16.34 10.68 -34.00 | -16.35 10.64 -33.99 | 0.034                     |
| 5.51 20.48 43.46    | 5.51 20.46 43.46    | 0.020                     |
| 11.75 25.78 11.49   | 11.75 25.80 11.49   | 0.018                     |
| $\vdots$            | $\vdots$            | $\vdots$                  |

5.7 PA4-G-output.txt

| $s_x, s_y, s_z$     | $c_x, c_y, c_z$     | $ \vec{s}_k - \vec{c}_k $ |
|---------------------|---------------------|---------------------------|
| -31.40 2.95 -13.76  | -31.40 2.95 -13.76  | 0.001                     |
| -41.16 -3.38 -17.37 | -41.15 -3.38 -17.37 | 0.002                     |
| -8.87 -20.46 -43.13 | -8.87 -20.46 -43.13 | 0.003                     |
| -7.55 8.85 -0.39    | -7.55 8.86 -0.39    | 0.005                     |
| $\vdots$            | $\vdots$            | $\vdots$                  |

5.8 PA4-H-output.txt

| $s_x, s_y, s_z$      | $c_x, c_y, c_z$      | $ \vec{s}_k - \vec{c}_k $ |
|----------------------|----------------------|---------------------------|
| -43.06 -6.12 -30.09  | -43.06 -6.12 -30.09  | 0.001                     |
| 5.44 23.97 20.28     | 5.44 23.98 20.28     | 0.006                     |
| -37.60 -14.73 -11.31 | -37.60 -14.73 -11.31 | 0.000                     |
| 35.47 10.44 -17.03   | 35.47 10.44 -17.04   | 0.003                     |
| $\vdots$             | $\vdots$             | $\vdots$                  |

5.9 PA4-J-output.txt

| $s_x, s_y, s_z$     | $c_x, c_y, c_z$     | $ \vec{s}_k - \vec{c}_k $ |
|---------------------|---------------------|---------------------------|
| 3.09 23.84 11.19    | 3.15 23.77 11.21    | 0.095                     |
| 8.91 20.81 52.54    | 8.85 20.99 52.56    | 0.185                     |
| -1.36 -11.02 11.50  | -1.35 -10.92 11.49  | 0.098                     |
| -2.70 -22.81 -19.86 | -2.62 -22.90 -19.83 | 0.129                     |
| $\vdots$            | $\vdots$            | $\vdots$                  |

5.10 PA4-K-output.txt

| $s_x, s_y, s_z$    | $c_x, c_y, c_z$    | $ \vec{s}_k - \vec{c}_k $ |
|--------------------|--------------------|---------------------------|
| -30.55 9.02 -20.14 | -30.56 9.06 -20.13 | 0.045                     |
| -9.10 3.57 25.65   | -9.01 3.59 25.62   | 0.100                     |
| 34.06 -6.45 -12.94 | 34.08 -6.46 -12.93 | 0.025                     |
| -7.67 5.64 -30.39  | -7.68 5.63 -30.39  | 0.018                     |
| $\vdots$           | $\vdots$           | $\vdots$                  |

6 Discussion

As described in the Verification section, we added a method to write the MSE of our ICP output coordinates into another output file, as well as the length of time taken to compute, and the MSE was always 0 or close to 0. In addition to this, we added print statements in our main method to check how long our ICP algorithm took. Overall, using a cov\_tree proved to be essential in the speed and accuracy of our ICP algorithm. While the first iteration took some time to run since it was searching for an initial set of points that were far from the closest set, it ended up allowing for an overall faster algorithm once the closest points were found and our bounding boxes were constrained. At the beginning of our assignment, we faced an issue where our covariance tree was accurate but took far too long to complete each iteration, mimicking a linear time ICP operation. This issue was not noticed in PA3, as the first iteration for both PA3 and PA4 ran slowly. However, subsequent iterations in PA4 should have been rapid. We found that our find\_closest\_point method was looking for points outside of the search bounds first, determining whether to search the right tree or the left. We fixed the issue by changing our code to check both the left and right trees and return the closest point found from each subtree. This drastically increased our speed by preventing a search outside of the bounds. As discussed in the verification, compromise between speed and accuracy was required. While we were easily able to create a highly accurate and extremely slow algorithm, or an inaccurate but speedy algorithm, finding a good medium, however, was more challenging.

7 Contributions

Ilana: created the initial files and program structure, wrote the frame.py and registration.py files, helped debug/conceptually work through the ICP algorithm, and worked on README.txt. Wrote the report and program structure diagram.

Arijit: created cov\_tree.py, pc\_io.py, icp.py, icp\_reg.py, thang.py, pa\_three.py, pa\_four.py, and worked on the README.txt. Also worked on debugging and did the program testing.

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