

# 1 The Radial Basis Function Kolmogorov-Arnold Network (RBF-KAN)

The RBF-KAN is a novel approach to function approximation, combining the theoretical foundations of the Kolmogorov-Arnold Theorem with the powerful approximation capabilities of radial basis functions (RBFs).

## 1.1 Kolmogorov-Arnold Theorem

The Kolmogorov-Arnold Theorem, also known as the Superposition Theorem or the Kolmogorov-Arnold Representation Theorem, states that any continuous multivariate function  $f : [0, 1]^{d_{\text{in}}} \rightarrow \mathbb{R}^{d_{\text{out}}}$  on the  $d_{\text{in}}$ -dimensional unit hypercube  $[0, 1]^{d_{\text{in}}}$  can be represented as a superposition (composition) of a limited number of one-variable (univariate) functions  $g_q$  and  $\psi_{p,q}$  such that:

$$f(\mathbf{x}) = \sum_{q=0}^{2d_{\text{in}}} g_q \left( \sum_{p=1}^{d_{\text{in}}} \psi_{p,q}(x_p) \right) \quad (1)$$

where  $\mathbf{x} = (x_1, x_2, \dots, x_{d_{\text{in}}})$ .

## 1.2 Radial Basis Functions (RBFs)

Radial basis functions (RBFs) are a class of functions that depend only on the distance (or radius) from a center point. A common choice for the RBF is the Gaussian function:

$$\phi(r) = e^{-(\beta r)^2} \quad (2)$$

where  $r$  is the Euclidean distance between the input sample and the RBF center, and  $\beta$  is a scaling factor.

## 1.3 The RBF-KAN Layer

In the RBF-KAN, the target multivariate function  $f(\mathbf{x})$  is approximated using a single layer of RBF interpolation:

$$\tilde{f}(\mathbf{x}) = \sum_{j=1}^{d_{\text{in}}} \sum_{k=1}^M \Theta_{j,k} \phi(r_{j,k}) \quad (3)$$

where  $\mathbf{x}$  is the normalized input tensor,  $M$  is the number of RBF centers,  $\phi(r_{j,k})$  is the RBF value computed using the distance between the  $j$ -th input dimension and the  $k$ -th RBF center, and  $\Theta \in \mathbb{R}^{d_{\text{in}} \times M}$  are the learnable coefficients for the RBF interpolation.

### 1.3.1 Input Normalization

Let  $\mathbf{X} \in \mathbb{R}^{N \times d_{\text{in}}}$  be the input tensor, where  $N$  is the batch size and  $d_{\text{in}}$  is the input dimension. The input tensor is normalized to the range  $[0, 1]$  using min-max normalization:

$$\mathbf{X}_{\text{norm}} = \frac{\mathbf{X} - \mathbf{X}_{\min}}{\mathbf{X}_{\max} - \mathbf{X}_{\min}} \quad (4)$$

where  $\mathbf{X}_{\min}$  and  $\mathbf{X}_{\max}$  are the minimum and maximum values of the input tensor, respectively.

### 1.3.2 RBF Computation

Define a set of  $M$  RBF centers  $\mathcal{C} = \{\mathbf{c}_1, \mathbf{c}_2, \dots, \mathbf{c}_M\}$ , where  $\mathbf{c}_j \in \mathbb{R}^{d_{\text{in}}}$  for  $j = 1, 2, \dots, M$ . These centers can be initialized randomly or using techniques like k-means clustering. Compute the RBF values for each input sample and RBF center using the Gaussian RBF:

$$\mathbf{R}_{i,j} = \phi(|\mathbf{X}_{\text{norm},i} - \mathbf{c}_j|) = e^{-\beta^2 |\mathbf{X}_{\text{norm},i} - \mathbf{c}_j|^2} \quad (5)$$

where  $\mathbf{R} \in \mathbb{R}^{N \times M}$  is the tensor containing the RBF values,  $\mathbf{X}_{\text{norm},i}$  is the  $i$ -th row of the normalized input tensor  $\mathbf{X}_{\text{norm}}$ , and  $|\cdot|$  denotes the Euclidean norm.

### 1.3.3 RBF Interpolation

Define a set of learnable weights  $\mathbf{W} \in \mathbb{R}^{M \times d_{\text{out}}}$ , where  $d_{\text{out}}$  is the output dimension. Compute the RBF interpolation as a weighted sum of the RBF values:

$$\mathbf{Y} = \mathbf{R}\mathbf{W} \quad (6)$$

where  $\mathbf{Y} \in \mathbb{R}^{N \times d_{\text{out}}}$  is the output tensor representing the approximation of the target function.

## 1.4 Training and Optimization

Train the RBF-KAN by optimizing the learnable weights  $\mathbf{W}$  and the RBF centers  $\mathcal{C}$  to minimize a loss function  $\mathcal{L}$  between the predicted output  $\mathbf{Y}$  and the true output  $\mathbf{Y}_{\text{true}}$ :

$$\mathcal{L}(\mathbf{W}, \mathcal{C}) = \frac{1}{N} \sum_{i=1}^N \ell(\mathbf{Y}_i, \mathbf{Y}_{\text{true},i}) \quad (7)$$

where  $\ell$  is a suitable loss function (e.g., mean squared error for regression tasks, cross-entropy for classification tasks). Compute the gradients of the loss with respect to the learnable weights  $\mathbf{W}$  and the RBF centers  $\mathcal{C}$  using backpropagation:

$$\frac{\partial \mathcal{L}}{\partial \mathbf{W}} = \frac{1}{N} \sum_{i=1}^N \frac{\partial \ell(\mathbf{Y}_i, \mathbf{Y}_{\text{true},i})}{\partial \mathbf{Y}_i} \cdot \mathbf{R}_i^\top \quad (8)$$

$$\frac{\partial \mathcal{L}}{\partial \mathcal{C}} = \frac{1}{N} \sum_{i=1}^N \frac{\partial \ell(\mathbf{Y}_i, \mathbf{Y}_{\text{true},i})}{\partial \mathbf{Y}_i} \cdot \mathbf{W}^\top \cdot \frac{\partial \mathbf{R}_i^\top}{\partial \mathcal{C}} \quad (9)$$

where  $\top$  denotes the transpose operation, and  $\frac{\partial \mathbf{R}_i^\top}{\partial \mathcal{C}}$  is the gradient of the RBF values with respect to the RBF centers, which can be computed using the chain rule and the derivative of the Gaussian RBF.