

Quantum thermometry with adaptive Bayesian strategies: a case study for release-recapture experiments

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Key works:

arXiv:2204.11816 (accepted in PRX Quantum)

Quantum Sci. Technol. (8) 015009, 2022

Phys. Rev. Lett. (127) 190402, 2021

with

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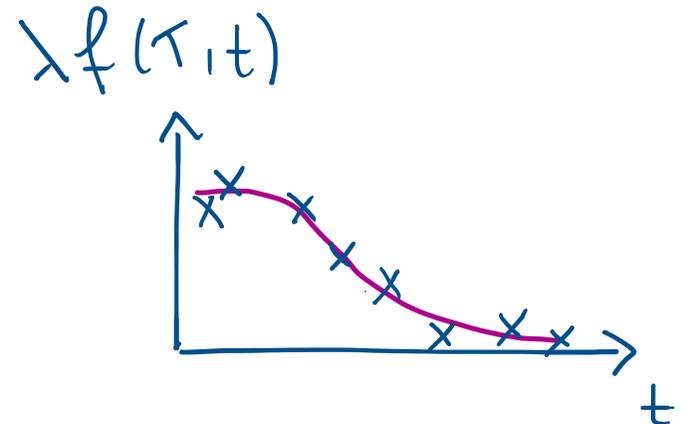
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BIRMINGHAM

JSPS London Symposium, Nottingham

15th Dec 2022

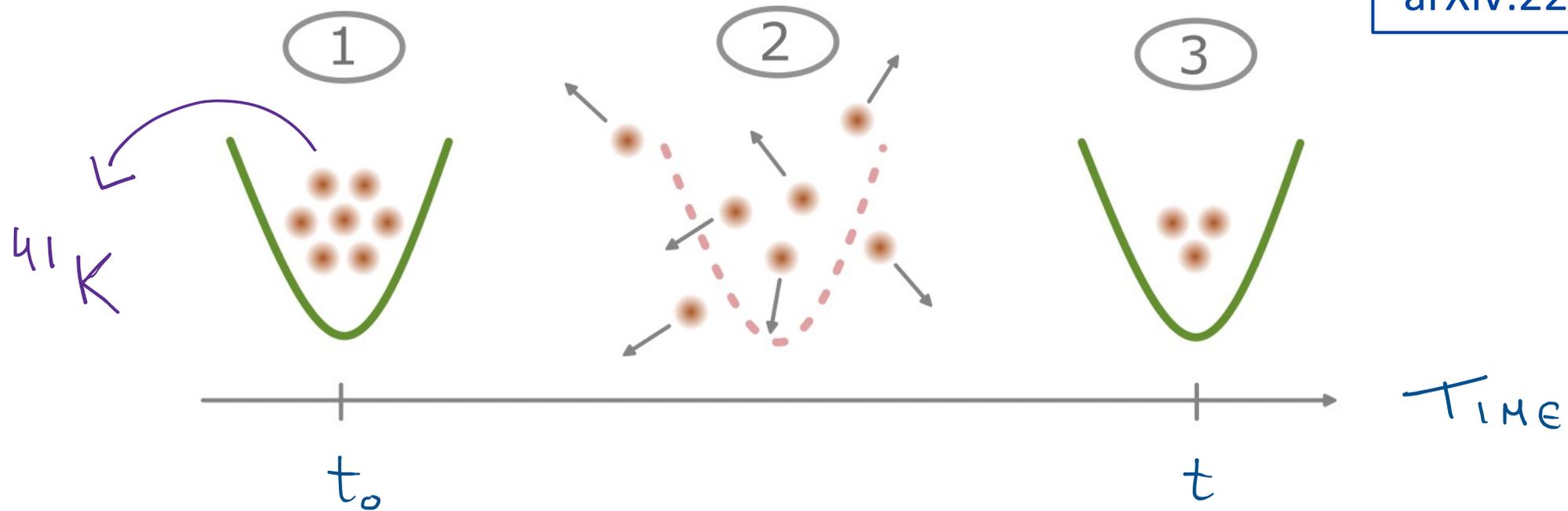
Our plan for today

- I. Release-recapture thermometry
- II. Bayes meets thermometry: a quick journey through the foundations of scale estimation
- III. Maximising information content in an adaptive fashion**
- IV. Conclusions and outlook



I. Release-recapture thermometry: an overview

arXiv:2204.11816



- 1) N_0 atoms at temperature T are trapped
- 2) Trap switched off \Rightarrow atoms expand ballistically
- 3) $n \leq N_0$ atoms are recaptured

I. Release-recapture thermometry: an overview

After \mathcal{N} trials:

$$\begin{aligned} \vec{n} &::= (n_1, n_2, \dots, n_N) \\ \vec{t} &::= (t_1, t_2, \dots, t_N) \end{aligned}$$

numbers of recaptured atoms

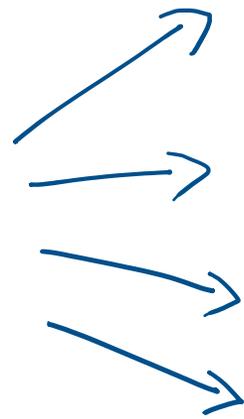
data

expansion times \equiv

controlled parameter

I. Release-recapture thermometry: an overview

PROTOCOLS FOR
TEMPERATURE
ESTIMATION



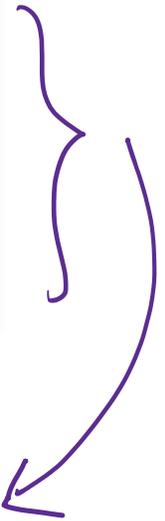
(i) Least squares

(ii) Unoptimised Bayes

(iii) A priori optimised

(iv) Fully adaptive

Standard approach



global quantum
thermometry

Least squares (standard approach)

$$\vec{n} = (\underbrace{n_{11}, \dots, n_{1v}}_{\downarrow}, \dots, \underbrace{n_{v1}, \dots, n_{vv}}_{\downarrow})$$

$$\left. \begin{array}{l} \langle \vec{n} \rangle = (\langle n_1 \rangle, \dots, \langle n_v \rangle) \\ \vec{t} = (t_1, \dots, t_v) \end{array} \right\} \text{data}$$

$$\left(\sum_{i=1}^v \alpha_i = \underbrace{M}_{\text{trials}} \right)$$

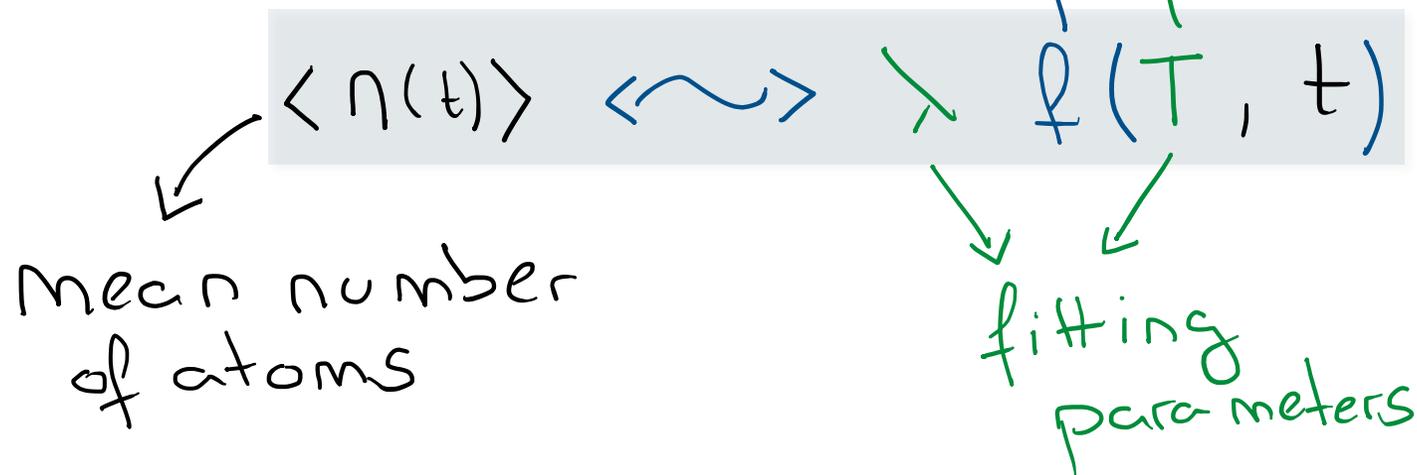
Least squares (standard approach)

$$\left. \begin{aligned} \langle \vec{n} \rangle &= (\langle n_1 \rangle, \dots, \langle n_\nu \rangle) \\ \vec{t} &= (t_1, \dots, t_\nu) \end{aligned} \right\} \text{data}$$

fraction of recaptured atoms

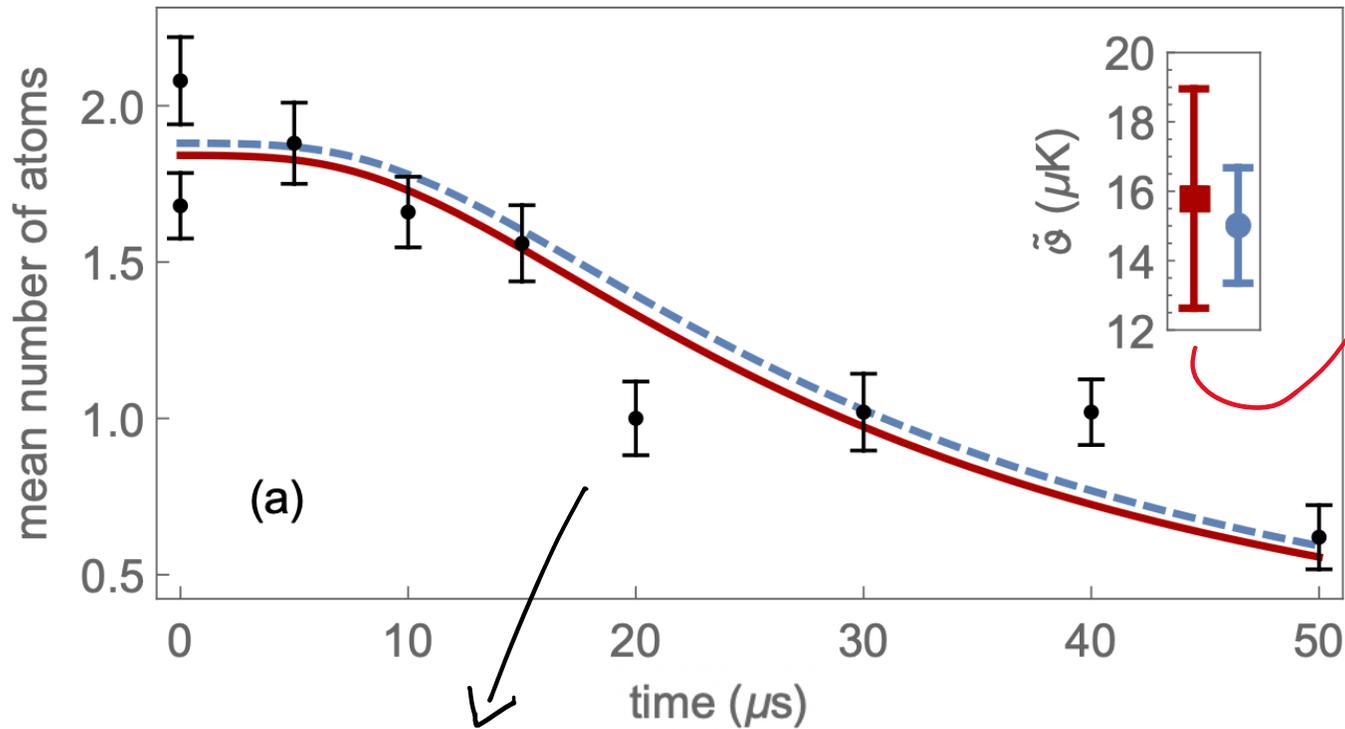
From statistical mechanics:

unknown temperature



Least squares (standard approach)

▣ Fitting to $\langle n(t) \rangle \leftrightarrow f(T, t)$, we get:

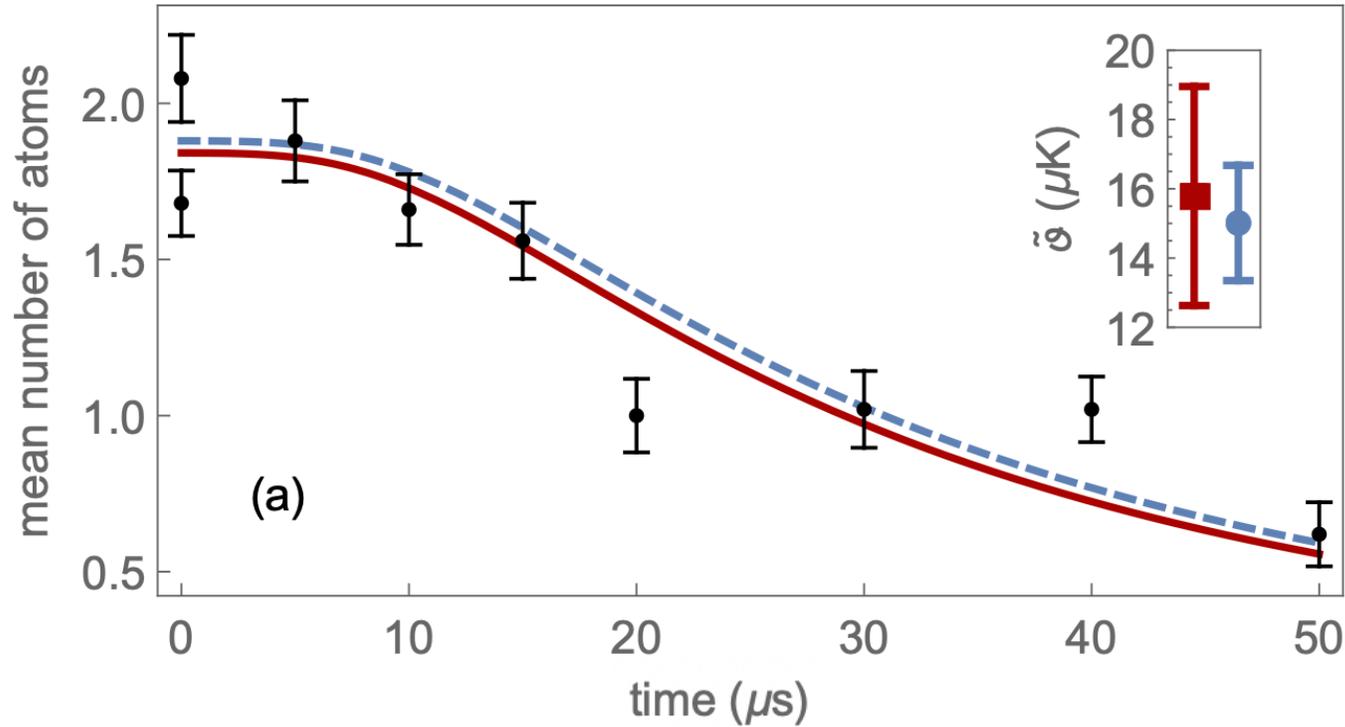


$\langle n_i \rangle$ at time t_i

$$\bar{T} = 15.8 \pm 3.2 \text{ MK}$$

FINAL TEMPERATURE
ESTIMATE

Least squares (standard approach)



$$\bar{T} = 15.8 \pm 3.2 \text{ } \mu\text{K}$$

• Why another method?

→ Fails to use all available info

→ Wastes resources

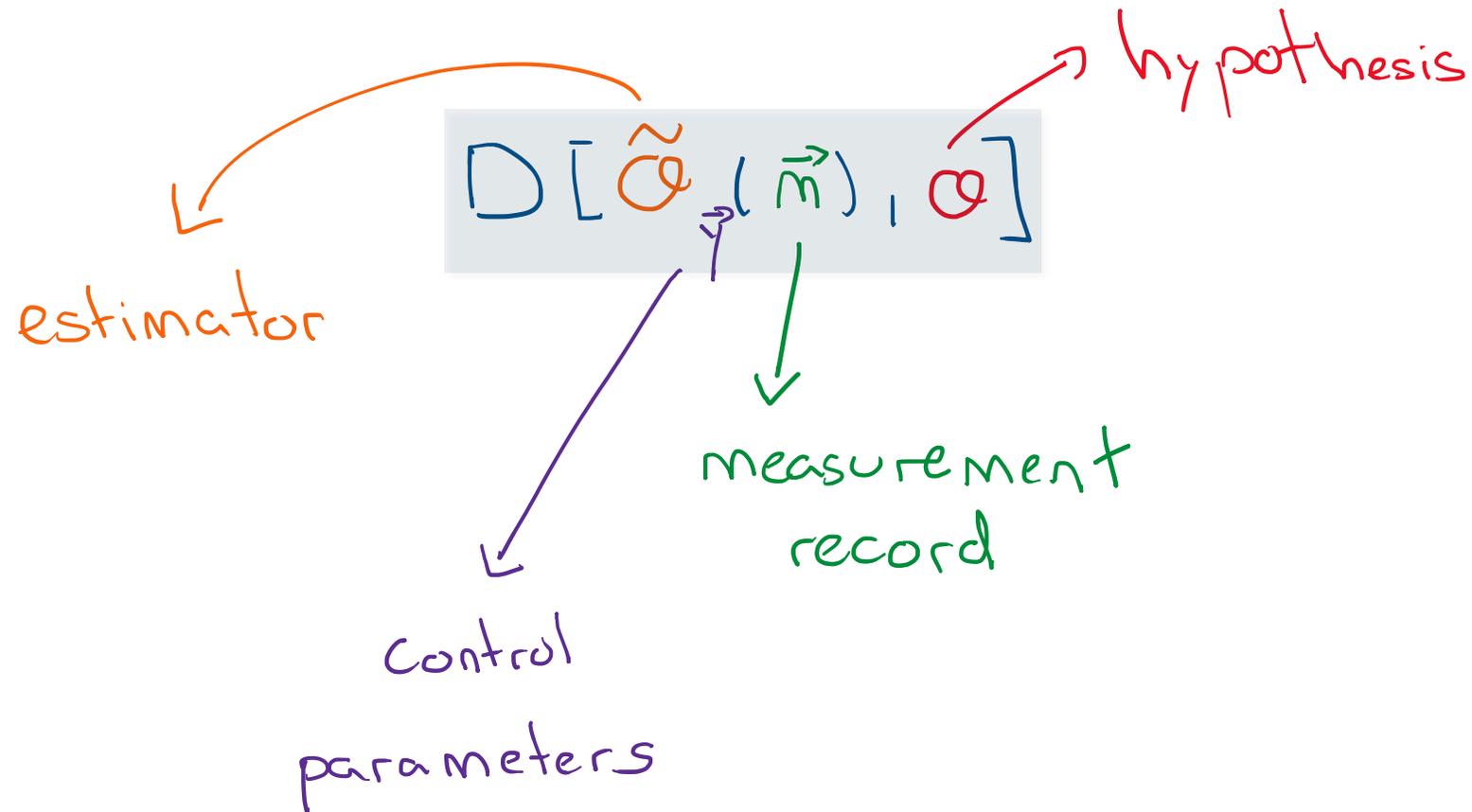
• With global quantum thermometry (this talk):

Twice as much precision with half of the measurement data

II. Bayes meets thermometry: a quick journey through the foundations of scale estimation



1 Construct a deviation function D :



II. Bayes meets thermometry: a quick journey through the foundations of scale estimation



1 Construct a deviation function $D[\tilde{\mathcal{Q}}_{\vec{y}}(\vec{m}), \vartheta]$

2 Construct an error functional $\bar{E}_{\vec{y}}$:

$$\bar{E}_{\vec{y}} = \int d\vartheta d\vec{m} p(\vartheta, \vec{m} | \vec{y}) D[\tilde{\mathcal{Q}}_{\vec{y}}(\vec{m}), \vartheta]$$

parameter independent

measurement independent

joint probability
(physical assumptions)

II. Bayes meets thermometry: a quick journey through the foundations of scale estimation



1 Construct a deviation function $D[\tilde{\Theta}_{\vec{\gamma}}(\vec{M}), \Theta]$

2 Construct an error functional

$$\bar{E}_{\vec{\gamma}} = \int d\Theta d\vec{M} p(\Theta, \vec{M} | \vec{\gamma}) D[\tilde{\Theta}_{\vec{\gamma}}(\vec{M}), \Theta]$$

3 Minimise $\bar{E}_{\vec{\gamma}}$ over:

- Estimator $\tilde{\Theta}_{\vec{\gamma}}(\vec{M})$
- Control parameters $\vec{\gamma}$
- POVM $\Pi_{\vec{\gamma}}(\vec{M})$

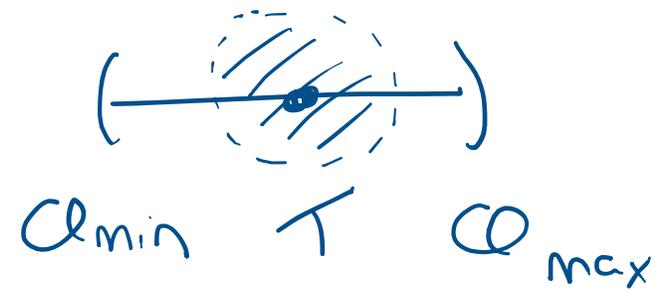
Optimisation problem

Release-recapture thermometry: unoptimised times



Physical assumptions:

$$p(\omega, \vec{n} | \vec{t}) = p(\omega) \underbrace{p(\vec{n} | \omega, \vec{t})}_{\text{likelihood}}$$



prior

$$\prod_{i=1}^n p(n_i | \omega, t_i)$$

independently estimated

$$P(n_i) \times f(\omega, t_i)$$

Poisson distribution

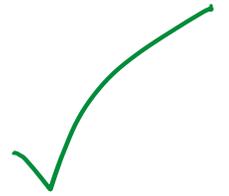
fraction of recaptured atoms

Release-recapture thermometry: unoptimised times



- Physical assumptions:

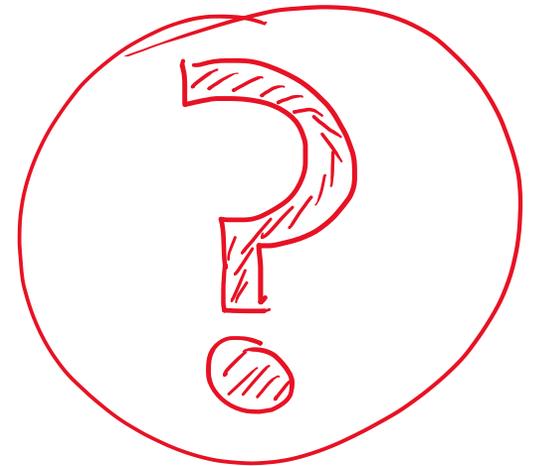
$$p(\alpha, \vec{n} | \vec{t}) = p(\alpha) \prod_{i=1}^n P[n_i | f(\alpha, t_i)]$$



- Missing:

→ prior $p(\alpha)$

→ deviation function $D[\hat{\alpha}(\vec{n}, \vec{t}), \alpha]$



Release-recapture thermometry: unoptimised times



fraction of recaptured atoms

expansion time

characteristic energy (trap parameters)

$$f(T, t) = \frac{g\{[E_k/(k_B T)] W(U_0/E_k)\}}{g[U_0/(k_B T)]}$$

temperature

trap depth (energy)

$\left\{ \begin{array}{l} g(\cdot) \\ W(\cdot) \equiv \text{Lambert function} \end{array} \right.$

Invariant under:

$$\left\{ \begin{array}{l} T \longrightarrow T' = \gamma T \\ U_0 \longrightarrow U_0' = \gamma U_0 \\ E_k \longrightarrow E_k' = \gamma E_k \end{array} \right.$$

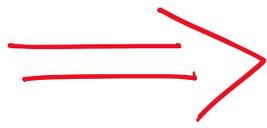
Release-recapture thermometry: unoptimised times



$$\begin{cases} T & \longrightarrow & T' = \gamma T \\ U_0 & \longrightarrow & U_0' = \gamma U_0 \\ E_{\nu} & \longrightarrow & E_{\nu}' = \gamma E_{\nu} \end{cases}$$

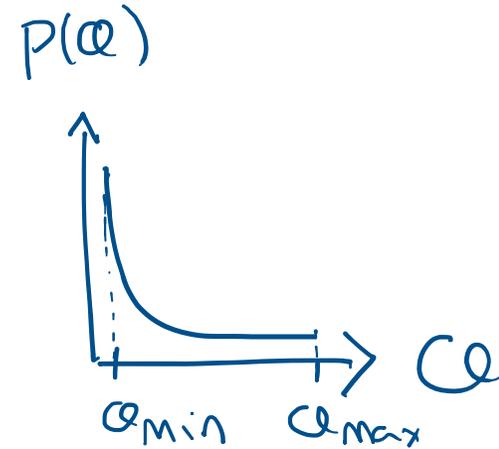
≡

SCALE
INVARIANCE



- **Jeffreys's prior:**

$$p(\theta) = \left[\theta \log \left(\frac{\theta_{\max}}{\theta_{\min}} \right) \right]^{-1}$$



- **Logarithmic error:**

$$D[\tilde{\theta}(\mathbf{n}, t), \theta] = \log^2[\tilde{\theta}(\mathbf{n}, t)/\theta]$$



IEEE Trans. Syst. Cybern. **4** 227–41 1968

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Release-recapture thermometry: unoptimised times



- Optimal rule to post-process measurements into a temperature reading:

$$\tilde{\vartheta}(\mathbf{n}, t) = \theta_u \exp \left[\int d\theta \underbrace{p(\theta | \mathbf{n}, t)} \log \left(\frac{\theta}{\theta_u} \right) \right]$$

UNIVERSALLY
VALID FOR
SCALE ESTIMATION

$$p(\theta | \mathbf{n}, t) \propto p(\theta) \prod_{i=1}^{\mu} p(n_i | \theta, t_i)$$

≡ BAYES THEOREM

Release-recapture thermometry: unoptimised times

How do we report temperature estimates in global quantum thermometry?

$$\tilde{\vartheta}(\mathbf{n}, t) \pm \Delta\tilde{\vartheta}(\mathbf{n}, t)$$

- Optimal **temperature estimator**:

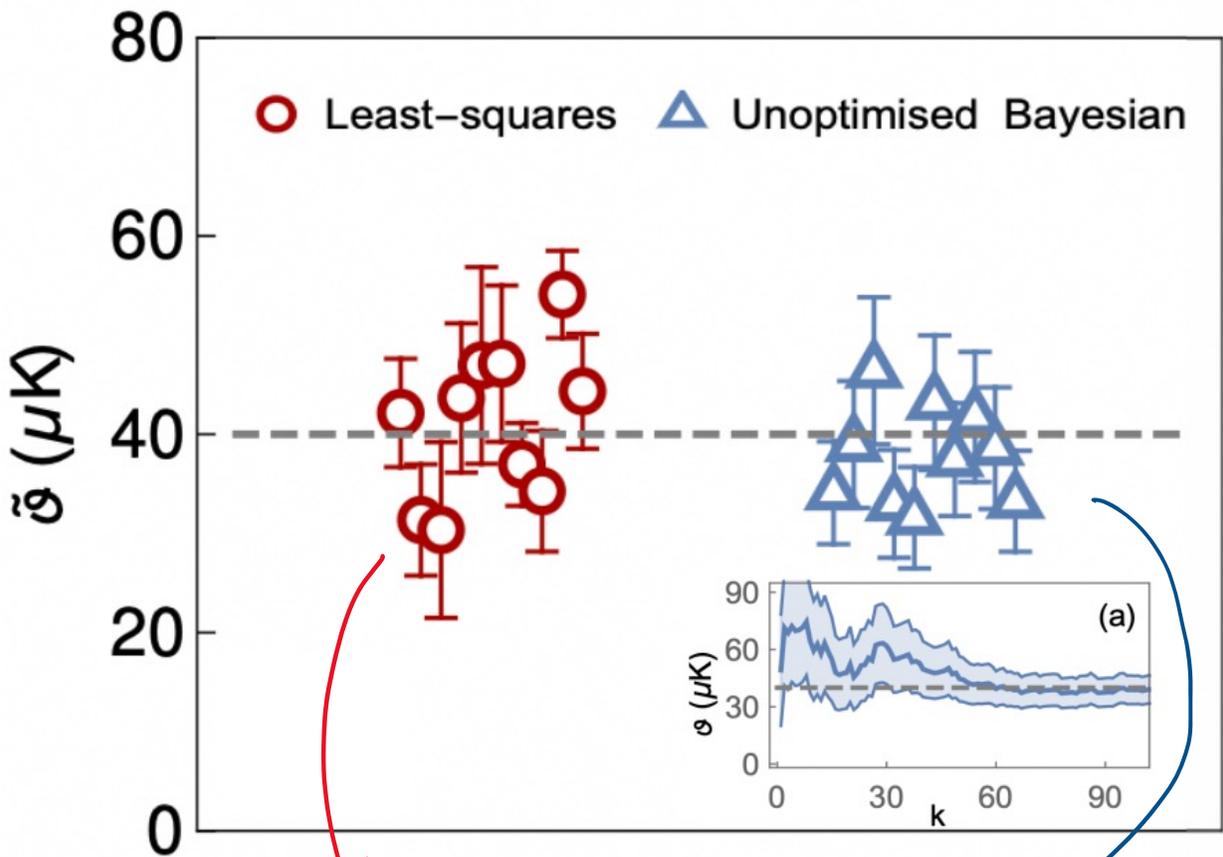
$$\tilde{\vartheta}(\mathbf{n}, t) = \theta_u \exp \left[\int d\theta p(\theta|\mathbf{n}, t) \log \left(\frac{\theta}{\theta_u} \right) \right]$$

- **Error bar**:

$$\Delta\tilde{\vartheta}(\mathbf{n}, t) = \tilde{\vartheta}(\mathbf{n}, t) \sqrt{\bar{\epsilon}_{\text{mle}}(\mathbf{n}, t)}$$

- Measurement-dependent **mean logarithmic error**:

$$\bar{\epsilon}_{\text{mle}}(\mathbf{n}, t) = \int d\theta p(\theta|\mathbf{n}, t) \log^2 \left[\frac{\tilde{\vartheta}(\mathbf{n}, t)}{\theta} \right]$$

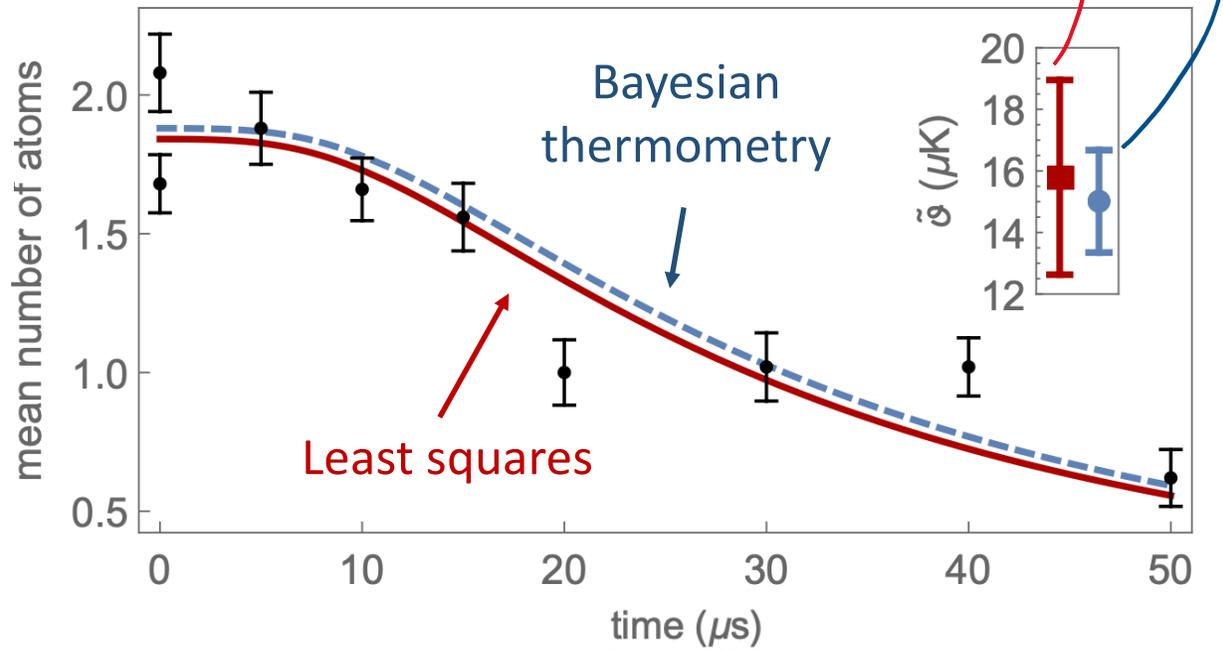


$$\hat{\tilde{\theta}} = 15.0 \pm 1.7 \text{ } \mu\text{K}$$

$$\hat{\tilde{\theta}} = 15.8 \pm 3.2 \text{ } \mu\text{K}$$

$$\frac{\Delta \hat{\tilde{\theta}}^2}{\hat{\tilde{\theta}}^2} = 0.064$$

$$\frac{\Delta \hat{\tilde{\theta}}^2}{\hat{\tilde{\theta}}^2} = 0.057$$



III. Maximising information content in an adaptive fashion

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Mean information gain for a single shot (supersedes the Fisher information):

$$\mathcal{K}(t) = \sum_n p(n|t) \log^2 \left[\frac{\tilde{v}(n, t)}{\tilde{v}_p} \right]$$

where

- optimal *a priori* estimate:

$$\tilde{v}_p = \theta_u \exp \left[\int d\theta p(\theta) \log \left(\frac{\theta}{\theta_u} \right) \right]$$

- evidence:

$$p(n|t) = \int d\theta p(\theta) p(n|\theta, t)$$

information provided by
the measurement w.r.t.
the optimal prior
estimate

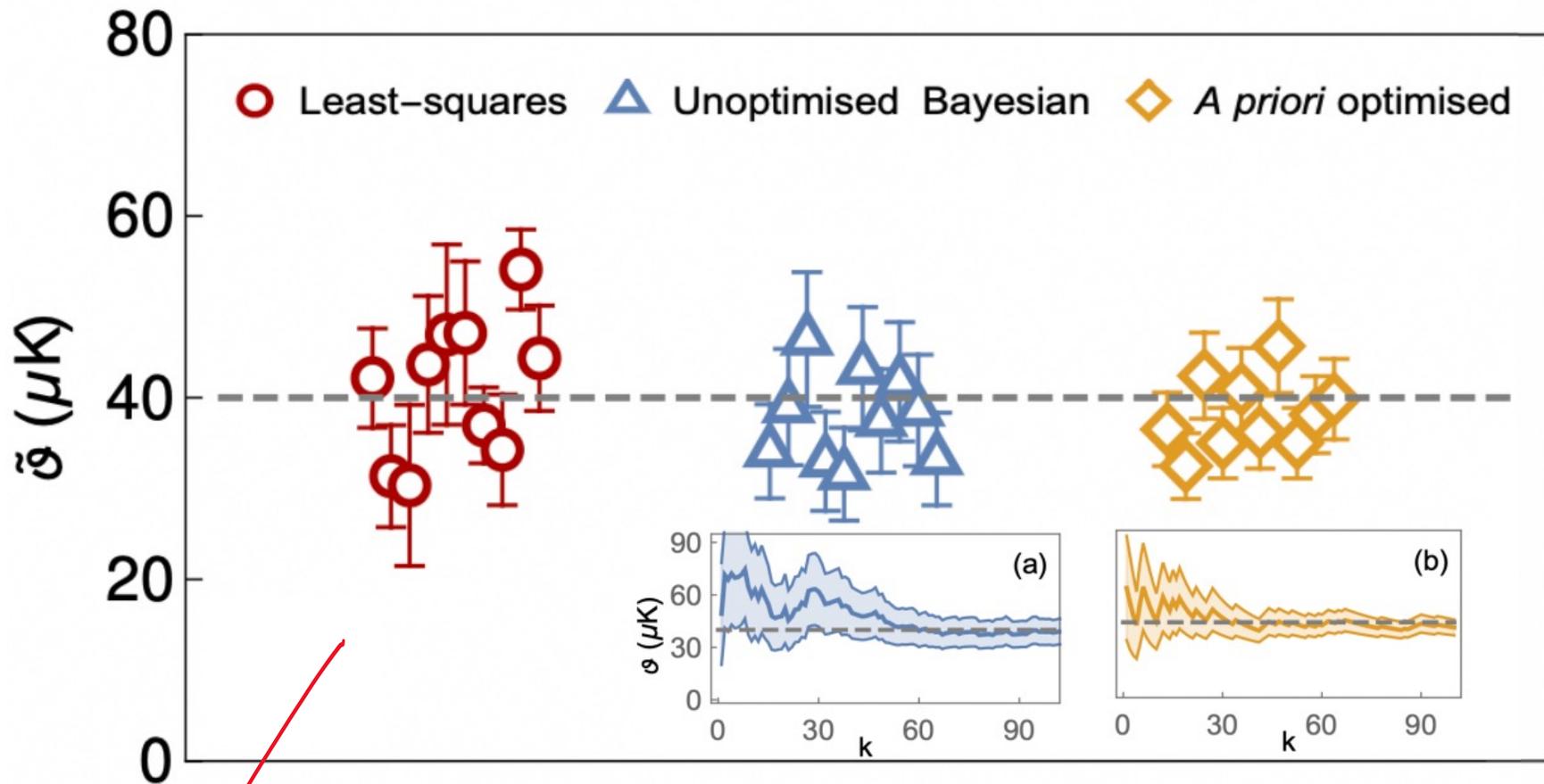
A priori optimised strategy

$$\mathbf{t} = \underbrace{(t_s, \dots, t_s)}_{\mu \text{ times}}$$

where t_s is solution to the optimisation problem:

$$\frac{dK(t)}{dt} = 0, \quad \frac{d^2K(t)}{dt^2} < 0.$$

Prescription first proposed in: *New J. Phys.* **21** 043037 2019



$$\frac{\Delta \hat{\theta}^2}{\hat{\theta}^2} = 0.064$$

$$\frac{\Delta \hat{\theta}^2}{\hat{\theta}^2} = 0.057$$

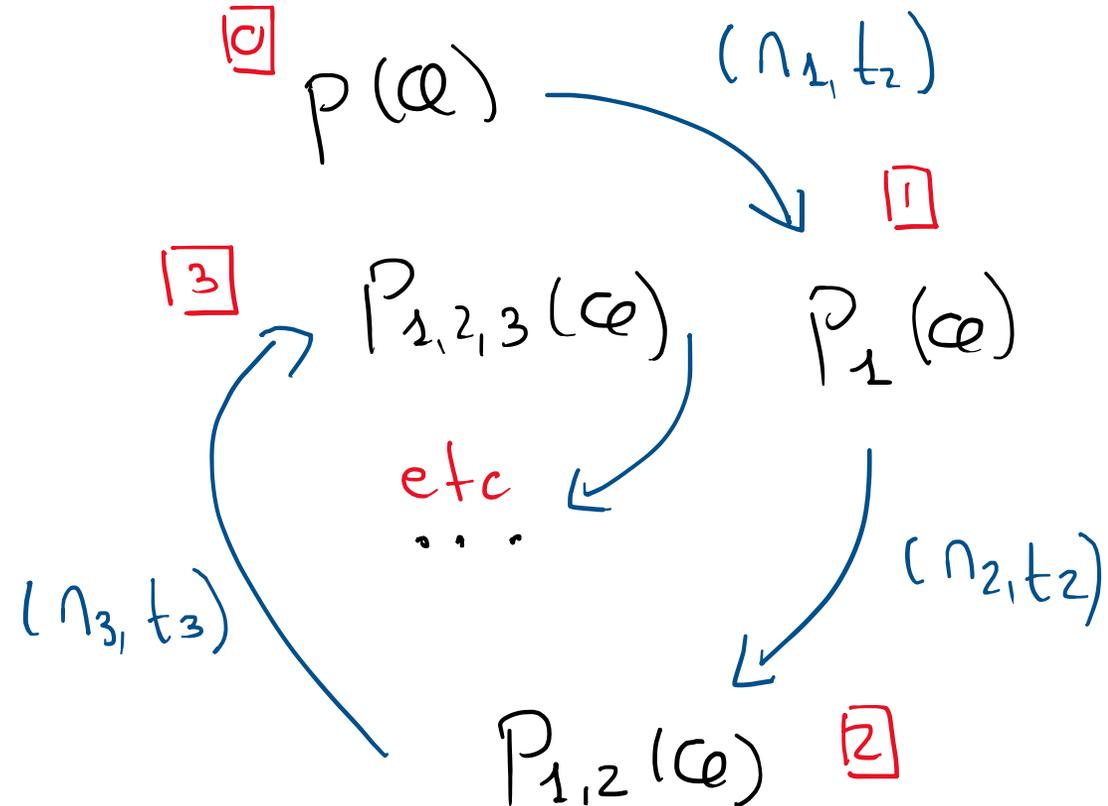
$$\frac{\Delta \hat{\theta}^2}{\hat{\theta}^2} = 0.034$$

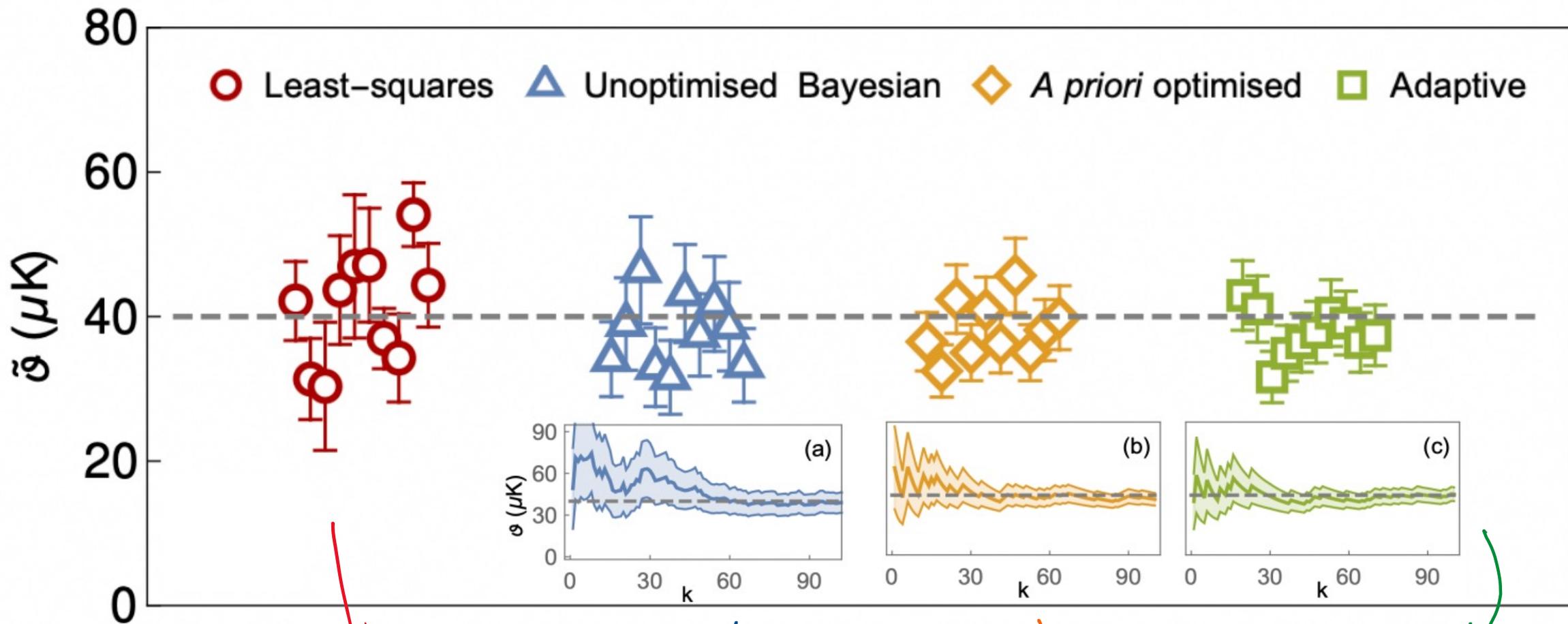
⇒

40% improvement

A fully adaptive approach

1. Given the prior $p(\theta)$ and the likelihood $p(n|\theta, t)$ for the first shot, maximise $\mathcal{K}(t)$ over t to find $t_1 = t_s$.
2. Perform a measurement at $t_1 = t_s$ and record n_1 .
3. Normalise $p(\theta) p(n_1|\theta, t_1)$ and use it as the new 'prior' for a second run [33, 34]. Then apply step 1 to find the optimal expansion time t_2 , and measure n_2 .
4. Iterate μ times. The resulting data can then be processed using Eqs. (3) and (4).





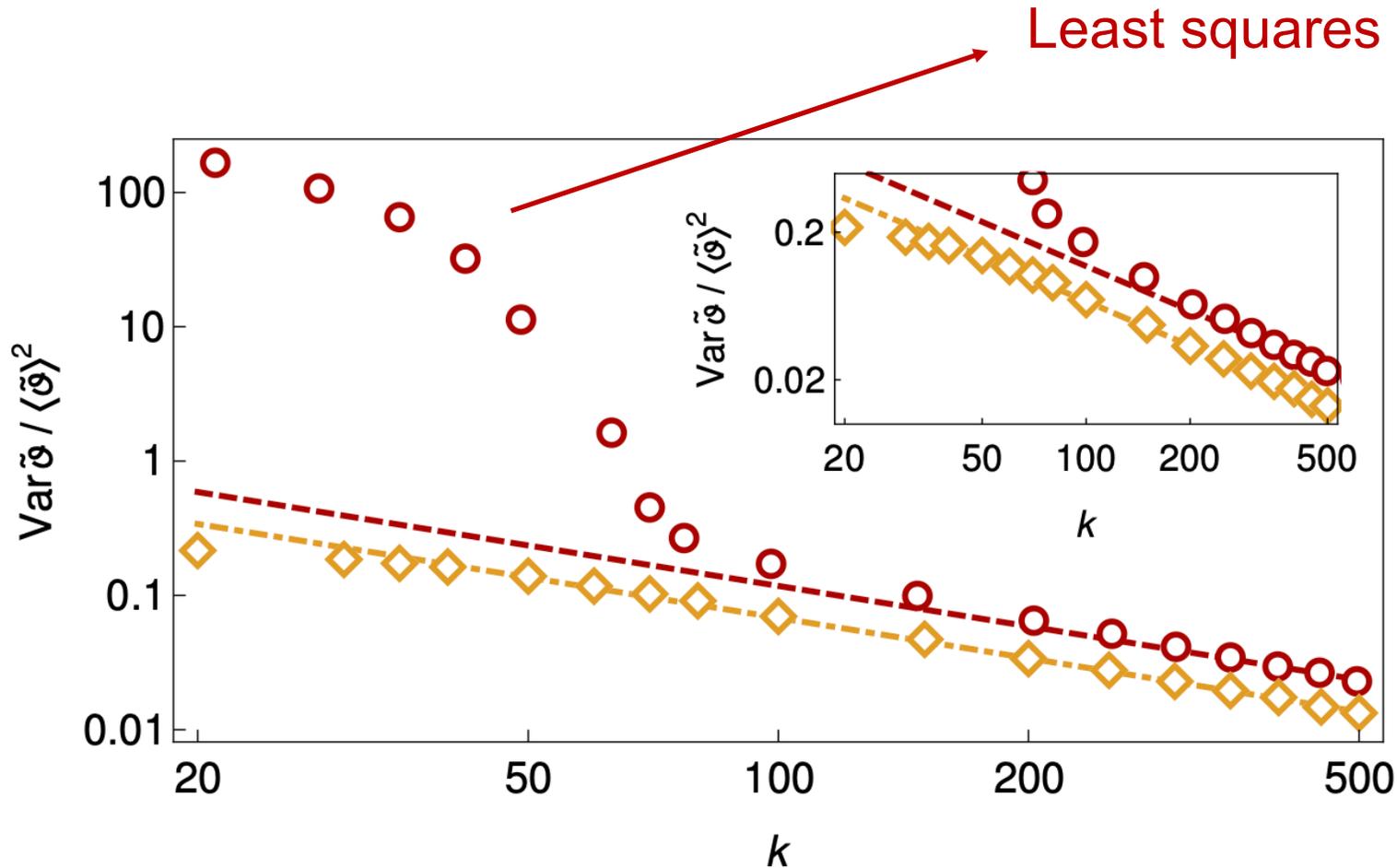
$$\frac{\Delta \tilde{\theta}^2}{\tilde{\theta}^2} = 0.064$$

$$\frac{\Delta \tilde{\theta}^2}{\tilde{\theta}^2} = 0.057$$

$$\frac{\Delta \tilde{\theta}^2}{\tilde{\theta}^2} = 0.034$$

$$\sim \frac{\Delta \tilde{\theta}^2}{\tilde{\theta}^2}$$

Precision and convergence



Least squares

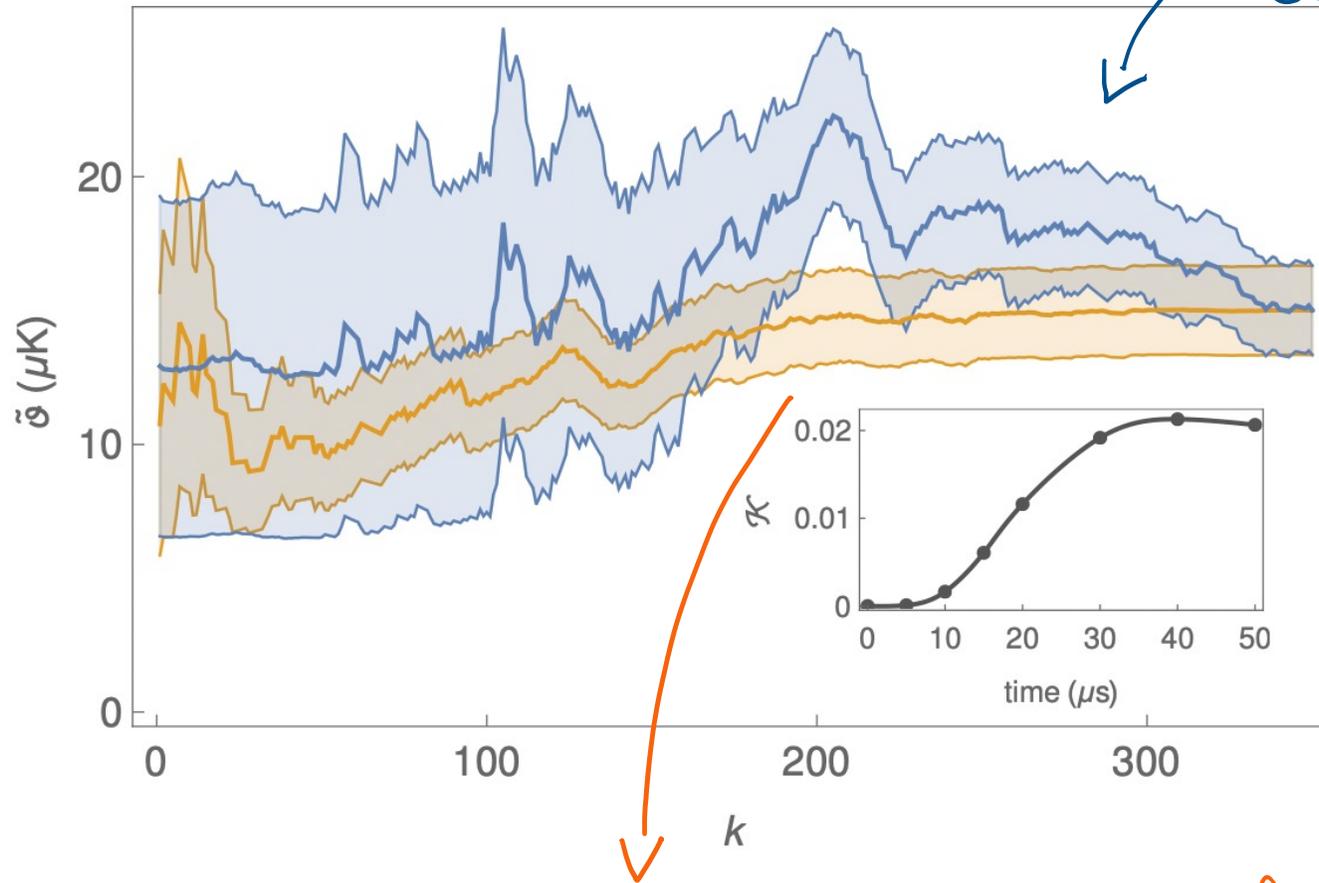
A priori optimised

Twice as much
precision with half of
the measurement data

arXiv:2204.11816

New J. Phys. **21**, 043037 (2019)

Why does it work?



mimicking the
a priori optimised
protocol with the
available
experimental
data

reordering events from most
to least informative

IV. Conclusions and outlook

Optimal cold atom thermometry using adaptive Bayesian strategies

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(Dated: 24th October 2022)

- *Standard* release-recapture thermometry is inefficient and wastes resources.
- **Global quantum thermometry can provide twice as much precision using half of the measurement data.**
- The global-Bayesian framework is applicable to *any* thermometric protocol where temperature plays the role of a scale parameter.
- Next steps: Bayesian formulation of non-equilibrium quantum thermometry under minimal assumptions.



Global Quantum Thermometry

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(Dated: 30th November 2020)

Thank you for
your attention