# CS6650: Smart Sensing for Internet of Things

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#### 1 Introduction

The Euclidean Distance Geometry problem, aims at reconstructing the configuration of points given the partial information on pairwise inter-point distances.

#### 1.0.1 Mathematical Setup

Consider a collection of n points in a r-dimensional Euclidean space, ascribed to the columns of the matrix  $P = [p_1, p_2, ..., p_n]^T \in \mathbb{R}^{n \times r}$ 

The Squared distance between any two points  $p_i$  and  $p_j$  is given by,  $d_{i,j}^2 = ||p_i - p_j||_2^2 = p_i^T p_i + p_i^T p_j - 2p_i^T p_j$ .

The Euclidean Distance Matrix  $D = [d_{i,j}^2]$ 

$$D = \mathbf{1} diag(\mathbf{P}\mathbf{P}^T)^T + diag(\mathbf{P}\mathbf{P}^T)\mathbf{1}^T - 2\mathbf{P}\mathbf{P}^T$$

 $X = \mathbf{PP}^T$  is the inner product matrix well known as the Gram matrix.

#### 1.0.2 Nuclear Norm Minimization

Assuming that the entries of D are sampled uniformly at random. By definition, **X** is **symmetric** and **positive semidefinite**.

Considering the constraint that the centroid of the points is located at the origin,  $\sum_{k=1}^{n}$ 

Since the nuclear norm of a positive semidefinite matrix equate to its trace the above optimization problem can be written as,

$$\min_{\bar{X} \in \mathbb{R}^{n \times n}} Trace\left(\bar{X}\right) subject \ to \ R_{\Omega}\left(\bar{X}\right) = R_{\Omega}\left(M\right)\bar{X} \cdot \mathbf{1} = \mathbf{0}; \ \ \bar{X} \geq \mathbf{0}$$

Let a matrix,  $\mathbf{C} \in \mathbb{R}^{n \times (n-1)}$  satisfying  $\mathbf{C}^T \mathbf{C} = I$  and  $\mathbf{C}^T \mathbf{1} = \mathbf{0}$ 

Rewriting the above optimization problem, by a change of variable  $\bar{\mathbf{X}} = \mathbf{C}\mathbf{X}\mathbf{C}^T$ .

$$\min_{\mathbf{X} \in \mathbb{R}^{(n-1)\times(n-1)}} \mathit{Trace}\left(\mathbf{C}\mathbf{X}\mathbf{C}^{T}\right) \mathit{subject to} \ R_{\Omega}\left(CXC\right) = R_{\Omega}\left(M\right)\mathbf{X} \geq \mathbf{0}$$

To enforce that  $\mathbf{X}$  is positive semidefinite, let  $\mathbf{X} = \mathbf{\bar{P}}\mathbf{\bar{P}}^T$  where  $\mathbf{\bar{P}} \in \mathbb{R}^{(n-1)\times q}$ . Since,  $\mathbf{\bar{P}}\mathbf{\bar{P}}^T$  has at most rank q, a good estimate of q should be a reasonable of the rank. The minimization problem now reduces to

$$\min_{\bar{P} \in \mathbb{R}^{(n-1) \times q}} Trace\left(\mathbf{C}\bar{\mathbf{P}}\bar{\mathbf{P}}^T\mathbf{C}^T\right) subject \ to \ R_{\Omega}\left(\mathbf{C}\bar{\mathbf{P}}\bar{\mathbf{P}}^T\mathbf{C}^T\right) = R_{\Omega}\left(M\right)$$

With  $P = C\bar{P}$ , the simplified minimization problem,

$$\min_{P \subseteq \mathbb{R}^{n \times q}} Trace\left(\mathbf{P}\mathbf{P}^{T}\right) \text{ subject to } R_{\Omega}\left(\mathbf{P}\mathbf{P}^{T}\right) = R_{\Omega}\left(\mathbf{M}\right)$$

Once  $\bar{X}$  is obtained, classical MDS can then be used to find the point coordinates. Therefore, solving for **P** above and then using MDS we can retrieve the topology.

We use the method of augmented Lagrangian to solve minimization problem. The constraint  $R_{\Omega}(\mathbf{PP}^T) = R_{\Omega}(\mathbf{M})$  can be written using the linear operator **A** defined as,

$$A: \mathbb{R}^{n \times n} \to \mathbb{R}^{|\Omega| \times 1}$$
 with,  $A(X) = f \in \mathbb{R}^{|\Omega| \times 1}$ ,  $f_i = \langle X, w_{\alpha_i} \rangle$  for  $\alpha_i \in \Omega$ 

For  $y \in \mathbb{R}^{|\Omega| \times 1}$ ,  $\langle AX, y \rangle = \sum_i \langle X, w_{\alpha_i} \rangle y_i = \langle X, \sum_i y_i w_{\alpha_i} \rangle$ . From this, the adjoint of A,A\* is as follows,  $A^*y = \sum_i y_i w_{\alpha_i}$ 

In practical settings, the available partial information is not exact but noisy. Considering an additive Gaussian noise **Z** with mean  $\mu$  and variance  $\sigma$ . The modified minimization problem can be written as

$$\min \|\mathbf{X}\|_{*}$$
 subject to  $\|R_{\Omega}(\mathbf{X}) - R_{\Omega}(\mathbf{M})\|_{F} \leq \delta$ 

Using the operator A introduced earlier, the above equation can be rewritten as,

$$\min_{P} Trace \left(\mathbf{P}\mathbf{P}^{T}\right) + \frac{\lambda}{2} \|A(\mathbf{P}\mathbf{P}^{T} - \mathbf{b})\|_{F}^{2}$$

where  $\mathbf{b} = A(\mathbf{M})$ . The augmented Lagrangian is given by

$$\mathbf{L}(\mathbf{P}; \Lambda) = Trace\left(\mathbf{P}\mathbf{P}^{T}\right) + \frac{\lambda}{2} \|A(\mathbf{P}\mathbf{P}^{T} - \mathbf{b})\|_{2}^{2}$$

where  $\Lambda \in \mathbb{R}^{n \times 1}$  denotes the Lagrangian multiplier and  $\lambda$  is the penalty term. **P** is computed using the Brazilai-Borwein steepest descent method with,

objective function  $Trace(\mathbf{PP}^T) + \frac{\lambda}{2} ||A(\mathbf{PP}^T - \mathbf{b})||_2^2$ 

and gradient  $2\mathbf{P} + 2\lambda A^* \left( A \left( \mathbf{P} \mathbf{P}^T \right) - \mathbf{b} \right) \mathbf{P}$ .

**Pseudo Code Initialization**: Set q = 10,  $\mathbf{P}^0 = rand(n,q)$ ,  $E^0_{Total} = 0$ , maxiterations, biterations,  $\lambda$ , Tol

for k=1:maxiterations do

Barzilai - Borwein(BB) descent for  $\mathbf{P}_k$ 

$$E_{Total}^{k} = Trace\left(\mathbf{P}^{k}\left(\mathbf{P}^{k}\right)^{T}\right) + \frac{\lambda}{2}\|A\left(\mathbf{P}^{k}\left(\mathbf{P}^{k}\right)^{T}\right) - \mathbf{b}\|_{2}^{2}$$

if  $E_{Total}^k < Tol$  then

# 2 Importing Libraries

```
[1]: import os
     import time
     import h5py
     import math
     import scipy
     import random
     import psutil
     import platform
     import scipy.io
     import numpy as np
     import pandas as pd
     import seaborn as sns
     from math import sin, cos
     import plotly.express as px
     import matplotlib.pyplot as plt
     import plotly.graph_objects as go
     from mpl_toolkits.mplot3d import Axes3D
     from plotly.subplots import make_subplots
     %matplotlib inline
     import plotly.io as pio
     pio.renderers.default='notebook'
     import warnings
     warnings.filterwarnings('ignore')
```

#### 2.0.1 Lat and Long to 3D Coordinates

```
[2]: def getXYZ(lat,lon,R):
    lat=(lat*math.pi/180)
    lon=(lon*math.pi/180)
    x = R * cos(lat) * cos(lon)
    y = R * cos(lat) * sin(lon)
    z = R *sin(lat)
    return (x, y, z)
```

#### 2.0.2 3D Coordinates to Lat and Long

```
[3]: def getLatLon(x,y,z,R) :
    lat = math.asin(z/R)
    lon = math.atan(y/x)
    return (lat/(math.pi)*180 , lon/(math.pi)*180)
```

## 3 EDM Solver Class

```
[4]: class EDMSolver():
         HHHH
         EDM SOlver Using Low rank Matrix Completion
         def __init__(self,points,opts,lsopts):
             Initializing Variables
             Parameters
             _____
             points : 3D array of shape (...,3)
                 3D Point Cloud
             opts : Params Class obj
                 general options
             lsopts - Params Class obj
                 options for BB gradient method
             self.points = points
             self.opts = opts
             self.lsopts = lsopts
             self.pts = self.points.T
         def getEDM(self,points):
             Computes the Euclidean Distance Matrix
             Parameters
             _____
             points: 3D array of shape (n,3)
                 3D point Cloud
             Returns
             An array of Distances of shape n \times n
```

```
num_pt = points.shape[0]
    pts = points.T
    PP = np.matmul(pts.T,pts)
    Dist = PP.diagonal().reshape(-1,1) + PP.diagonal().reshape(-1,1).T - 2*PP
    for i in range(num_pt):
        Dist[i,i] = 0
    return Dist
def getWeight(self):
    Generating Weight for random missing distances
    Parameters
    _____
    None
    Returns
        Mask/Weight matrix, Weight == 1 means available distance
    num_pt = self.points.shape[0]
    rate = self.opts.rate
    Weight = np.random.rand(num_pt,num_pt)
    Weight[Weight > 1-rate] = 1
    Weight[Weight < 1] = 0
    Weight[Weight > 0] = 1
    for i in range(num_pt):
        Weight[i,i] = 1
        for j in range(i+1,num_pt):
            Weight[i,j] = Weight[j,i]
    self.ratio = ((len(np.where(Weight == 1)[0])-num_pt)/(num_pt**2))*100
    print('%f percent of the EDM Distances are known\n'%(self.ratio))
    return Weight
def AOp(self,X):
    HHHH
    The linear operator A for:
    R_{\{ \setminus Omega\} (X) = R_{\{ \setminus Omega\} (M) \}}
    Parameters
```

```
X: array_like
           The array input to the linear operator
      Returns
       _____
      y : array_like
          y = A(X)
      # diagonal of XX'
      X_diag = np.sum(np.multiply(X,X),axis=1,keepdims=True)
       # off diagonal of XX'
      X_offdiag = np.sum(np.multiply(X[self.I,:],X[self.J,:
→]),axis=1,keepdims=True)
      Y = X_diag[self.I] + X_diag[self.J] - 2*X_offdiag;
      return Y
  def AOpstar(self,y):
       The Adjoint operator of the linear operator A, A*
      A*(y) = \sum_{i=1}^{n} y^{i} \{alpha_{i}\} w_{i}
      Parameters
      y : array\_like
      Returns
       _____
      X : array_like
      X = np.zeros((self.dim,self.dim))
      X = X.reshape(-1,1)
      X[self.edgeind] = -2*y
      X_copy = X.reshape(self.dim,self.dim)
      X[self.diagind] = -np.sum(X_copy,1,keepdims=True)
      X = X.reshape((self.dim,self.dim))
      return X
  def gradient(self,P):
      11 11 11
       The gradient function for the BBGradient
      Parameters
       -----
      P : array_like
           Input to the BB Gradient Descent
      Returns
```

```
F : int
           The value of the objective function
       G: array_like
           The gradient
       tmp1 = self.AOp(P) - self.b;
       # The Objective Function
       F = np.sum(np.sum(np.multiply(P,P),axis=1)) + 0.5*self.lambda_*np.linalg.
→norm(tmp1, 'fro')**2
       # The Gradient
       G = 2*P + np.matmul(2.0*self.lambda_*self.AOpstar(tmp1),P)
       return F,G
  def addNoise(self):
       Helper function to add noise to the distance matrix
       # Gram Matrix(Inner Product Matrix)
       self.lambda_ = self.opts.lambda_
       self.Rk = self.opts.rank
       self.dim,_ = self.Dist.shape
       # Gram Matrix(Inner Product Matrix)
       self.Gram_Truth = self.Dist - np.matmul(np.mean(self.
→Dist,axis=1,keepdims=True),np.ones((1,self.dim)))
       self.Gram_Truth = -0.5*(self.Gram_Truth - np.matmul(np.ones((self.
→dim,1)),np.mean(self.Gram_Truth,axis=0,keepdims=True)))
       # Obtaining Omega i.e., indices of randomly choosen entries of D
       self.I,self.J = np.where(self.Weight == 1)
       # Diagonal indices
       self.diagind = np.array(range(1,self.dim+1)) + self.dim*np.
→array(range(self.dim)) - 1
       # Edge indices of choosen entries of D (includes symmetrical indices)
       self.edgeind = self.I + (self.J)*self.dim
       M_noisefree = self.Dist.reshape(-1,1)[self.edgeind]
       # Minimum sampled distance
       mindist = np.min(M_noisefree[M_noisefree > 0])
       # Adding noise in the range of the minimum measured distance
       #and ensuring all distances are positive
```

```
M = M_noisefree + 4*mindist + mindist*np.random.
→standard_normal(M_noisefree.shape);
      M[M_noisefree == 0] = 0
      self.b = M
  def BBGradient(self,x):
       The Brazilai-Borwein steepest descent method
       # The Objective and the gradient initialization
       [f,g] = self.gradient(x)
      # Norm of the Gradient
      gnorm = np.linalg.norm(g,'fro')
      # Initialize Q, C and alpha
      Q = 1;
      C = f;
      alpha = self.lsopts.alpha
      # Initializing xmin and fmin
      xmin = x
      fmin = f
      n = x.shape[0]
       # Maximum and Minimum values of alpha
      MAXalpha = 1e16
      MINalpha = 1e-16
      for it in range(self.lsopts.maxit):
           xold = x
           fold = f
           gold = g
           numlinesearch = 1
           wolfe_factor = self.lsopts.rho*gnorm**2
           while True:
              x = xold - alpha*gold
               [f,g] = self.gradient(x)
               if f <= C - alpha*wolfe_factor or numlinesearch >= 5:
                   break
               alpha = self.lsopts.sigma*alpha
               numlinesearch = numlinesearch + 1
```

```
if f < fmin:</pre>
               xmin = x
               fmin = f
           # Stopping Conditions
           gnorm = np.linalg.norm(g, 'fro')
           s = x - xold
           xstop = np.linalg.norm(s,'fro')/np.sqrt(n)
           fstop = abs(fold - f)/(abs(fold) + 1)
           if (xstop < self.lsopts.xtol and self.lsopts.ftol) or (gnorm < self.
→lsopts.gtol):
               break
           # Updates
           y = g - gold
           normy = np.linalg.norm(y,'fro')
           sy = abs(np.trace(np.matmul(s.T,y)))
           if (it+1)\%2 == 0:
               alpha = (np.linalg.norm(s, 'fro')**2)/sy
           else:
               alpha = sy/(normy**2)
           # To make sure alpha lies in bound
           alpha = min(alpha,MAXalpha)
           alpha = max(alpha,MINalpha)
           # Updating Cost
           Qold = Q
           Q = self.lsopts.eta*Qold + 1
           C = (self.lsopts.eta*Qold*C + f)/Q
       x = xmin
       return x
  def matrixCompletion(self):
       Performs the BB method to solve for P, getting the gram matrix and
       reconstructed Gram matrix and performing canonical Multidimensional
       scaling(MDS)
       HHHH
       ## Initializing Values
       # P, lagrangian Multipliers
       P = np.random.random((self.dim,self.Rk))
       # Energies
```

```
E = np.zeros((self.opts.maxit,1))
       # Iteration counter, relative error
      num_it = 0
      cre = 1
      # Barzilai-Borwein(BB) descent to solve for P
      for i in range(self.opts.maxit):
          num_it = num_it + 1
          # Gradient descent for P
          P = self.BBGradient(P)
           # Calculating total energy
          E[i] = np.sum(np.multiply(P,P)) + 0.5*self.lambda_*np.linalg.
→norm(self.AOp(P)-self.b,'fro')**2
           # To print energy
          if self.opts.printenergy == 1:
               print("Iteration %d, TotalE = %f\n" % (i,E[i]))
           # Calculating relative error
          if i > 0:
               cre = abs(E[i] - E[i-1])/E[i]
           # stopping condition
          if cre < 1e-5:
              break
       # Plotting Energies with respect to iteration number
      if self.opts.plotting == 1:
          plt.figure(figsize=(12,8))
          plt.semilogy(range(num_it), E[:num_it], linewidth=2)
          plt.xlabel("Iteration number",fontsize=18)
          plt.ylabel("$\log(E)$",fontsize=18)
          plt.grid()
          plt.show()
       # Obtaining the Gram Matrix
      Gram = np.matmul(P,P.T)
       # Reconstructing the Gram Matrix and Multidimensional Scaling (MDS)
      self.Gram_Recon = Gram - (1/self.dim)*np.tile(np.
→sum(Gram,axis=0,keepdims=True),(self.dim,1)) + (1/(self.dim**2))*np.tile(np.
→sum(Gram),(self.dim,self.dim))
      self.Gram_Recon = (self.Gram_Recon + self.Gram_Recon.T)/2
      self.Gram_err = np.linalg.norm(self.Gram_Truth - self.Gram_Recon, 'fro')/
→np.linalg.norm(self.Gram_Truth, 'fro')
```

```
D,V = scipy.sparse.linalg.eigs(self.Gram_Recon,self.opts.rank,which='LM')
      D = D.reshape(-1,1)
      D,IJ = -np.sort(-D,axis=0),np.argsort(-D,axis=0)
      V = V[:,IJ]
      temp = np.real(V[:,0:3])
      temp = temp.reshape(temp.shape[0],temp.shape[1]*temp.shape[2])
      Global_Coordinate = np.matmul(temp,np.diag(np.sqrt(D[0:3]).reshape(-1)))
      output = dict()
      output['E'] = E[num_it-1]
      output['ReconError'] = self.Gram_err
      output['numit'] = num_it
      return Global_Coordinate, self.Gram_Recon, output
  # helper function 1
  def plot3d(self,coords,title=None):
      Helper function to plot the 3D plots
      Parameters
       coords : array_like
           The 3D point cloud locations (x, y, z)
      x, y, z = coords[:,0],coords[:,1],coords[:,2]
      data = go.Scatter3d(x=x,y=y,z=z,mode='markers',
              marker=dict(size=5,color=z,colorscale='Viridis')) # set color to_
→an array/list of desired values
      fig = go.Figure(data=[data])
       # tight layout
      fig.update_layout(margin=dict(l=0, r=0, b=0, t=0),title=title)
      fig.show()
  # helper function 2
  def rigid_transform_3D(self,A, B):
      Helper function to fix the orientation of the reconstructed poit cloud
      Parameters
      A : array\_like of shape (3,...)
      B : array_like of shape (3,...)
```

```
Returns
       _____
       R : array_like of shape (3,3)
           Rotation Matrix
       t : array_like of shape(3,1)
           Translation Matrix
       Refrence
       This function has been directly used from the below github repository
       https://github.com/nghiaho12/rigid_transform_3D
       assert A.shape == B.shape
       num_rows, num_cols = A.shape
       if num_rows != 3:
           raise Exception(f"matrix A is not 3xN, it is {num_rows}x{num_cols}")
       num_rows, num_cols = B.shape
       if num_rows != 3:
           raise Exception(f"matrix B is not 3xN, it is {num_rows}x{num_cols}")
       # find mean column wise
       centroid_A = np.mean(A, axis=1)
       centroid_B = np.mean(B, axis=1)
       # ensure centroids are 3x1
       centroid_A = centroid_A.reshape(-1, 1)
       centroid_B = centroid_B.reshape(-1, 1)
       # subtract mean
       Am = A - centroid_A
       Bm = B - centroid_B
       H = Am @ np.transpose(Bm)
       # sanity check
       #if linalg.matrix_rank(H) < 3:</pre>
           raise ValueError("rank of H = {}, expecting 3".format(linalg.
\rightarrow matrix\_rank(H)))
       # find rotation
       U, S, Vt = np.linalg.svd(H)
       R = Vt.T @ U.T
```

```
# special reflection case
    if np.linalg.det(R) < 0:
        print("det(R) < R, reflection detected!, correcting for it ...")</pre>
        Vt[2,:] *= -1
        R = Vt.T @ U.T
    t = -R @ centroid_A + centroid_B
    return R, t
# helper function 4
def run(self):
    Helper function to combine all the above functions and run it in a
    sequential manner
    Returns
    _____
    output : dict
        keys:
                      : smapling rate: actual 3D point cloud
            rate
           points
            tpoints
                       : transformed reconstructed point cloud
            system
                       : system parameters
            release
                      : system parameters
            Version
                       : system parameters
            Machine : system parameters
           Processor : system parameters
            error
                        : Obtained error
                       : Number of BB iterations
            numit
            timeelapsed : Total time elapsed in computing
    H H H
    # Recording starting time
    start = time.time()
    # Obtaining the EDM Matrix
    self.Dist = self.getEDM(self.points)
    self.scaler = np.max(self.Dist)
    # Normalizing the entries of EDM matrix
    self.Dist = self.Dist/self.scaler
    # Getting the mask
    self.Weight = self.getWeight()
    # Adding noise to the randomly choosen distances
    self.addNoise()
```

```
# Matrix Completion and Multi Dimensional Scaling(MDS)
      self.GCor, ipm, _output = self.matrixCompletion()
      self.GCor = np.real(self.GCor)
       # Rescaling the points to original ranges
      self.GCor = self.GCor*np.sqrt(self.scaler)
      if self.opts.plotting:
           # 3D View of Constructed points
          self.plot3d(self.GCor, 'Constructed Points')
           # 3D View of Actual points
          self.plot3d(self.pts.T,'Actual Points')
       # 3D transformation to original orientation
      R,t = self.rigid_transform_3D(self.GCor.T,self.points.T)
      self.transformed_points = (R @ self.GCor.T) + t
      self.transformed_points = self.transformed_points.T
      if self.opts.plotting:
           # plotting the 3D transformed points
          self.plot3d(self.transformed_points, 'Transformed Reconstructed_
→Points')
       # Record finish time
      end = time.time()
       # Error Valuation
      error = np.linalg.norm(self.points-self.transformed_points, 'fro')/np.
→linalg.norm(self.points,'fro') # relative error
      output = dict()
      uname = platform.uname()
      output['ratio'] = self.ratio
      output['rate'] = self.opts.rate
      output['points'] = self.points
      output['tpoints'] = self.transformed_points
      output['System'] = uname.system
      output['Release'] = uname.release
      output['Version'] = uname.version
      output['Machine'] = uname.machine
      output['Processor'] = uname.processor
      output['error'] = error
      output['numit'] = _output['numit']
      output['timeelapsed'] = end - start
```

```
return output
```

#### 3.1 Parameters Class

```
[5]: class params():
         lambda_ = None
         printenergy = None
         plotting = None
         rank = None
         rate = None
         maxit = None
         tol = None
         maxit = None
         xtol = None
         gtol = None
         ftol = None
         alpha = None
         rho = None
         sigma = None
         eta = None
```

## 4 Dataset Generation

# 4.1 Dataset 1: Mapping the major cities of a country

## 4.1.1 Loading Countries Location Data

```
[6]: WorldData = pd.read_csv('./data/worldcitiespop.csv')
WorldData
```

```
AccentCity Region Population
[6]:
              Country
                                City
                                                                           Latitude
                                                                          42.483333
     0
                    ad
                               aixas
                                             Aixàs
                                                          6
                                                                     NaN
                                        Aixirivali
     1
                    ad
                         aixirivali
                                                          6
                                                                     {\tt NaN}
                                                                          42.466667
     2
                                        Aixirivall
                                                          6
                                                                     NaN
                                                                          42.466667
                    ad
                         aixirivall
     3
                                                          6
                    ad
                          aixirvall
                                         Aixirvall
                                                                     NaN
                                                                          42.466667
     4
                                                          6
                    ad
                           aixovall
                                          Aixovall
                                                                     {\tt NaN}
                                                                          42.466667
                                 . . .
                                                        . . .
                                                                     . . .
     . . .
                   . . .
     3173953
                         zimre park
                                        Zimre Park
                                                         4
                                                                     NaN -17.866111
                   ZW
     3173954
                        ziyakamanas Ziyakamanas
                                                         0
                                                                     NaN -18.216667
                   zw
     3173955
                         zizalisari
                                        Zizalisari
                                                          4
                                                                     NaN -17.758889
                    zw
                                           Zuzumba
                                                          6
                                                                     NaN -20.033333
     3173956
                             zuzumba
                    ZW
     3173957
                    zw
                         zvishavane
                                        Zvishavane
                                                                79876.0 -20.333333
```

```
Longitude
0 1.466667
1 1.500000
```

```
2 1.500000

3 1.500000

4 1.483333

... ... ...

3173953 31.213611

3173954 27.950000

3173955 31.010556

3173956 27.933333

3173957 30.033333

[3173958 rows x 7 columns]
```

#### 4.1.2 Selecting A Country: India

```
[7]: Data = WorldData[WorldData['Country'] == 'in']
Data = Data.reset_index(drop = True)
Data.shape
```

[7]: (39813, 7)

#### 4.1.3 Extracting and saving 3D Point Cloud

```
[8]: points1 = []
for i in random.sample(range(0, Data.shape[0]), 2500):
    points1.append(getXYZ(Data['Latitude'][i],Data['Longitude'][i],6371))

points1 = np.array(points1)
    print(points1.shape)

np.savetxt('./data/india_latlon.csv', points1, delimiter=',')

(2500, 3)
```

#### 4.1.4 Viewing the 3D point Cloud

#### 4.2 Dataset 2: Sewage Pipes Data

#### 4.2.1 Extracting and saving 3D Point Cloud

```
[10]: path = './data/training_pointcloud_hdf5_real.h5'
with h5py.File(path, 'r') as hdf:
    Data = np.asarray(hdf['Training/PointClouds'][:])
Data = np.array(Data)
print(Data.shape)

np.savetxt('./data/pipe.csv',Data[0,:,:],delimiter=',')

points2 = Data[0,:,:]
(274, 1024, 3)
```

## 4.2.2 Viewing the 3D point Cloud

# 5 Dataset 3: Location data collected from our peers

## 5.1 Visualisation

```
fig.show()
```

# 5.2 Dataset 4: Intramolecular Data to identify protein structure

```
[14]: f = open("./data/atom.txt", "r")
Lines = f.readlines()
count = 0
points4 = []
for line in Lines:
    count += 1
    x=line.split()
    try:
        t=[float(x[6]),float(x[7]),float(x[8])]
        points4.append(t)
    except:
        continue
points4 = np.array(points4)
```

#### 5.3 Visualisation

## 5.4 Running the Solver

## 5.4.1 Defining EDMSolver Class Parameters

```
[16]: opts = params()
lsopts = params()

# Matrix Completion parameters
opts.lambda_ = 5
opts.rate = 0.08
opts.printenergy = 0
opts.plotting = 0
opts.rank = 10
opts.maxit = 30
opts.tol = 1e-5
```

```
# BBGradient parameters
lsopts.maxit = 20
lsopts.xtol = 1e-8
lsopts.gtol = 1e-8
lsopts.ftol = 1e-10
lsopts.alpha = 1e-3
lsopts.rho = 1e-4
lsopts.sigma = 0.1
lsopts.eta = 0.8
```

#### 5.4.2 Defining and Running the Solver

```
[17]: solver1 = EDMSolver(points1,opts,lsopts)
    solver2 = EDMSolver(points2,opts,lsopts)
    solver3 = EDMSolver(points3,opts,lsopts)
    solver4 = EDMSolver(points4,opts,lsopts)

output1 = solver1.run()
    output2 = solver2.run()
    output3 = solver3.run()
    output4 = solver4.run()

output5 = [output1,output2,output3,output4]
    solvers = [solver1,solver2,solver3,solver4]
    for i in range(len(outputs)):
        outputs[i]['Dataset'] = 'Dataset '+str(i+1)
```

8.003296 percent of the EDM Distances are known

7.968140 percent of the EDM Distances are known

10.666667 percent of the EDM Distances are known

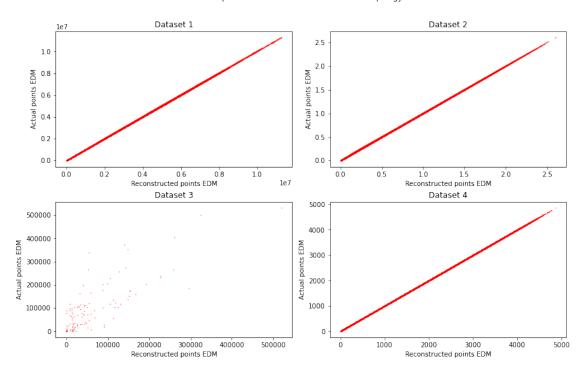
det(R) < R, reflection detected!, correcting for it ...
7.986740 percent of the EDM Distances are known</pre>

```
fig.add_trace(go.Scatter3d(x=points1[:,0], y=points1[:,1], z=points1[:
 →,2],mode='markers',marker=dict(size=2,color=points1[:
 →,2],colorscale='Viridis')),
             row=1, col=1)
fig.add_trace(go.Scatter3d(x=points2[:,0], y=points2[:,1], z=points2[:
 →,2],mode='markers',marker=dict(size=2,color=points2[:
 →,2],colorscale='Viridis')),
              row=1, col=2)
fig.add_trace(go.Scatter3d(x=points3[:,0], y=points3[:,1], z=points3[:
 →,2],mode='markers',marker=dict(size=2,color=points3[:
 →,2],colorscale='Viridis')),
              row=2, col=1)
fig.add_trace(go.Scatter3d(x=points4[:,0], y=points4[:,1], z=points4[:
 →,2],mode='markers',marker=dict(size=2,color=points4[:
 →,2],colorscale='Viridis')),
             row=2, col=2)
fig.update_layout(height=1000, showlegend=False)
fig.show()
```

```
[19]: # Transformed points
      tpoints1,tpoints2,tpoints3,tpoints4 =
       -output1['tpoints'],output2['tpoints'],output3['tpoints'],output4['tpoints']
      fig = make_subplots(
          rows=2, cols=2,
          specs=[[{"type": "scene"}, {"type": "scene"}],
                 [{"type": "scene"}, {"type": "scene"}]],
      )
      fig.add_trace(go.Scatter3d(x=tpoints1[:,0], y=tpoints1[:,1], z=tpoints1[:
       →,2],mode='markers',marker=dict(size=2,color=tpoints1[:
       →,2],colorscale='Viridis')),
                    row=1, col=1)
      fig.add_trace(go.Scatter3d(x=tpoints2[:,0], y=tpoints2[:,1], z=tpoints2[:
       →,2],mode='markers',marker=dict(size=2,color=tpoints2[:
       →,2],colorscale='Viridis')),
                    row=1, col=2)
```

```
fig,ax = plt.subplots(2,2,figsize=(12,8))
for i in range(len(outputs)):
    solver,output = solvers[i],outputs[i]
    dact = np.tril(solver.getEDM(output['points'])).ravel()
    d0bs = np.tril(solver.getEDM(output['tpoints'])).ravel()
    ax[int(i/2),int(i%2)].scatter(d0bs,dact,color='r',s=0.1)
    ax[int(i/2),int(i%2)].set_title('{}'.format(output['Dataset']))
    ax[int(i/2),int(i%2)].set_xlabel('Reconstructed points EDM')
    ax[int(i/2),int(i%2)].set_ylabel('Actual points EDM')

plt.tight_layout()
fig.subplots_adjust(top=0.88)
fig.suptitle('EDM comparision of actual and reconstructed topology')
plt.show()
```



Here we have seen different datasets and sampled them to obtain our dataset. The EDM after construction has been sampled at certain rate and the following errors and properties have been noted.

```
[21]: print(' Dataset
                                           Error(%)
                          Rate
                                 Ratio
                                                      Num Iterations
                                                                      Num Points Time
       →Elapsed(s)')
      for output in outputs:
          print(output['Dataset']," %0.2f
                                               %0.2f
                                                       %2.6f
                                                                     %0.2f
                                                                                   %5d
       →4f"%(output['rate'],output['ratio'],output['error']*100,output['numit'],output['points'].
       →shape[0],output['timeelapsed']))
      Dataset
                  Rate
                         Ratio
                                  Error(%)
                                              Num Iterations
                                                              Num Points Time
     Elapsed(s)
     Dataset 1
                                                 26.00
                                                                 2500
                  0.08
                         8.00
                                0.189246
                                                                         79.2980
     Dataset 2
                  0.08
                         7.97
                                0.311351
                                                 17.00
                                                                 1024
                                                                         8.7557
                                                  6.00
                                                                         0.0190
     Dataset 3
                  0.08
                         10.67
                                 2.217214
                                                                   15
                                                 27.00
                                                                         258.3121
     Dataset 4
                  0.08
                         7.99
                                0.027376
                                                                 4384
```

#### 5.5 Error Analysis: Dataset 1

#### 5.5.1 Defining EDMSolver Class Parameters

```
[22]: opts = params()
      lsopts = params()
      # Matrix Completion parameters
      opts.lambda_ = 5
      # opts.rate = 0.02
      opts.printenergy = 0
      opts.plotting = 0
      opts.rank = 10
      opts.maxit = 30
      opts.tol = 1e-5
      # BBGradient parameters
      lsopts.maxit = 20
      lsopts.xtol = 1e-8
      lsopts.gtol = 1e-8
      lsopts.ftol = 1e-10
      lsopts.alpha = 1e-3
      lsopts.rho = 1e-4
      lsopts.sigma = 0.1
      lsopts.eta = 0.8
```

#### 5.5.2 Defining and Running the Solver

```
coutputs1 = []
rates1 = [0.01,0.02,0.03,0.05]

points = np.genfromtxt('./data/india_latlon.csv', delimiter=',')

for rate in rates1:
    opts.rate = rate
    solver = EDMSolver(points,opts,lsopts)
    output = solver.run()
    outputs1.append(output)

0.994080 percent of the EDM Distances are known

1.981792 percent of the EDM Distances are known

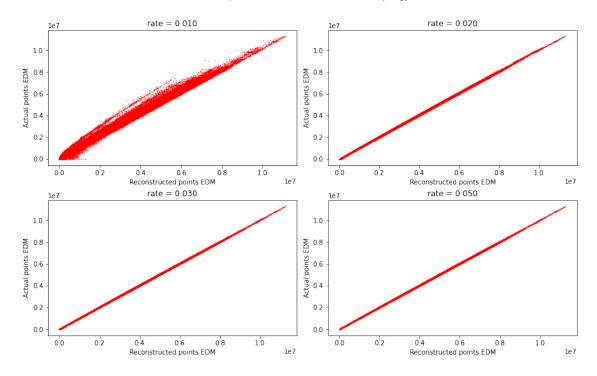
det(R) < R, reflection detected!, correcting for it ...
2.994368 percent of the EDM Distances are known

det(R) < R, reflection detected!, correcting for it ...
5.016000 percent of the EDM Distances are known</pre>
```

```
fig,ax = plt.subplots(2,2,figsize=(12,8))
for i in range(len(outputs1)):
    output = outputs1[i]
    dact = np.tril(solver.getEDM(output['points'])).ravel()
    d0bs = np.tril(solver.getEDM(output['tpoints'])).ravel()
    ax[int(i/2),int(i%2)].scatter(d0bs,dact,color='r',s=0.1)
    ax[int(i/2),int(i%2)].set_title('rate = %.3f'%output['rate'])
    ax[int(i/2),int(i%2)].set_xlabel('Reconstructed points EDM')
    ax[int(i/2),int(i%2)].set_ylabel('Actual points EDM')

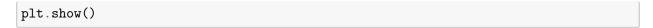
plt.tight_layout()
fig.subplots_adjust(top=0.88)
fig.suptitle('EDM comparision of actual and reconstructed topology')
plt.show()
```

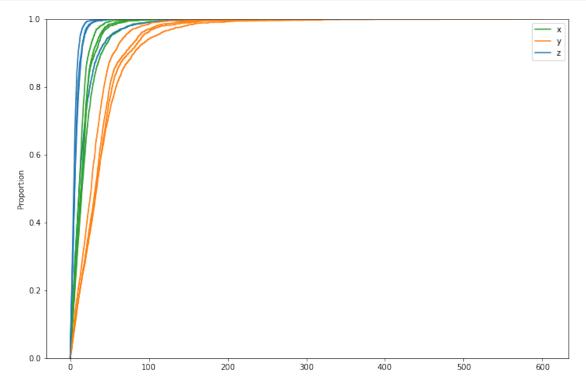
#### EDM comparision of actual and reconstructed topology



#### 5.5.3 Coordinate Error CDF

```
[25]: plt.figure(figsize=(12,8))
for i in range(len(outputs1)):
        sns.ecdfplot((abs(outputs1[i]['points'] - outputs1[i]['tpoints'])))
plt.legend(['x','y','z'])
```





Here we have taken India cities Latitudes and Longitudes and sampled them to obtain our dataset. The EDM after construction has been sampled at various rates and the following errors and properties have been noted.

```
[26]: print('Rate Ratio Error(%) Num Iterations Time Elapsed(s)')
for output in outputs1:
    print("%0.2f %0.2f %0.7f %0.2f %0.

4f"%(output['rate'],output['ratio'],output['error']*100,output['numit'],output['timeelapsed']

Rate Ratio Error(%) Num Iterations Time Elapsed(s)
0.01 0.99 1.0994404 19.00 29.2304
```

Rate	Ratio	ELLOL(%)	Num Iterations	lime Elapsed(s)
0.01	0.99	1.0994404	19.00	29.2304
0.02	1.98	0.7978092	16.00	32.7008
0.03	2.99	0.8518338	16.00	34.3210
0.05	5.02	0.6650077	19.00	57.8823

# 5.6 Error Analysis: Dataset 2

## 5.6.1 Defining EDMSolver Class Parameters

```
[27]: opts = params()
lsopts = params()
```

```
# Matrix Completion parameters
opts.lambda_ = 5
# opts.rate = 0.02
opts.printenergy = 0
opts.plotting = 0
opts.rank = 10
opts.maxit = 30
opts.tol = 1e-5
# BBGradient parameters
lsopts.maxit = 20
lsopts.xtol = 1e-8
lsopts.gtol = 1e-8
lsopts.ftol = 1e-10
lsopts.alpha = 1e-3
lsopts.rho = 1e-4
lsopts.sigma = 0.1
lsopts.eta = 0.8
```

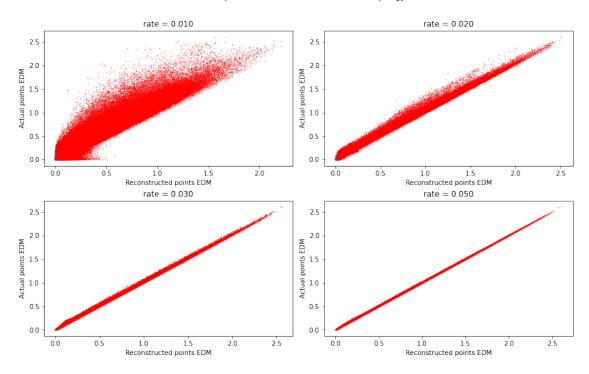
#### 5.6.2 Defining and Running the Solver

```
[28]: outputs2 = []
      rates2 = [0.01, 0.02, 0.03, 0.05]
      points = np.genfromtxt('./data/pipe.csv', delimiter=',')
      for rate in rates2:
          opts.rate = rate
          solver = EDMSolver(points,opts,lsopts)
          output = solver.run()
          outputs2.append(output)
     1.002312 percent of the EDM Distances are known
     det(R) < R, reflection detected!, correcting for it ...
     1.998138 percent of the EDM Distances are known
     det(R) < R, reflection detected!, correcting for it ...</pre>
     3.017235 percent of the EDM Distances are known
     5.010223 percent of the EDM Distances are known
[29]: fig,ax = plt.subplots(2,2,figsize=(12,8))
      for i in range(len(outputs2)):
          output = outputs2[i]
          dact = np.tril(solver.getEDM(output['points'])).ravel()
```

```
d0bs = np.tril(solver.getEDM(output['tpoints'])).ravel()
   ax[int(i/2),int(i%2)].scatter(d0bs,dact,color='r',s=0.1)
   ax[int(i/2),int(i%2)].set_title('rate = %.3f'%output['rate'])
   ax[int(i/2),int(i%2)].set_xlabel('Reconstructed points EDM')
   ax[int(i/2),int(i%2)].set_ylabel('Actual points EDM')

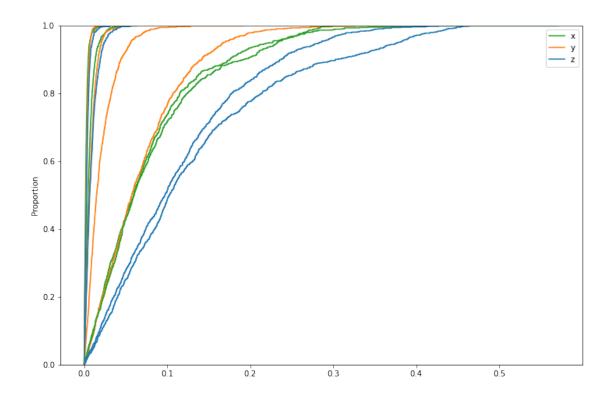
plt.tight_layout()
fig.subplots_adjust(top=0.88)
fig.suptitle('EDM comparision of actual and reconstructed topology')
plt.show()
```

EDM comparision of actual and reconstructed topology



#### 5.6.3 Coordinate Error CDF

```
[30]: plt.figure(figsize=(12,8))
   for output in outputs2:
        sns.ecdfplot((abs(output['points'] - output['tpoints'])))
   plt.legend(['x','y','z'])
   plt.show()
```



Rate	Ratio	Error(%)	Num Iterations	Time Elapsed(s)
0.01	1.00	21.3511732	30.00	7.6700
0.02	2.00	22.1612297	28.00	8.1524
0.03	3.02	2.0046797	19.00	6.7776
0.05	5.01	0.7753925	15.00	6.2720

# 5.7 Error Analysis: Dataset 3

# 5.7.1 Defining EDMSolver Class Parameters

```
[32]: opts = params()
    lsopts = params()

    opts.lambda_ = 5
    # opts.rate = 0.4
    opts.printenergy = 0
    opts.plotting = 0
    opts.rank = 10
    opts.maxit = 30
```

```
opts.tol = 1e-5

lsopts.maxit = 20
lsopts.xtol = 1e-8
lsopts.gtol = 1e-8
lsopts.ftol = 1e-10
lsopts.alpha = 1e-3
lsopts.rho = 1e-4
lsopts.sigma = 0.1
lsopts.eta = 0.8
```

#### 5.7.2 Defining and Running the Solver

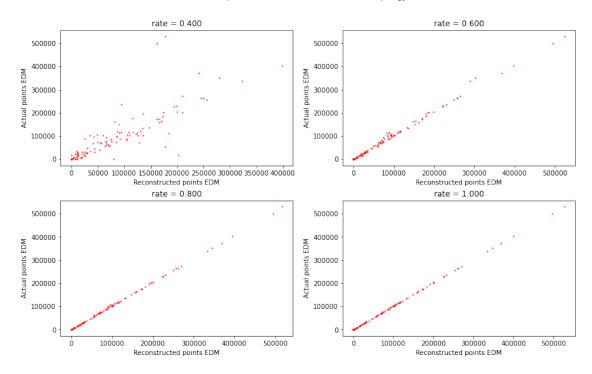
```
[33]: outputs3 = []
  outputdata = []
  rates3 = [0.4,0.6,0.8,1]

for rate in rates3:
    opts.rate = rate
    solver = EDMSolver(points3,opts,lsopts)
    output = solver.run()
    outputs3.append(output)
```

41.777778 percent of the EDM Distances are known 58.666667 percent of the EDM Distances are known 74.666667 percent of the EDM Distances are known 93.333333 percent of the EDM Distances are known

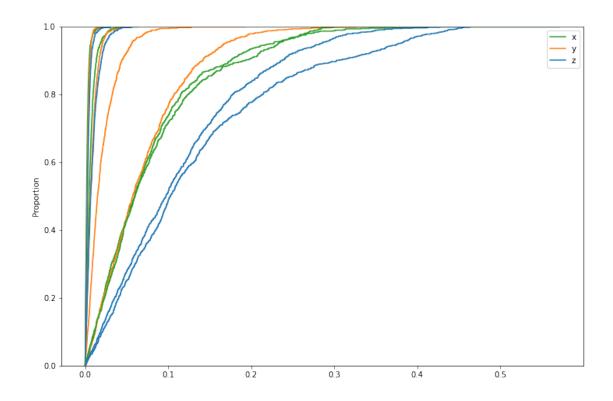
```
fig,ax = plt.subplots(2,2,figsize=(12,8))
for i in range(len(outputs3)):
    output = outputs3[i]
    dact = np.tril(solver.getEDM(output['points'])).ravel()
    d0bs = np.tril(solver.getEDM(output['tpoints'])).ravel()
    ax[int(i/2),int(i%2)].scatter(d0bs,dact,color='r',s=1)
    ax[int(i/2),int(i%2)].set_title('rate = %.3f'%output['rate'])
    ax[int(i/2),int(i%2)].set_xlabel('Reconstructed points EDM')
    ax[int(i/2),int(i%2)].set_ylabel('Actual points EDM')

plt.tight_layout()
fig.subplots_adjust(top=0.88)
fig.suptitle('EDM comparision of actual and reconstructed topology')
plt.show()
```



#### 5.7.3 Coordinate Error CDF

```
[35]: plt.figure(figsize=(12,8))
    for output in outputs2:
        sns.ecdfplot((abs(output['points'] - output['tpoints'])))
    plt.legend(['x','y','z'])
    plt.show()
```



```
[36]: print('Rate
                    Ratio
                             Error(%)
                                        Num Iterations Time Elapsed(s)')
      for output in outputs3:
          print("%0.2f %0.2f
                                 %0.7f
                                             %0.2f
                                                           %0.
       →4f"%(output['rate'],output['ratio'],output['error']*100,output['numit'],output['timeelapsed']
                                Num Iterations Time Elapsed(s)
     Rate
            Ratio
                     Error(%)
     0.40
            41.78
                    2.2919492
                                    5.00
                                                 0.0140
                    0.3994958
                                    11.00
                                                  0.0299
     0.60
            58.67
```

0.0219

0.0150

# 5.8 Error Analysis: Dataset 4

74.67

93.33

0.80

1.00

# 5.8.1 Defining EDMSolver Class Parameters

0.1712731

0.0687895

```
[37]: opts = params()
    lsopts = params()

    opts.lambda_ = 5
    # opts.rate = 0.4
    opts.printenergy = 0
    opts.plotting = 0
    opts.rank = 10
    opts.maxit = 30
```

8.00

5.00

```
opts.tol = 1e-5

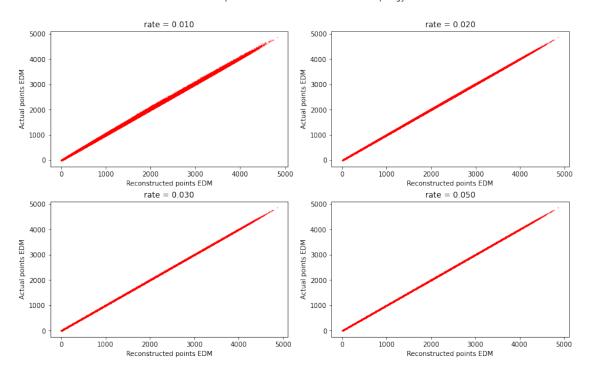
lsopts.maxit = 20
lsopts.xtol = 1e-8
lsopts.gtol = 1e-8
lsopts.ftol = 1e-10
lsopts.alpha = 1e-3
lsopts.rho = 1e-4
lsopts.sigma = 0.1
lsopts.eta = 0.8
```

#### 5.8.2 Defining and Running the Solver

```
[38]: outputs4 = []
      rates4 = [0.01, 0.02, 0.03, 0.05]
      for rate in rates4:
          opts.rate = rate
          solver = EDMSolver(points4,opts,lsopts)
          output = solver.run()
          outputs4.append(output)
     0.998259 percent of the EDM Distances are known
     det(R) < R, reflection detected!, correcting for it ...</pre>
     2.001472 percent of the EDM Distances are known
     det(R) < R, reflection detected!, correcting for it ...</pre>
     3.002135 percent of the EDM Distances are known
     4.999413 percent of the EDM Distances are known
     det(R) < R, reflection detected!, correcting for it ...</pre>
[39]: fig,ax = plt.subplots(2,2,figsize=(12,8))
      fig.suptitle('EDM comparision of actual and reconstructed topology')
      for i in range(len(outputs4)):
          output = outputs4[i]
          dact = np.tril(solver.getEDM(output['points'])).ravel()
          dObs = np.tril(solver.getEDM(output['tpoints'])).ravel()
          ax[int(i/2),int(i%2)].scatter(d0bs,dact,color='r',s=0.1)
          ax[int(i/2),int(i%2)].set_title('rate = %.3f'%output['rate'])
          ax[int(i/2),int(i\(\frac{1}{2}\)].set_xlabel('Reconstructed points EDM')
          ax[int(i/2),int(i%2)].set_ylabel('Actual points EDM')
      plt.tight_layout()
      fig.subplots_adjust(top=0.88)
```

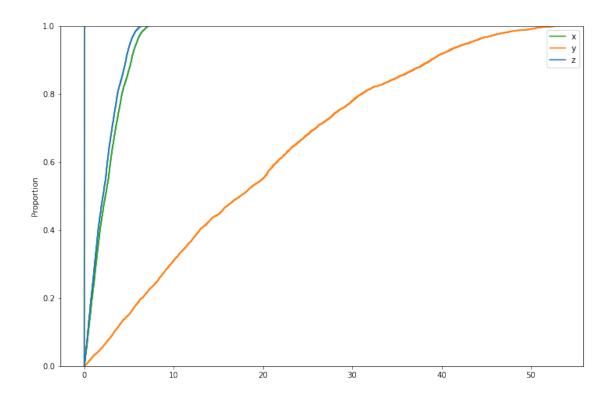
# plt.show()

EDM comparision of actual and reconstructed topology



## 5.8.3 Coordinate Error CDF

```
[40]: plt.figure(figsize=(12,8))
    for output in outputs4:
        sns.ecdfplot((abs(output['points'] - output['tpoints'])))
    plt.legend(['x','y','z'])
    plt.show()
```



```
[41]: print('Rate
                             Error(%)
                                        Num Iterations Time Elapsed(s)')
                    Ratio
      for output in outputs4:
          print("%0.2f %0.2f
                                 %0.7f
                                             %0.2f
                                                           %0.
       →4f"%(output['rate'],output['ratio'],output['error']*100,output['numit'],output['timeelapsed']
                     Error(%)
                                Num Iterations Time Elapsed(s)
     Rate
            Ratio
     0.01
            1.00
                   47.7290209
                                                   66.8547
                                     14.00
```

69.5083

103.2109

179.7898

# 6 observations

2.00

3.00

5.00

47.8567700

0.0577544

47.8917195

0.02

0.03

0.05

We can see that the errors plotted for the 4 datasets in the above run is tabularised as follows

12.00

22.00

15.00

Table 1: Dataset 1

Rate	Ratio	Error(%)	Num Iterations	Time Elapsed(s)
0.01	0.99	1.0994404	19.00	29.2304
0.02	1.98	0.7978092	16.00	32.7008
0.03	2.99	0.8518338	16.00	34.3210
0.05	5.02	0.6650077	19.00	57.8823

It is evident that as the sampling rate, the number of entries from the EDM matrix increased , the performance of our algorithm has increased considerably. However there is a improvement in error as we move from 0.03 to 0.05 sampling rate. This could be due to randomised algorithm which we are using to label some our EDM matrix entries as missing. It could be possible that the data available was not representative enough of the spatial orientation. This introduces another problem of identifying the crucial edges of our topology which play an important role in our algorithm.

Table 2 : Dataset 2

Rate	Ratio	Error(%)	Num Iterations	Time Elapsed(s)
0.01	1.00	21.3511732	30.00	7.6700
0.02	2.00	22.1612297	28.00	8.1524
0.03	3.02	2.0046797	19.00	6.7776
0.05	5.01	0.7753925	15.00	6.2720

We see a perfect decrease in error as the sample rating increases, almost near ideal accuracies as we approach a sampling rate of 5%. This is in line with our expected observations.

Table 3 : Dataset 3

Rate	Ratio	Error(%)	Num Iterations	Time Elapsed(s)
0.40	41.78	2.2919492	5.00	0.0140
0.60	58.67	0.3994958	11.00	0.0299
0.80	74.67	0.1712731	8.00	0.0219
1.00	93.33	0.0687895	5.00	0.0150

Since this is a very small dataset in comparison to our previous ones, we had to use 10-50 times the sampling rate in the previous datasets. This is because as the number of points decreases, the total number of connections also decrease and thus our solver has much lesser data to work with. Reasonable accuracies were obtained with about 40% of the data which is acceptable in most real life applications.

Table 4 : Dataset 4

Rate	Ratio	Error(%)	Num Iterations	Time Elapsed(s)
0.01	1.00	47.7290209	14.00	66.8547
0.02	2.00	47.8567700	12.00	69.5083
0.03	3.00	0.0577544	15.00	103.2109
0.05	5.00	47.8917195	22.00	179.7898

We found this dataset interesting because it is one of the state-of-the-art applications of our algorithm. From NMR spectral data we derived the molecular conformation of "HAEMOGLOBIN IN THE DEOXY QUATERNARY STATE WITH LIGAND BOUND AT THE ALPHA HAEMS". However, we can see that the error is not that fancy, this could be because the intra molecular distance is very very small ( $\sim 10^{-10}$ ) and any small deviations due to our calculations of the euclidean

distance will results in huge margin for our relative error that is being calculated in this case.

## 7 Conclusions

Our error that is being calculated is not a precise measure of our algorithm because of a few inherent flaws in it.

One of the major issues is that it is always possible that some set of distance matrix entries are more important and representative our of data than another one. Hence we are leaving the performance of our algorithm at random when we start from a point cloud (All these datasets). However, in most real life applications we start with an EDM matrix and hence this should not be a problem for our algorithm.

Another important issue we are dealing with is the normalising of EDM matrix and subsequent scaling of the output point matrix obtained. These rounding off may seem small to the naked eye but we can see that we are using the kilometers as a basic unit for our algorithm (For datasets 1 and 3). A small variation in this about  $\sim 10^{-2}$  could results in deviations of about 10 metres.

# 8 Future extensions for our project

Identify the crucial edges for our topology by generating the performance data of our algorithm with different sets of entries in our EDM matrix and its parameters and training an ML model using this data to assign an importance factor to each of these EDM matrix entries

Dynamic reconstruction of the topology of moving nodes using IMU data obtained.

#### 9 References

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Rigid Transform 3D - https://github.com/nghiaho12/rigid\_transform\_3D

Dataset 1 - https://www.kaggle.com/max-mind/world-cities-database

Dataset 2 - https://www.kaggle.com/aalborguniversity/sewerpointclouds

Dataset 4 - https://www.rcsb.org/structure/1COH