

Introduction to computational models

Lab assignment 3. Radial basis function neural networks

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Module “Introduction to computational models”
4th year of “Grado en Ingeniería Informática”
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- 1 Contents
- 2 Introduction
- 3 Architecture of an RBF neural network
- 4 Training of an RBF network
 - Phase 1: clustering
 - Phase 2: adjustment of the radii
 - Phase 3: output layer weights



Objectives of the lab assignment

- To familiarise the student with the concept of the neuronal network of radially based functions (RBF).
- To implement such a network.
- To familiarise the learner with the use of [scikit-learn](#) as an environment for the creation of automatic learning models.



Radial basis function neural network

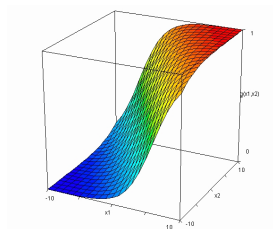
- Radial Basis Function Neural Networks (**RBFNNs**): they are based on a local approach.
 - The hidden layer neurons are radial basis functions (RBFs): **local functions**.
 - Opposite to MLP networks, where the hidden layer neurons are of the sigmoid type: **projection functions**.
- **Projection functions**: high value, not zero, over a wide region of the input space.
- **Local functions**: high value, not zero, only on a localised region of the input space.



Local Vs global functions

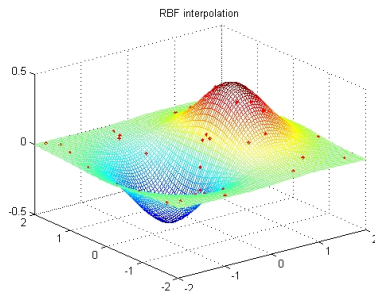
Sigmoidal units

- Additive projection model.

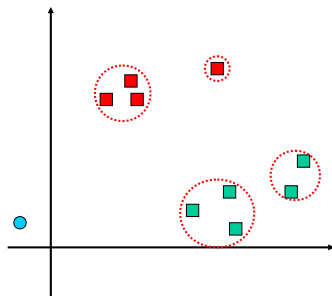
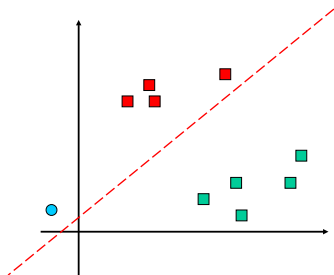


Radial basis function

- Local model.



Local Vs global functions



RBFs

- A function is said to be RBF if its output depends on the distance between the input vector and a vector stored in it (**centre**).
- Each RBF stores a **centre** as a reference, and every time a new pattern is presented to it, the distance to this centre is calculated.
 - If the distance is small, the RBF **is activated** (its output is 1).
 - If the distance is high, the RBF **is not activated** (its output is 0).
- What do we consider as large or small? \Rightarrow For this, we incorporate an additional element: the **radius**.
 - If the radius is small, activation will only occur when the pattern is very close to the centre.
 - If the radius is large, the activation will take place at a greater distance.



RBFs

- There are many functions that fulfil these properties: Cauchy RBF, inverse multi-quadratic...
- But the most common is the Gaussian function:

$$\varphi(\mathbf{x}, \mathbf{c}, \sigma) = \exp\left(-\frac{\|\mathbf{x} - \mathbf{c}\|^2}{2(\sigma)^2}\right)$$

where \mathbf{c} is the centre of the RBF, σ is the radius and \mathbf{x} is the pattern we are evaluating.

- $\|\mathbf{x} - \mathbf{c}\|$ is the norm of the difference vector between the centre and the pattern, or what is the same as the Euclidean distance:

$$\|\mathbf{x} - \mathbf{c}\| = \sqrt{\sum_{i=1}^n (x_i - c_i)^2}$$



RBFs

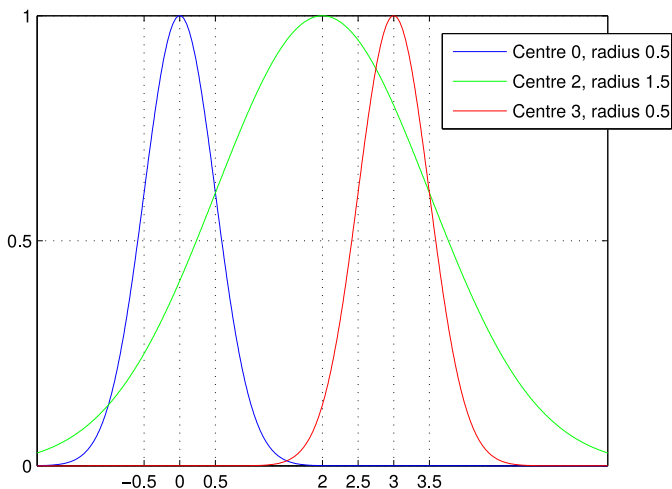
- By linking the two expressions and simplifying:

$$\varphi(\mathbf{x}, \mathbf{c}, \sigma) = \exp \left(- \frac{\left(\sqrt{\sum_{i=1}^n (x_i - c_i)^2} \right)^2}{2(\sigma)^2} \right)$$

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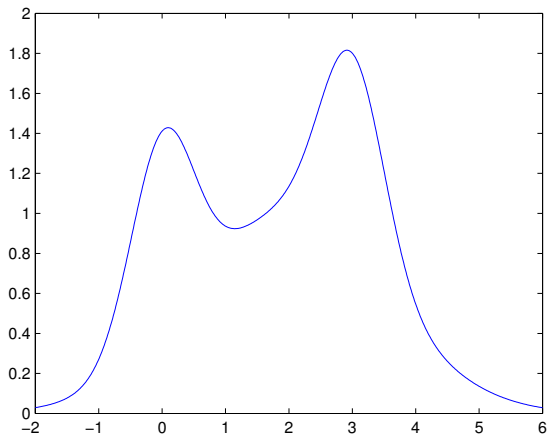


RBFs: effect of the centre and the radius



RBF neural network

Sum of the previous two RBFs:

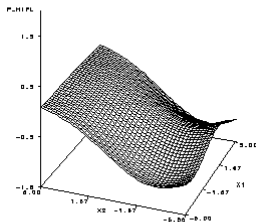


Example in 2D:

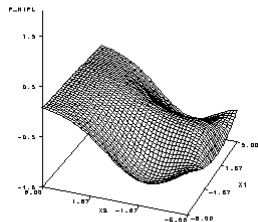
Neural Network: Ordinary RBF Network with Equal Widths and Heights

2 Hidden Units: 8 Parameters

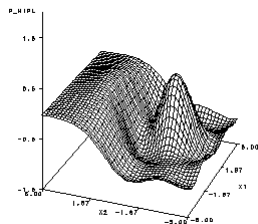
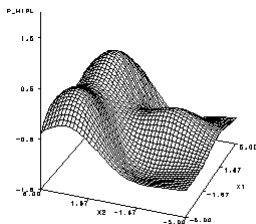
3 Hidden Units: 11 Parameters



5 Hidden Units: 17 Parameters



9 Hidden Units: 29 Parameters



Architecture of an RBFNN

- Three layers:
 - Input layer.
 - Hidden layer (**RBF** functions).
 - Output layer:
 - **Regression**: linear function (weighted sum of the RBF outputs).
 - **Classification**: *softmax* function.



Architecture of an RBFNN

- Following the notation we used for the MLP:

- RBF neurons:

- $net_j^h = \sum_{i=1}^{n_{h-1}} (w_{ji}^h - out_i^{h-1})^2$

- $out_j^h = \exp\left(-\frac{net_j^h}{2 (w_{j0}^h)^2}\right)$

- Linear type neurons:

- $out_j^H = net_j^H = w_{j0}^H + \sum_{i=1}^{n_{H-1}} w_{ji}^H out_i^{H-1}$

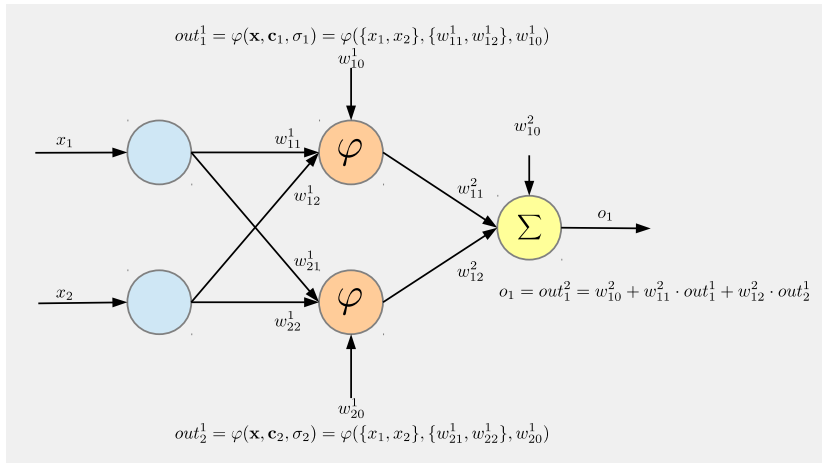
- *softmax* neurons:

- $net_j^H = w_{j0}^H + \sum_{i=1}^{n_{H-1}} w_{ji}^H out_i^{H-1}$

- $out_j^H = \frac{\exp(net_j^H)}{\sum_{i=1}^{n_H} \exp(net_i^H)}$



Architecture of an RBFNN (regression)



Training of an RBFNN

- How do we adjust the parameters?
 - RBF functions are derivable → Apply the error backpropagation algorithm (**fully supervised** training).
 - Derivatives would have to be calculated with respect to radius and centres.
 - They are complex and have a higher computational cost than for the MLP.
 - **Hybrid** training: **unsupervised** part (*clustering*) and **supervised** part (logistic regression or matrix inversion).
 - The local properties of the RBF networks are better exploited.
 - The computational cost is generally lower than using the gradient descent algorithm.



Training of an RBFNN

- We have to obtain three things:

- ① Coordinates of the RBF centres: $\mathbf{c}_1, \mathbf{c}_2, \dots, \mathbf{c}_{n_1}$.

- Weights from input layer to hidden layer:

$$\mathbf{c}_1 = \{w_{11}^1, w_{12}^1, \dots, w_{1n}^1\}$$

$$\mathbf{c}_2 = \{w_{21}^1, w_{22}^1, \dots, w_{2n}^1\}$$

...

$$\mathbf{c}_{n_1} = \{w_{n_1 1}^1, w_{n_1 2}^1, \dots, w_{n_1 n}^1\}.$$

- ② Width of the RBF: $\sigma_1, \sigma_2, \dots, \sigma_{n_1}$.

- We will use the bias position:

$$\sigma_1 = w_{10}^1, \sigma_2 = w_{20}^1, \dots, \sigma_{n_1} = w_{n_1 0}^1$$

- ③ Weights from hidden layer to output layer (with bias):

$$w_{10}^2, w_{11}^2, w_{12}^2, \dots, w_{1n_1}^2$$

$$w_{20}^2, w_{21}^2, w_{22}^2, \dots, w_{2n_1}^2$$

...

$$w_{k0}^2, w_{k1}^2, w_{k2}^2, \dots, w_{kn_1}^2$$

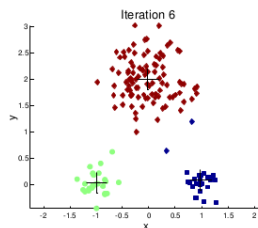
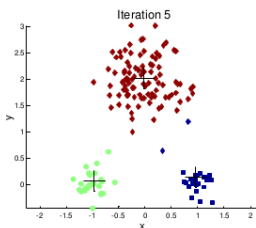
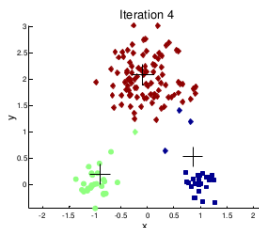
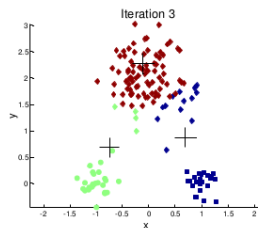
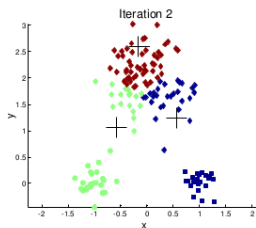
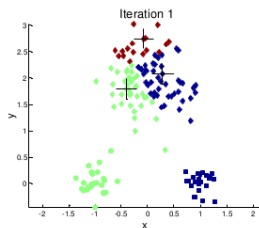


Training of an RBFNN: phase 1 (clustering)

- The adjustment of the centres in the network can be done by a *clustering* procedure.
- The idea is to detect the group of patterns (or *clusters*) in the input space and locate one RBF in each *cluster*.
- We are going to use the most popular clustering method, the *K*-means.
- The coordinates of the centroids of each *cluster* will be the centres of the RBFs.



Training of an RBFNN: phase 1 (clustering)



Training of an RBFNN: phase 1 (clustering)

- *K*-means: partitional clustering method.
 - The number of *clusters* must be specified, in our case, it will be the number of hidden neurons of our RBFNN (n_1).
 - Each *cluster* has a centroid (geometric mean of the *cluster*).
 - The points are assigned to the *cluster* with the closest centroid (using any distance metric).
 - Iteratively, we update the centroids as a function of the assignments of the points to the *clusters*, until the centroids no longer change.
 - The results depend on the **initialization of the centroids**:
 - For **classification**, we select randomly, and in a **stratified manner**, n_1 patterns.
 - For **regression**, we randomly select n_1 patterns.



Training of an RBFNN: phase 2 (adjustment of the radii)

- More complex procedures (density estimation) can be used to fine tune the radii of the RBFs.
- However, we will adjust the radii in a very simple way, we will take half (because it is a radius and not a diameter) of the average distance to the rest of centroids.
- That is, the radius of the neuron j will be:

$$\sigma_j = w_{j0}^1 = \frac{1}{2 \cdot (n_1 - 1)} \sum_{i \neq j} \|c_j - c_i\| = \quad (1)$$

$$= \frac{1}{2 \cdot (n_1 - 1)} \sum_{i \neq j} \sqrt{\sum_{d=1}^n (c_{jd} - c_{id})^2} \quad (2)$$



Training of an RBFNN: phase 3 (output layer weights, case 1 classification)

- We will adjust the weights of the output layer in two ways, depending on whether we are facing a classification or regression problem.
 - For **classification**, the weights will be adjusted using **logistic regression**.
 - For **regression**, the weights will be adjusted using the **pseudo-inverse**.



Training of an RBFNN: phase 3 (output layer weights, case 1 classification)

- In both cases, we will need the RBF output matrix, which we will call \mathbf{R} :

$$\mathbf{R} = \begin{pmatrix} out_1^1(\mathbf{x}_1) & out_2^1(\mathbf{x}_1) & \dots & out_{n_1}^1(\mathbf{x}_1) & 1 \\ out_1^1(\mathbf{x}_2) & out_2^1(\mathbf{x}_2) & \dots & out_{n_1}^1(\mathbf{x}_2) & 1 \\ \dots & \dots & \dots & \dots & \dots \\ out_1^1(\mathbf{x}_N) & out_2^1(\mathbf{x}_N) & \dots & out_{n_1}^1(\mathbf{x}_N) & 1 \end{pmatrix} \quad (3)$$

where $out_j^1(\mathbf{x}_i)$ is the output of the j -th RBF neuron when it is feed using the training pattern \mathbf{x}_i . To simulate the bias we have included a constant column equal to 1.



Training of an RBFNN: phase 3 (output layer weights, case 1 classification)

- Once the matrix has been constructed, in the case of classification, we will apply **logistic regression**.
 - We will use a function to which we will pass the **R** matrix as if it were the input matrix of my database.
 - Logistic regression is a linear classification model, which approximates the probability of belonging to a class in the following way (*softmax* function):

$$P(\mathbf{x} \in C_j) = o_j = \frac{\exp(\beta_{j0} + \sum_{i=1}^n \beta_{ji} x_i)}{\sum_{l=1}^k \exp(\beta_{l0} + \sum_{i=1}^n \beta_{li} x_i)} \quad (4)$$

- The objective of the logistic regression is to obtain the values of β_{ji} that maximize the cross entropy:

$$L = \frac{1}{N} \sum_{p=1}^N \left(\frac{1}{k} \sum_{o=1}^k d_j \ln(o_j) \right) \quad (5)$$



Training of an RBFNN: phase 3 (output layer weights, case 1 classification)

- ...we will apply **regresión logística**.
 - Logistic regression can include regularization, which is a mechanism for making the maximum number of parameters β_{ji} tend to zero (or almost zero).

- L2 regularization:

$$L = \left(\frac{1}{N} \sum_{p=1}^N \left(\frac{1}{k} \sum_{o=1}^k d_j \ln(o_j) \right) \right) - \eta \left(\sum_{j=1}^k \sum_{i=0}^n \beta_{ji}^2 \right) \quad (6)$$

- L1 regularization:

$$L = \left(\frac{1}{N} \sum_{p=1}^N \left(\frac{1}{k} \sum_{o=1}^k d_j \ln(o_j) \right) \right) - \eta \left(\sum_{j=1}^k \sum_{i=0}^n |\beta_{ji}| \right) \quad (7)$$

- The η parameter must be set by the user and establishes the importance given to regularization.



Training of an RBFNN: phase 3 (output layer weights, case 1 classification)

- ...we will apply **regresión logística**.
 - Regularisation provides simpler models that are less prone to over-fitting.
 - Difference between L2 and L1:
 - The L2 regularisation tends to provide smaller weights (although not necessarily equal to zero).
 - L1 regularization tends to prune more variables, making many weights equal to zero (although non-zero weights are not necessarily small in absolute value).



Training of an RBFNN: phase 3 (output layer weights, case 2 regression)

- Once the matrix has been constructed, in the case of regression, we will apply the **pseudo-inverse**.
 - From the point of view of linear algebra, the output of the network can be written as:

$$\mathbf{R}_{(N \times (n_1+1))} \times \boldsymbol{\beta}_{((n_1+1) \times k)}^T = \hat{\mathbf{Y}}_{(N \times k)} \quad (8)$$

where \mathbf{R} is the matrix containing the outputs of the RBF neurons, $\boldsymbol{\beta}$ is a matrix containing a vector of parameters for each output to be predicted and $\hat{\mathbf{Y}}$ is a matrix with all the estimated outputs.

$$\begin{pmatrix} out_{11}^1 & \dots & out_{n_1 1}^1 & 1 \\ out_{12}^1 & \dots & out_{n_1 2}^1 & 1 \\ \dots & \dots & \dots & \dots \\ out_{1N}^1 & \dots & out_{n_1 N}^1 & 1 \end{pmatrix} \begin{pmatrix} \beta_{11} & \beta_{21} & \dots & \beta_{k1} \\ \dots & \dots & \dots & \dots \\ \beta_{1n_1} & \beta_{2n_1} & \dots & \beta_{kn_1} \\ \beta_{10} & \beta_{20} & \dots & \beta_{k0} \end{pmatrix} \quad (9)$$



Training of an RBFNN: phase 3 (output layer weights, case 2 regression)

- ...we will apply the **pseudo-inverse**.
 - If we want to obtain the best possible values for the parameters, we use the following equation:

$$\mathbf{R}_{(N \times (n_1+1))} \times \boldsymbol{\beta}_{((n_1+1) \times k)}^T = \mathbf{Y}_{(N \times k)} \quad (10)$$

where \mathbf{Y} is the matrix with desired outputs:

$$\mathbf{Y} = \begin{pmatrix} d_{11} & \dots & d_{1k} \\ d_{21} & \dots & d_{2k} \\ \dots & \dots & \dots \\ d_{N1} & \dots & d_{Nk} \end{pmatrix} \quad (11)$$



Training of an RBFNN: phase 3 (output layer weights, case 2 regression)

- ...we will apply the **pseudo-inverse**.
 - If \mathbf{R} is square ($N = (n_1 + 1)$), then \mathbf{R} has an inverse and we can clear it directly:

$$\beta_{((n_1+1) \times k)}^T = (\mathbf{R}_{(N \times N)})^{-1} \mathbf{Y}_{(N \times k)} \quad (12)$$

- If $(n_1 + 1) > N$, then there are many solutions, and some kind of feature selection algorithm must be used to reduce the value of n_1 .
- If $(n_1 + 1) < N$ (most common case), there is one single solution, but, given that \mathbf{R} is not square, we have to use the Moore Penrose pseudo-inverse.



Training of an RBFNN: phase 3 (output layer weights, case 2 regression)

- ...we will apply the **pseudo-inverse**.
 - Moore Penrose pseudo-inverse:

$$\beta_{((n_1+1) \times k)}^T = (\mathbf{R}^+)_{((n_1+1) \times N)} \mathbf{Y}_{(N \times k)} \quad (13)$$

$$(\mathbf{R}^+)_{((n_1+1) \times N)} = \left(\mathbf{R}_{((n_1+1) \times N)}^T \times \mathbf{R}_{(N \times (n_1+1))} \right)^{-1} \mathbf{R}_{((n_1+1) \times N)}^T \quad (14)$$

- We will use a matrix library to do these operations and obtain $\beta_{((n_1+1) \times k)}^T$.



RBF training algorithm *off-line*

Start

- ① $\text{initialCentroids} \leftarrow$ randomly select n_1 patterns (**regression**) or n_1 patterns in a stratified way (**classification**).
- ② $\text{centroids} \leftarrow \text{K-means}(\mathbf{X}, n_1, \text{initialCentroids})$ // \mathbf{X} is the input matrix with the inputs of all patterns
- ③ $\sigma_j \leftarrow$ (average of the distance from j to the rest of centroids)/2.
- ④ Construct the $\mathbf{R}_{(N \times (n_1 + 1))}$ matrix, where $\mathbf{R}_{ij} = \text{out}_j^1(\mathbf{x}_i)$ for $j \neq (n_1 + 1)$, and $\mathbf{R}_{ij} = 1$ for $j = (n_1 + 1)$.
- ⑤ **If classification**
 - ① $\text{outputWeights} \leftarrow \text{applyLogisticRegression}(\mathbf{R}, \text{eta})$
- ⑥ **Si regresión**
 - ① $\text{outputWeights} \leftarrow \text{calculatePseudoinverse}(\mathbf{R})$

End



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