

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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Thomas Hewitt
& Giovanni Barontini

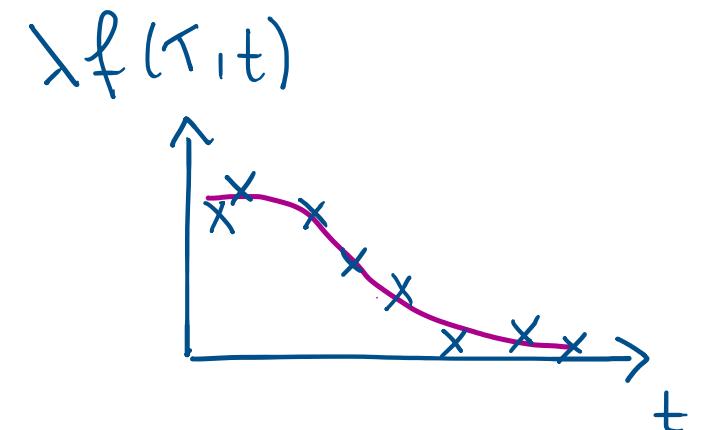


JSPS London Symposium, Nottingham

15th Dec 2022

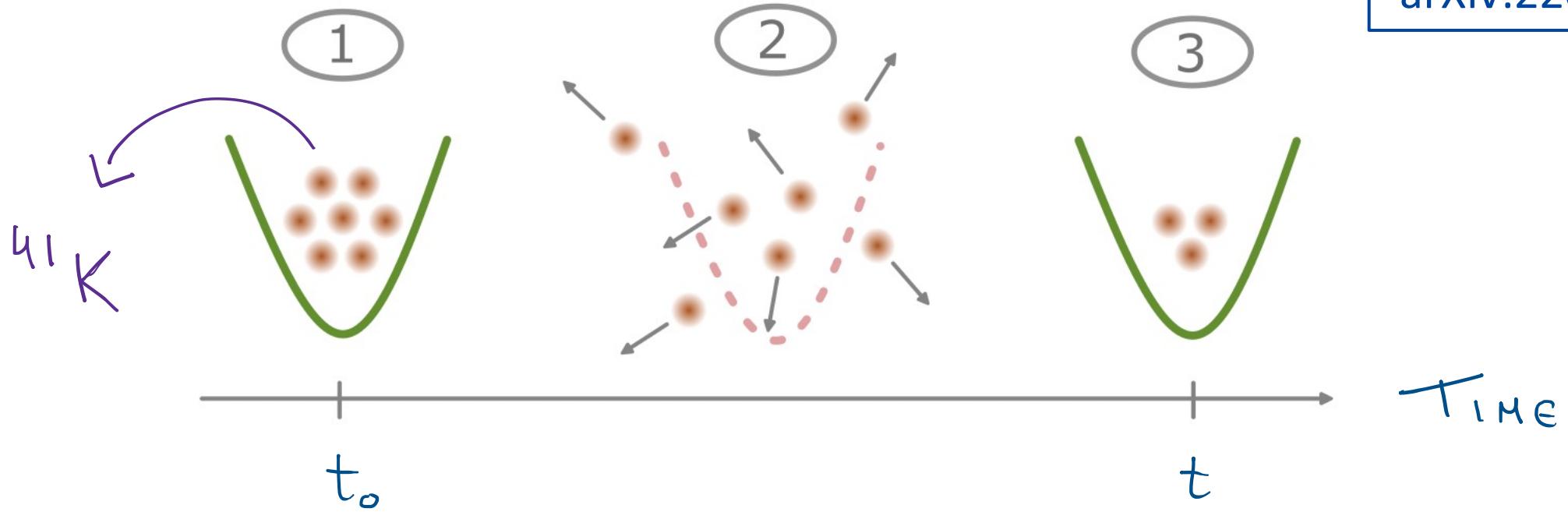
Our plan for today

- 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 $\textcircled{1}$ trials:

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

numbers of recaptured atoms

data

expansion times =

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)

$$\vec{n} = (\underbrace{n_{11}, \dots, n_{1d_1}}_{\text{data}}, \dots, \underbrace{n_{v1}, \dots, n_{vd_v}}_{\text{data}})$$

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

$$\left(\sum_{i=1}^v \alpha_i = \mu + \text{rials} \right)$$

Least squares (standard approach)

$$\begin{aligned}\langle \vec{n} \rangle &= (\langle n_1 \rangle, \dots, \langle n_v \rangle) \\ \vec{t} &= (t_1, \dots, t_v)\end{aligned}\left.\right\}$$

data

fraction of
recaptured atoms

From statistical mechanics:

$$\langle n(t) \rangle \sim f(T, t)$$

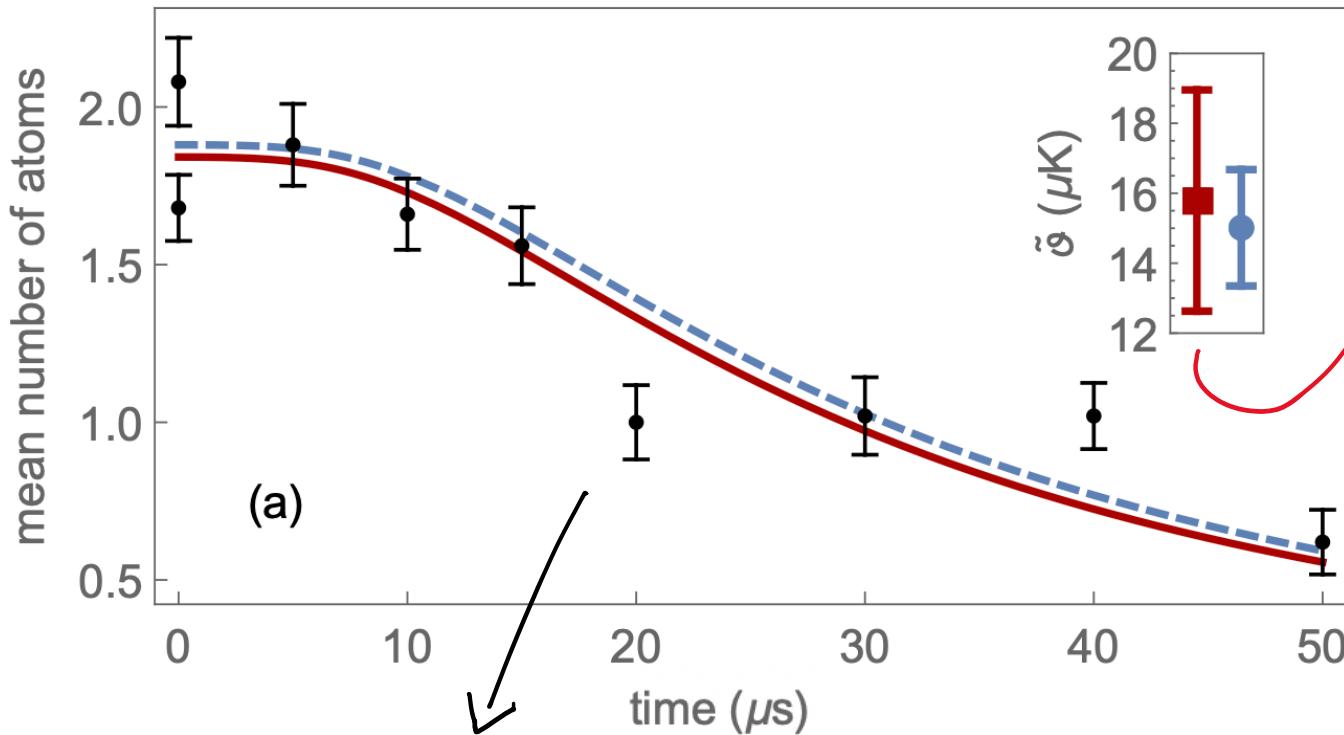
mean number
of atoms

unknown temperature

fitting
parameters

Least squares (standard approach)

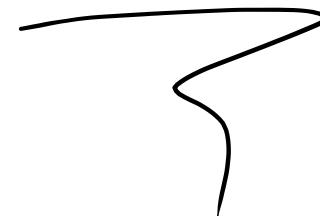
█ Fitting to $\langle n(t) \rangle \curvearrowright \propto f(T, t)$, we get:



$$\tilde{T} = 15.8 \pm 3.2 \text{ mK}$$

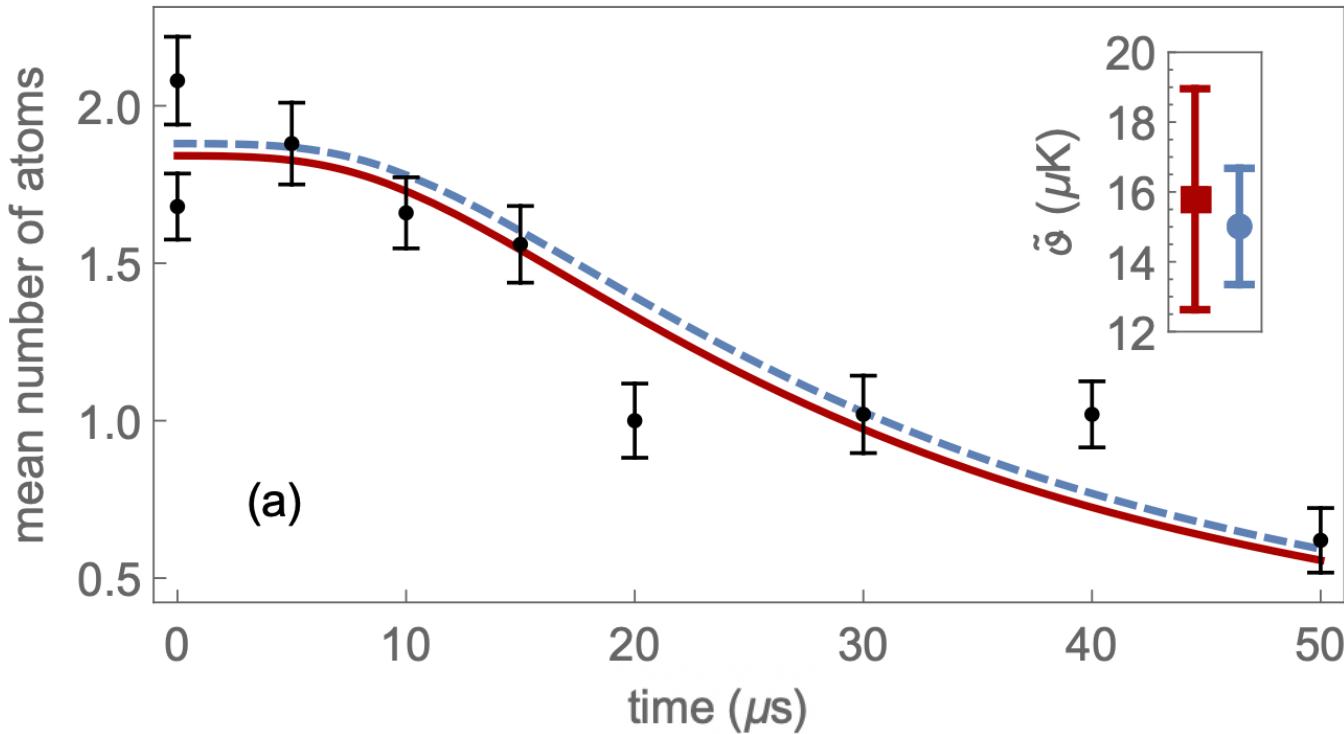
FINAL TEMPERATURE

ESTIMATE



$\langle n_i \rangle$ at time t_i

Least squares (standard approach)



$$\tilde{T} = 15.8 \pm 3.2 \text{ mK}$$

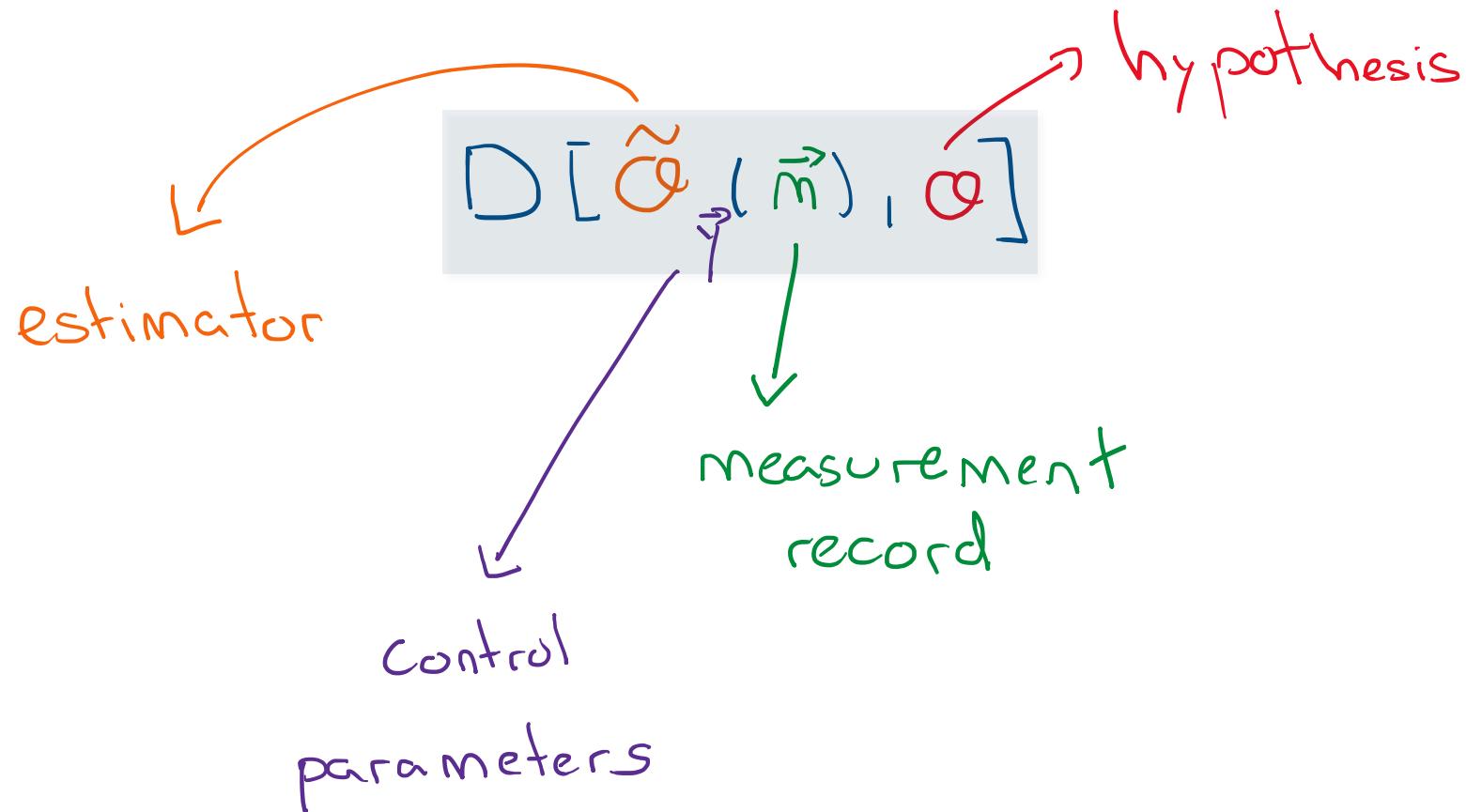
- Why another method?
 - Fails to use all available info
 - Wastes resources
- With global quantum Thermometry (this talk):

Twice as much precision with half of the measurement data

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



I Construct a deviation function D :



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



[1] Construct a deviation function $D[\tilde{\alpha}_{\vec{r}}(\vec{m}), \alpha]$

[2] Construct an error functional $\bar{E}_{\vec{r}}$:

$$\bar{E}_{\vec{r}} = \int d\alpha d\vec{m} p(\alpha, \vec{m} | \vec{r}) D[\tilde{\alpha}_{\vec{r}}(\vec{m}), \alpha]$$

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 $D[\tilde{\phi}_{\vec{\gamma}}(\vec{m}), \phi]$

[2] Construct an error functional

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

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

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

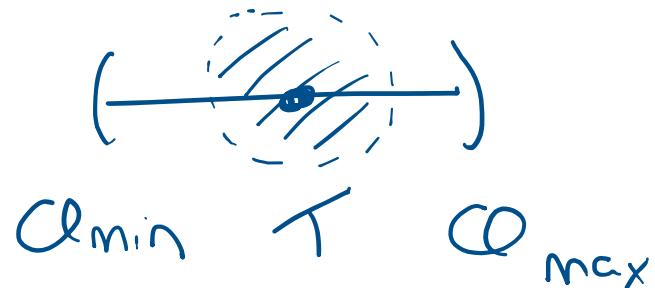
optimisation
problem

Release-recapture thermometry: unoptimised times



- Physical assumptions:

$$p(\alpha, \vec{n} | \vec{t}) = p(\alpha) p(\vec{n} | \alpha, \vec{t})$$



prior

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

$P[n_i | f(\alpha, t_i)]$

Poisson distribution

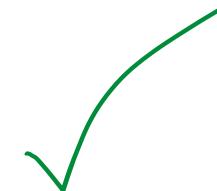
independently estimated
fraction of recaptured atoms

Release-recapture thermometry: unoptimised times



- Physical assumptions:

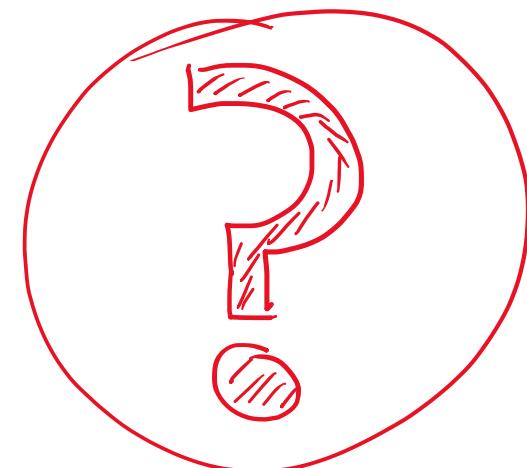
$$p(\alpha, \vec{n} | \vec{t}) = p(\alpha) \prod_{i=1}^n P[n_i | f(\alpha, t_i)]$$



- Missing:

→ prior $p(\alpha)$

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



Release-recapture thermometry: unoptimised times



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

fraction of recaptured atoms

expansion time

temperature

characteristic energy (trap parameters)

trap depth (energy)

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

Invariant under:

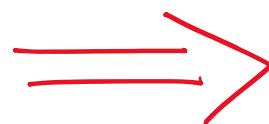
$$\begin{cases} T \rightarrow T' = \gamma T \\ U_0 \rightarrow U'_0 = \gamma U_0 \\ E_k \rightarrow E'_k = \gamma E_k \end{cases}$$

Release-recapture thermometry: unoptimised times



$$\begin{cases} T \rightarrow T' = \gamma T \\ U_0 \rightarrow U'_0 = \gamma U_0 \\ E_K \rightarrow E'_K = \gamma E_K \end{cases} \equiv$$

SCALE
INvariance

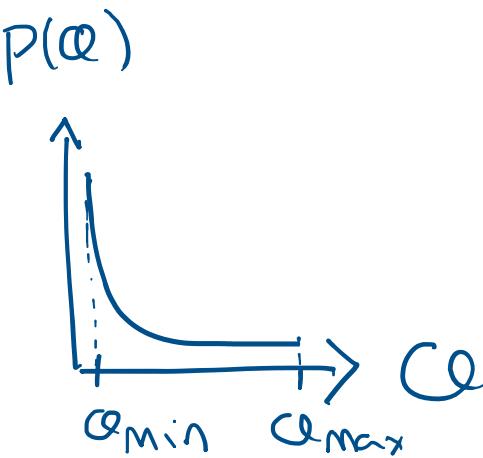


- **Jeffreys's prior:**

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

- **Logarithmic error:**

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



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:

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

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

= BAYES THEOREM

UNIVERSALLY
VALID FOR
SCALE ESTIMATION

Release-recapture thermometry: unoptimised times

How do we report temperature estimates in global quantum thermometry?

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

- Optimal **temperature estimator**:

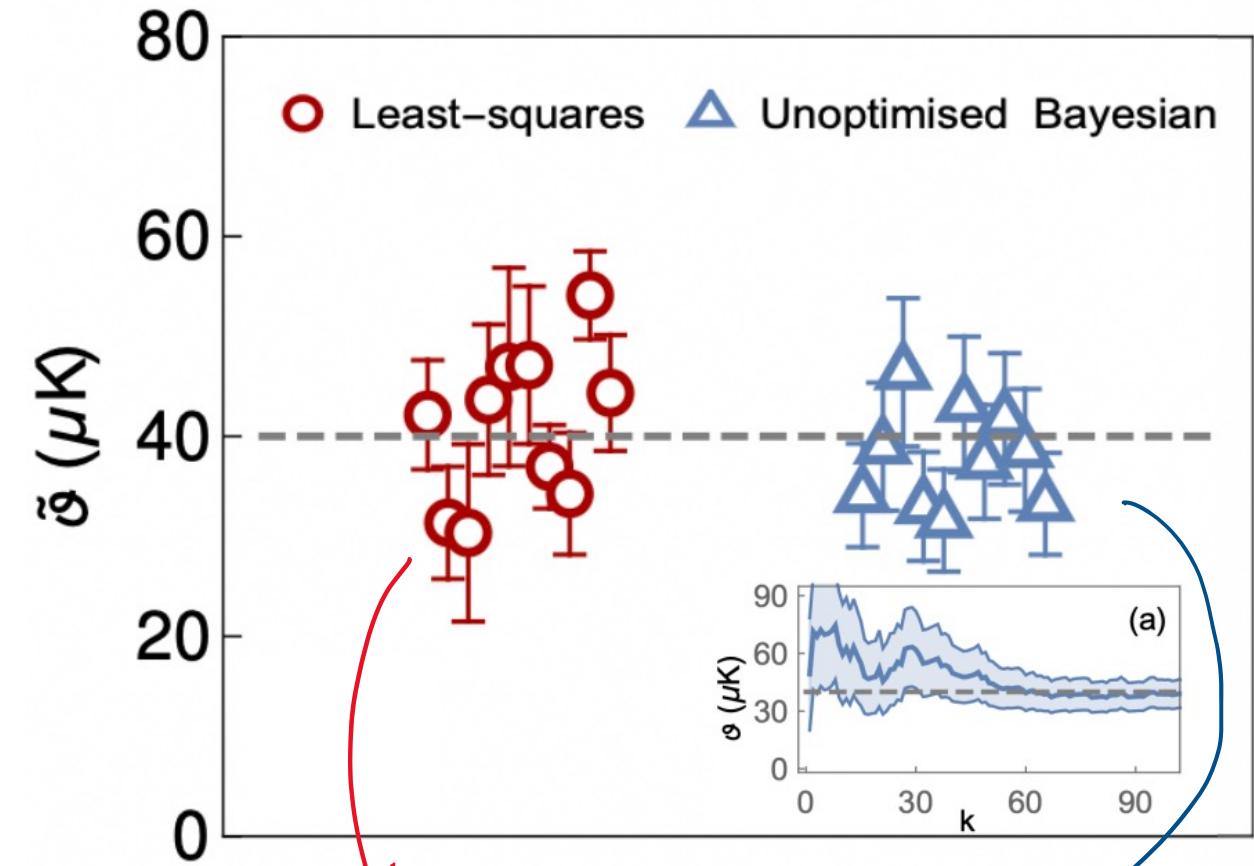
$$\tilde{\vartheta}(\mathbf{n}, \mathbf{t}) = \theta_u \exp \left[\int d\theta p(\theta | \mathbf{n}, \mathbf{t}) \log \left(\frac{\theta}{\theta_u} \right) \right]$$

- **Error bar**:

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

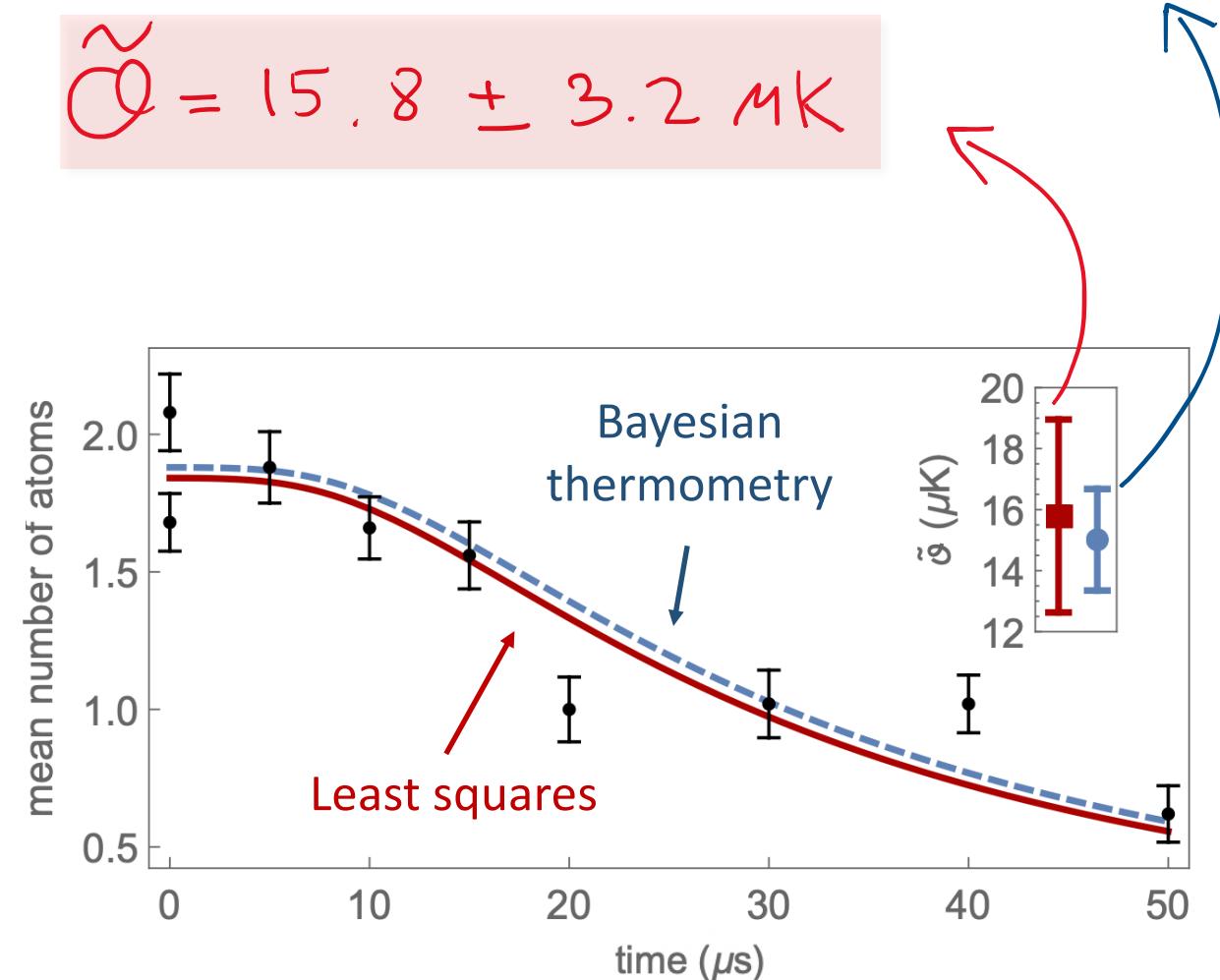
- Measurement-dependent **mean logarithmic error**:

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



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

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



III. Maximising information content in an adaptive fashion

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

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



where

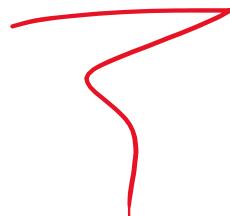
- optimal *a priori* estimate:

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

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

- evidence:

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



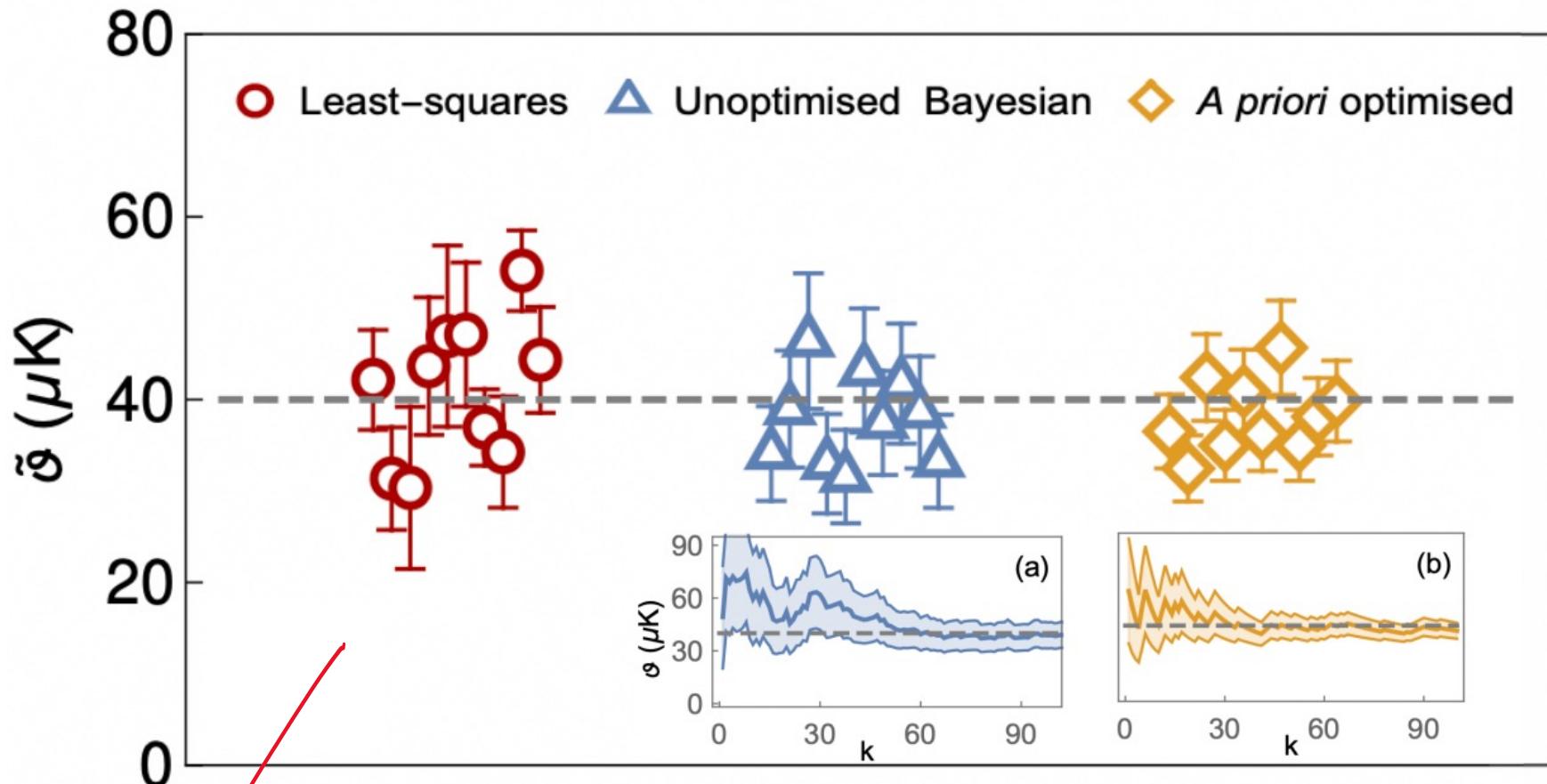
A priori optimised strategy

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

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



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

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

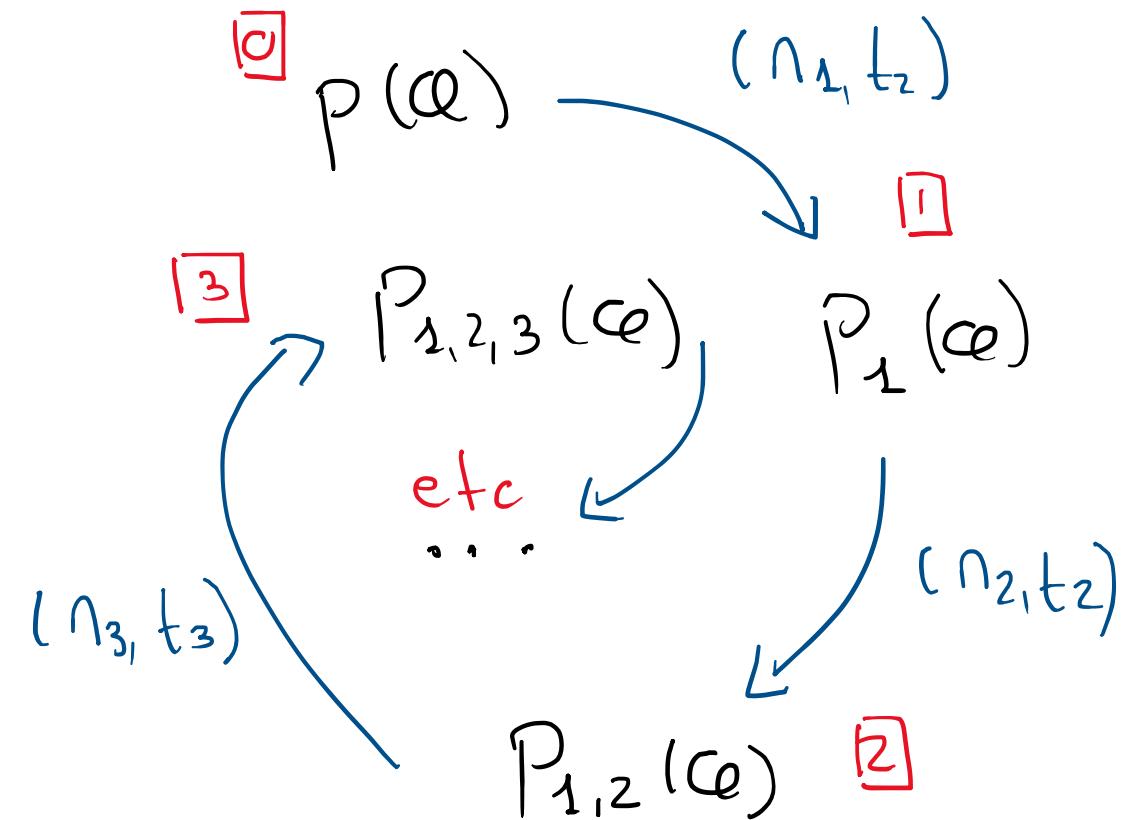
$$\frac{\Delta \tilde{\sigma}^2}{\tilde{\sigma}^2} = 0.034$$

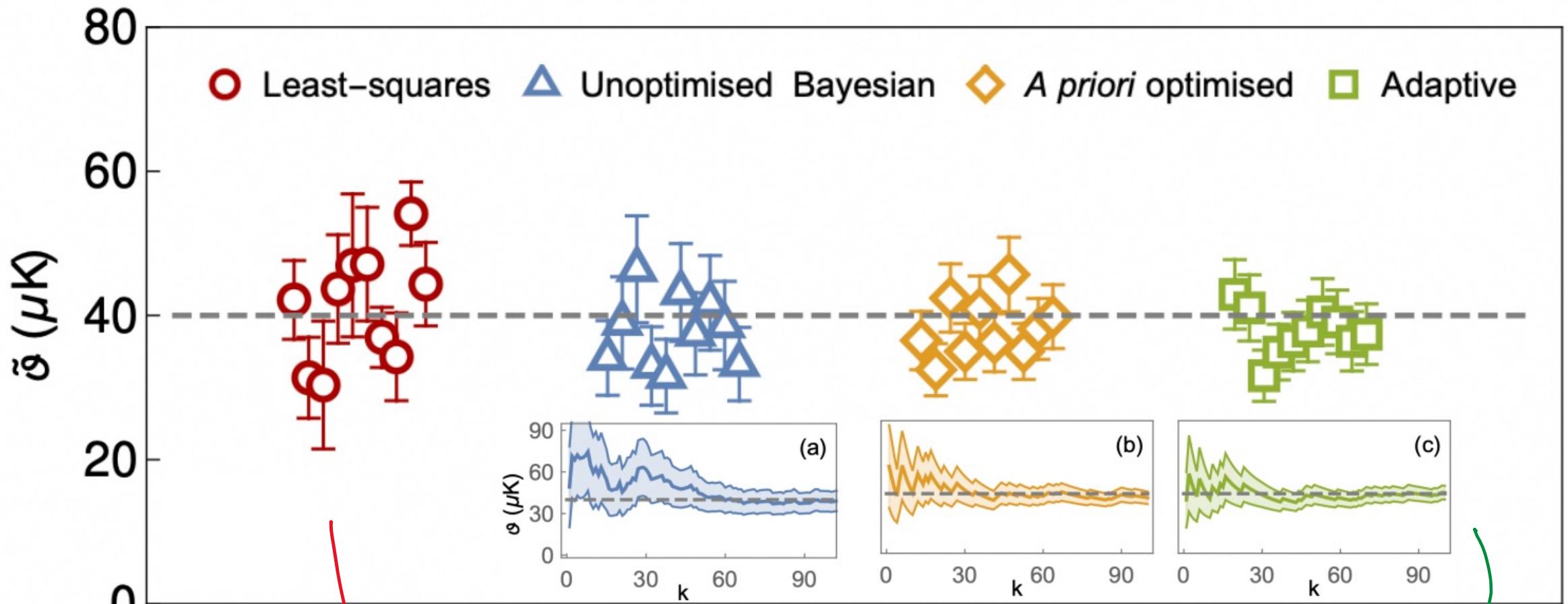
\Rightarrow

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).





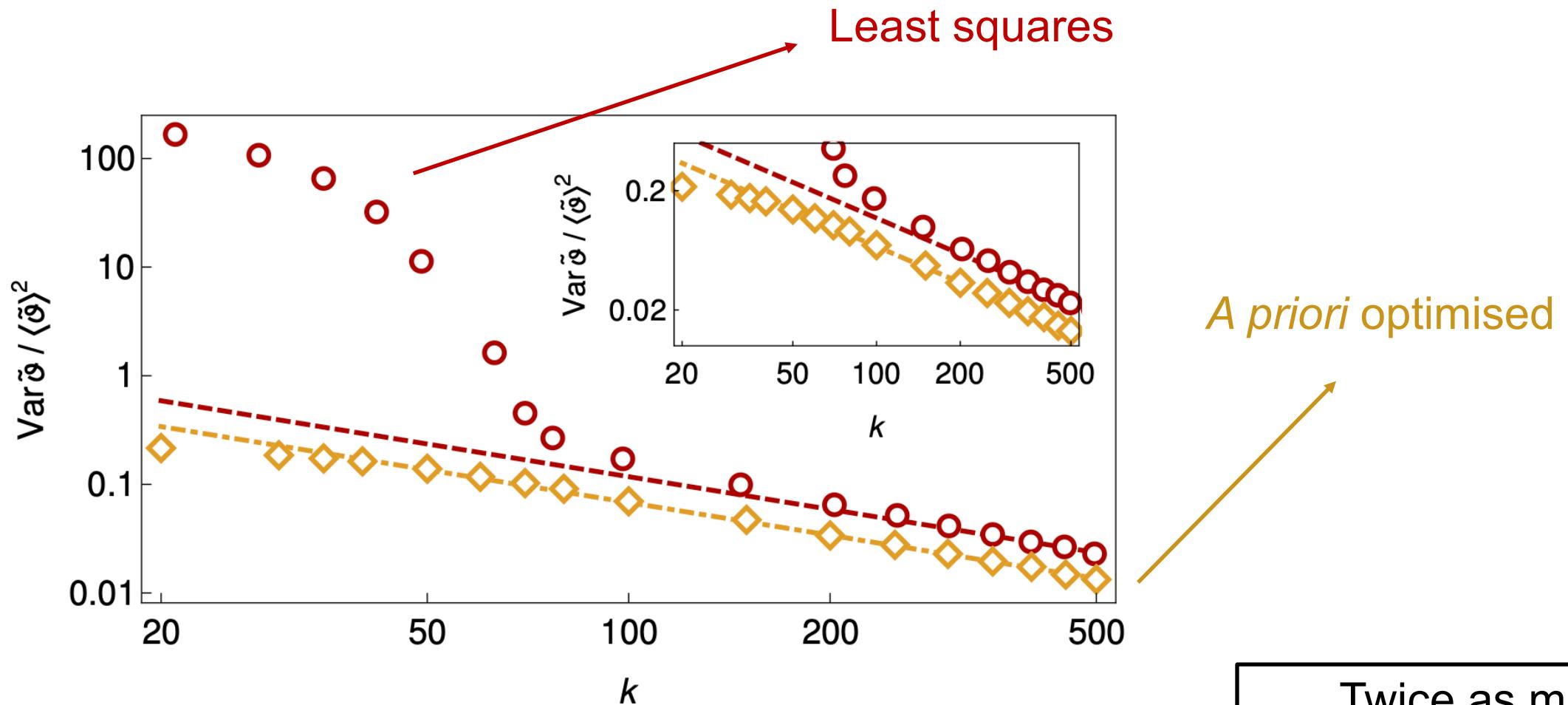
$$\frac{\Delta \tilde{\sigma}^2}{\tilde{\sigma}^2} = 0.064$$

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

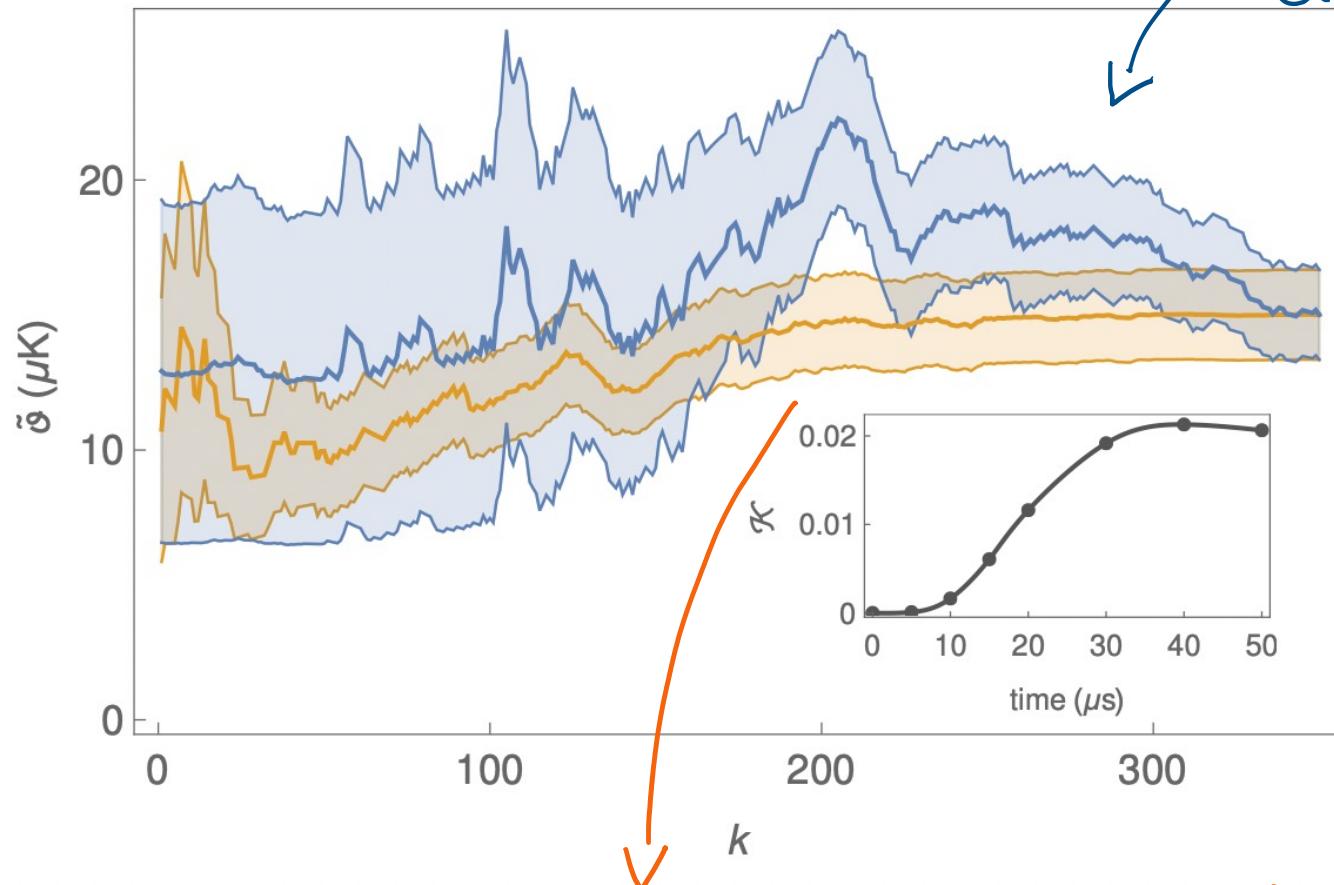
$$\frac{\Delta \tilde{\sigma}^2}{\tilde{\sigma}^2} = 0.034$$

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

Precision and convergence



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.

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PAPER

Quantum scale estimation

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Global Quantum Thermometry

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(Dated: 30th November 2020)

Thank you for
your attention