- 1. Classification
 - (a) K-Nearest Neighbor
 - (b) Adaboost
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 - (d) C4.5
 - (e) Naive Bayes
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- 2. Neural Networks
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 - (a) Gradient Descent
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Algorithm 1 k-Nearest Neighbor [Tay et al., 2014] link:36

Input: X: training data, Y:Class labels of X, x: unknown sample

Output: Class label of unknown sample

- 1: **function** CLASSIFY(X, Y, x)
- 2: **for** i = 1 to m **do**
- 3: Compute distance $d(X_i, x)$
- 4: end for
- 5: Compute set I containing indices for the k smallest distances $d(X_i, x)$
- 6: Return majority label $\{Y_i \text{ where } i \in I\}$
- 7: end function

Algorithm 2 Adaboost [Schapire, 2014]

Input:

Training data $\{(x_i, y_i)_{i=1}^N \text{ where } x_i \in \mathbb{R}^k \text{ and } y_i \in \{-1, 1\}\}$ Large number of classifiers denoted by $f_m(x) \in \{-1, 1\}$ 0-1 loss function I defined as

$$I(f_m(x,y)) = \begin{cases} 0, & \text{if } f_m(x_i) = y_i \\ 1, & \text{if } f_m(x_i) \neq y_i \end{cases}$$
 (1)

```
Output: The final classifier
```

```
1: for i = 1 to N do
               for i = 1 to M do
 2:
                       Fit weak classifier m to minimize the objective function:
 3:
                       where I(f_m(x_i)) \neq y_i and 0 otherwise \alpha_m = \ln \frac{\sum_{i=1}^N w_i^m I(f_m(x_i)) \neq y_i}{x^2 + 2x + 1} where I(f_m(x_i) \neq y_i) = 1 if f_m(x_i) \neq y_i and 0 otherwise \alpha_m = \ln \frac{1 - \epsilon_m}{\epsilon_m}
 4:
 5:
 6:
               end for
 7:
                \begin{aligned} & \mathbf{for} \ \text{all} \ i \ \ \mathbf{do} \\ & w_i^{m+1} = w_i^{(m)} e^{\alpha_{mI(f_m(x_i) \neq y_i)}} \end{aligned} 
 8:
 9:
               end for
10:
11: end for
```

Algorithm 3 Adaboost [Hertzmann et al., 2015]

Input:

```
Training data \{(x_i, y_i)_{i=1}^N \text{ where } x_i \in \mathbb{R}^k \text{ and } y_i \in \{-1, 1\}\}

Output: Weighted sum that represents the final output of the boosted classifier

1: Given Training data \{(x_i, y_i) \text{ where } y_i \in \{-1, 1\}\}

2: initialize D_1 = \text{uniform distribution on training examples}

3: for t = 1 to T do

4: Train weak classifier h_t on D_t

5: choose \alpha_t > 0

6: compute new distribution D_{t+1}:

7: for all i do

8: multiply D_t(x) by
```

$$\begin{cases}
e^{-\alpha_t}, & (<1) \text{ if } y_i = h_t(x_i) \\
e^{\alpha_t}, & (>1) \text{ if } y_i \neq h_t(x_i)
\end{cases} \tag{3}$$

```
9: renormalize

10: end for

11: output final classifier H_{final(x)} = sign(\sum \alpha_t h_t(x))

12: end for
```

Algorithm 4 Random forest [Bernstein, 2016] Link:39

```
Input: S: training set, F:Features and number of trees in forest B
Output: Constructed tree
 1: function RANDOMFOREST(S, F)
       H \leftarrow \emptyset
 2:
       for i \in 1, ....B do
 3:
           S^{(i)} \leftarrow A bootstrap sample from S
 4:
           h_i \leftarrow RANDOMIZEDTREELEARN(S^i, F)
 5:
 6:
           H \leftarrow H \bigcup \{h_i\}
       end for
 7:
       return H
 9: end function
10: function RANDOMIZEDTREELEARN(S, F)
       At each node:
11:
12:
        f \leftarrow a very small subset of F
13:
       Split on best feature in f
       return The learned tree
14:
15: end function
```

Algorithm 5 Iterative Dichotomiser 3 [., 2015] Link:40

```
Input: D: Training Data, X: Set of Input Attributes
Output: A decision tree
 1: function ID3(D, X)
       Let T be a new tree
       if all instances in D have the same class c then
 3:
           Label (T) = c; Return T
 4:
       end if
 5:
       if X = \emptyset or no attribute has positive information gain then
 6:
 7:
           Label (T) = most common class in D; Return T
       end if
 8:
 9:
       X \leftarrow attribute with highest information gain
       Label(T) = X
10:
       for each value x of X do
11:
           D_x \leftarrow \text{ instances in } D \text{ with } X = x
12:
           if D_x is empty then
13:
               Let T_x be a new tree
14:
               Label(T_x) = most common class in D
15:
           else
16:
               T_x = \mathrm{ID3}(D_x, X - \{x\})
17:
18:
       Add a branch from T to T_x labeled by x
19:
20: end for
21: return T
22nd function
```

```
Algorithm 6 Perceptron [Brownlee, 2015d] Link:65
```

```
Input: ProblemSize, InputPatterns, iterations_max, learn_rate

Output: Weights

1: for i = 1 to iterations_{max} do

2: Pattern_i \leftarrow SelectInputPattern(InputPatterns)

3: Activation_i \leftarrow ActivateNetwork(Pattern_i, Weights)

4: Output_i \leftarrow TransferActivation(Activation_i)

5: UpdateWeights(Pattern_i, Output_i, learn_{rate})

6: end for

7: Return Weights
```

```
Algorithm 7 Back-propagation [Brownlee, 2015a]
Input: ProblemSize, InputPatterns, iterations_{max}, learn_{rate}
Output: Network
 1: Network \leftarrow ConstructNetworkLayers()
 2: Network_w eights \leftarrow InitializeWeights(Network, Problem Size)
 3: for i = 1 to iterations_{max} do
       Pattern_i \leftarrow SelectInputPattern(InputPatterns)
       Output_i \leftarrow ForwardPropagate(Pattern_i, Network)
 6:
       BackwardPropagateError(Pattern_i, Output_i, Network)
       UpdateWeights(Pattern_i, Output_i, Network, learn_{rate})
 7:
 8: end for
 9: Return Network
```

Algorithm 8 Back-propagation1

```
Input:
       Training Set x^{(1)}, y^{(1)}, \dots, (x^{(m)}, y^{(m)})
Output:
       Gradient of the cost function
  1: \Delta_{ij}^{(l)} = 0(for all l,i,j)
  2: for i = 1 to m do
            Set a^{(1)} = x^{(i)}
            Perform forward propagation to compute a^{(l)} for l = 2, 3, ....L
  4:
            Using y^{(i)}, compute \delta^{(L)} = a^{(L)} - y^{(i)}
            Compute \delta^{(L-1)}, \delta^{(L-2)}, .... \delta^{(2)}
 7: \Delta_{ij}^{(l)} := \Delta_{ij}^{(l)} + a_j^{(l)} \delta_i^{(l+1)}
8: end for
8: end for
9: D_{ij}^{(l)} := \frac{1}{m} \Delta_{ij} + \lambda \theta_{ij}^{(l)} if j \neq 0
10: D_{ij}^{(l)} := \frac{1}{m} \Delta_{ij} if j = 0
```

```
Input: ProblemSize, InputPatterns, iterations_{max}, CodebookVectors_{num}, learn_{rate}
Output: CodebookVectors
 1: \ Codebook Vectors \leftarrow Initialize Codebook Vectors (Codebook Vectors_{num}, Problem Size)
 2: for i = 1 to iterations_{max} do
         Pattern_i \leftarrow SelectInputPattern(InputPatterns)
 3:
         Bmu_i \leftarrow SelectBestMatchingUnit(Pattern_i, CodebookVectors)
 4:
        for Bmu_i^{attribute} \in Bmu_i do
 5:
             if Bmu_i^{class} \equiv Pattern_i^{class} then
 6:
                 Bmu_i^{attribute} \leftarrow Bmu_i^{attribute} + learn_{rate} \times (Pattern_i^{attribute} - Bmu_i^{attribute})
 7:
 8:
                 Bmu_i^{attribute} \leftarrow Bmu_i^{attribute} - learn_{rate} \times (Pattern_i^{attribute} - Bmu_i^{attribute})
 9:
10:
        end if
```

Algorithm 9 Learning Vector Quantization [Brownlee, 2015c]

Return CodebookVectors

11: end for end for

```
Algorithm 10 Self Organizing Map [Brownlee, 2015b]
Input: InputPatterns, iterations_{max}, learn_{rate}, Grid_width, Grid_height
Output: CodebookVectors
 1: \ Codebook Vectors \leftarrow Initialize Codebook Vectors (Grid_{width}, Grid_{height}, Input Patterns)
 2: for i = 1 to iterations_{max} do
        Learn_{rate}^{i} \leftarrow CalculateLearningRate(i, learn_{rate}^{init})
 3:
        neighborhood_{size}^{i} \leftarrow CalculateNeighborhoodSize(i, neighborhood_{init}^{size})
 4:
        Pattern_i \leftarrow SelectInputPattern(InputPatterns)
 5:
        Bmu_i \leftarrow SelectBestMatchingUnit(Pattern_i, CodebookVectors)
 6:
        Neighborhood \leftarrow Bmu_i
 7:
        Neighborhood \leftarrow SelectNeighbors(Bmu_i, CodebookVectors, neighborhood_{size}^i)
 8:
        for Vector_i \in Neighborhood do
 9:
            for Vector_i^{attribute} \in Vector_i do
10:
                 Vector_i^{attribute} \leftarrow Vector_i^{attribute} + learn_{rate} \times (Pattern_i^{attribute} - Vector_i^{attribute})
11:
12:
            end for
        end for
13:
14: end for
15: Return CodebookVectors
```

```
Algorithm 11 Hierarchial Agglomerative Algorithm [Stein, 2016a]
Input:
     \langle V, E, w \rangle. Weighted graph
     d_c. Distance measure for two clusters
Output: \langle V_T, E_T \rangle. Cluster hierarchy or dendogram
 1: C = \{\{v \mid v \in V\}\}
                                                                                                                                  ▶ Initial Clustering
 2: V_t = \{v_C \mid C \in C\}, E_T = \emptyset
                                                                                                                                {\,\vartriangleright\,} \operatorname{Initial\,Dendogram}
 3: while |C| > 1 do
 4:
          update\_distance\_matrix(C, G, d_c)
          \{C, C'\} = \underset{\{C_i, C_j\} \in C: C_i \neq C_j}{argmin} d_c(C)
C = (C \setminus \{C, C'\}) \cup \{C \cup C'\}
 5:
                                              d_c(C_i,C_i)
                                                                                                                                               ▶ Merging
          V_T = V_T \cup \{v_{C,C'}\}, E_T = E_T \cup \{\{v_{C,C'}, v_C\}, \{v_{C,C'}, v_C\}\}\
                                                                                                                                          ▶ Dendogram
 8: end while
 9: Return T
```

```
Algorithm 12 Hierarchial Divisive Algorithm [Stein, 2016b]
Input:
     \langle V, E, w \rangle. Weighted graph
     d_c. Distance measure for two clusters
Output: \langle V_T, E_T \rangle. Cluster hierarchy or dendogram
                                                                                                                       ▶ Initial Clustering
 1: C = \{V\}
 2: V_t = \{v_C \mid C \in C\}, E_T = \emptyset
                                                                                                                     ▷ Initial Dendogram
 3: while \exists C_x : (C_x \in C \land |C| > 1) do
         update\_distance\_matrix(C, G, d_c)
 4:
                      \underset{\{C_i,C_j\}:C_i\cup C_j=C_x\wedge\ C_i\cap C_j=\emptyset}{argmax}
                                                        d_c(C_i, C_j)
 5:
         C = (C \backslash \{C, C'\}) \cup \{C \cup C'\}
                                                                                                                                  ▶ Merging
         V_T = V_T \cup \{v_{C,C'}\}, E_T = E_T \cup \{\{v_{C,C'}, v_C\}, \{v_{C,C'}, v_C\}\}\
                                                                                                                              ▶ Dendogram
 8: end while
 9: Return T
```

Algorithm 13 C4.5 [Dai and Ji, 2014]

```
Input:
    T: Training dataset
    S: Attributes
Output: decision tree Tree
 1: function C4.5(T)
 2:
        if T is NULL then
 3:
            return failure
        end if
 4:
        if S is NULL then
 5:
            return Tree as a single node with most frequent class label in T
 6:
 7:
        end if
        if |S| = 1 then
 8:
            return Tree as a single node S
 9:
        end if
10:
        set Tree = \{\}
11:
12:
        for a \in S do
            set Info(a,T) = 0 and SplitInfo(a,T) = 0
13:
            compute Entropy(a)
14:
            for v \in values(a, T) do
15:
               set T_{a,v} as the subset of T with attribute a=v
16:
               Info(a,T) + = \frac{|T_{a,v}|}{|T_a|} Entropy(a)
17:
               SplitInfo(a,T) + = -\frac{|T_{a,v}|}{|T_a|} \log \frac{|T_{a,v}|}{|T_a|}
18:
            end for
19:
            Gain(a, T) = Entropy(a) - Info(a, T)
20:
           GainRatio(a,T) = \frac{Gain(a,T)}{SplitInfo(a,T)}
21:
        end for
22:
        set a_{best} = argmax\{GainRatio(a, T)\}
23:
        a_{best}into Tree
24:
        for v \in values(a_{best}, T) do call C4.5(T_{a,v})
25:
        end for
26:
        return Tree
27:
28: end function
```

Algorithm 14 Gradient Descent

```
Input:
    f
    starting value x_1
    termination tolerances
Output: x_{maxIters}
 1: for i = 1 to maxIters do
 2:
        Compute the search direction d_t = -\delta f(x_t)
        if |d_T| < \epsilon_g then
 3:
            return "Converged to critical point", output x_t
 4:
            Find \alpha_t so that f(x_t + \alpha_t d_t) < f(x_t)
 5:
        end if
 6:
        if |\alpha_t d_T| < \epsilon_x then
 7:
            return "Converged in x", output x_t
 8:
            Find \alpha_t so that f(x_t + \alpha_t d_t) < f(x_t)
 9:
10:
        end if
        Let x_{t+1} = x_t + \alpha_t d_t
11:
12: end for
13: Return "Max number of iterations reached", output x_{maxIters}
```

```
Algorithm 15 Naive Bayes
```

```
Input:
    C: A  fixed set of classes
    D: Documents
Output: Category(Class) of the Documents
 1: function TrainMultinomialNB(C, D)
       V \leftarrow EXTRACTVOCABULARY(D)
       N \leftarrow COUNTDOCS(D)
 3:
       for each c \in C do
 4:
           N_c \leftarrow COUNTDOCSINCLASS(D, c)
 5:
 6:
           prior|c| \leftarrow N_c/N
           text_c \leftarrow CONCATENATETEXTOFALLDOCSINCLASS(D,C)
 7:
           for each t \in V do
 8:
              condprob|t||c| \leftarrow \frac{T_{ct}+1}{\sum_{t'}(T_{ct'+1})}
 9:
           end for
10:
       end for
11:
       return V, prior, condprob
12:
13: end function
14: function APPLYMULTINOMIALNB(C, D, prior, condprob, d)
       W \leftarrow EXTRACTTOKENSFROMDOC(V, d)
15:
       for each c \in C do
16:
           score|c| \leftarrow \log \ prior|c|
17:
           for each t \in W do
18:
              score|c| + = \log condprob|t||c|
19:
20:
           end for
       end for
21:
       return arg\ max_{c \in C} score|c|
22:
23: end function
```

Algorithm 16 Lasso Regression

```
Input:
     ipy: Inner product vector, ipy_i = \langle y, X_{i} \rangle
     ipx: Inner product matrix, ipx_{ij} = \langle X_{\cdot i}, X_{\cdot j} \rangle
     \lambda: Penalty parameter
     N: Number of samples
Output: beta: Regression parameter vector
 1: function FastLasso(ipy, ipx, \lambda, N)
                                                                                                   ▶ Threshold for stopping iteration
 2:
         stop_thr
 3:
         p \leftarrow length(ipy)
 4:
         beta \leftarrow 0 with length p
         gc \leftarrow 0 with length p
 5:
         while difBeta_{max} \geq \text{stop\_thr do}
 6:
              difBeta_{max} \leftarrow 0
 7:
              for j = 1 \leftarrow p do
 8:
                  z \leftarrow (ipy|j| - gc|j|)/N + beta|j|
 9:
                  \texttt{beta\_tmp} \leftarrow \max(0, z - \lambda) - \max(0, -z - \lambda)
10:
                  difBeta \leftarrow \mathtt{beta\_tmp} - beta|j|
11:
                  difabs \leftarrow abs(difBeta)
12:
                  if difabs > 0 then
13:
                      beta|j| \leftarrow \mathtt{beta\_tmp}
15:
                      gc \leftarrow gc + ipx|j| \times difBeta
                      difBeta_{max} = max(difBeta_{max}, difabs)
16:
                  end if
17:
              end for
         end while
20: end function
```

Algorithm 17 Bagging

Input:

B: the number of bags or base hypotheses

L: Base Learning Algorithm

Output: New Training Sets

- 1: function Bagging(examples, B, L)
- 2: **for** i = 1 to B **do**
- 3: $examples_i \leftarrow a \text{ bootstrap sample of } examples$
- 4: end for
- 5: Compute set I containing indices for the k smallest distances $d(X_i, x)$
- 6: $h_i \leftarrow \text{apply}Ltoexamples_i$
- 7: Return $h_1, h_2, ...h_B$
- 8: end function

Algorithm 18 Deep Q-Learning with Experience Replay [Mnih et al., 2013]

Input:

D: data set

Q: Action-Value Function

Output: New Training Sets

- 1: for i = 1 to M do
- 2: Initialise sequence $s_1 = \{x_1\}$ and preprocessed sequenced $\phi = \phi(s_1)$
- 3: **for** i = 1 to T **do**
- 4: With probability ϵ select a random action a_t otherwise select $a_t = max_aQ * (\phi(s_t).a:\theta)$
- 5: Execute action a_t in emulator and observe reward r and image x_{t+1}
- 6: Set $s_{t+1} = s_t, a_t, x_{t+1}$ and preprocess $\phi_{t+1} = \phi(s_{t+1})$
- 7: Store transition $(\phi_t, a_t, r_t, \phi_{t+1})$ in D
- 8: Set $y_j =$

$$\begin{cases}
r_j, & \text{for terminal } \phi_{j+1} \\
r_j + \gamma \max_{a'} Q(\phi_{j+1,a';\theta}), & \text{for terminal } \phi_{j+1}
\end{cases}$$
(5)

9: Perform a gradient descent step on $(y_j - Q(\phi_j, a_j; \theta))^2$ according to the following equation

10:

$$\Delta_{\theta}L_{i}(\theta_{i}) = \mathbb{E}_{s,a \sim \rho(.);s' \sim \epsilon[(r + \gamma max_{a'}Q(s',a';\theta_{i-1}) - Q(s,a;\theta_{i}))\Delta_{\theta_{i}}Q(s,a;\theta_{i})]}$$

11: end for

12: end for

Algorithm 19 PageRank

```
Input:
    G: inlink file
    iteration: Number of iteration
Output: PageRank
 1: function PageRank(G, iteration)
        d \leftarrow 0.85
                                                                                                          ⊳ damping factor: 0.85
 3:
        oh \leftarrow G
                                                                                                      ⊳ get outlink hash from G
        ih \leftarrow G
                                                                                                       4:
        N \leftarrow G
 5:
                                                                                                 ⊳ get number of pages from G
        for all p in the graph do
 6:
             opg[p] \leftarrow \tfrac{1}{N}
 7:
        end for
 8:
        while iteration > 0 do
 9:
             dp \leftarrow 0
10:
11:
             for all p that has no out-links do
                 dp \leftarrow dp + d * \frac{opg[p]}{N}
12:
             end for
13:
             for all p in the graph do
14:
                 npg[p] \leftarrow dp + \frac{[1-d]}{N}
                 for all ip in ih[p] do
16:
                     npg[p] \leftarrow dp + \frac{d*opg[ip]}{oh[ip]}
17:
                 end for
18:
19:
             end for
20:
             opg \leftarrow npg
             iteration \leftarrow iteration - 1
21:
        end while
22:
23: end function
```

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Algorithm 20 DBSCAN link:42

```
Input:
   D: Data
    \epsilon:Threshold distance
    MinPts: Minimum number of points required to form a cluster
Output: Clustered Data
 1: function DBSCAN(D, \epsilon, minPts)
       C = 0
 2:
       for each point Pin datasetD do
 3:
          if P is visited then
 4:
              continue next point
 5:
          end if
 6:
          mark P as visited
 7:
          NeighborPts = regionQuery(P, \epsilon)
 8:
          if \ size of(Neighbor Pts) < Min Pts \ then
 9:
10:
              mark P as NOISE
          else
11:
              C = \text{next cluster}
12:
              expandCluster(P, NeighborPts, C, \epsilon, MinPts)
13:
       end if
15: end for
and function
function EXPANDCLUSTER(P, NeighborPts, C, \epsilon, MinPts)
18: add P to Cluster C
19: for each point P'inNeighborPts do
       if P' is not visited then
20:
          mark P' as visited
21:
          NeighborPts' = regionQuery(P', eps)
22:
          if sizeof(NeighborPts) >= MinPts then
23:
              NeighborPts = NeighborPts joined with NeighborPts'
24:
          end if
25:
       end if
26:
       if P' is not yet member of any cluster then
27:
          add P' to cluster C
28:
       end if
29:
30: end for
and function
function REGIONQUERY(P, \epsilon)
33: return all points within P's \in neighborhood
end function
```

Algorithm 21 Logistic Regression

Input:

Training data of the form $\{(x_1, 1), (x_2, 0), ...\}$

x: unknown sample

Output: The output is a probability that the given input point belongs to a certain class

- 1: $0 \leftarrow \beta$
- 2: Compute y by setting its elements to

$$y = \begin{cases} 1, & \text{if } g_i = 1\\ 0, & \text{if } g_i = 2 \end{cases}$$
 (7)

i = 1,2,...N

3: Compute p by setting its elements to

$$p(x_i, \beta) = \frac{e^{\beta^T x_i}}{1 + e^{\beta^T x_i}}$$

i = 1,2,...N

- 4: Compute the diagonal matrix W. The ith diagonal element is $p(x_i, \beta)(1 p(x_i; \beta))$
- 5: $z \leftarrow X\beta + W^{-1}(y-p)$ 6: $\beta \leftarrow (X^TWX)^{-1}X^TWz$
- 7: If the stopping criteria, stop; otherwise go back to step 3

Algorithm 22 Gaussian Process

Input:

$$X = \begin{bmatrix} x_1^T \\ \dots \\ x_n^T \end{bmatrix} \in \mathbb{R}^{n \times D}, \text{m training inputs}$$

$$y = \begin{bmatrix} y_1^T \\ \dots \\ y_n^T \end{bmatrix} \in \mathbb{R}^n$$

$$k(\dots) : \mathbb{R}^{D \times D}$$

 x_* test input

 σ^2 noise level on the observations

$$[y(x) = f(x) + \epsilon, \epsilon \sim N(0, \sigma^2)]$$

Output:

$$f_* \\ cov(f_*)$$

 $cov(f_*)$ 1: $K \in \mathbb{R}^{n \times n}$ Gram matrix. $K_{ij} = k(x_i, x_j)$

$$k(x_*) = k_* = k(X, x_*) = \begin{bmatrix} k(x_1, x_*) \\ \dots \\ k(x_n, x_*) \end{bmatrix} \in \mathbb{R}$$

2:
$$\alpha = (K + \sigma_2 \mathbb{I}_n)^{-1} y$$

3:
$$f_* = k_*^T \alpha \in \mathbb{R}$$

4: $cov(f_*) = k(x_*, x_*) - k_*^T [K + \sigma_2 \mathbb{I} n]^{-1} k_*$