

# Distributed and Secure Kernel-Based Quantum Machine Learning

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## Abstract

Quantum computing promises to revolutionize machine learning, offering significant efficiency gains in tasks such as clustering and distance estimation. Additionally, it provides enhanced security through fundamental principles like the measurement postulate and the no-cloning theorem, enabling secure protocols such as quantum teleportation and quantum key distribution. While advancements in secure quantum machine learning are notable, the development of secure and distributed quantum analogues of kernel-based machine learning techniques remains underexplored.

In this work, we present a novel approach for securely computing common kernels, including polynomial, radial basis function (RBF), and Laplacian kernels, when data is distributed, using quantum feature maps. Our methodology introduces a robust framework that leverages quantum teleportation to ensure secure and distributed kernel learning. The proposed architecture is validated using IBM's Qiskit Aer Simulator on various public datasets.

## 1 Introduction

Quantum computing is set to revolutionize machine learning (ML), by leveraging its capability to encode high dimensional data into quantum bits, or qubits. These qubits exist in a superposition of states, enabling quantum data to represent data exponentially more efficiently than classical computing; data represented using  $N$  classical bits can equivalently be represented by  $\log_2 N$  qubits. Although practical quantum computers are still in their infancy, various quantum machine learning (QML) techniques have been and are being proposed.

Notably, quantum computing has exhibited a substantial efficiency gain in some computational tasks as compared to classical computing (Schuld & Petruccione, 2018): estimating distances and inner products between post-processed  $N$ -dimensional vectors is achieved in  $\mathcal{O}(\log(N))$  as compared to the  $\mathcal{O}(N)$ . Similarly, clustering  $N$ -dimensional vectors into  $M$  clusters is expedited to  $\mathcal{O}(\log(MN))$  using quantum data, as compared to  $\mathcal{O}(\text{poly}(MN))$ .

Quantum computers however are not only effective because they handle high dimensional data better, but also offer strengthened security due to two fundamental principles of quantum mechanics - the measurement postulate and the no-cloning theorem (Wootters & Zurek, 1982). Quantum data collapses upon measurement, and can't be copied without destroying the original data, offering absolutely secure communication. Secure quantum computing is well studied, and comprises of protocols such as quantum teleportation (Bennett et al., 1993; Bouwmeester et al., 1997), quantum key distribution (Bennett & Brassard, 2014; Bennett et al., 1992), and quantum secure direct communication (Long & Liu, 2002; Deng et al., 2003; Wang et al., 2005; Zhang et al., 2006). Additionally, quantum computers utilizing various technologies, such as trapped ions, photons, superconducting circuits and so forth, are actively being developed, enhancing the practical implementation of these secure communication protocols.

Meanwhile, kernel-based ML comprises a set of techniques that are particularly effective for classification and regression tasks. These methods assess the similarity between data points in higher-dimensional spaces, which is essential for learning from data when data is not trivially separable. In contrast to more advanced

counterparts such as deep-learning, many kernel-based ML techniques often offer greater interpretability (Morocho-Cayamcela et al., 2019; Ponte & Melko, 2017), and provide better accuracy and models when the high-dimensional data is limited which is often the case in many real-world applications (Ding et al., 2021; Montesinos-López et al., 2021). While much of the focus in QML has been on developing quantum or hybrid (quantum-classical) deep learning and neural networks (Garg & Ramakrishnan, 2020; Kwak et al., 2023), quantum analogues of kernel-based ML techniques are important alternatives with landmark studies that work on centralized data (Havlíček et al., 2019; Schuld & Killoran, 2019).

In real-world scenarios, data is often distributed amongst various parties, who would want to collaboratively train a model but ensure privacy of their data. A critical challenge in QML is designing methods that work with distributed data in a secure manner. Current research in the field of QML in the context of distributed kernel-based techniques (Yu et al., 2006; Hannemann et al., 2023) is limited, with only one notable study addressing it (Sheng & Zhou, 2017). Although it is not described as a kernel method originally, Schuld & Killoran (2019) later laid the mathematical framework to describe the work to be a special case of the linear kernel applied in a 2-qubit system.

Our work addresses the gaps left by previous research by introducing a novel approach to securely compute some commonly used kernels. We achieve this by encoding classical data into quantum states using Random Fourier Features (RFF) (Rahimi & Recht, 2007). We provide a robust architecture for secure and distributed kernel-based learning, utilizing a centralized semi-honest server to compute kernels and train machine learning models. This approach leverages quantum teleportation to ensure data security during transmission.

To validate our proposed architecture empirically, we use various publicly available datasets and Qiskit’s Aer Simulator (Wille et al., 2019) to simulate feasible kernel-based machine learning techniques. Our results demonstrate that this method not only ensures data security but also achieves performance comparable to both centralized classical and quantum scenarios. Although quantum analogs of classical algorithms provide promising avenues for secure and efficient computations in high-dimensional data settings, it is widely recognized that these methods exhibit lower accuracy (Bharti et al., 2022). This is primarily due to inherent quantum noise, approximations in quantum state preparation, and the current limitations of quantum simulators and hardware.

We make the following three contributions:

1. **Introduction of Quantum Feature Maps:** We introduce quantum feature maps for the polynomial, Radial Basis Function (RBF), and Laplacian kernels, and theoretically prove the correctness of these feature maps.
2. **Architecture for Secure Kernel Computation:** We propose a secure architecture to securely compute the linear, polynomial, RBF, and Laplacian kernels in a federated manner on distributed datasets using quantum encoding.
3. **Implementation and Validation:** We theoretically validate our architecture, and for empirical validation, we implement the architecture for linear kernels on publicly available datasets using the capabilities of Qiskit’s Aer Simulator. Due to the limitations of the simulator and our lack of access to a real quantum computer, we were unable to test other kernels at this stage.

## 2 Background

### 2.1 Kernel-based Machine Learning

In machine learning, one typically works with a dataset  $\mathcal{X}$  consisting of data points  $\{x_1, x_2, \dots, x_N\} \in \mathcal{X}$ , where the goal is to identify patterns to evaluate previously unseen data. Kernel-based techniques employ a similarity measure called the kernel function, between two inputs to construct models that effectively capture the underlying properties of the data distribution. This kernel function is often an inner product in a feature space, typically of higher dimensionality, where non-linear relationships between data points becomes more apparent. Various kernel functions are used in practice, such as Linear, RBF, Polynomial, and Laplacian Kernels. These functions are designed to accommodate diverse data characteristics, making them suitable

for various applications. Besides many applications, these methods have a rich theoretical foundation that we briefly explore below.

**Definition 1.** *Kernel function (Aizerman, 1964)*

A kernel function  $K$  is a map  $K : \mathcal{X} \times \mathcal{X} \rightarrow \mathbb{C}$  that satisfies  $K(x, y) = \langle \phi(x), \phi(y) \rangle$ , where  $\phi : \mathcal{X} \rightarrow \mathcal{H}$  is a map from the input space  $\mathcal{X}$  to a Hilbert space  $(\mathcal{H}, \langle \cdot, \cdot \rangle)$ .

One refers to  $\phi$  as a feature map. Since for any unitary operator  $U : \mathcal{H} \rightarrow \mathcal{H}$ ,  $\langle \phi(x), \phi(y) \rangle = \langle U\phi(x), U\phi(y) \rangle$ , a kernel can be related to many different feature maps. However, kernel theory defines a unique Hilbert space associated with a kernel, called the Reproducing Kernel Hilbert space (RKHS) as follows.

**Definition 2.** *Reproducing Kernel Hilbert Space (RKHS) (Aronszajn, 1950)*

Let  $\mathcal{X}$  be an input space, and  $(\mathcal{R}, \langle \cdot, \cdot \rangle)$  the Hilbert space of functions  $f : \mathcal{X} \rightarrow \mathbb{C}$ . Then  $\mathcal{R}$  is an RKHS if there exists a function  $K : \mathcal{X} \times \mathcal{X} \rightarrow \mathbb{C}$  such that for all  $x \in \mathcal{X}$  and  $f \in \mathcal{R}$ , the following holds true:

$$f(x) = \langle f, K(x, \cdot) \rangle.$$

Alternatively, considering an associated feature map,  $\phi : \mathcal{X} \rightarrow \mathcal{H}$ , then  $\mathcal{R}$  is the space of functions  $f : \mathcal{X} \rightarrow \mathbb{C}$  such that for all  $x \in \mathcal{X}$  and  $\nu \in \mathcal{H}$ ,

$$f(x) = \langle \nu, \phi(x) \rangle_{\mathcal{H}}.$$

Typically, a large family of machine learning problems aim to compute a prediction function  $f : \mathcal{X} \rightarrow \mathbb{C}$  that takes training or test data and predicts the corresponding label. This is often formulated as the solution to the following optimization problem:

$$\min_{f \in \mathcal{R}} \left( \sum_{j=1}^n L(y_j, f(x_j)) + \lambda \|f\|^2 \right), \quad (1)$$

where  $L$  is a loss function,  $x_j$  are training data points,  $y_j$  the corresponding labels, and  $\lambda$  a regularization term. This prediction function generally lives in an RKHS. The representer theorem (Schölkopf & Smola, 2002) then states that the solution to this optimization problem can be formulated as follows.

$$f^*(x) = \sum_{j=1}^n \alpha_j K(x, x_j),$$

where  $K$  is the corresponding kernel in the RKHS. Hence, the optimization in the infinite-dimensional space is reduced to a finite-dimensional problem of solving for  $\alpha_j$  by computing the kernel values at the training data points.

In summary, kernel functions are fundamental in machine learning as they enable the transformation of data into higher-dimensional spaces where non-trivial relationships between data can be studied. By leveraging the theoretical framework of RKHS and the representer theorem, kernels facilitate efficient computations for a large class of machine learning models, such as Support Vector Machines (SVM), Gaussian Processes, and Principal Component Analysis (PCA).

## 2.2 Quantum Encoding

Quantum encoding techniques are crucial for translating classical data into quantum states. There are various methods to do so, such as basis encoding, angle encoding, amplitude encoding, and Hamiltonian evolution ansatz encoding - each with its distinct advantages and disadvantages. For instance, one defines one such encoding below.

**Definition 3.** *Amplitude Encoding (Schuld et al., 2015)*

Given classical data  $x = (x_1, x_2, \dots, x_N)^T$ , where  $N = 2^n$ , one defines the amplitude encoding of the data as the quantum state:

$$|\psi(x)\rangle := \sum_{j=1}^N \frac{x_j}{\|x\|} |j\rangle,$$

where  $|j\rangle$  represents the computational basis states of an  $n$ -qubit system.

Hence, encoding classical information into qubits is achieved by mapping an input  $x \in \mathcal{X}$  to a quantum state  $|\psi(x)\rangle$  in a separable complex Hilbert space  $\mathcal{H}$ . This encoding  $x \mapsto |\psi(x)\rangle$  can then just be viewed as a *quantum feature map* (Schuld & Killoran, 2019), inducing a kernel and an associated RKHS. Utilizing the representer theorem, this connection enables the application of the rich theory of kernel methods in optimizing machine learning (ML) algorithms. For example, Schuld & Killoran (2019) detailed that the amplitude encoding for example, corresponds to the linear kernel (Vapnik, 2013):

$$k(x, y) = x^T y.$$

Further, Schuld & Killoran (2019) also introduce the quantum feature map *Copies of Quantum States*:

$$x = (x_1, \dots, x_N) \mapsto \left( \sum_j^N \frac{x_j}{\|x_j\|} |j\rangle \right)^{\otimes d}.$$

This map is then associated with the homogeneous polynomial kernel, expressed as

$$k(x, y) = (x^T y)^d.$$

As discussed earlier, a large family of ML algorithms optimize a functional (1) to obtain a prediction function. There are two primary approaches to this optimization in the context of QML: the implicit approach and the explicit approach. The implicit approach uses the representer theorem and computes kernels, while offloading the remaining tasks to classical computing, as demonstrated by Rebentrost et al. (2014); Schuld & Killoran (2019); Schuld (2021). The explicit approach uses variational circuits to solve the optimization problem in the infinite-dimensional RKHS, as discussed by Havlíček et al. (2019); Schuld & Killoran (2019); Cerezo et al. (2021). Our work follows the implicit approach, in a distributed setting where quantum states are used for kernel computation, and the modeling is offloaded to classical computing.

### 2.3 Random Fourier Features

Kernel methods often face significant computational challenges, particularly with large datasets. To address this issue, Rahimi & Recht (2007) introduced Random Fourier Features (RFF) as an effective approach to estimate kernel functions using finite-dimensional feature maps. RFF enable efficient computation of kernel approximations by leveraging the Fourier transform properties of shift-invariant kernels. One defines RFF as below:

**Definition 4.** *Random Fourier Features (RFF) (Rahimi & Recht, 2007)*

*Given a shift invariant kernel  $k(x - y)$  that is the Fourier transform of a probability distribution  $\chi$ , the corresponding lower dimensional feature map  $z : \mathbb{R}^D \rightarrow \mathbb{R}^d$  defined by*

$$z(x) := \left( \sqrt{2} \cos(w_1 x + b_1), \dots, \sqrt{2} \cos(w_d x + b_d) \right),$$

*with  $w_i \sim \chi((w_1, \dots, w_d))$  and  $b_i$  are independent samples from the uniform distribution  $U[0, 2\pi]$  satisfies the following inequality for all  $\epsilon$ :*

$$\mathbb{P} \left( |z^T(x)z(y) - k(x, y)| \geq \epsilon \right) \leq 2 \exp \left( \frac{-D\epsilon^2}{4} \right).$$

*The feature map  $z$  is called an RFF.*

In the context of our work where we want to compute the quantum feature maps associated with widely used kernels, we adopt RFF to determine the quantum encoding necessary for computing the RBF and Laplacian kernels.

### 3 Related Work

To the best of our knowledge, only one study (Sheng & Zhou, 2017) has implemented kernel-based techniques using quantum computing within a distributed framework. It primarily devised a method to compute distances between 2D vectors using a polarization based 1-qubit system. The method although not described as a kernel computation originally, can be seen as a special case of the linear kernel, later proposed by Schuld & Killoran (2019). However, the work did not identify or utilize the implicit relationship between quantum encoding and kernels. Consequently, it overlooked the broader kernel framework that can be leveraged for various supervised and unsupervised machine learning tasks across different types of data, including images, text, and numeric data.

In contrast, our research substantially broadens these initial concepts by facilitating the computation of encoding-induced kernels and other standard kernels such as the polynomial, RBF, and Laplacian kernels for data of any dimensionality. Our approach offers a comprehensive solution by clearly defining the roles of participants in the pipeline and establishing a well-defined adversary model. This generalization to kernel based techniques is much stronger, and important for future study.

### 4 Quantum Feature Maps

Although Schuld & Killoran (2019) pointed out the implicit connection between quantum encoding techniques and feature maps, they only devised the feature maps associated with the linear kernel, the homogeneous polynomial kernel and the cosine kernel. We extend this by defining quantum feature maps associated with three very commonly used kernels in the context of ML - the polynomial, RBF and the Laplacian kernel.

#### 4.1 Polynomial Kernel

Given classical data  $x = (x_1, x_2, \dots, x_N)^T$ , we define the following quantum feature map:

$$x \mapsto \psi(x) = \bigotimes_{j=1}^{(N+d)} \frac{\sqrt{a}\sqrt{d!}}{\sqrt{k_1!k_2! \dots k_{N+1}!}} x_1^{k_1} \dots x_N^{k_N} \sqrt{c}^{k_{N+1}} |j-1\rangle,$$

where the multi-index  $k = (k_1, \dots, k_{N+1})$  runs over all combinations such that  $\sum_{l=1}^{N+1} k_l = d$ , and  $c = 1 - a\|x\|$  if  $d \in \mathbb{N}$ , or  $c = -1 - a\|x\|$  if  $d \in 2\mathbb{N}$ .

**Theorem 1.** *The quantum feature map above is a well-defined quantum state.*

*Proof.* To be well defined, we require the map to be normalizable. Consider the map  $x \mapsto \psi(x)$ . Then, using the multinomial theorem (Aizerman, 1964; Boser et al., 1992), we have that

$$\begin{aligned} \|\psi(x)\|^2 &= \sum_{\sum_l k_l = d} \left( \frac{\sqrt{a}\sqrt{d!}}{\sqrt{k_1!k_2! \dots k_{N+1}!}} \right)^2 (x_1^{k_1})^2 \dots (x_N^{k_N})^2 c^{k_{N+1}}, \\ &= (a\|x\| + c)^d = 1. \end{aligned}$$

This completes the proof. □

**Theorem 2.** *The quantum feature map defined above yields the polynomial kernel (Schölkopf & Smola, 2002),*

$$K_{poly} = (ax^T y + c)^d,$$

*under an inner product.*

*Proof.* This follows directly from the multinomial theorem. Take  $\phi(x)$  and  $\phi(y)$  to be two quantum feature maps of classical data  $x$  and  $y$ , defined as above. Then,

$$\begin{aligned} \langle \phi(x) | \phi(y) \rangle &= \sum_{\sum_l k_l = d} \left( \frac{\sqrt{a} \sqrt{d!}}{\sqrt{k_1! k_2! \dots k_{N+1}!}} \right)^2 x_1^{k_1} \dots x_N^{k_N} y_1^{k_1} \dots y_N^{k_N} c^{k_{N+1}}, \\ &= (ax^T y + c)^d, \\ &= K_{poly}(x, y). \end{aligned}$$

This completes the proof.  $\square$

## 4.2 RBF Kernel

Using RFF, given classical data  $x = (x_1, x_2, \dots, x_N)^T$ , we define the following quantum feature map :

$$x \mapsto \psi(x) = \frac{1}{\sqrt{D}} \sum_{j=1}^D (\cos(w_j^T x) |2j-2\rangle + \sin(w_j^T x) |2j-1\rangle),$$

where  $\lceil \log_2(2D) \rceil$  determines the number of qubits used and the approximation quality, and  $w_i$  are independent samples from the normal distribution  $\mathcal{N}(0, \sigma^{-2}I)$ .

**Theorem 3.** *The quantum feature map above is a well-defined quantum state.*

*Proof.* To be well defined, we require the map to be normalizable. Consider the map  $x \mapsto \psi(x)$ . Then, we have that

$$\|\psi(x)\|^2 = \frac{1}{D} \sum_{j=1}^D \cos^2(w_j^T x) + \sin^2(w_j^T x) = 1.$$

This completes the proof.  $\square$

**Theorem 4.** *The quantum feature map defined above yields the RBF kernel (Broomhead & Lowe, 1988),*

$$K_{RBF}(x, y) = \exp\left(-\frac{\|x - y\|^2}{2\sigma^2}\right),$$

*under an inner product.*

*Proof.* Take  $\phi(x)$  and  $\phi(y)$  to be two quantum feature maps of classical data  $x$  and  $y$ , defined as above. It follows that

$$\begin{aligned} \mathbb{E}[\langle \phi(x) | \phi(y) \rangle] &= \frac{1}{D} \sum_{j=1}^D \mathbb{E}[\cos(w_j^T x) \cos(w_j^T y) + \sin(w_j^T x) \sin(w_j^T y)], \\ &= \frac{1}{D} \sum_{j=1}^D \mathbb{E}[\cos(w_j^T (x - y))]. \end{aligned}$$

Using Euler's formula, we can rewrite this as

$$\mathbb{E}[\langle \phi(x) | \phi(y) \rangle] = \frac{1}{2D} \sum_{j=1}^D (\mathbb{E}[\exp(iw_j^T (x - y))] + \mathbb{E}[\exp(-iw_j^T (x - y))]). \quad (2)$$

Since normal distributions are closed under linear transformations (Wackerly et al., 2008),

$$w_j^T (x - y) = \sum_{k=1}^N w_{jk} (x_k - y_k) \sim N\left(0, \frac{1}{\sigma^2} \sum_{k=1}^N (x_k - y_k)^2\right) \sim \mathcal{N}\left(0, \frac{1}{\sigma^2} \|x - y\|^2\right).$$

Since  $w_j^T(x - y)$  is a normal distribution, we rewrite (2) as

$$\mathbb{E}[\langle \phi(x) | \phi(y) \rangle] = \frac{1}{2D} \sum_{j=1}^D \left( M_{w_j^T(x-y)}(i) + M_{w_j^T(x-y)}(-i) \right),$$

where  $M_Z(t) = \mathbb{E}[\exp(tZ)]$  is the moment generating function of a random variable  $Z$ . Hence, since the moment generating function of a normal distribution  $Z \sim \mathcal{N}(\mu, \gamma^2)$  is given by  $M_Z(t) = \exp(t\mu + \frac{1}{2}\gamma^2 t^2)$ , we have

$$\begin{aligned} \mathbb{E}[\langle \phi(x) | \phi(y) \rangle] &= \frac{1}{2D} \sum_{j=1}^D \left[ \exp\left(-\frac{1}{2\sigma^2}\|x - y\|^2\right) + \exp\left(-\frac{1}{2\sigma^2}\|x - y\|^2\right) \right], \\ &= \exp\left(-\frac{1}{2\sigma^2}\|x - y\|^2\right) = K_{RBF}(x, y). \end{aligned}$$

This completes the proof.  $\square$

### 4.3 Laplacian Kernel

Using RFF, given classical data  $x = (x_1, x_2, \dots, x_N)^T$ , we define the following quantum feature map:

$$x \mapsto \psi(x) = \frac{1}{\sqrt{D}} \sum_{j=1}^D (\cos(w_j^T x + \alpha_j) |2j - 2\rangle + \sin(w_j^T x + \alpha_j) |2j - 1\rangle),$$

where  $\lceil \log_2(2D) \rceil$  determines the number of qubits used and the approximation quality,  $w_j$  are independent samples from the Cauchy distribution  $\mathcal{C}(0, \alpha^{-1}I)$ , and  $\alpha_j$  are independent samples from the Uniform distribution  $\mathcal{U}(0, 2\pi)$ .

**Theorem 5.** *The quantum feature map above is a well-defined quantum state.*

*Proof.* To be well defined, we require the map to be normalizable. Consider the map  $x \mapsto \psi(x)$ . Then, we have that

$$\|\psi(x)\|^2 = \frac{1}{D} \sum_{j=1}^D \cos^2(w_j^T x + \alpha_j) + \sin^2(w_j^T x + \alpha_j) = 1.$$

This completes the proof.  $\square$

**Theorem 6.** *The quantum feature map defined above yields the Laplacian kernel (Smola & Kondor, 2003),*

$$K_L(x, y) = \exp\left(-\frac{\|x - y\|_1}{\alpha}\right),$$

*under an inner product.*

*Proof.* Take  $\phi(x)$  and  $\phi(y)$  to be two quantum feature maps of classical data  $x$  and  $y$ , defined as above. It follows like in Theorem 4 that

$$\mathbb{E}[\langle \phi(x) | \phi(y) \rangle] = \frac{1}{2D} \sum_{j=1}^D (\mathbb{E}[\exp(iw_j^T(x - y))] + \mathbb{E}[\exp(-iw_j^T(x - y))]). \quad (3)$$

Since Cauchy distributions are closed under linear transformations (Nolan, 2012),

$$w_j^T(x - y) = \sum_{k=1}^N w_{jk}(x_k - y_k) \sim \mathcal{C}\left(0, \frac{1}{\alpha} \sum_{k=1}^N |x_k - y_k|^2\right) \sim \mathcal{C}\left(0, \frac{\|x - y\|_1}{\alpha}\right).$$

Since,  $w_j^T(x - y)$  is a Cauchy distribution, we rewrite (3) as

$$\mathbb{E}[\langle \phi(x) | \phi(y) \rangle] = \frac{1}{2D} \sum_{j=1}^D \left( \phi_{w_j^T(x-y)}(1) + \phi_{w_j^T(x-y)}(-1) \right),$$

where  $\phi_Z(t) = \mathbb{E}[\exp(itZ)]$  is the characteristic function of a random variable  $Z$ . Hence, since the characteristic function of a Cauchy distribution  $Z \sim \mathcal{C}(\mu, \gamma)$  is given by  $\phi_Z(t) = \exp(it\mu - \gamma|t|)$ , we have

$$\begin{aligned} \mathbb{E}[\langle \phi(x) | \phi(y) \rangle] &= \frac{1}{2D} \sum_{j=1}^D \left[ \exp\left(-\frac{\|x - y\|_1}{\alpha}\right) + \exp\left(-\frac{\|x - y\|_1}{\alpha}\right) \right], \\ &= \exp\left(-\frac{\|x - y\|_1}{\alpha}\right) = K_L(x, y). \end{aligned}$$

This completes the proof. □

## 5 Distributed Secure Computation of Kernels

### 5.1 Architecture

Our architecture comprises multiple clients, a central server, and a helper entity. The clients hold sensitive data that they privately and collaboratively want to learn from. The central server is tasked with computing the kernel securely and privately. The helper prepares entangled quantum states to facilitate quantum communication. All entities in this setup are capable of performing the necessary quantum operations. The architecture is depicted in Figure 1.

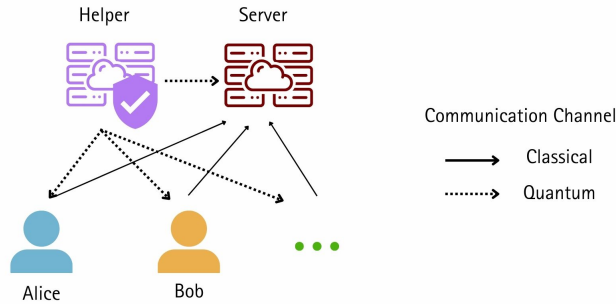


Figure 1: Visualization of our architecture consisting of a helper, a server and multiple clients.

We employ an infrastructure in which clients are initially provided with their shared seeds securely, using established cryptographic primitives. The helper party ensures the fair distribution of seeds, adhering to standard privacy-preserving protocols.

### 5.2 Protocol Description

Without loss of generality, we describe our protocol with two participants. Our method naturally extends to any number of participants. To start, Alice and Bob, declare the size of their classical data in bits, denoted by  $N$ . The helper entity then computes the number of qubits,  $n$ , needed to encode the data for a single participant based on the chosen encoding technique. Then the protocol is established within a total system of  $(6n + 1)$  qubits.

The protocol’s circuit diagram is detailed in Figure 2 below. The correctness of the protocol is theoretically shown in Appendix A.



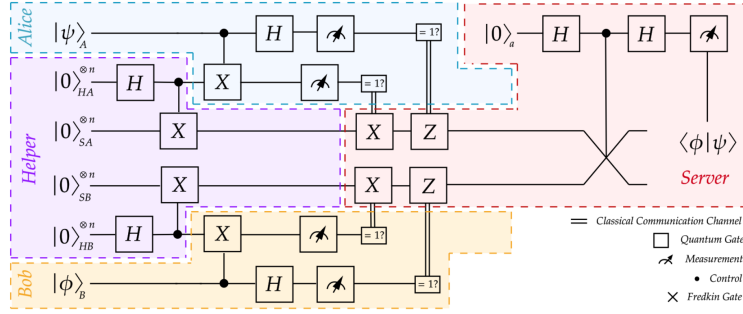


Figure 2: Quantum circuit diagram associated with our secure and distributed quantum-based kernel computation architecture.

### 5.2.1 Helper: Quantum State Preparation for Teleportation

The helper generates  $2n$  Bell states, distributing the qubits between Alice, Bob, and the server as follows:

1. In the first set of  $n$  Bell states, one qubit from each entangled pair represented by  $|0\rangle_{HA}$  is sent to Alice, and the other represented by  $|0\rangle_{SA}$  to the server.
2. In the remaining  $n$  Bell states, one qubit from each entangled pair represented by  $|0\rangle_{HB}$  is sent to Bob, and the other represented by  $|0\rangle_{SB}$  to the server.

This then enables us to quantum teleport Alice’s and Bob’s encoded data to the server for secure computation.

### 5.2.2 Clients: Data Encoding and Measurement

Alice and Bob determine the encoding of their data represented by  $|\psi\rangle_A$  and  $|\phi\rangle_B$  respectively, with multiple encodings for every data point based on the required model accuracy. The encoding sequence is derived from the initial shared seed. Subsequently, Alice and Bob execute the following steps:

1. Apply a Controlled-X gate to the qubits they received from the helper using their original quantum data as control.
2. Perform a Hadamard gate on their original data.
3. Measure their data and the received qubits in the computational basis.
4. Communicate the results to the server through an encrypted classical communication channel.

Upon receiving the measurements, the server adjusts the qubits it holds by applying appropriate X and Z gates.

### 5.2.3 Server: Inner Product Measurement

The server then executes a standard swap test (Barenco et al., 1997). It prepares an ancilla qubit in the zero state, applies a Hadamard gate, and uses it to conditionally swap the two sets of qubits received from Alice and Bob. After reverting the ancilla qubit with another Hadamard and measuring it, the output helps determine the required inner product (Buhrman et al., 2001). This measurement process is repeated  $p$  times to enhance accuracy.

## 5.3 Security of Protocol

In our proposed adversarial model, clients such as Alice and Bob, along with the server, are semi-honest. They adhere to the defined protocol yet may attempt to infer additional information from the data they

handle. A helper entity, deemed a semi-honest third party, guarantees the integrity of the quantum states used in communication, similar to protocols implementing secure multi-party computation (SMPC) (Yao, 1982). The server is explicitly characterized as non-colluding with the clients, consistent with established norms in distributed and federated architectures (Hannemann et al., 2024).

In our setup, each client processes exclusively their own data and cannot access information from other clients, effectively mitigating the risk of adversarial clients learning the data from honest input parties. The non-colluding server, does not know the series of encodings applied on the original data, and hence can't reconstruct the original classical data from the quantum data it receives, since it doesn't know how to measure it. It only learns the kernel matrix reflecting similarities between participants' data. However, since the labels are obfuscated and are irrelevant to model training, the server gains no knowledge beyond the similarity distribution pertaining to obscured labels.

An adversarial third party attempting to eavesdrop on the quantum data would face significant challenges due to the no-cloning theorem (Wootters & Zurek, 1982), which prohibits the duplication of quantum information without destroying the original information. In the event of interception, the malicious entity would need to generate and transmit its own quantum data to the server. This can be effectively detected if clients and the server intermittently exchange predetermined random quantum states, enabling the server to check for any discrepancies indicative of interference (Sheng & Zhou, 2017). Additionally, the utility of intercepted data is limited for the third party, as the encoding of data for transmission is randomized, and only known to the clients through the pre-shared seed.

## 6 Experimental Evaluation

All the proof-of-concept experiments in our following evaluation were conducted using classical computing resources on a High-Performance Computing (HPC) cluster. Each node within this HPC environment was equipped with an Intel XEON CPU E5-2650 v4, complemented by 256 GB of memory and a 2 TB SSD storage capacity. We used the Qiskit Aer Simulator to run the program offline due to limited access to IBM's quantum resources. Due to this, we were restricted to simulating only 31 qubits in our environment. Note that we don't report any timings since the experiments are run on a simulator.

Our experiments focused on computing the linear kernel. Given the limitation of simulating only 31 qubits, which confines us to  $2^7$  features, we adopted this approach and assigned  $n = 7$  qubits to each party in our distributed setup. While implementing other kernels, such as encoding-induced kernels, RBF kernels, polynomial kernels, and Laplacian kernels, would require more qubits than available, our primary goal is to validate the architecture rather than exhaustively test every encoding. Since the validity of these encodings has already been theoretically established, our focus is on demonstrating that our architecture functions as expected within this framework.

Additionally, we tested our methodology in a two-party configuration. This is trivially expandable, as the data can be redistributed to additional participants, thereby maintaining consistent results. Due to the constraints on qubit simulation, a two-party configuration is employed, allocating 14 qubits for the two data providers, 14 for the helper, and 1 for the server.

### 6.1 Accuracy Analysis

We present a comparative analysis of our distributed quantum kernel learning setup against centralized quantum kernel computation and centralized classical kernel computation. Centralized quantum kernel computation only performs the swap test, and doesn't constitute quantum teleportation. The datasets used for this analysis are widely used, and publicly available, and include the Wine dataset (178 samples, 13 features) (Asuncion et al., 2007), the Parkinson's disease dataset (197 samples, 23 features) (Sakar et al., 2019), and the Framingham Heart Study dataset (4238 samples, 15 features) (Bhardwaj, 2022). Kernel-based training was performed using SVM for all datasets, and PCA was applied to the binary datasets (Parkinson's and Framingham Heart Study) to reduce dimensionality. After applying PCA, SVM was used on the transformed data to obtain accuracy metrics. All SVM training and evaluation were performed using

stratified 5-fold cross-validation to ensure unbiased accuracy metrics. The accuracies of the different models on the datasets are summarized in Table 1 below.

Table 1: Comparison of accuracies across different methods and datasets.

Dataset (Samples $\times$ Features)	Method	Accuracy		
		Centralised Classical	Centralised Quantum	Distributed Quantum
Wine (178 $\times$ 13)	kernel-SVM	0.9860 $\pm$ 0.0172	0.8805	0.8874 $\pm$ 0.0259
Parkinsons (197 $\times$ 23)	kernel-SVM	0.8196 $\pm$ 0.0644	0.7875	0.7983 $\pm$ 0.0798
	kernel-PCA	0.7872 $\pm$ 0.0716	0.7451	0.7660 $\pm$ 0.0744
Framingham Heart Study (4238 $\times$ 15)	kernel-SVM	0.6788 $\pm$ 0.0108	0.6308	0.6340 $\pm$ 0.0143
	kernel-PCA	0.6788 $\pm$ 0.0095	0.6249	0.6422 $\pm$ 0.0092

As expected centralized classical methods generally achieve the highest accuracy, serving as a baseline. Centralized quantum methods show competitive performance, although slightly lower than their classical counterparts, due to the inherent characteristics of quantum data and quantum simulators. Our distributed quantum architecture exhibits comparable but not the same accuracy as centralized quantum architecture, due to the complexity introduced by additional gates in the quantum circuit. All experiments were conducted with 1024 shots of the quantum circuit to ensure reliable accuracy. Here, shots refers to the number of times,  $p$ , that a circuit is repeated.

## 6.2 Effect of Noise

Quantum computing is susceptible to various types of errors due to environmental interactions and imperfections in quantum gate implementations. Our objective is to evaluate the performance of distributed kernel-based QML under different noise conditions on Qiskit and compare it with a classical SVM. We employed three noise models:

**No Noise:** This model assumes an ideal environment without any noise. It serves as a baseline to evaluate the performance of the our protocol in the absence of errors.

**Noise Level 1:** This model introduces a depolarizing error with a 0.1% error rate for single-qubit gates and two-qubit gates. Depolarizing error is a type of quantum error where a qubit, with a certain probability, is replaced by a completely mixed state, losing all its original information.

**Noise Level 2:** This model simulates a more challenging environment with a depolarizing error rate of 1%.

Our results reported in Figure 3 show that increasing noise had a negative impact on model performance.

## 6.3 Effect of Shots

Here, we detail the impact of varying the number of shots used in Qiskit to repeat a quantum circuit on the performance of our proposed algorithm. We used a subset of the Digits dataset containing 100 samples (Pedregosa et al., 2011). The objective was to classify these samples into 10 labels (0-9) and evaluate the classification accuracy using linear kernel-based SVM.

We varied the number of shots, specifically using 128, 256, 512, and 1024 shots, to observe the effect on the classification accuracy. The results, depicted in Figure 4, indicate improved performance with increased amount of shots.

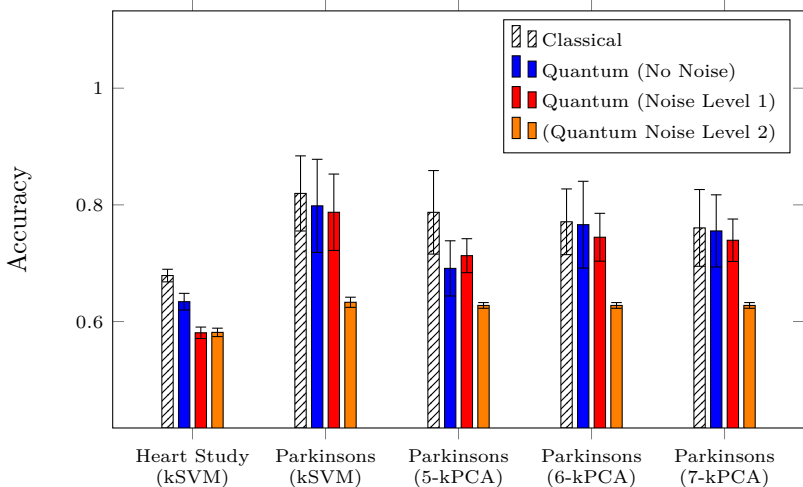


Figure 3: Comparison of accuracy scores across different noise levels. The baseline includes centralized classical kernel computation and our distributed quantum kernel computation with no noise. We incrementally introduce noise, using depolarizing error at Level 1 and Level 2, to evaluate and report the corresponding accuracy loss.

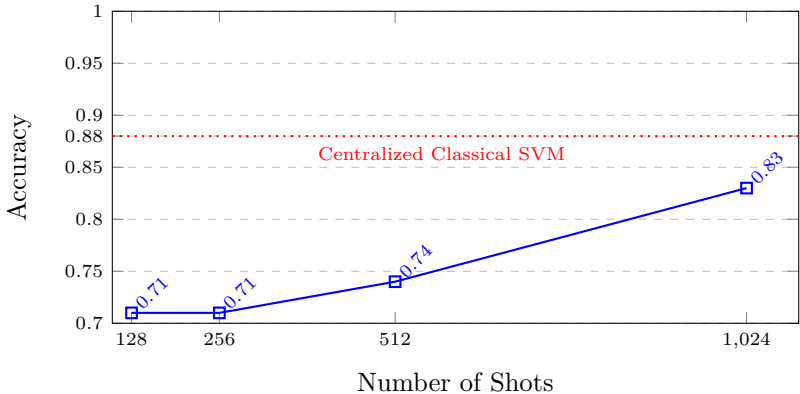


Figure 4: Accuracy of linear kernel-based SVM on a subset of the Digits dataset featuring 100 samples and 10 labels, compared against the number of shots run by the simulator.

## 7 Conclusion and Future Work

In this study, we introduced a novel kernel-based QML algorithm that operates within a distributed and secure framework. By utilizing the implicit connection between quantum encoding and kernel computation, our method supports the calculation of encoding-induced as well as traditional kernels. To that end, we extended the existing body of knowledge by introducing three novel quantum feature maps, designed to compute the polynomial, RBF, and the Laplacian kernels, building upon previous studies that mainly focused on linear and homogeneous polynomial kernels. Furthermore, we have theoretically validated that the proposed quantum feature maps help compute the concerned kernels.

Using a hybrid quantum-classical architecture, our approach functions under a distributed environment where a central server aids data providers in processing their data collaboratively, akin to a centralized model. This setup primarily addresses the case of a semi-honest scenario—a common consideration in studies involving distributed architectures. The architecture is further enhanced by the inclusion of a trusted third party helper, that ensures the integrity of the Bell pairs necessary for quantum teleportation within our framework.

We have demonstrated that our proposed framework upholds security against semi-honest parties as well as external eavesdroppers.

The application of our architecture to compute the linear kernel for publicly available datasets using Qiskit's Aer Simulator validates our distributed framework, yielding accuracies comparable to those achieved in centralized classical and quantum frameworks.

In the future, we aim to adapt our methodology to scenarios involving actively malicious entities. Additionally, provided the rich potential of kernel theory, further theoretical and practical explorations into quantum feature maps represent a promising direction for future research in the field.

## Supplementary information

Our code and results are available at the following URL: <https://anonymous.4open.science/r/distributed-secure-kernel-based-QML-480F/>.

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## A Correctness of Architecture

In this section, we provide a theoretical proof of correctness for the proposed circuit architecture, demonstrating that it accurately computes the kernel matrix for given input data. Without loss of generality, consider an  $n$ -qubit system, where Alice’s encoded data is represented by  $|\psi\rangle_A$  and Bob’s data by  $|\psi\rangle_B$ . As illustrated in Figure 1, the Helper initializes the system by preparing  $2n$  Bell states. The initial state consists of  $|0\rangle_{HA}^{\otimes n}$ ,  $|0\rangle_{SA}^{\otimes n}$ ,  $|0\rangle_{SB}^{\otimes n}$ , and  $|0\rangle_{HB}^{\otimes n}$ , where the superscript denotes qubits in each subsystem, e.g., the  $i$ -th qubit of Alice’s data is represented as  $|\psi\rangle_A^i$ .

Let  $|\psi\rangle$  in the computational basis be written as  $\alpha|0\rangle + \beta|1\rangle$  and  $|\phi\rangle$  as  $\delta|0\rangle + \gamma|1\rangle$ . Initially, the entire system is in the state:

$$|\psi\rangle_A^{\otimes n} \otimes |0\rangle_{HA}^{\otimes n} \otimes |0\rangle_{SA}^{\otimes n} \otimes |0\rangle_{SB}^{\otimes n} \otimes |0\rangle_{HB}^{\otimes n} \otimes |\phi\rangle_B^{\otimes n}$$

For simplicity, we track only the  $i$ -th qubit of each  $n$ -qubit state:

$$|\psi\rangle_A^i \otimes |0\rangle_{HA}^i \otimes |0\rangle_{SA}^i \otimes |0\rangle_{SB}^i \otimes |0\rangle_{HB}^i \otimes |\phi\rangle_B^i$$

After applying Hadamard gates to the Helper qubits, the system evolves to the following state:

$$|\psi\rangle_A^i \otimes \frac{1}{\sqrt{2}} \left( |0\rangle_{HA}^i + |1\rangle_{HA}^i \right) \otimes |0\rangle_{SA}^i \otimes |0\rangle_{SB}^i \otimes \frac{1}{\sqrt{2}} \left( |0\rangle_{HB}^i + |1\rangle_{HB}^i \right) \otimes |\phi\rangle_B^i.$$

Upon applying the Controlled-X gates, the system is entangled, preparing it for quantum teleportation:

$$\frac{1}{2} \left( |\psi\rangle_A^i \otimes \left( |00\rangle_{HA,SA}^i + |11\rangle_{HA,SA}^i \right) \otimes \left( |00\rangle_{SB,HB}^i + |11\rangle_{SB,HB}^i \right) \otimes |\phi\rangle_B^i \right).$$

Next, Alice and Bob perform Controlled-X gates, resulting in the state:

$$\begin{aligned} & \frac{1}{2} \left( \alpha |000\rangle_{A,HA,SA}^i + \beta |110\rangle_{A,HA,SA}^i + \alpha |011\rangle_{A,HA,SA}^i + \beta |101\rangle_{A,HA,SA}^i \right) \\ & \otimes \frac{1}{2} \left( \gamma |000\rangle_{SB,HB,B}^i + \delta |011\rangle_{SB,HB,B}^i + \gamma |110\rangle_{SB,HB,B}^i + \delta |101\rangle_{SB,HB,B}^i \right). \end{aligned}$$

After applying Hadamard gates, the system evolves to:

$$\begin{aligned} & \frac{1}{4} \left( \alpha |000\rangle_{A,HA,SA}^i + \alpha |100\rangle_{A,HA,SA}^i + \beta |010\rangle_{A,HA,SA}^i - \beta |110\rangle_{A,HA,SA}^i \right. \\ & \quad \left. + \alpha |011\rangle_{A,HA,SA}^i + \alpha |111\rangle_{A,HA,SA}^i + \beta |001\rangle_{A,HA,SA}^i - \beta |101\rangle_{A,HA,SA}^i \right) \\ & \otimes \frac{1}{4} \left( \left( \gamma |000\rangle_{SB,HB,B}^i + \gamma |001\rangle_{SB,HB,B}^i - \delta |011\rangle_{SB,HB,B}^i + \delta |010\rangle_{SB,HB,B}^i \right. \right. \\ & \quad \left. \left. + \gamma |110\rangle_{SB,HB,B}^i + \gamma |111\rangle_{SB,HB,B}^i - \delta |101\rangle_{SB,HB,B}^i + \delta |100\rangle_{SB,HB,B}^i \right) \right). \end{aligned}$$

Once the classical bits are communicated to the server, the server applies the appropriate X and Z gates, resulting in the system state:

$$|0\rangle_a \left( \alpha |0\rangle_{SA} + \beta |1\rangle_{SA} \right) \otimes \left( \gamma |0\rangle_{SB} + \delta |1\rangle_{SB} \right).$$

After applying a Hadamard gate, the server obtains:

$$\frac{1}{\sqrt{2}} \left( |0\rangle_a + |1\rangle_a \right) \left( |\psi\rangle_{SA} \right) \otimes \left( |\phi\rangle_{SB} \right).$$

The final step involves applying Fredkin gates, with the ancilla qubit as the control:

$$\frac{1}{\sqrt{2}} \left( |0\psi\phi\rangle_{a,SA,SB} + |1\phi\psi\rangle_{a,SA,SB} \right).$$

Upon applying a Hadamard gate to the ancilla qubit, we obtain:

$$\frac{1}{2} \left( |0\rangle_a \otimes \left( |\psi\phi\rangle_{SA,SB} + |\phi\psi\rangle_{SA,SB} \right) + |1\rangle_a \otimes \left( |\psi\phi\rangle_{SA,SB} - |\phi\psi\rangle_{SA,SB} \right) \right).$$

Measuring the ancilla qubit along the computational basis yields:

$$Pr(0)_a = \frac{1}{4} \left( \langle\psi|\langle\phi| + \langle\phi|\langle\psi| \right) \left( |\psi\rangle|\phi\rangle + |\phi\rangle|\psi\rangle \right) = \frac{1}{2} + \frac{1}{2} \|\langle\psi|\phi\rangle\|^2,$$

where,  $P(0)_a$  is the probability that the ancilla qubit is in the  $|0\rangle$  state. Rearranging, this then determines the inner product of  $|\psi\rangle$  and  $|\phi\rangle$ .

$$\|\langle\psi|\phi\rangle\| = \sqrt{2Pr(0)_a - 1},$$