

# Looking at Non-Linear Dimension Reduction as Models in the Data Space

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## Abstract

Nonlinear dimension reduction (NLDR) techniques such as tSNE, and UMAP provide a low-dimensional representation of high-dimensional (high-D) data using non-linear transformation. The methods and parameter choices can create wildly different representations, making it difficult to decide which is best, or whether any or all are accurate or misleading. NLDR often exaggerates random patterns, sometimes due to the samples observed. But NLDR views have an important role in data analysis because, if done well, they provide a concise visual (and conceptual) summary of high-D distributions. To help evaluate the NLDR we have developed an algorithm to show the 2D NLDR model in the high-D space, viewed with a tour. One can see if the model fits everywhere or better in some subspaces, or completely mismatches the data. It is used to evaluate which 2D layout is the best representation of the high-D distribution and see how different methods may have similar summaries or quirks.

*Keywords:* high-dimensional data, dimension reduction, hexagon binning, low-dimensional manifold, tour, data visualization, model in the data space

# 1 Introduction

Non-linear dimension reduction (NLDR) is popular for making a convenient low-dimensional ( $k$ -D) representation of high-dimensional ( $p$ -D) data. Recently developed methods include t-distributed stochastic neighbor embedding (tSNE) (van der Maaten & Hinton 2008), uniform manifold approximation and projection (UMAP) (McInnes & Healy 2018), potential of heat-diffusion for affinity-based trajectory embedding (PHATE) algorithm (Moon et al. 2019), large-scale dimensionality reduction Using triplets (TriMAP) (Amid & Warmuth 2022), and pairwise controlled manifold approximation (PaCMAP) (Wang et al. 2021). However, the representation generated can vary dramatically from method to method, and with different choices of parameters or random seeds made using the same method (Figure 1). The dilemma for the analyst is then, **which representation to use**. The choice might result in different procedures used in the downstream analysis, or different inferential conclusions. The research described here provides new visual tools to aid with this decision.

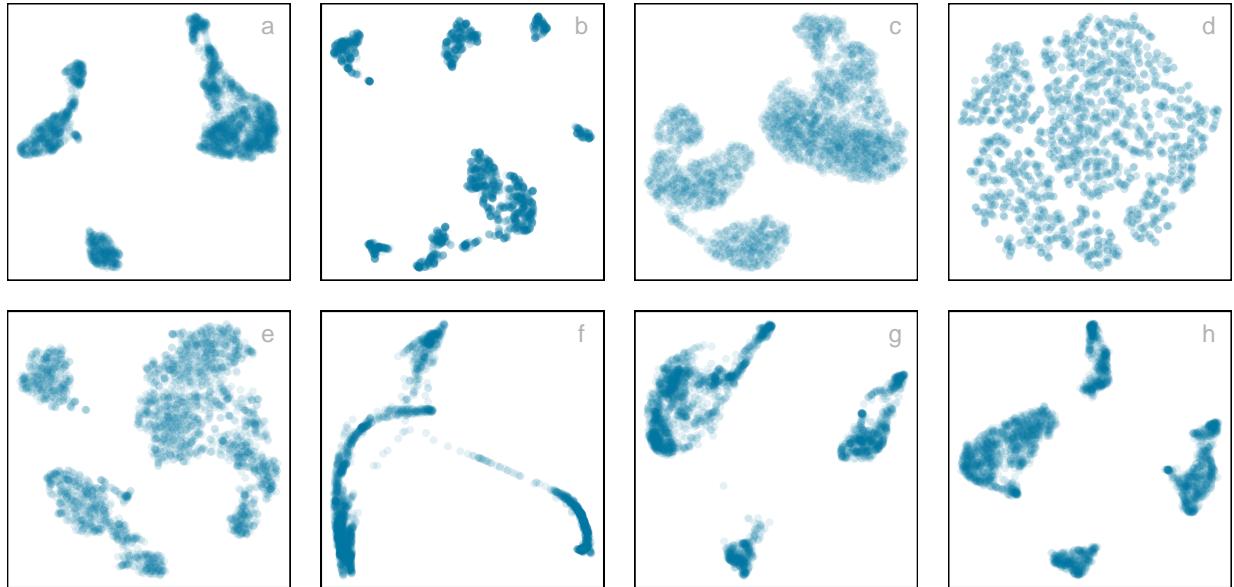


Figure 1: Six different NLDR representations of the same data. Different techniques and different parameter choices are used. Researchers may have seen any of these in their analysis of this data, depending on their choice of method, or typical parameter choice. Would they make different decisions downstream in the analysis depending on which version seen? Which is the most accurate representation of the structure in high dimensions?

The paper is organised as follows. Section 2 provides a summary of the literature on NLDR, and high-dimensional data visualization methods. Section 3 contains the details of the new methodology, including simulated data examples. Two applications illustrating the use of the new methodology for bioinformatics and image classification are in Section 6. Limitations and future directions are provided in Section 7.

## 2 Background

Historically,  $k$ - $D$  representations of  $p$ - $D$  data have been computed using multidimensional scaling (MDS) (Borg & Groenen 2005), which includes principal components analysis (PCA) (Jolliffe 2011) as a special case. The  $k$ - $D$  representation can be considered to be a layout of points in  $k$ - $D$  produced by an embedding procedure that maps the data from  $p$ - $D$ . In MDS, the  $k$ - $D$  layout is constructed by minimizing a stress function that differences distances between points in  $p$ - $D$  with potential distances between points in  $k$ - $D$ . Various formulations of the stress function result in non-metric scaling (Saeed et al. 2018) and isomap (Silva & Tenenbaum 2002). Challenges in working with high-dimensional data, including visualization, are outlined in Johnstone & Titterington (2009).

Many new methods for NLDR have emerged in recent years, all designed to better capture specific structures potentially existing in  $p$ - $D$ . Here we focus on five currently popular techniques, tSNE, UMAP, PHATE, TriMAP and PaCMAP. tNSE and UMAP can be considered to produce the  $k$ - $D$  minimizing the divergence between two distributions, where the distributions are modeling the inter-point distances. PHATE, TriMAP and PaCMAP are examples of diffusion processes (Coifman et al. 2005) spreading to capture geometric shapes, that include both global and local structure.

The array of layouts in Figure 1 illustrate what can emerge from the choices of method and parameters, and the random seed that initiates the computation. Key structures interpreted from these views suggest: (1) highly **separated clusters** (a, b, e, g, h) with the number ranging from 3-6; (2) **stringy branches** (f), and (3) **barely separated clusters** (c, d) which would **contradict** the other representations.

It happens because these methods and parameter choices provide different lenses on the interpoint distances in the data.

The alternative approach to visualizing the high-dimensional data is to use linear projections. PCA is the classical approach, resulting in a set of new variables which are linear combinations of the original variables. Tours, defined by Lee et al. (2021), broaden the scope by providing movies of linear projections, that provide views the data from all directions. Lee et al. (2021) provides an review of the main developments in tours. There are many tour algorithms implemented, with many available in the R package `tourr` (Wickham et al. 2011), and versions enabling better interactivity in `langevitour` (Harrison 2023) and `detourr` (Hart & Wang 2022). Linear projections are a safe way to view high-dimensional data, because they do not warp the space, so they are more faithful representations of the structure. However, linear projections can be cluttered, and global patterns can obscure local structure. The simple activity of projecting data from  $p$ - $D$  suffers from piling (Laa et al. 2022), where data concentrates in the center of projections. NLDR is designed to escape these issues, to exaggerate structure so that it can be observed. But as a result NLDR can hallucinate wildly, to suggest patterns that are not actually present in the data.

The solution is to use the tour to examine how the NLDR is warping the space. This approach follows what Wickham et al. (2015) describes as *model-in-the-data-space*. The fitted model should be overlaid on the data, to examine the fit relative the spread of the observations. While this is straightforward, and commonly done when data is 2D, it is also possible in  $p$ - $D$ , for many models, when a tour is used.

[Wickham et al. \(2015\)](#) provides several examples of models overlaid on the data in  $p$ - $D$ . In hierarchical clustering, a representation of the dendrogram using points and lines can be constructed by augmenting the data with points marking merging of clusters. Showing the movie of linear projections reveals shows how the algorithm sequentially fitted the cluster model to the data. For linear discriminant analysis or model-based clustering the model can be indicated by  $(p - 1)$ - $D$  ellipses. It is possible to see whether the elliptical shapes appropriately matches the variance of the relevant clusters, and to compare and contrast different fits. For PCA, one can display the  $k$ - $D$  plane of the reduced dimension using wireframes of transformed cubes. Using a wireframe is the approach we take here, to represent the NLDR model in  $p$ - $D$ .

## 3 Method

### 3.1 What is the NLDR model?

At first glance, thinking of NLDR as a modeling technique might seem strange. It is a simplified representation or abstraction of a system, process, or phenomenon in the real world. The  $p$ - $D$  observations are the realization of the phenomenon, and the  $k$ - $D$  NLDR layout is the simplified representation. From a statistical perspective we can consider the distances between points in the  $k$ - $D$  layout to be variance that the model explains, and the (relative) difference with their distances in  $p$ - $D$  is the error, or unexplained variance. We can also imagine that the positioning of points in 2- $D$  represent the fitted values, that will have some prescribed position in  $p$ - $D$  that can be compared with their observed values. This is the conceptual framework underlying the more formal versions of factor analysis ([Jöreskog 1969](#)) and multidimensional scaling (MDS) ([Borg & Groenen 2005](#)). (Note that, for this thinking the full  $p$ - $D$  data needs to be available, not just the interpoint distances.)

We define the NLDR as a function  $g: \mathbb{R}^{n \times p} \rightarrow \mathbb{R}^{n \times k}$ , with (hyper-)parameters  $\boldsymbol{\theta}$ . The parameters,  $\boldsymbol{\theta}$ , depend on the choice of  $g$ , and can be considered part of model fitting in the traditional sense. Common choices for  $g$  include functions used in tSNE, UMAP, PHATE, TriMAP, PaCMAP, or MDS, although in theory any function that does this mapping is suitable.

With our goal being to make a representation of this 2- $D$  layout that can be lifted into high-dimensional space, the layout needs to be augmented to include neighbour information. A simple approach would be to triangulate the points and add edges. A more stable approach is to first hexagonally bin the data, reducing it from  $n$  to  $m \leq n$  observations, and connect the bin centroids. This process serves to reduce some noisiness in the resulting surface shown in  $p$ - $D$ . The steps in this process are shown in Figure 2, and documented below.

To illustrate the method, we use 7- $D$  simulated data, which we call the “S-curve”. It is constructed by simulating  $n = 750$  observations from  $\theta \sim U(-3\pi/2, 3\pi/2)$ ,  $X_1 = \sin(\theta)$ ,  $X_2 \sim U(0, 2)$  (adding thickness to the S),  $X_3 = \text{sign}(\theta) \times (\cos(\theta) - 1)$ . The remaining variables  $X_4, X_5, X_6, X_7$  are all uniform error, with small variance. We would consider  $T = (X_1, X_2, X_3)$  to be the geometric structure (true model) that we hope to capture.

Notation	Description
$n, p, k$	number of observations, variables, embedding dimension, respectively
$\mathbf{X}, \mathbf{x}$	$p$ -dimensional data (population, sample)
$\mathbf{y}$	$k$ -dimensional layout
$P$	orthonormal basis, generating a $d$ -dimensional linear projection of $p$ -dimensional data
$T$	true model
$g$	functional mapping from $p$ -D to $k$ -D, especially as prescribed by NLDR
$\theta$	(Hyper-) parameters for NLDR method
$r$	ranges of the embedding components
$C^{(j)}$	$j$ -dimensional bin centers
$(b_1, b_2)$	number of bins in each direction
$(a_1, a_2)$	binwidths, distance between centroids in each direction
$(s_1, s_2)$	starting coordinates of the hexagonal grid
$q$	buffer to ensure hexgrid covers data, proportion of data range, 0-1
$m$	number of non-empty bins
$b$	number of hexagons in the grid
$h$	hexagonal id

Table 1: Summary of notation for describing new methodology.

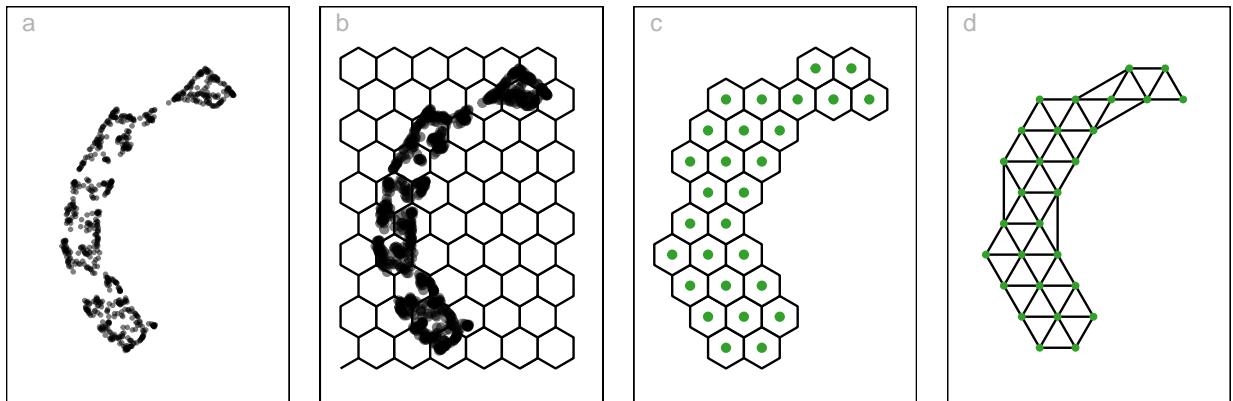


Figure 2: Key steps for constructing the model on the UMAP layout ( $k = 2$ ): (a) data, (b) hexagon bins, (c) bin centroids, and (d) triangulated centroids. The S-curve data is shown.

### 3.1.1 Scaling the data

Because we are working with distances between points, starting with data having a standard scale, e.g.  $[0, 1]$ , is recommended. When the aspect ratio of the NLDR  $(r_1, r_2, \dots, r_k)$  is meaningful, the data can be scaled to take this into account. When  $k = 2$ , as in hexagon binning, the default range is  $[0, y_{i,\max}]$ ,  $i = 1, 2$ , where  $y_{1,\max} = 1$  and  $y_{2,\max} = \frac{r_2}{r_1}$  (Figure 2).

### 3.1.2 Computing hexagon grid configuration

The 2-D hexagon grid is defined by the number of bins in each direction  $(b_1, b_2)$ , giving total number of bins as  $b = b_1 \times b_2$ , and a unique hexagon id,  $h = 1, \dots, b$ . Each hexagon,  $H_h$  is uniquely described by centroid,  $C_h^{(2)} = (c_{h1}, c_{h2})$ . (Note that, we detail the algorithm for constructing the grid and binning the data because we found that there was no readily available implementation that could produce what was needed for this project.)

To make the grid, a buffer parameter ( $q$ ) is set as a proportion of the plot space. By default this is 0.1. A value should be chosen to provide at buffer of a full hexagon width ( $a_1$ ) and height ( $a_2$ ) for the grid to extend beyond the data. The lower left position where the grid starts is defined as  $(s_1, s_2)$ , corresponding to the centroid of the lowest left hexagon, which should be below the minimum data value. This is computed as one buffer unit,  $q$  below the minimum data value, by default. The user provides the preferred number of bins in the horizontal direction,  $b_1$ , and the algorithm computes the appropriate number of bins in the vertical direction,  $b_2$ .

The value for  $b_2$  is computed by fixing  $b_1$ . Considering the lower bound of the NLDR,  $a_1 > -2q$ , and  $a_1 > \frac{1+q}{b_1-1}$ . Similarly, according to the upper bound of the NLDR,  $a_1 > \frac{2r_2(1+q)}{\sqrt{3}(b_2-1)}$ , because  $a_2 = \frac{\sqrt{3}}{2}a_1$  for regular hexagons. Therefore,  $b_2 = \left\lceil 1 + \frac{2r_2(b_1-1)}{\sqrt{3}} \right\rceil$ .

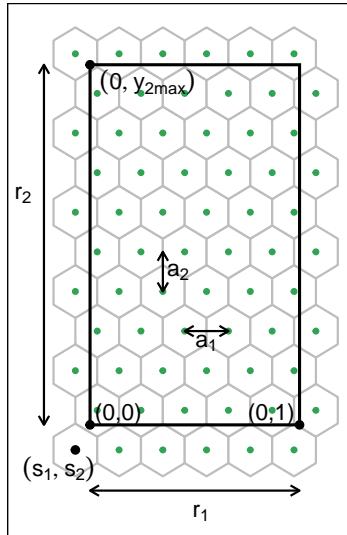


Figure 3: The components of the hexagon grid illustrating notation.

### 3.1.3 Binning the data

Observations are grouped into bins based on their nearest centroid. This produces a reduction in size of the data from  $n$  to  $m$ , where  $m \leq b$  (total number of bins). This can be defined using the function  $u : \mathbb{R}^{n \times 2} \rightarrow \mathbb{R}^{m \times 2}$ , where  $u(i) = \arg \min_{j=1, \dots, b} \sqrt{(y_{i1} - C_{j1}^{(2)})^2 + (y_{i2} - C_{j2}^{(2)})^2}$ , mapping observation  $i$  into  $H_h = \{i | u(i) = h\}$ .

By default, the bin centroid is used for describing a hexagon (as done in Figure 2 (c)), but any measure of center, such as a mean or weighted mean of the points within each hexagon, could be used. The bin centers, and the binned data, are the two important components needed to render the model representation in high dimensions.

### 3.1.4 Indicating neighborhood

Delaunay triangulation (Lee & Schachter 1980) is used to connect neighboring centroids, which is needed to preserve neighborhood information when the model is lifted into  $p$ -D.

It can also happen that distant centroids can be connected, which can result in long line segments. In order to generate a smooth surface in 2-D, these long line segments must be removed (Figure 2 (d)).

## 3.2 Rendering the model in $p$ -D

The last step is to lift the  $k$ -D model into  $p$ -D by computing  $p$ -D vectors that represent bin centroids. We use the  $p$ -D mean of the points in  $H_h$  to map the centroid  $C_h^{(2)} = (c_{h1}, c_{h2})$  to a point in  $p$ -D. Let the  $p$ -D mean be

$$C_h^{(p)} = \frac{1}{n_h} \sum_{i=1}^{n_h} x_i, h = 1, \dots, b; n_h > 0.$$

Furthermore, line segments that exist in the  $k$ -D model generate line segments in  $p$ -D by connecting the  $p$ -D means of the corresponding  $k$ -D bin centroids.

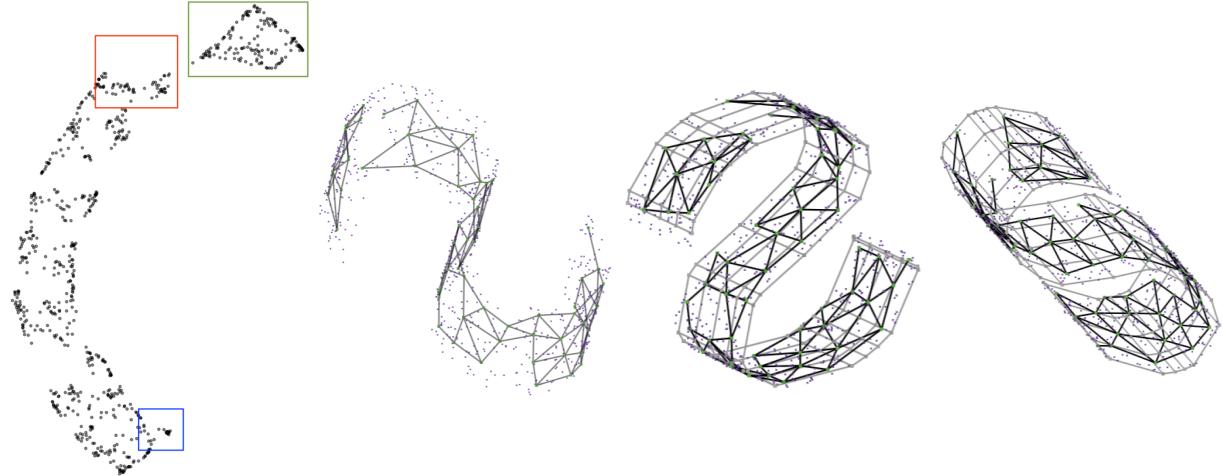


Figure 4: Projections of the 7-D S-curve data (purple points) overlaid with the model (black grid with green points indicating bin means) constructed from the UMAP layout, with 60 non-empty bins (using the **langevitour** software). The model closely fits the shape of this data. A video of the full tour is available at (<https://youtu.be/G1m4q9k--v4>).

### 3.3 Measuring the fit

#### 3.3.1 Fitted values

The fitted values of the points in  $H_h$  refers to the  $p$ - $D$  mapping  $C_h^{(p)}$  of the corresponding  $k$ - $D$  model point  $C_h^{(2)}$ .

#### 3.3.2 Residuals and error

To assess how well our method captures and represents the underlying structure of the  $p$ - $D$  data, residuals are important. Residuals are computed by taking the squared  $p$ - $D$  Euclidean distance,  $\sum_{h=1}^b \sum_{i=1}^{n_h} \sum_{j=1}^P (x_{hij} - C_{hj}^{(p)})^2$ . Additionally, Mean Squared Error (MSE),  $\frac{1}{n} \sum_{h=1}^b \sum_{i=1}^{n_h} \sum_{j=1}^P (x_{hij} - C_{hj}^{(p)})^2$ , is computed based on residuals.

### 3.4 Prediction

The prediction approach involves finding the nearest  $k$ - $D$  model point for a new  $p$ - $D$  point. We define the function  $z : \mathbb{R}^{n \times p} \rightarrow \mathbb{R}^{m \times p}$ , where  $z(i) = \arg \min_{j=1, \dots, b} \sqrt{\sum_{v=1}^P (x_{iv} - C_{jv}^{(p)})^2}$  maps each  $p$ - $D$  point to its nearest  $p$ - $D$  mapping of the model. Therefore, the new observation  $i$  falls in the hexagon,  $H_h = \{i | z(i) = h\}$  and the corresponding  $k$ - $D$  bin centroids,  $C_h^{(2)} = (c_{h1}, c_{h2})$  be the predicted values.

### 3.5 Tuning

The performance and robustness of our model depend on three key parameters: (i) the total number of bins ( $b$ ), (ii) a benchmark value used to remove low-density hexagons, and (iii) a benchmark value used to remove long edges. However, there is no analytical formula to calculate an appropriate value for these parameters. The selection of these parameter values depends on the model performance computed by MSE (see Section 3.3).

#### 3.5.1 Choice of bins

The number of hexagon bins in the hexagon grid has a considerable impact on the construction of the 2-D model, serving as the initial step. The chosen total number of bins must effectively capture the structure of the NLDR data. If the number of bins is too low, the model may not be able to capture the structure of the NLDR data effectively (see Figure 5 (a)), while if there are too many bins, it may result in over-fitting the individual points of the NLDR data (see Figure 5 (c)). Therefore, it is important to determine an effective number of bins to construct a reasonable model (see Figure 5 (b)).

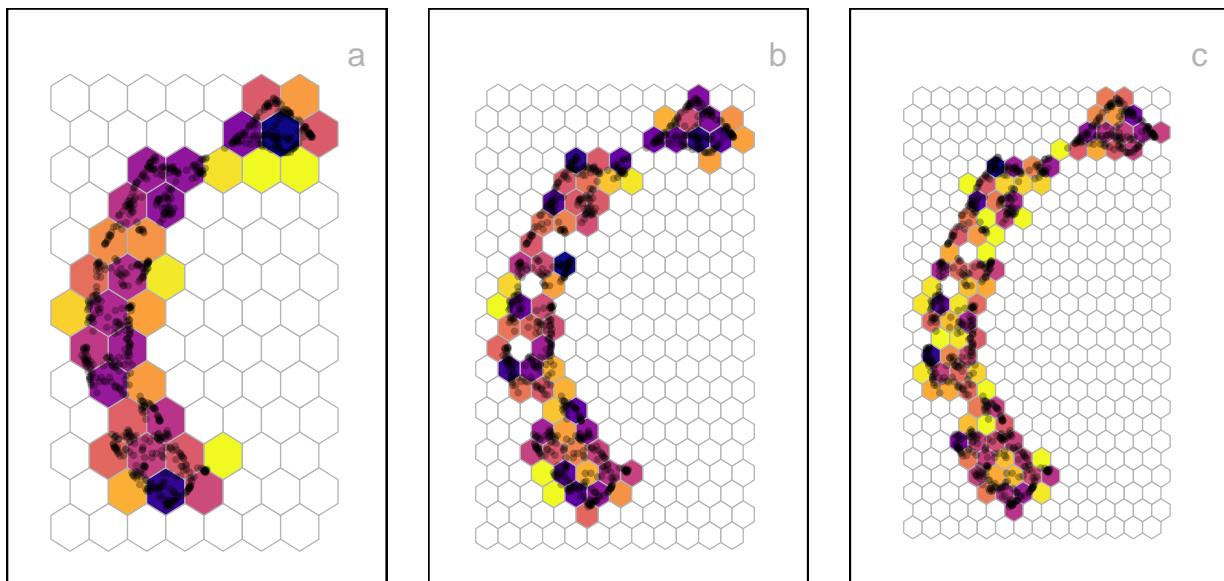


Figure 5: Hexbin density plots of UMAP layout of the S-curve data, using three different bin inputs: (a)  $b = 91$  (7, 13), (b)  $b = 190$  (10, 19), and (c)  $b = 364$  (14, 26). Color indicates standardized counts, dark indicating high count and light indicates low count. At the smallest bin size the data segregates into two separate groups, suggesting this is too many bins. Using the MSE of the model fit in  $p - D$  helps decide on a useful choice of number of bins.

To determine the effective  $b$ , candidate values are selected based on the range between the minimum and approximate maximum  $b_1$ , because  $b_2$  is computed from  $b_1$ . The minimum  $b_1$  is set to 2, while the maximum number is estimated by taking the square root of  $\frac{n}{2}$ . The analysis evaluates the MSE across varying  $b$  within this range, covering the minimum to maximum values along both axes (see [?@fig-mse-scurve-b](#)).

### 3.5.2 Removal of low density bins

Once setting up the hexagon grid with an appropriate number of bins, some hexagon bins may have few or no data points within them (see Figure 5 (b)). To ensure comprehensive coverage of the NLDR data, it is necessary to select hexagon bins with a considerable number of data points. This involves calculating the number of points within each hexagon. Then, the standard count is computed by dividing the number of points within each hexagon by the maximum number of points in the grid. Next, bins with a standard count less than a benchmark value are removed (see Figure 6 (a)). There is no specific rule for selecting a benchmark value. However, the following steps can help determine a suitable value for removing low-density hexagons:

1. Plot the distribution of the standardized counts (see `?@fig-stdcts-scurve`).
2. Examine the distribution of counts.
3. Select the first quantile value if the distribution is skewed.

The benchmark value for removing low-density hexagons ranges between 0 and 1. When analyzing how these benchmark values influence model performance, it's essential to observe the change in MSE as the benchmark value increases (see `?@fig-mse-scurve-lwd`). The MSE shows a gradual increase as the benchmark value progresses from 0 to 1. Evaluating this rate of increase is important. If the increment is not considerable, the decision might lean towards retaining low-density hexagons.

Furthermore, selecting the benchmark value for removing low-density hexagons is important. Removing unnecessary bins may lead to the formation of long edges and an uneven 2-D model. Hence, rather than solely relying on the benchmark value to identify hexagons for removal, it's essential to consider the standard number of points in the neighboring hexagons of the identified low-density bins (see Figure 6 (b)). If neighboring bins also show low counts, only those bins will be removed. The remaining bins are used to construct the 2-D model.

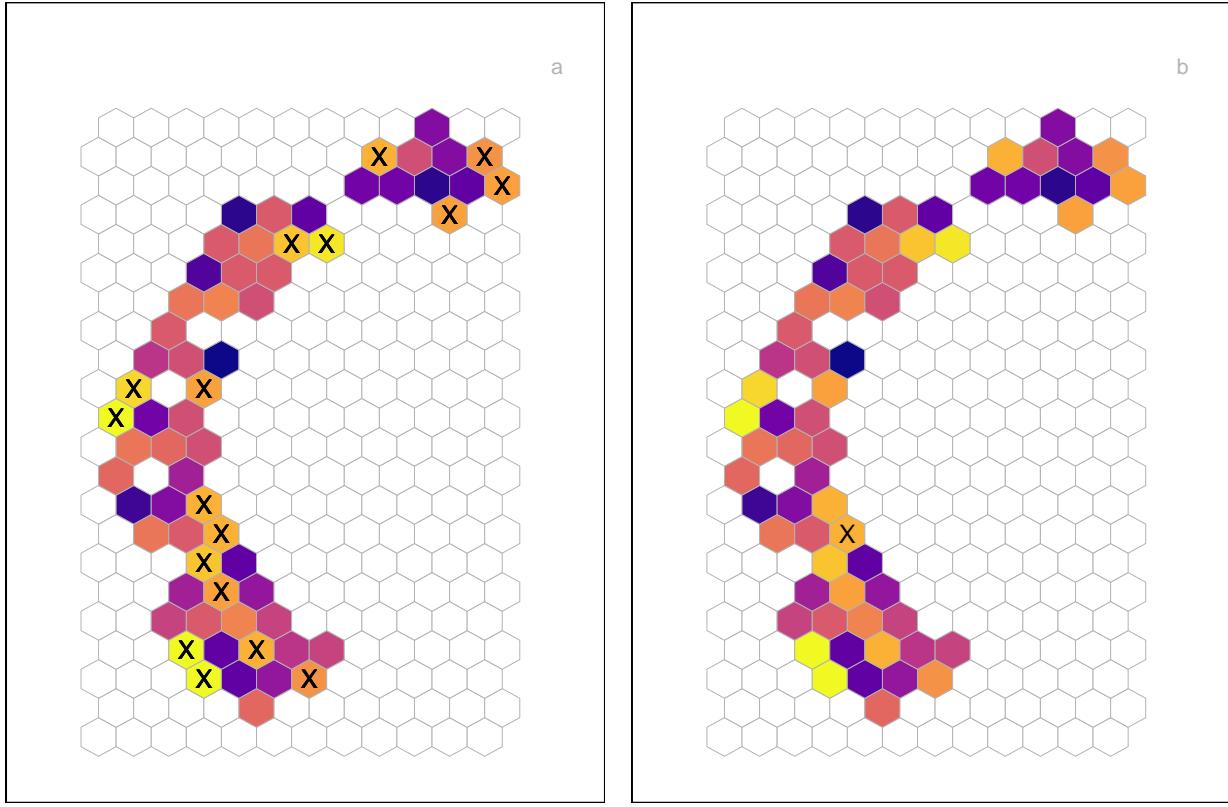


Figure 6: Two approaches to identify low-density hexagons. Hexagon grid (a) marks hexagons with a standardized count less than 0.25. Hexagon grid (b) marks hexagons by considering the densities of neighboring bins. Removing hexagons by also investigating the densities of neighboring bins is more reliable for preserving the structure.

### 3.5.3 Removing long edges

To create a smooth 2-D representation (see Figure 2 (d)), it is necessary to remove edges that connect distant bin centroids in the triangular mesh. These edges only exist in the 2-D model and do not extend into  $p$ -D, so their removal does not impact the model in  $p$ -D. Although there are no specific criteria for determining the benchmark value to remove long edges, the following steps provide an approach to identifying a suitable threshold:

1. Plot the distribution of the 2D Euclidean distances (see [?@fig-dist-scurve](#)).
2. Identify the first largest difference between consecutive distance values.
3. Take the distance value corresponding to this difference as the benchmark value.

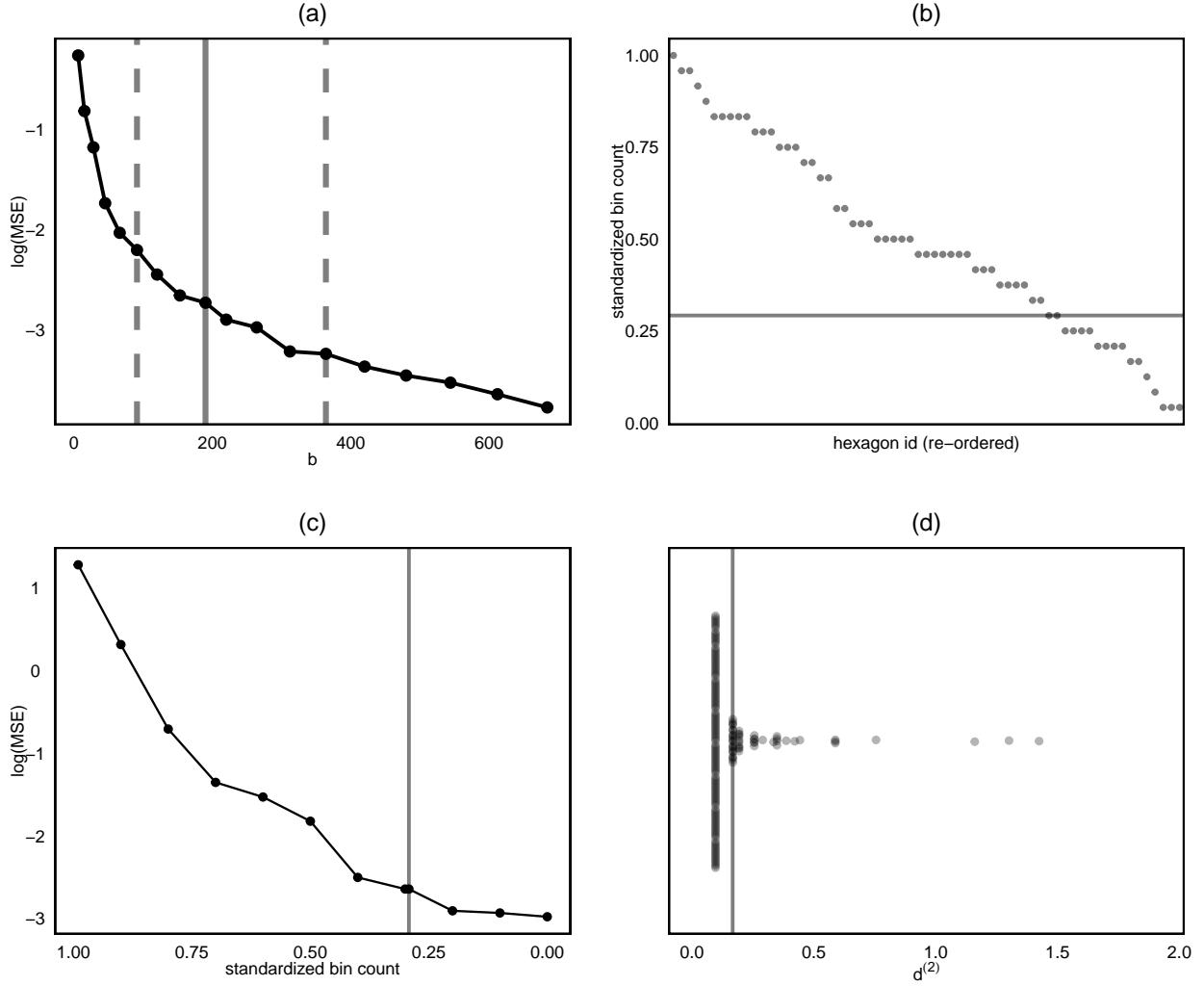


Figure 7: Various plots to help assess best number of bins, low density bin and large edge removal. Both (a) and (c) show MSE, against number of bins and standardised count. A good benchmark value for these parameters is when the MSE drops and then flattens out. Plot (b) shows the distribution of standardised counts of hexagons. Plot (c) shows the distribution of 2-D Euclidean distances between bin centroids, with a good benchmark value for removing large edges would being the distance that shows the first large increase.

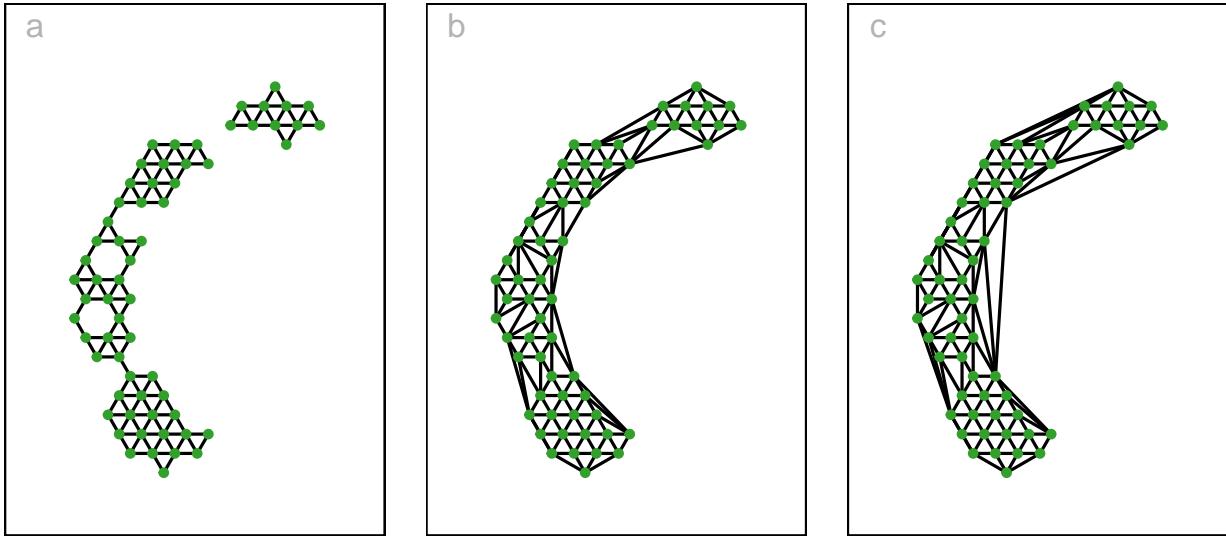


Figure 8: 2-D model generated for different benchmark values to remove long edges: (a) benchmark = 0.201 (default), (b) benchmark = 0.35, and (c) benchmark = 1.1. What is the effective benchmark value to remove long edges?

## 4 Deciding on the most suitable fit

The appropriate model for the S-curve using UMAP (n\_neighbors: 15) was created with 420 bins. It was determined to be the best model after removing edges with a length greater than 0.134. The decision was made considering that the model accurately captures the geometry of the S-curve. But better in some places only ([?@fig-scurve\\_sc\\_best](#)).

In contrast, a poor model for the S-curve with PHATE was created using only 190 bins and removing edges with a length greater than 0.3. This model is inadequate as it has a higher MSE compared to the appropriate model. Visually, the model squeezed to the middle of the geometry of the S-curve ([?@fig-scurve-sc-bad](#)).

## 5 Illustration on simulated data

In this section, the effectiveness of the algorithm is described using a simulated dataset. The dataset consists of five spherical Gaussian clusters in 4-D, with each cluster containing an equal number of points and the same within-cluster variation.

In the 2D layouts generated by various NLDR techniques, as shown in Figure 9, five well-separated clusters are shown. In tSNE (see Figure 9 (a)), these clusters appear closely. UMAP arranges all clusters in a parallel manner, with three aligned in one line and the other two in a separate line (see Figure 9 (b)). In contrast, PHATE shows two closely positioned clusters and three more distant ones (see Figure 9 (c)). In TriMAP, two clusters are close, though not as tightly as PHATE, while the other three are well-separated (see Figure 9 (d)). Finally, PaCMAP shows one central cluster and the remaining four spread out in different directions (see Figure 9 (e)).

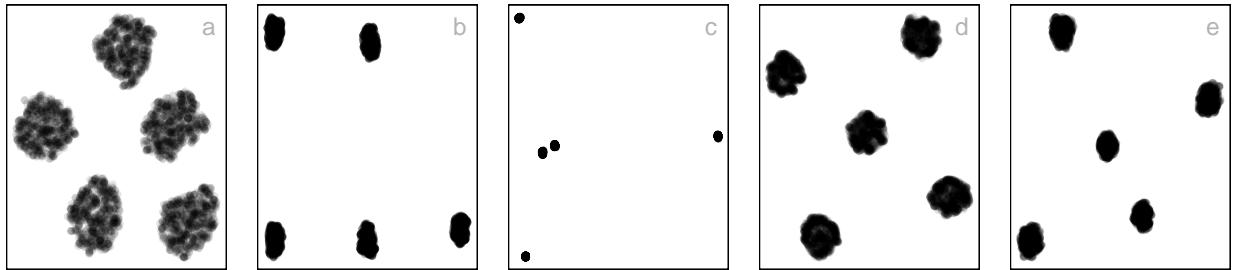


Figure 9: Five different NLDR representations of the same data. Different techniques and different parameter choices are used. Is there a best representation of the original data or are they all providing equivalent information?

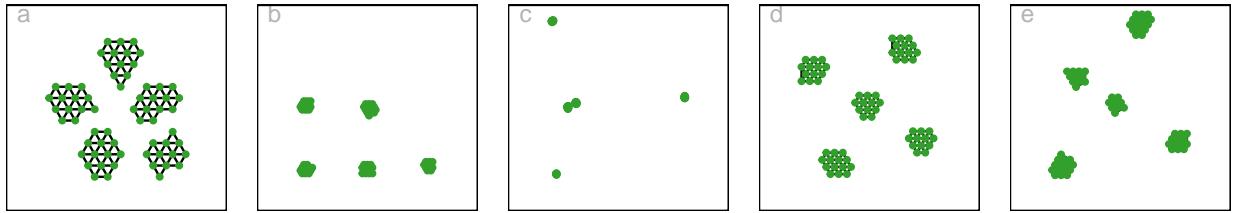


Figure 10: Model generated with five different NLDR methods in 2-D with approximately 65 non-empty bins in each.

To investigate which is the reasonable representation to visualize the five spherical Gaussian cluster data or all NLDR methods provide equivalent information, we visualize all the models in  $p$ - $D$  space. Models from all NLDR methods show five well-separated clusters (see Figure 11, Figure 12, Figure 13, Figure 14, and Figure 15). This suggests that for the five Gaussian cluster dataset, all NLDR methods effectively preserve the global structure. tSNE displays clusters with varying densities, indicating their ability to capture within-cluster variation (see Figure 11). On the other hand, both UMAP, PHATE, PaCMAP and TriMAP show clusters with flat surfaces, suggesting a failure to capture within-cluster variation (see Figure 12, Figure 13, Figure 14 and Figure 15). Therefore, UMAP, PHATE, PaCMAP and TriMAP do not capture the local structure as effectively as other methods.

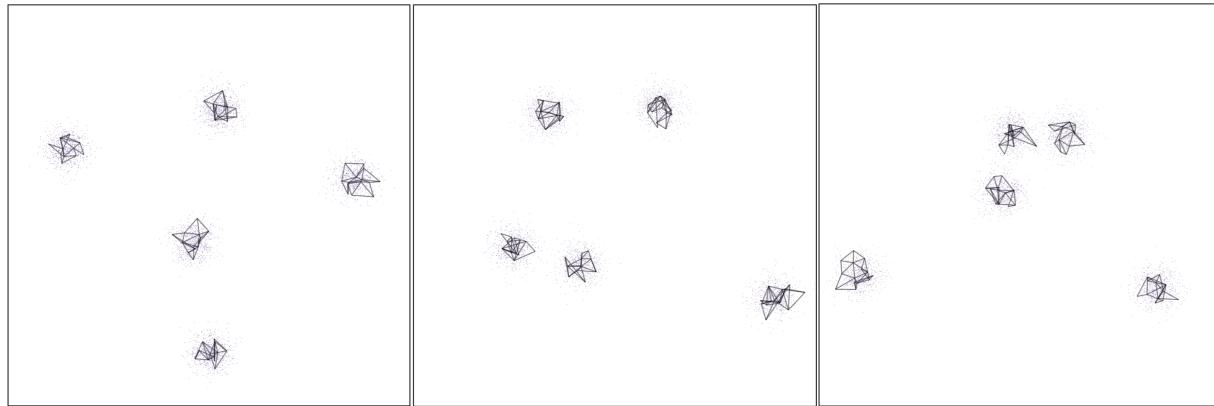


Figure 11: Screen shots of the **langevitour** of the five Gaussian clusters dataset, shows the model with tSNE in  $p$ - $D$ , a video of the tour animation is available at (<https://youtu.be/RASEE7N5MbM>).

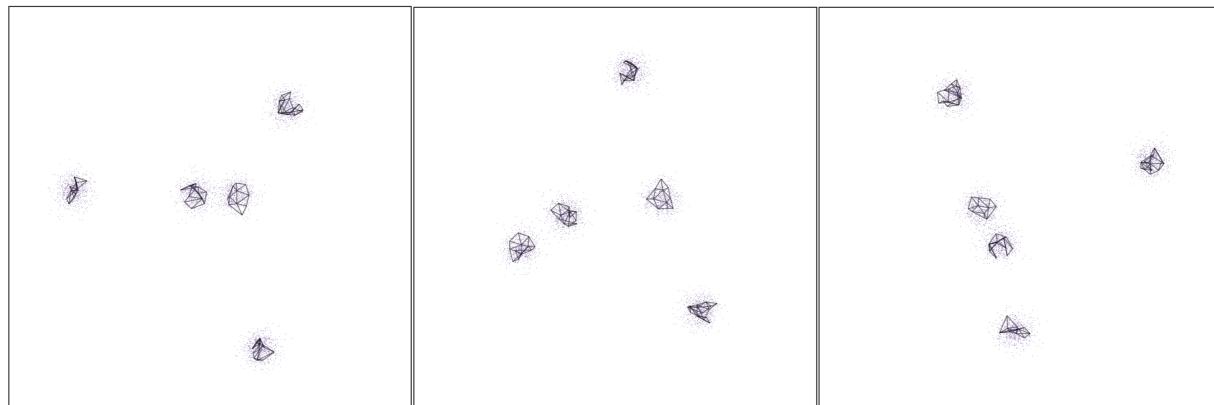


Figure 12: Screen shots of the **langevitour** of the five Gaussian clusters dataset, shows the model with UMAP in  $p$ - $D$ , a video of the tour animation is available at (<https://youtu.be/iG4bCPkJlw>).

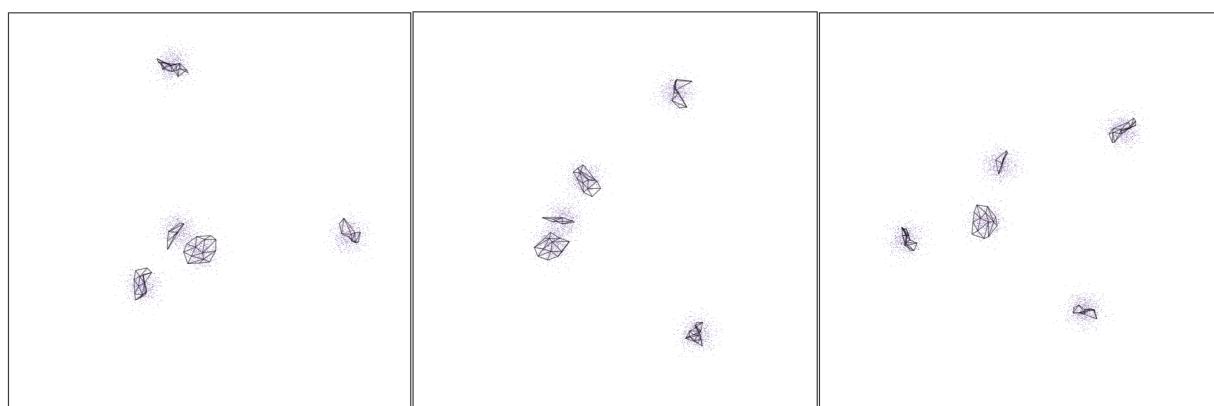


Figure 13: Screen shots of the **langevitour** of the five Gaussian clusters dataset, shows the model with PHATE in  $p$ - $D$ , a video of the tour animation is available at ([https://youtu.be/L\\_PVLGwfOS0](https://youtu.be/L_PVLGwfOS0)).

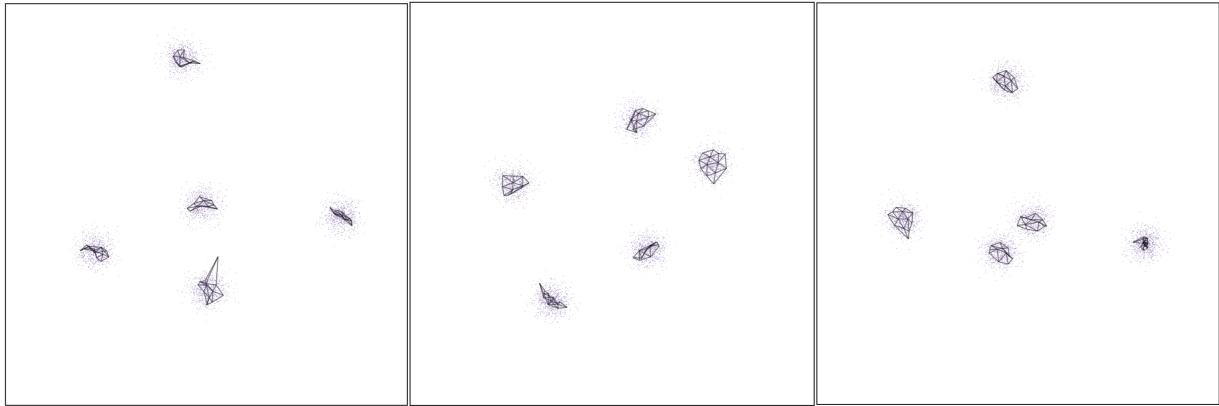


Figure 14: Screen shots of the **langevitour** of the five Gaussian clusters dataset, shows the model with PaCMAP in  $p$ - $D$ , a video of the tour animation is available at (<https://youtu.be/z07cKXi8EJQ>).

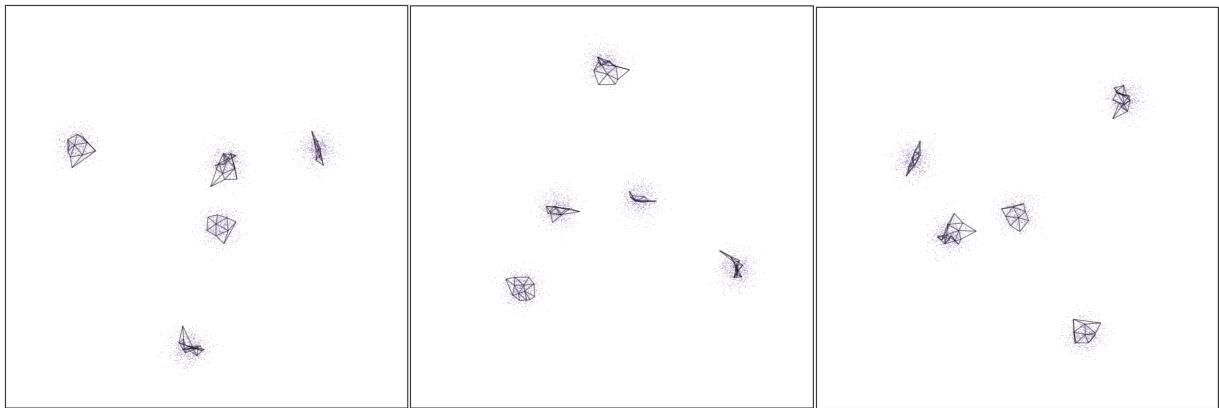


Figure 15: Screen shots of the **langevitour** of the five Gaussian clusters dataset, shows the model with TriMAP in  $p$ - $D$ , a video of the tour animation is available at (<https://youtu.be/Chs1lYAoX2w>).

When compare the NLDR representations and generated models, tSNE with perplexity 61 appears to be a reasonable representation for visualizing the five Gaussian cluster dataset. This is supported by investigating the model generated with tSNE in the data space, which provides evidence that it preserves both local and global structures. Also, the NLDR representation with tSNE shows five well-separated clusters.

## 6 Applications

### 6.1 Single-cell gene expression

In the field of single-cell studies, a common analytical task involves clustering to identify groups of cells with similar expression profiles. Analysts often turn to NLDR techniques to verify and identify these clusters and explore developmental trajectories. To illustrate the

importance of NLDR techniques and parameter selection in identifying clusters, Human Peripheral Blood Mononuclear Cells (PBMC3k) dataset (Chen et al. 2023) is used. In a study by Chen et al. (2023), this dataset was used to demonstrate how UMAP represents clusters (see Figure 16). As shown in Figure 16, there are three distant and well-separated clusters.



Figure 16: 2- $D$  layout from UMAP applied for the PBMC3k dataset. Is this a best representation of the original data? The parameter setting is  $n\_neighbors = 30$ ,  $\text{min\_dist} = 0.3$ .

To determine whether the UMAP representation with the parameter choice suggested by Chen et al. (2023) preserves the original data structure, we visualize the model constructed with UMAP overlaid on the  $p$ - $D$  data. The figures in Figure 18 show three well-separated clusters, indicating that the suggested UMAP representation preserves the global structure (see Figure 16). However, as shown in Figure 18, these clusters are close to each other in  $p$ - $D$ . Also, non-linear continuity patterns and high-density patches within the clusters are observed (see Figure 18). Therefore, the suggested UMAP representation (see Figure 16) does not accurately preserve the local structure of the PBMC3k dataset.

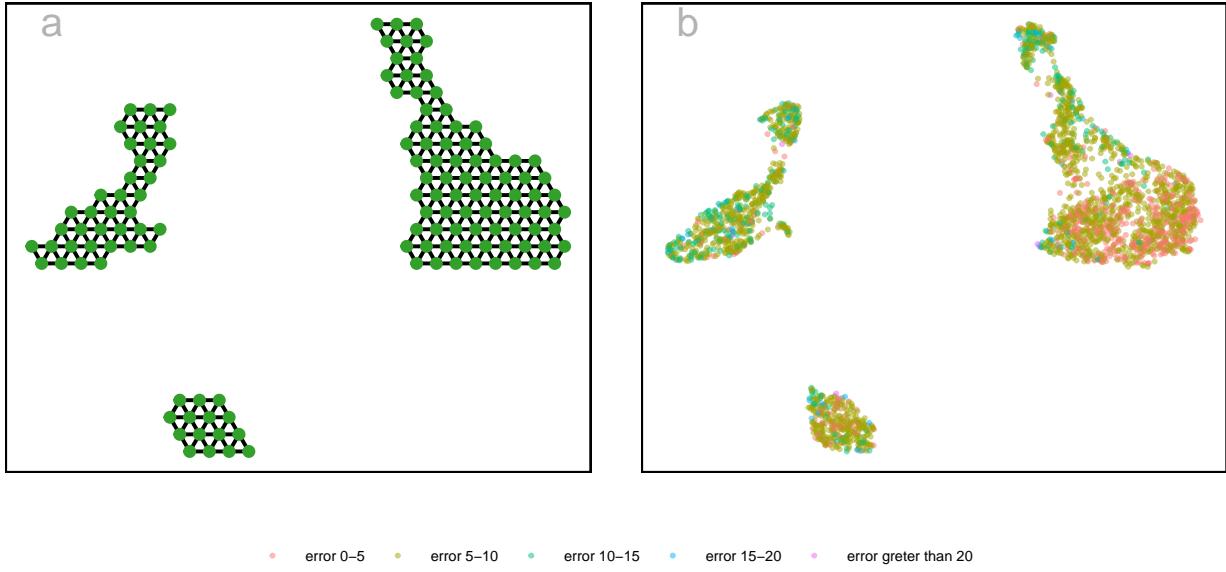


Figure 17: (a) Model generated in 2-D with UMAP, and (b)  $p$ -D model error in 2-D. The 2-D model shows three well-separated distant clusters. The  $p$ -D model errors are distributed along clusters.

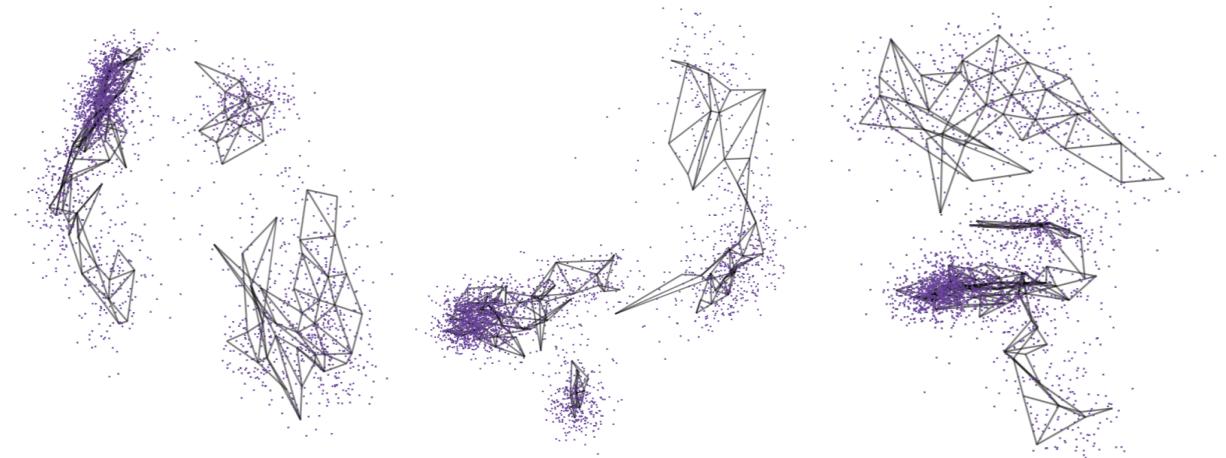


Figure 18: Screen shots of the **langevitour** of the PBMC3k data set, shows the model-in-data space, a video of the tour animation is available at (<https://youtu.be/VqqWuE0Jj6A>).

In order to find a reasonable NLDR representation for the PBMC3k dataset, we calculated the absolute error for different numbers of non-empty bins using various NLDR techniques and different parameter settings (see Figure 19). After analyzing the results, we found that tSNE with a perplexity set to 30 had the lowest error when the number of non-empty bins was 137. Therefore, tSNE with a perplexity of 30, which is the default parameter setting, is considered as a reasonable representation for the PBMC3k dataset.

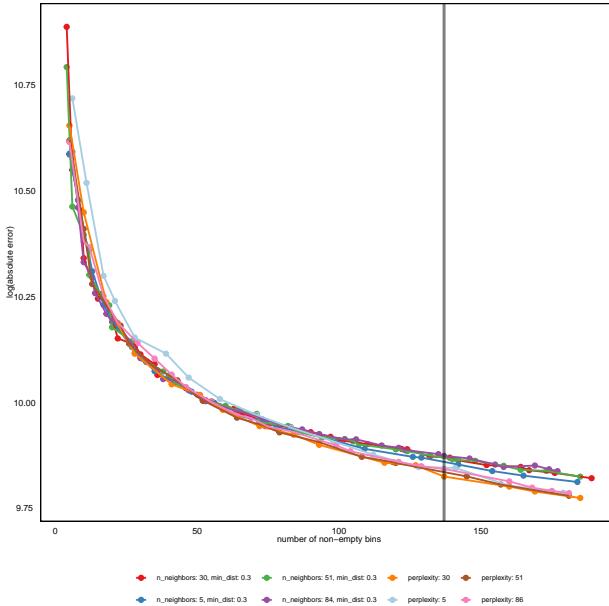


Figure 19: Absolute error from UMAP and tSNE applied to training PBMC3k dataset with different parameter choices. What is the best parameter choice to create the model? The residual plot have a steep slope at the beginning, indicating that a smaller number of non-empty bins causes a larger amount of error. Then, the slope gradually declines or level off, indicating that a higher number of non-empty bins generates a smaller error. Using the elbow method, it was observed that when the number of non-empty bins is set to 137, the lowest error occurred with the parameters perplexity: 30.

As shown in Figure 20, there are three well-separated clusters, although they are located close to each other. Additionally, non-linear structures can also be observed within the clusters (see Figure 21 (a)). In this manner, tSNE was able to capture the data structure for the PBMC3k dataset that UMAP failed to do.

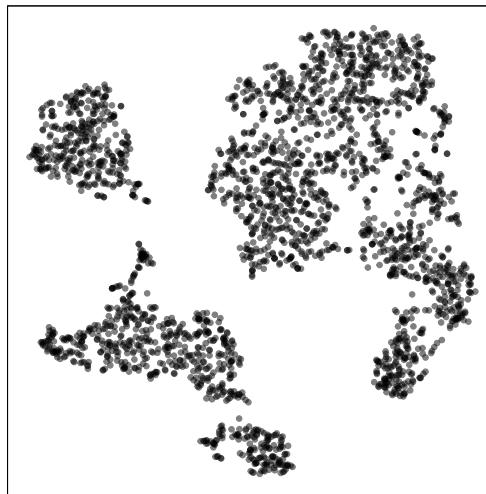


Figure 20: 2-D layout from tSNE applied for the PBMC3k dataset. Is this a best representation of the original data? The parameter setting is perplexity=30.

We then fit the model for tSNE, and visualize the resultant model in the  $p$ - $D$  data space. The model shows a quirk, as shown in Figure 22. All three clusters are connected by an edge except the small and large clusters. Because the clusters are so close in 2-D, they attempt to maintain the structure in  $p$ - $D$  as well. This is evident that tSNE with perplexity 30 provides a reasonable representation of PBMC3k data.

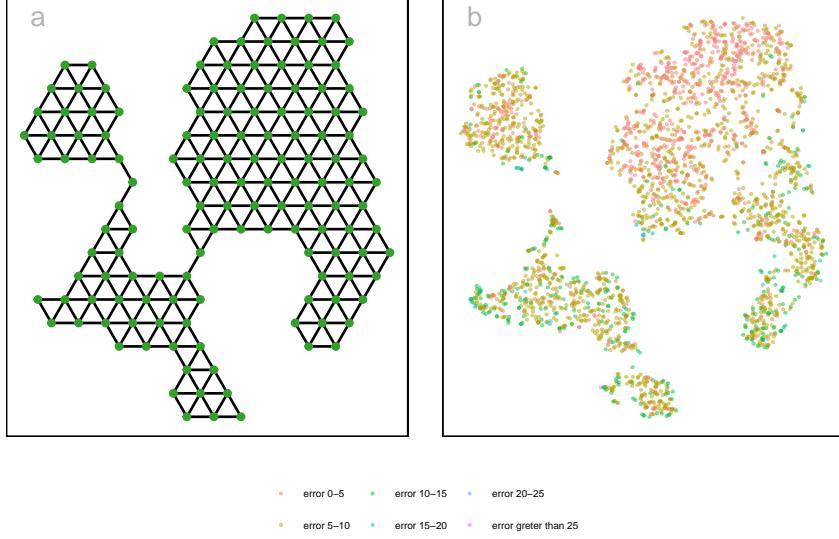


Figure 21: (a) Model generated in 2-D with tSNE, and (b)  $p$ - $D$  model error in 2-D. The 2-D model shows three well-separated distant clusters. The  $p$ - $D$  model errors are distributed along clusters, but most low  $p$ - $D$  model errors present in the large cluster.

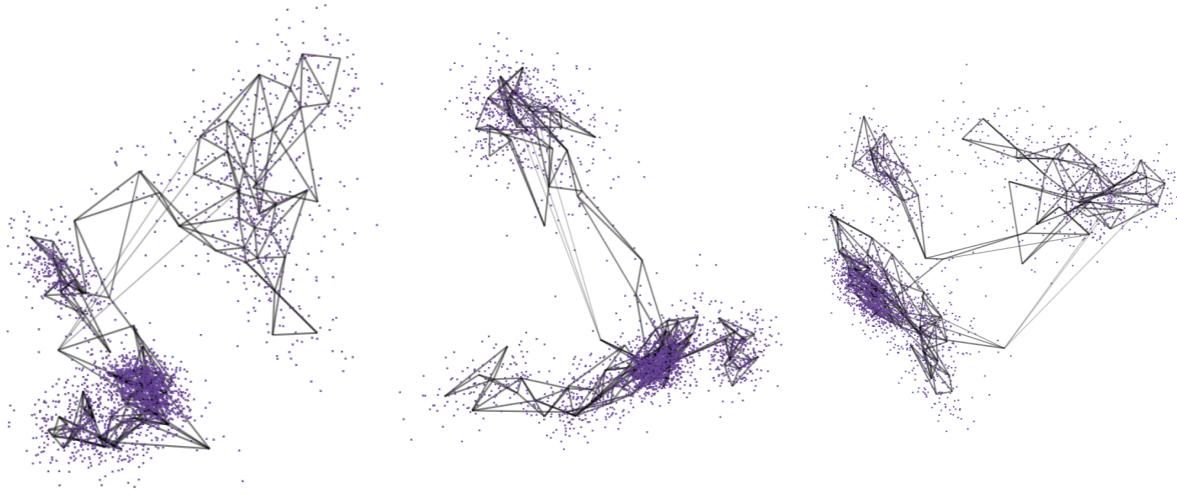


Figure 22: Screen shots of the **langevitour** of the PBMC3k data set, shows the model in high-D, a video of the tour animation is available at (<https://youtu.be/5Y1hE4i7N2k>).

## 6.2 Hand-written digits

The MNIST dataset consists of grayscale images of handwritten digits (LeCun & Cortes 2010). Wang et al. (2021) used this dataset to demonstrate how PaCMAP preserves local

structure. We selected the 2-*D* embedding of PaCMAP for the handwritten digit 1 to assess whether this is a reasonable representation using our method. As shown in Figure 24, the angle of the digit 1 images varies along the 2-*D* structure.

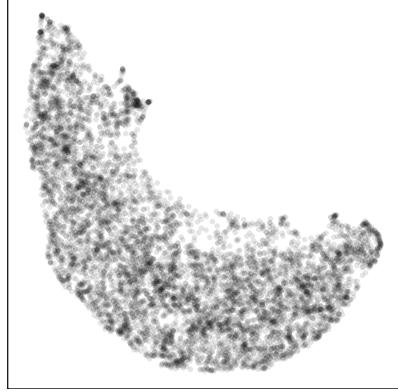


Figure 23: 2-*D* layout from PaCMAP applied for the digit 1 of the MNIST dataset. Is this the best representation of the digit 1? The parameter setting is n\_components=2, n\_neighbors=10, init=random, MN\_ratio=0.9, FP\_ratio=2.0.

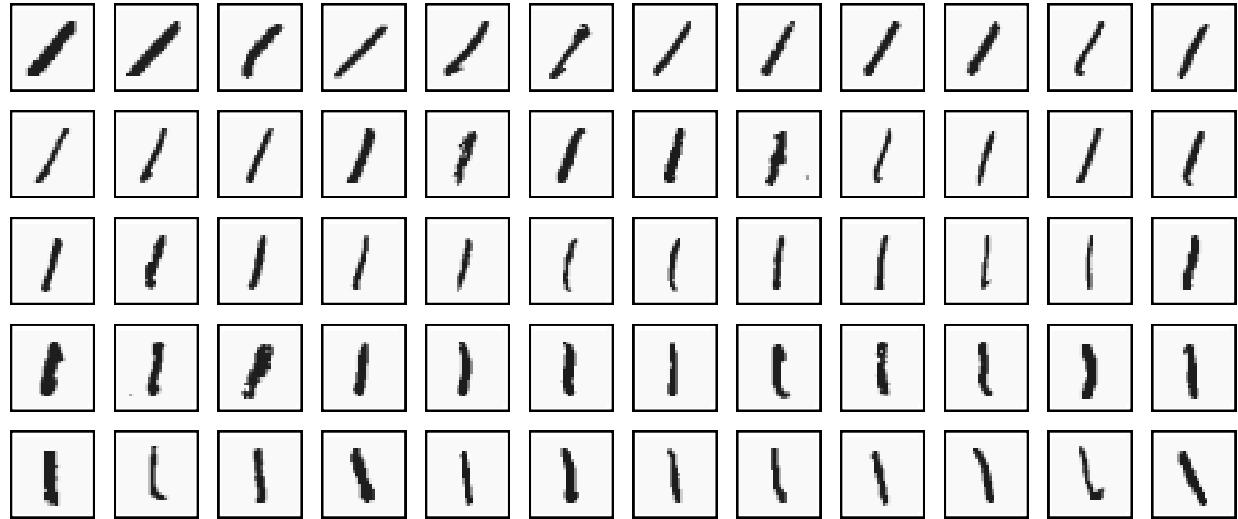


Figure 24: Images of the handwritten digit 1 are ordered from the bottom-right to the top-left of the 2-*D* structure. The angle of the digit varies along this structure. Images at the bottom-right of the 2-*D* layout show the digit 1 angled more to the right, while images at the top-left show the digit 1 angled more to the left. This demonstrates how the angle changes from right to left along the 2-*D* structure.

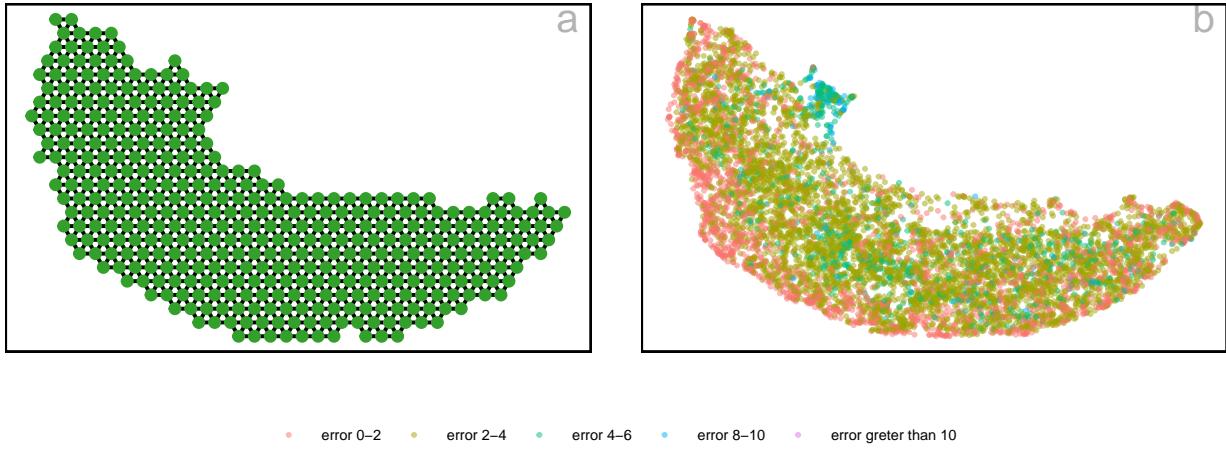


Figure 25: (a) Model generated in 2-D, and (b)  $p$ -D model error in 2-D. The 2-D model shows a non-linear continuous structure. Most low  $p$ -D model errors are distributed along the lower edge of the 2-D structure, while most high  $p$ -D model errors are concentrated along the upper edge.

According to Figure 26a, the non-linear continuous structure observed in the 2-D representation of PaCMAP (see Figure 23) is also visible when visualizing the model overlaid on the data space. This indicates that PaCMAP accurately captures the structure of the  $p$ -D data. Additionally, the model shows a twisted pattern within the non-linear structure in  $p$ -D space (see Figure 26b), which is an additional pattern not visible in the 2-D representation (see Figure 23). Furthermore, as shown in Figure 26c, some long edges exist in the  $p$ -D space that are not recognized as long edges in the 2-D representation. However, PaCMAP is a reasonable 2-D representation of MNIST digit 1 data. Because PaCMAP preserves the local structure.

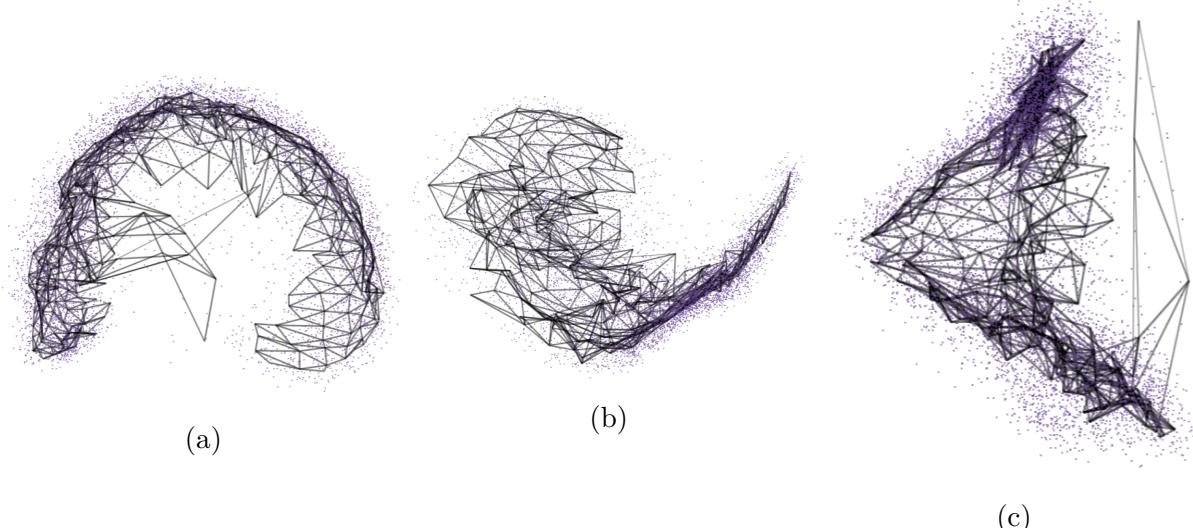


Figure 26: Screen shots of the **langevitour** of the MNIST digit 1 data set, shows the model-in-data space, a video of the tour animation is available at ([https://youtu.be/zcg\\_GXBmjqA](https://youtu.be/zcg_GXBmjqA)).

There are certain data points that exhibit high error rates due to their deviation from the usual  $p$ - $D$  data structure, which makes them anomalies (see Figure 25 (b)). These anomalies can be classified into two types: those that are anomalies within the non-linear structure and those that lie outside of it. The images associated with high model error points within the non-linear structure display different patterns of the digit 1, as shown in Figure 27 (a). However, when comparing these images to the ones found outside of the non-linear structure, it becomes evident that the latter display different patterns of the digit 1 (see Figure 27 (b)).

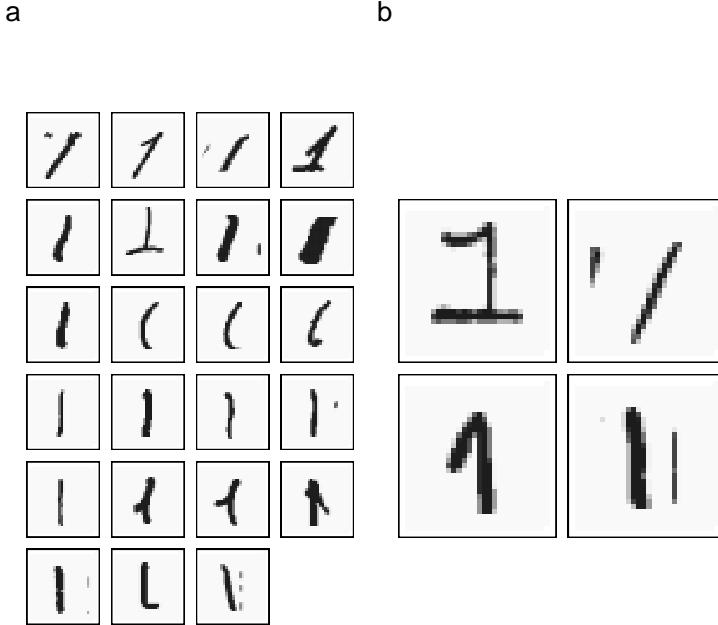


Figure 27: Some images of handwritten digit 1 which occur high model error (a) within the non-linear structure, and (b) outside the non-linear structure. The images shows different patterns of digit 1.

## 7 Discussion

This study makes several important contributions to the field of NLDR. We have developed an algorithm to evaluate the most useful NLDR method and (hyper-)parameter choices for creating a reasonable 2- $D$  layout of high-dimensional data. Our objective is to fit a model for the 2- $D$  layout that preserves the relationships between neighboring points and turns it into a high-dimensional wireframe, which can be overlaid on the data and visualized using a tour. This approach is defined as *model-in-data-space*. Viewing a model in the data space is an ideal way to examine the fit.

The effectiveness of this approach is illustrated through various examples. For instance, the S-curve example demonstrates how the model accurately fits the points, capturing both local and global structures in high-dimensional space. Our simulation case study further, five Gaussian cluster example shows that while all observed NLDR methods preserve the global structure, only tSNE effectively maintains the local structure, highlighting the specific strengths and quirks of different methods.

Human behavior often shows a desire for more certainty and a tendency to prefer well-separated views. This emphasizes the importance of clear and distinct clusters. For example, in the UMAP layout of the **pbmc** dataset suggested by Chen et al. (2023), three distant, well-separated clusters are shown. However, our model reveals that these clusters are actually close to each other in  $p$ - $D$ . Additionally, the model discovers non-uniform data distribution and non-linear structures within the clusters that are not visible in the UMAP layout, demonstrating the ability of our model in uncovering hidden data characteristics.

Evaluating the error or unexplained variance is important for assessing how well the model fits the data. By examining the error for different numbers of bins, we found that tSNE with a perplexity of 30 provides a reasonable representation for the **pbmc** dataset. Connecting the closest clusters with line segments in the fitted model further supports the preservation of neighborhood relationships.

The **digit: 1** example further illustrates the model’s ability to accurately capture non-linear structures and provide additional information. Key findings include a twisted pattern that compresses the structure in some projections and long line segments that detect anomalies.

Predicting new observations in  $k$ - $D$  is particularly valuable due to the limitations of some NLDR methods, like tSNE, which don’t provide a straightforward method for prediction. As a result, our approach offers a solution that capable of generating predicted  $k$ - $D$  embedding regardless of the NLDR method employed, effectively addressing this functional gap.

In conclusion, while our method effectively captures and represents high-dimensional data structures, further enhancements could involve introducing approaches to bind the data, indicate line segments beyond 2- $D$ , and diagnose the fitted model. These improvements would help in creating a more accurate representation of the data when 2- $D$  layout is inadequate.

## 8 Supplementary Materials

Code, and data for reproducing this paper are available at <https://github.com/JayaniLakshika/paper-nldr-vis-algorithm>.

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