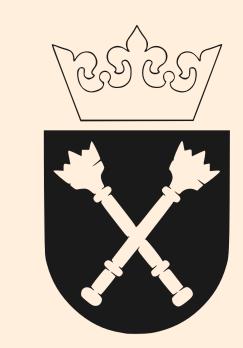
# Machine classification for probe-based quantum thermometry



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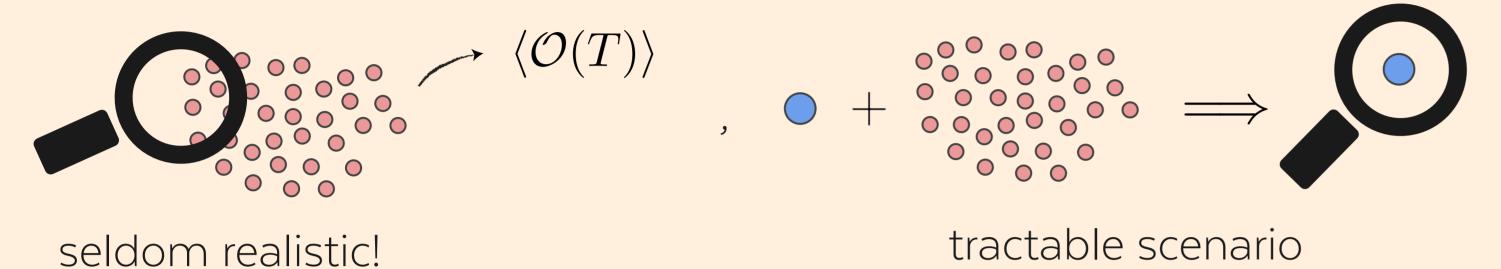
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#### Introduction

Motivation: Quantum thermometry are crucial for experimental applications. But known strategies are highly model-dependent.

This work: introduces machine classification for quantum thermometry and show that it provides reliable and entirely model-independent predictions.

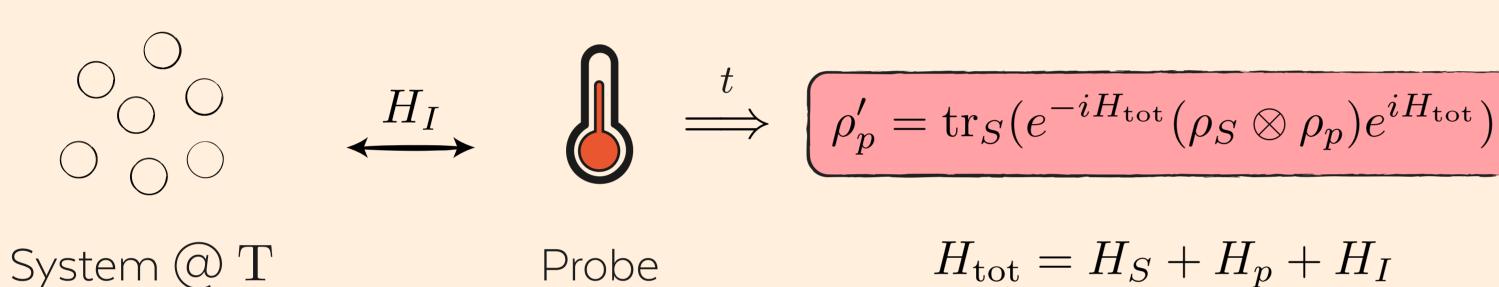
#### ■ Setting the scene

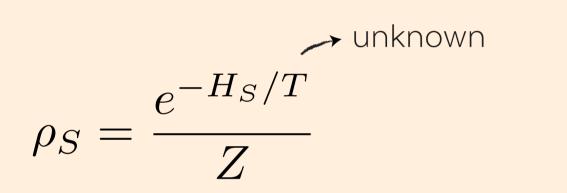


### Probe-based thermometry

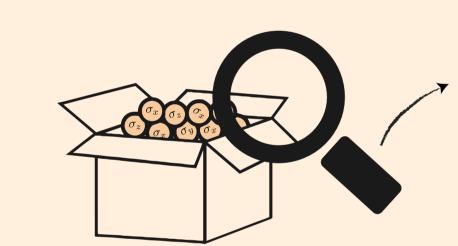
The temperature of the system is estimated by coupling it to a probe, which is subsequently **measured**. The protocol can be summarised as follow:

#### 1.Interaction





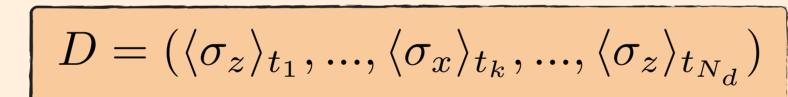
Impurities in ultra-cold gases, phonon occupation number of trapped ion, or a mechanical resonator represents a **prototypical example** of probe-based thermometry.



Experimentalist

 $ho_p$ 

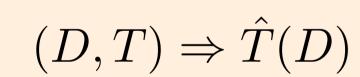
dataset

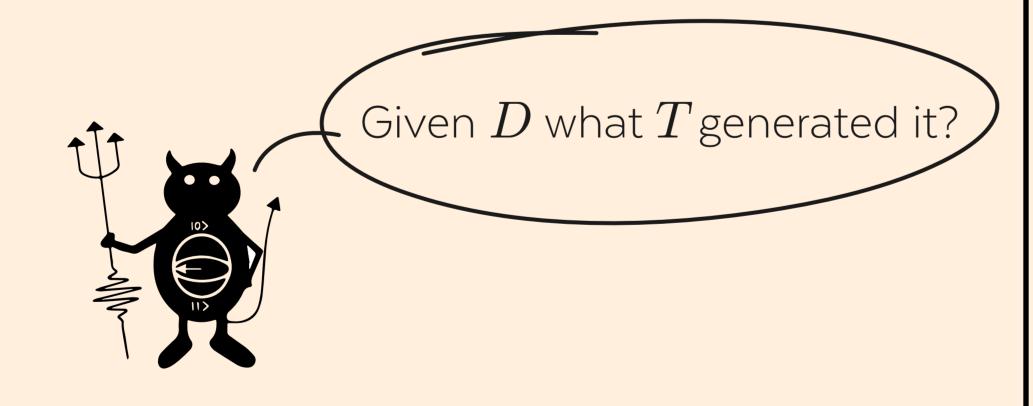


Probe observables

2. Measurement







### The k-nearest-neighbours (KNN) algorithm

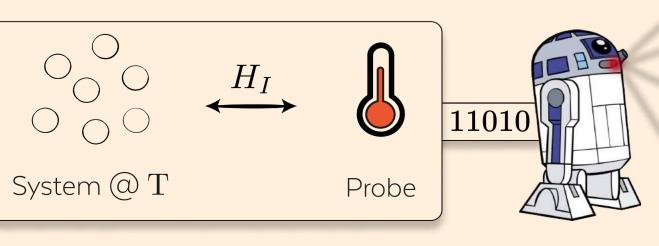
Classification is a pattern recognition method that can be employed as a concrete estimation strategy.

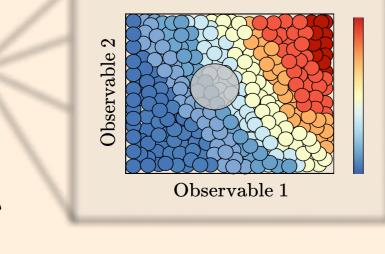
#### KNN in a nutshell

The green dot should be classified either to blue squares or to red triangles. If k=3 (solid circle) it is assigned to pink triangles. However, if k=5(dashed circle), it is assigned to the blue squares.

#### Probe-based thermometry and machine classification

Prior information:  $T \in [T_{\min}, T_{\max}]$ 





- 1. Discretize T2. Train using  $(D_i, T_i)$
- 3. KNN:  $\hat{T}$

## Data: training (70%) and validation (30%) sets

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Rzeczpospolita Polska

MSE

Jaynes-Cummings (JC) model - We illustrate the idea using the JC model. The probe is described by a qubit and the system by a bosonic mode. The total Hamiltonian is

Results

$$H = \omega a^{\dagger} a + \frac{\Omega}{2} \sigma_z + \gamma (a^{\dagger} \sigma_- + a \sigma_+)$$

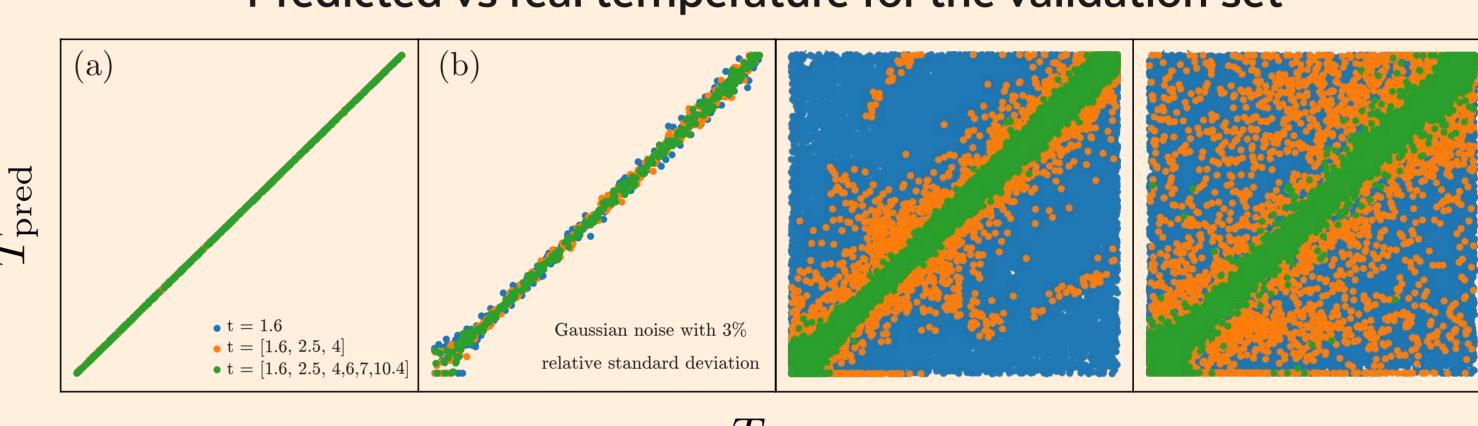
■ The probe is taken to be resonant with the system  $(\Omega = \omega)$  and start in the pure state:

 $|\psi_P\rangle = |+\rangle = \frac{1}{\sqrt{2}}(|0\rangle + |1\rangle)$ 

■ Setting:  $T, \gamma \in [0.1, 2]$ 

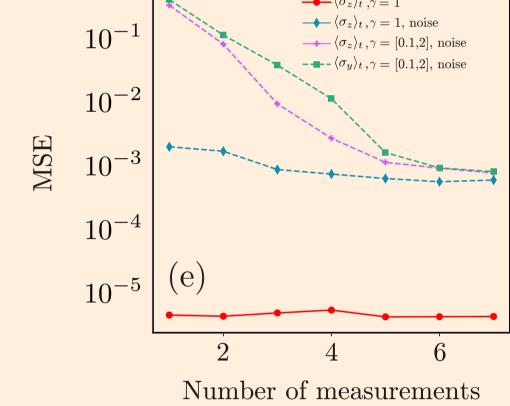
Free parameters:  $\gamma, T$ 

#### Predicted vs real temperature for the validation set



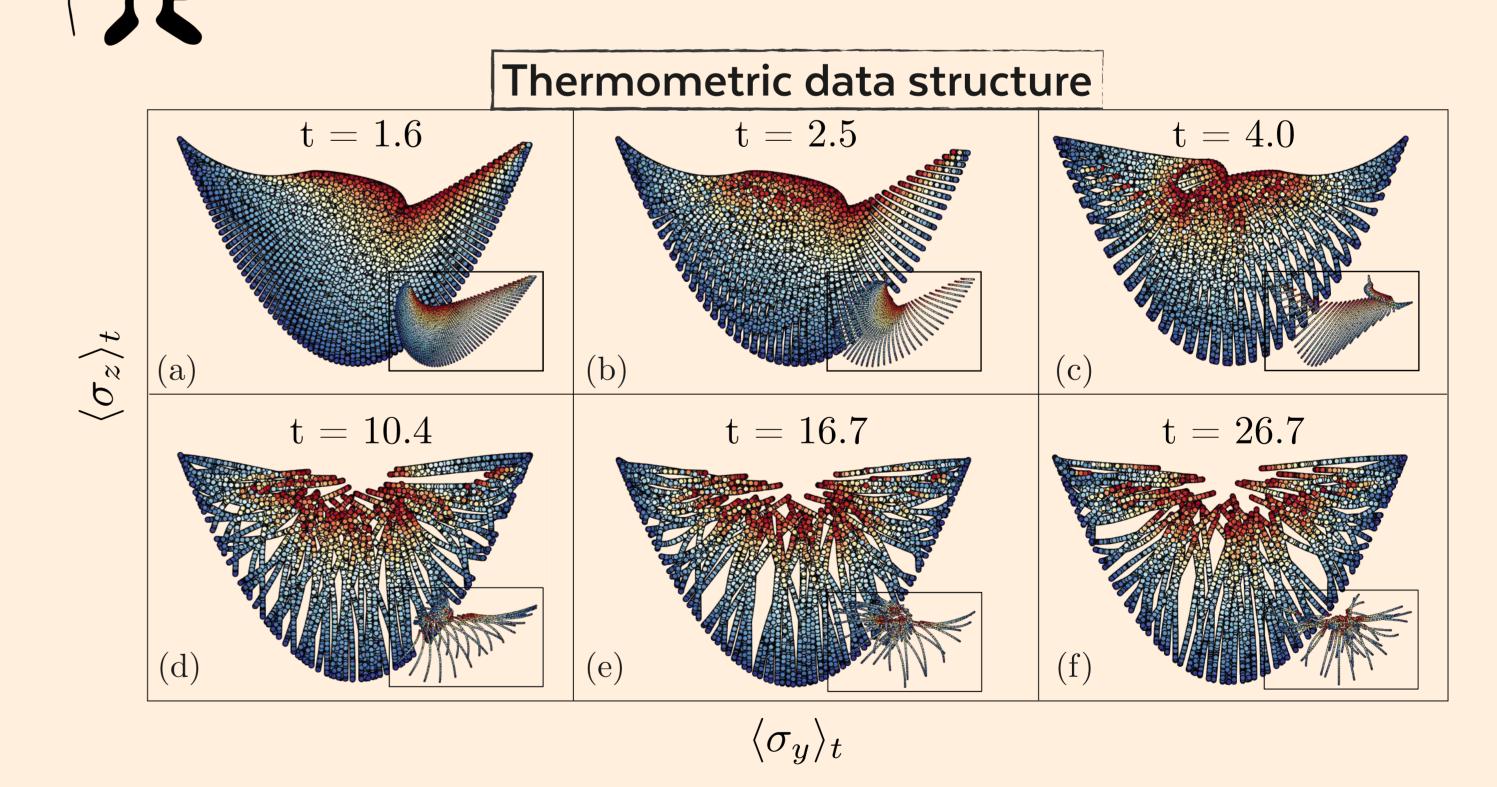
 $T_{\rm real}$ 

- (a) Using only data from  $\langle \sigma_z \rangle_t$  with  $\gamma = 1$ .
- (b) Same, but with noise.
- (c) Similar to (b) but with  $\gamma \in [0.1, 2]$ .
- (d) Same, but using  $\langle \sigma_y \rangle_t$  instead.
- (e) Net mean-squared error as function of the number of measurement times.



 $MSE = \frac{1}{N_{\text{val}}} \sum_{\text{val} \text{ set}} (T_{\text{pred}} - T_{\text{real}})^2$ 

# Why does KNN lead to higher precision?



- The dataset is segmented into well-defined regions, e.g., the change from hot to cold regions is smooth.
- (a)-(f)  $\langle \sigma_z \rangle_t$  vs  $\langle \sigma_y \rangle_t$  JC model at different times, for  $T \in [0.1, 2]$ and  $\gamma \in [0.1, 2]$ . The colors represent the temperature of the correspoding data point. The insets are similar, but for the Rabi model instead.
- We also have explored other systems, such as qudits and spin chains. Also, performed a variety of parameter choices: resonant vs non-resonant energy gaps, different probe states, and so on.

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Number of measurements