ML 813: Topics on Dimensionality Reduction and Manifold Learning

Uniform Manifold Approximation and Projection (UMAP)

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Uniform Manifold Approximation and Projection: UMAP

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What is UMAP? (1)

Definition

UMAP (Uniform Manifold Approximation and Projection) is a powerful probabilistic-based technique in dimensionality reduction within machine learning. It tackles the challenge of simplifying complex, high-dimensional data into a more digestible, low-dimensional space, preserving crucial information.

Major Aspects

- Reduces dimensionality: Imagine a dataset with dozens of features. UMAP helps you
 condense it into a few key dimensions, making it easier to visualize, analyze, and train
 machine learning models.
- Preserves relationships: Unlike some techniques, UMAP excels at capturing the
 underlying structure of your data. It prioritizes relationships between points, ensuring
 similar data points stay close in the low-dimensional space, even if they're far apart in the
 original high-dimensional space.

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What is UMAP? (2)

UMAP Importance

- Speed and Scalability: UMAP is significantly faster than some popular alternatives like t-SNE, making it a practical choice for large datasets.
- Flexibility: UMAP offers various customizable parameters, allowing you to tailor the
 results to your specific data and needs.
- Global and local behavior: UMAP balances the "big picture" (global structure) with the fine details (local relationships), offering a comprehensive view of your data.

Major Aspects

- Visualization: UMAP's low-dimensional representations are ideal for creating informative and interpretable visualizations of complex data.
- Clustering: Identifying groups within your data becomes easier with UMAP's well-preserved relationships.
- Feature engineering: Extracted UMAP features can be used as inputs to other machine learning tasks, boosting their performance.

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Data Graph in the Input Space

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Data Graph in the Input Space

- Uniform Manifold Approximation and Projection (UMAP) was proposed in 2018 [1].
- Consider a training dataset $\mathbf{X} = [\mathbf{x}_1, \dots, \mathbf{x}_n] \in \mathbb{R}^{d \times n}$ where n is the sample size and d is the dimensionality.
- We construct a k-Nearest Neighbors (kNN) graph for this dataset. It has been empirically observed that UMAP requires fewer number of neighbors than t-SNE [2]. Its default value is k=15. We denote the j-th neighbor of x_i by $x_{i,j}$. Let \mathcal{N}_i denote the set of neighbor points for the point x_i , i.e., $\mathcal{N}_i := \{x_{i,1}, \ldots, x_{i,k}\}$.
- We treat neighborhood relationship between points stochastically. Inspired by SNE [3] and t-SNE [4, 5], we use the **Gaussian or Radial Basis Function (RBF)** kernel for the measure of similarity between points in the **input space**. The probability that a point x_i has the point x_j as its neighbor can be computed by the similarity of these points:

$$p_{j|i} := \begin{cases} \exp\left(-\frac{\|\mathbf{x}_i - \mathbf{x}_j\|_2 - \rho_i}{\sigma_i}\right) & \text{if } \mathbf{x}_j \in \mathcal{N}_i \\ 0 & \text{Otherwise,} \end{cases}$$
 (1)

where $\|.\|_2$ denotes the ℓ_2 norm.

• The ρ_i is the distance from x_i to its nearest neighbor:

$$\rho_i := \min\{\|\mathbf{x}_i - \mathbf{x}_{i,j}\|_2 \mid 1 \le j \le k\}. \tag{2}$$

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Data Graph in the Input Space (1)

• The σ_i is the scale parameter which is calculated such that the total similarity of point x_i to its k nearest neighbors is normalized. By binary search, we find σ_i to satisfy:

$$\sum_{j=1}^{k} \exp\left(-\frac{\|\mathbf{x}_{i} - \mathbf{x}_{i,j}\|_{2} - \rho_{i}}{\sigma_{i}}\right) = \log_{2}(k).$$
 (3)

• Note that t-SNE [4] has a similar search for its scale using entropy as perplexity. These searches make the neighborhoods of various points behave similarly because the scale for a point in a dense region of dataset becomes small while the scale of a point in a sparse region of data becomes large.

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Data Graph in the Input Space (2)

- In other words, UMAP and t-SNE both assume (or approximate) that points are uniformly distributed on an underlying low-dimensional manifold. This approximation is also included in the name of UMAP.
- Eq. (1):

$$p_{j|i} := \left\{ egin{array}{ll} \exp\left(-rac{\|oldsymbol{x}_i - oldsymbol{x}_j\|_2 -
ho_i}{\sigma_i}
ight) & ext{if } oldsymbol{x}_j \in \mathcal{N}_i \ 0 & ext{Otherwise,} \end{array}
ight.$$

is a **directional similarity** measure. To have a **symmetric** measure with respect to i and j, we symmetrize it as:

$$\mathbb{R} \ni p_{ij} := p_{j|i} + p_{i|j} - p_{j|i} p_{i|j}. \tag{4}$$

This is a symmetric measure of similarity between points x_i and x_j in the input space.

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Data Graph in the Embedding Space

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Data Graph in the Embedding Space

- Let the embeddings of points be $\mathbf{Y} = [\mathbf{y}_1, \dots, \mathbf{y}_n] \in \mathbb{R}^{p \times n}$ where p is the dimensionality of embedding space and is smaller than input dimensionality, i.e., $p \ll d$. Note that \mathbf{y}_i is the embedding corresponding to \mathbf{x}_i .
- In the embedding space, the probability that a point y_i has the point y_j as its neighbor can be computed by the similarity of these points:

$$\mathbb{R} \ni q_{ij} := (1 + a \| \mathbf{y}_i - \mathbf{y}_j \|_2^{2b})^{-1}, \tag{5}$$

which is **symmetric** with respect to i and j.

• The variables a>0 and b>0 are hyperparameters determined by the user. By default, we have $a\approx 1.929$ and $b\approx 0.7915$ [1], although it has been empirically seen that setting a=b=1 does not qualitatively impact the results [6].

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 UMAP aims to make the data graph in the low-dimensional embedding space similar to the data graph in the high-dimensional embedding space. In other words, we treat Eqs. (4) and (5) as probability distributions and minimize the difference of these distributions to make similarities of points in the embedding space as the similarities of points in the input space. A measure for the difference of these similarities of graphs is the fuzzy cross-entropy defined as:

$$c_1 := \sum_{i=1}^n \sum_{j=1, j \neq i}^n \left(p_{ij} \ln(\frac{p_{ij}}{q_{ij}}) + (1 - p_{ij}) \ln(\frac{1 - p_{ij}}{1 - q_{ij}}) \right), \tag{6}$$

where ln(.) is the natural logarithm. The definition of this cross-entropy is in the field of fuzzy category theory (see our tutorial paper [7] for more information).

- The first term in Eq. (6) is the attractive force which attracts the embeddings of neighbor points toward each other. This term should only appear when $p_{ii} \neq 0$ which means either x_i is a neighbor of x_i , or x_i is a neighbor of x_i , or both (see Eq. (4), $\mathbb{R}\ni p_{ij}:=p_{i|j}+p_{i|j}-p_{i|i}p_{i|j}).$
- The second term in Eq. (6) is the repulsive force which repulses the embeddings of non-neighbor points away from each other.

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- As the number of all permutations of non-neighbor points is very large, computation of the second term is non-tractable in big data. Inspired by Word2Vec [8] and LargeVis [9], UMAP uses negative sampling where, for every point x_i, m points are sampled randomly from the training dataset and treat them as non-negative (negative) points for x_i.
- As the dataset is usually large, i.e. $m \ll n$, the sampled points will be actual negative points with high probability.
- We had:

$$c_1 := \sum_{i=1}^n \sum_{j=1, j \neq i}^n \Big(p_{ij} \ln(\frac{p_{ij}}{q_{ij}}) + (1-p_{ij}) \ln(\frac{1-p_{ij}}{1-q_{ij}}) \Big).$$

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- The summation over the second term in Eq. (6) is computed only over these negative samples rather than all negative points.
- UMAP changes the data graph in the embedding space to make it similar to the data graph in the input space.
- Eq. (6) is the cost function which is minimized in UMAP where the optimization variables are {y_i}ⁿ_{i-1}:

$$\begin{split} \min_{\left\{ \mathbf{y}_{i} \right\}_{i=1}^{n}} c_{1} &:= \min_{\left\{ \mathbf{y}_{i} \right\}_{i=1}^{n}} \sum_{j=1}^{n} \sum_{j=1, j \neq i}^{n} \left(p_{ij} \ln(p_{ij}) - p_{ij} \ln(q_{ij}) \right. \\ &+ \left. \left(1 - p_{ij} \right) \ln(1 - p_{ij}) - \left(1 - p_{ij} \right) \ln(1 - q_{ij}) \right). \end{split}$$

• According to Eqs. (1), (4), and (5), in contrast to q_{ij} , the p_{ij} is independent of the optimization variables $\{y_i\}_{i=1}^n$. Hence, we can drop the constant terms to revise the cost function:

$$c_2 := -\sum_{i=1}^n \sum_{j=1, j \neq i}^n \left(p_{ij} \ln(q_{ij}) + (1 - p_{ij}) \ln(1 - q_{ij}) \right), \tag{7}$$

which should be minimized.

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We had:

$$c_2 := -\sum_{i=1}^n \sum_{j=1, j
eq i}^n \Big(
ho_{ij} \ln(q_{ij}) + (1-
ho_{ij}) \ln(1-q_{ij}) \Big).$$

Two important terms in this cost function are:

$$c_{i,j}^{\mathfrak{a}} := -\ln(q_{ij}), \tag{8}$$

$$c_{i,j}^r := -\ln(1-q_{ij}),$$
 (9)

and we can write:

$$c_2 := \sum_{i=1}^n \sum_{j=1, j \neq i}^n \left(p_{ij} \, c_{i,j}^{a} + (1 - p_{ij}) \, c_{i,j}^{r} \right) \tag{10}$$

$$\stackrel{\text{(a)}}{=} 2 \sum_{i=1}^{n} \sum_{j=i+1}^{n} \left(p_{ij} c_{i,j}^{a} + (1 - p_{ij}) c_{i,j}^{r} \right), \tag{11}$$

where (a) is because $p_{ij} = p_{ji}$, $c_{i,j}^a = c_{i,i}^a$, and $c_{i,j}^r = c_{i,i}^r$ are symmetric.

- The Eqs. (8) and (9) are the attractive and repulsive forces in Eq. (7), respectively. The attractive force attracts the neighbor points toward each other in the embedding space while the repulsive force pushes the non-neighbor points (i.e., points with low probability of being neighbors) away from each other in the embedding space.
- According to Eq. (10), $c_{i,j}^a$ and $c_{i,j}^r$ occur with probability p_{ij} and $(1 p_{ij})$, respectively. For every point, we call it the **anchor** point and we call its neighbor and non-neighbor points, with large and small p_{ii} , as the **positive** and **negative** points, respectively.

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The Training Algorithm of UMAP

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The Training Algorithm of UMAP

```
1 Input: Training data \{x_i\}_{i=1}^n
 2 Construct kNN graph
 \mathfrak{z} Initialize \{oldsymbol{y}_i\}_{i=1}^n by Laplacian eigenmap
 4 Calculate p_{ij} and q_{ij} for \forall i, j \in \{1, ..., n\} by
       Eqs. (4) and (5)
 5 n \leftarrow 1, \nu \leftarrow 0
 6 while not converged do
           \nu \leftarrow \nu + 1 // epoch index
           for i from 1 to n do
                  for j from 1 to n do
                        u \sim U(0,1)
10
                        if u \leq p_{ij} then
11
12
13
                              for m iterations do
14
                                  \begin{vmatrix} l \sim U\{1, \dots, n\} \\ \mathbf{y}_i \leftarrow \mathbf{y}_i - \eta \frac{\partial c_{i,l}^r}{\partial \mathbf{y}_i} \end{vmatrix} 
15
16
                                     // The next line does not exist
17
                                     in original UMAP:
                                    \boldsymbol{y}_l \leftarrow \boldsymbol{y}_l - \eta \frac{\partial c_{i,l}^r}{\partial \boldsymbol{y}_l}
18
        \eta \leftarrow 1 - \frac{\nu}{\nu_{max}}
20 Return \{y_i\}_{i=1}^n
```

Algorithm 1: UMAP algorithm

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The Training Algorithm of UMAP(1)

- As this algorithm shows, a kNN graph is constructed from the training data $\{x_i\}_{i=1}^n$.
- UMAP uses Laplacian eigenmap [10, 11], also called spectral embedding, for initializing the embeddings of points denoted by $\{y_i\}_{i=1}^n$.
- Using Eqs. (4) and (5), p_{ij} and q_{ij} are calculated for all points.
- Stochastic Gradient Descent (SGD) is used for optimization where optimization is performed iteratively.
- In every iteration (epoch), we iterate over points twice with indices i and j where the i-th point is called the **anchor**. For every pair of points x_i and x_j , we update their embeddings x_i and x_j with probability p_{ij} (recall Eq. (7)).
- If p_{ij} is large, it means that the points x_i and x_j are probably neighbors (in this case, the j-th point is called the **positive** point) and their embeddings are highly likely to be updated to become close in the embedding space based on the attractive force. For implementing it, we can **sample a uniform** value from the continuous uniform distribution U(0,1) and if that is less than p_{ij} , we update the embeddings.

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The Training Algorithm of UMAP (2)

- We update the embeddings \mathbf{y}_i and \mathbf{y}_j by gradients $\partial c_{i,j}^a/\partial \mathbf{y}_i$ and $\partial c_{i,j}^a/\partial \mathbf{y}_j$, respectively, where η is the learning rate.
- For repulsive forces, we use negative sampling as was explained before. If m denotes the size of negative sample, we sample m indices from the discrete uniform distribution $U\{1,\ldots,n\}$. These are the indices of points which are considered as negative samples $\{y_i\}$ where $|\{y_i\}|=m$. As the size of dataset is usually large enough to satisfy $n\gg m$, these negative points are probably valid because many of the points are non-neighbors of the considered anchor.
- In negative sampling, the **original UMAP** [1] updates only the embedding of anchor y_i by gradient of the repulsive force $\partial c_{i,j}^a/\partial y_i$. One can **additionally update the embedding of negative point** y_i by gradient of the repulsive force $\partial c_{i,j}^a/\partial y_i$ [12].

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The Training Algorithm of UMAP

We had:

$$egin{aligned} c_{i,j}^{a} &:= -\ln(q_{ij}), \ c_{i,j}^{r} &:= -\ln(1-q_{ij}). \end{aligned}$$

• The gradients of attractive and repulsive cost functions in UMAP are:

$$\frac{\partial c_{i,j}^{a}}{\partial \mathbf{y}_{i}} = \frac{2ab\|\mathbf{y}_{i} - \mathbf{y}_{j}\|_{2}^{2(b-1)}}{(1 + a\|\mathbf{y}_{i} - \mathbf{y}_{i}\|_{2}^{2b})}(\mathbf{y}_{i} - \mathbf{y}_{j}), \tag{12}$$

$$\frac{\partial c_{i,j}^r}{\partial \mathbf{y}_i} = \frac{-2b}{(\varepsilon + \|\mathbf{y}_i - \mathbf{y}_j\|_2^2)(1 + a\|\mathbf{y}_i - \mathbf{y}_j\|_2^{2b})}(\mathbf{y}_i - \mathbf{y}_j), \tag{13}$$

where ε is a small positive number, e.g. $\varepsilon=0.001$, for stability to prevent division by zero when $\mathbf{y}_i\approx\mathbf{y}_i$.

Likewise. we have:

$$\begin{split} \frac{\partial c_{i,j}^{a}}{\partial \mathbf{y}_{j}} &= \frac{2ab\|\mathbf{y}_{i} - \mathbf{y}_{j}\|_{2}^{2(b-1)}}{(1 + a\|\mathbf{y}_{i} - \mathbf{y}_{j}\|_{2}^{2b})}(\mathbf{y}_{j} - \mathbf{y}_{i}), \\ \frac{\partial c_{i,j}^{r}}{\partial \mathbf{y}_{i}} &= \frac{-2b}{(\varepsilon + \|\mathbf{y}_{i} - \mathbf{y}_{i}\|_{2}^{2})(1 + a\|\mathbf{y}_{i} - \mathbf{y}_{i}\|_{2}^{2b})}(\mathbf{y}_{j} - \mathbf{y}_{i}). \end{split}$$

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The Training Algorithm of UMAP (1)

Proof:

We had:

$$egin{aligned} c_{i,j}^{a} &:= -\ln(q_{ij}), \ c_{i,j}^{r} &:= -\ln(1-q_{ij}), \ \mathbb{R} &\ni q_{ij} &:= (1+a\|oldsymbol{y}_{i}-oldsymbol{y}_{j}\|_{2}^{2b})^{-1}. \end{aligned}$$

For the first equation, we have:

$$\frac{\partial c_{i,j}^{a}}{\partial \mathbf{y}_{i}} = \frac{\partial c_{i,j}^{a}}{\partial q_{ij}} \times \frac{\partial q_{ij}}{\partial \mathbf{y}_{i}} = \frac{-1}{q_{ij}} \times \left(\frac{-1}{(1+a\|\mathbf{y}_{i}-\mathbf{y}_{j}\|_{2}^{2b})^{2}} \times 2ab(\mathbf{y}_{i}-\mathbf{y}_{j}) \times \|\mathbf{y}_{i}-\mathbf{y}_{j}\|_{2}^{2(b-1)}\right)$$

$$\stackrel{(5)}{=} \frac{2ab\|\mathbf{y}_{i}-\mathbf{y}_{j}\|_{2}^{2(b-1)}}{(1+a\|\mathbf{y}_{i}-\mathbf{y}_{j}\|_{2}^{2b})} (\mathbf{y}_{i}-\mathbf{y}_{j}).$$
(14)

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The Training Algorithm of UMAP (2)

• For the second equation, we have:

$$\frac{\partial c_{i,j}^{r}}{\partial \mathbf{y}_{i}} = \frac{\partial c_{i,j}^{r}}{\partial q_{ij}} \times \frac{\partial q_{ij}}{\partial \mathbf{y}_{i}}$$

$$= \frac{1}{1 - q_{ij}} \times \left(\frac{-1}{(1 + a \|\mathbf{y}_{i} - \mathbf{y}_{j}\|_{2}^{2b})^{2}} \times 2ab(\mathbf{y}_{i} - \mathbf{y}_{j}) \times \|\mathbf{y}_{i} - \mathbf{y}_{j}\|_{2}^{2(b-1)} \right)$$

$$= \frac{-2ab\|\mathbf{y}_{i} - \mathbf{y}_{j}\|_{2}^{2(b-1)}}{(1 - q_{ij})(1 + a \|\mathbf{y}_{i} - \mathbf{y}_{j}\|_{2}^{2b})^{2}} (\mathbf{y}_{i} - \mathbf{y}_{j}). \tag{15}$$

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The Training Algorithm of UMAP(3)

- Eq. (5) was: $\mathbb{R} \ni q_{ij} := (1 + a \| \mathbf{y}_i \mathbf{y}_i \|_2^{2b})^{-1}$.
- The term in the numerator can be simplified as:

$$-2ab\|\mathbf{y}_{i} - \mathbf{y}_{j}\|_{2}^{2(b-1)} = -2b(a\|\mathbf{y}_{i} - \mathbf{y}_{j}\|_{2}^{2b})\|\mathbf{y}_{i} - \mathbf{y}_{j}\|_{2}^{-2}$$

$$\stackrel{(5)}{=} -2b(q_{ij}^{-1} - 1)\|\mathbf{y}_{i} - \mathbf{y}_{j}\|_{2}^{-2}.$$

• The term in the denominator can be simplified as:

$$(1-q_{ij})(1+a\|\boldsymbol{y}_i-\boldsymbol{y}_j\|_2^{2b})^2 \stackrel{(5)}{=} (1-q_{ij})q_{ij}^{-2} = q_{ij}^{-2}-q_{ij}^{-1} = q_{ij}^{-1}(q_{ij}^{-1}-1).$$

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The Training Algorithm of UMAP (4)

With these simplification, Eq. (15) becomes:

$$\frac{\partial c_{i,j}^r}{\partial \mathbf{y}_i} = \frac{-2ab\|\mathbf{y}_i - \mathbf{y}_j\|_2^{2(b-1)}}{(1 - q_{ij})(1 + a\|\mathbf{y}_i - \mathbf{y}_j\|_2^{2b})^2} (\mathbf{y}_i - \mathbf{y}_j),$$

Which can be simplified further as:

$$\frac{\partial c_{i,j}^{r}}{\partial \mathbf{y}_{i}} = \frac{-2b (q_{ij}^{-1} - 1) \|\mathbf{y}_{i} - \mathbf{y}_{j}\|_{2}^{-2}}{q_{ij}^{-1} (q_{ij}^{-1} - 1)} (\mathbf{y}_{i} - \mathbf{y}_{j}) = \frac{-2b}{\|\mathbf{y}_{i} - \mathbf{y}_{j}\|_{2}^{2} q_{ij}^{-1}} (\mathbf{y}_{i} - \mathbf{y}_{j})$$

$$\stackrel{(5)}{=} \frac{-2b}{\|\mathbf{y}_{i} - \mathbf{y}_{j}\|_{2}^{2} (1 + a \|\mathbf{y}_{i} - \mathbf{y}_{j}\|_{2}^{2b})} (\mathbf{y}_{i} - \mathbf{y}_{j}).$$

• If we add ε for stability to the squared distance in the denominator, the equation is obtained. Q.E.D.

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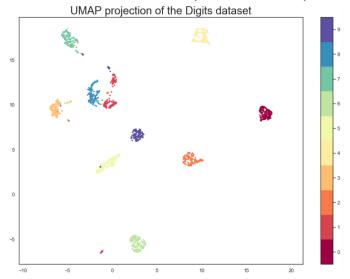
Example of UMAP Embedding (Digit Dataset)

5 9 9 0 8792 4625 26

Credit of image: https://umap-learn.readthedocs.io/en/latest/basic_usage.html

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Example of UMAP Embedding (Digit Dataset)



Credit of image: https://umap-learn.readthedocs.io/en/latest/basic_usage.html

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Acknowledgment

- Some slides are based on our tutorial paper: "Uniform Manifold approximation and projection (UMAP) and its variants: tutorial and survey" [7]
- For more information on UMAP, refer to our tutorial paper [7].
- UMAP library: https://umap-learn.readthedocs.io/en/latest/basic_usage.html

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