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Weak-value amplification and optimal parameter estimation in the presence of correlated noise

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We analytically and numerically investigate the performance of weak-value amplification (WVA) and related parameter estimation methods in the presence of temporally correlated noise. WVA is a special instance of a general measurement strategy that involves sorting data into separate subsets based on the outcome of a second “partitioning” measurement. Using a simplified correlated noise model that can be analyzed exactly together with optimal statistical estimators, we compare WVA to a conventional measurement method. We find that WVA indeed yields a much lower variance of the parameter of interest than the conventional technique does, optimized in the absence of any partitioning measurements. In contrast, a statistically optimal analysis that employs partitioning measurements, incorporating all partitioned results and their known correlations, is found to yield an improvement—typically slight—over the noise reduction achieved by WVA. This result occurs because the simple WVA technique is not tailored to any specific noise environment and therefore does not make use of correlations between the different partitions. We also compare WVA to traditional background subtraction, a familiar technique where measurement outcomes are partitioned to eliminate unknown offsets or errors in calibration. Surprisingly, for the cases we consider, background subtraction turns out to be a special case of the optimal partitioning approach, possessing a similar typically slight advantage over WVA. These results give deeper insight into the role of partitioning measurements (with or without postselection) in enhancing measurement precision, which some have found puzzling. They also resolve previously made conflicting claims about the usefulness of weak-value amplification to precision measurement in the presence of correlated noise. We finish by presenting numerical results to model a more realistic laboratory situation of time-decaying correlations, showing that our conclusions hold for a wide range of statistical models.

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I. INTRODUCTION

Weak-value amplification (WVA) [1] is a technique that has been used in a variety of experimental settings to permit the precise measurement of small parameters [2–22]. Over the past few years, WVA has been the subject of an ongoing debate over whether or not it can provide an actual advantage in terms of the resulting measurement precision [20,23–29]. Experimental scenarios with temporally correlated noise are one area where there have been claims about an advantage for WVA. In this paper, we delineate the situations under which WVA does, in fact, improve measurement precision in the presence of correlated noise, and we compare it with competing approaches. This investigation leads to interesting connections between WVA and other techniques, such as background subtraction and lock-in amplification, which elucidate the technical advantages that have already been experimentally observed.

Alongside arguments regarding the usefulness of WVA to precision measurement, there has continued to be a great deal of work extending and improving the basic technique of WVA in other situations. Some recent advances in the field include the incorporation of photon recycling of discarded events [30–32], the observation that WVA can improve measurement precision in cases with detector saturation [33], the optimization of the shape of the meter probe [34], and a generalized approach to probabilistic quantum metrology [35]. Of particular interest is weak-value amplification with entanglement [29,36], squeezing [37], and the observation that weak-value amplification can suppress systematic errors [38], which is closely related to the present work. We refer the reader to recent reviews for a wider overview of the field [39,40].

Weak-value amplification is an experimental technique that involves two measurements. In the first, a system observable is measured via a weak interaction with a measurement apparatus. The effect of this weak interaction is to induce a small shift in the pointer of the measurement apparatus. The size of this shift is determined by the observable (which is typically known) and the coupling strength, which we are interested in estimating. In weak-value amplification this coupling strength is usually very weak, so that very little information is gained about the state being measured, and the corresponding measurement disturbance is minimized. The second measurement is a strong projective measurement in a different basis on the system. Because the system and measurement device are left weakly entangled by the first measurement, there are interesting correlations between the two systems. These correlations can be seen by repeating the two-measurement procedure identically many times and dividing the data set compiled from the outcomes of the first measurement into partitions based on the results of the second projective measurement, and then averaging the different partition with different weights. In conventional WVA, the weights are 0 and 1. This corresponds to discarding certain
measurement outcomes based on the result of the second projective measurement, and is called postselection. The mean shift of the pointer conditioned upon the postselection succeeding is proportional to the weak value. It is defined in terms of the initial state of the system (|i⟩), the state the system is found to be in if the postselection succeeds, ⟨|f⟩⟩, and the operator associated with the first measurement, \( \mathbf{A} \),

\[
A_w = \frac{\langle f | \mathbf{A} | i \rangle}{\langle f | i \rangle}.
\]

The weak value can become quite large when the overlap of the initial and final state becomes very small. This “amplification” of what would be a small shift in our measurement pointer seems to open up the possibility for improving the measurement precision. The catch is that when the initial and final state have small overlap, the postselection occurs very rarely and so the retained data set becomes quite small. The improvement in precision which might be expected from an amplified signal is in fact often perfectly canceled by the reduction in the size of the data set in experiments where the measurements are uncorrelated. In spite of this fact, there have been several experimental metrological works, such as Refs. [2,3], which argued that WVA could offer an advantage in specific experimental situations. These claims ignited a fierce debate over proper resource counting and whether or not the advantage reported was a feature of the experimental design or of the technique itself. In this paper, we focus on the case of additive correlated noise, and detail how WVA can yield an advantage over other common experimental techniques.

This topic was raised in 2011, when Feizpour, Xing, and Steinberg (FXS) studied WVA in the presence of additive, time-correlated, Gaussian random noise and claimed it yielded an improvement in the signal-to-noise ratio of the measured parameter over other common techniques [41]. In the scenario considered by FXS, measurements are performed sequentially in time under equivalent circumstances, but are not independent due to temporal correlations in the noise. In contrast in WVA, postselection events naturally occur more rarely, helping the postselected data set remain uncorrelated. WVA is therefore robust to the detrimental effects of time-decaying temporal correlations and, FXS claim, superior to other approaches. More recently, however, others have argued using Fisher Information methods that the WVA method is inherently suboptimal because it involves discarding a portion of the measurement outcomes [24]. Using the same Fisher information based approach, WVA has since been shown to be capable of capturing nearly all of the information, becoming asymptotically optimal. The optimal limit is possible to reach because the information contained in the discarded measurement outcomes is a tiny fraction of the total information (despite the discarded outcomes making up the vast majority of the total number of outcomes) [20,27–29,42].

The case with correlated noise is more complex, however, because the correlations within the postselected partition are sensitive to controlled parameters like the probability of postselection, and because there are correlations (and therefore information) between the postselected and nonpostselected partitions which must be considered. We will revisit the case of correlated noise and show that while WVA is much better than the conventional approach, in the temporally correlated noise case it is (typically) slightly inferior to a statistical analysis which optimally utilizes all partitioned outcomes and the correlations between them. See Fig. 1, for a schematic representation of the three measurement strategies. We will also show that WVA and the optimal partitioning approach are closely related to background subtraction and lock-in amplification techniques which are well known.

This paper is organized as follows. In Sec. II, we introduce the tools used in estimation theory: estimators; variance; the Cramér-Rao Bound; our metric of choice, the Fisher information; and a model for correlated Gaussian noise. In Sec. III, we explore the Fisher information as an information metric using a simple two-measurement-outcome example. In Sec. IV, we give an eigenvalue analysis of the Fisher information, and show that it may be expressed as a weighted average of the eigenvalues of the covariance matrix. The weak-value amplification effect is introduced in Sec. V. An optimal partitioning measurement approach is introduced in Sec. VI, which improves slightly on the advantage achieved by WVA over the direct method by including all partitions and the correlations between them. This physics is illustrated in Sec. VII, where an exactly solvable model is introduced, and the variance of all the estimation strategies is given explicitly and compared. A numerical investigation of these issues is presented in Sec. IX, where an experimentally motivated correlated noise model is given, and analyzed. Our conclusions are given in Sec. X.

II. PARAMETER ESTIMATION

Consider a common scenario in the natural sciences in which the goal is the measurement of some unknown parameter. If the measurement is noisy, this can be accomplished by performing the same measurement procedure many times
under equivalent circumstances. Each measurement yields an outcome, which is compiled into a data set. We represent this data set as \( \{s_i\} \), where \( i \) identifies the \( i \)th measurement, \( i = 1, 2, \ldots, N \). Measurement outcomes can be scaled and shifted so the measured signal \( s_i \) can be modeled as the parameter of interest \( d \) and noise \( x_i \),

\[
s_i = d + x_i,
\]

where \( x_i \) are zero mean Gaussian random variables [41]. In the most generic case, these variables may be correlated. For example, this could occur if the source of noise possessed temporal correlations which lasted for several measurements. We define the correlation function \( C_{i,j}(x) = \langle x_i x_j \rangle \)—this is also called the covariance matrix. Our task is to estimate the unknown parameter \( d \) from the data set and our knowledge of the covariance matrix. In this paper, we consider the covariance matrix known, and the detailed knowledge able to be applied to implement the optimal estimator of the unknown quantity \( d \). The extra resources required to estimate this matrix and implement the optimal estimator are not considered [27], but this would amount to completely characterizing the sources of noise in our experiment. By choice of estimator, we simply mean some algorithm which maps our data set to an estimate of \( d \). Taking the arithmetic average corresponds to the most conventional and straightforward estimator. Because the noise has zero mean, this estimator is unbiased, which means that in the limit of \( N \to \infty \), the estimate of the parameter converges to \( d \). There are many other unbiased estimators which could be constructed depending on how the measurement is designed and implemented.

We will discuss two different classes of measurement design. The first class, which most conventional measurements fall into, we term “direct.” Direct measurements only involve measurements on the parameter of interest. The second class, which we term “partitioning,” introduces a second measurement that is used to sort the first measurement’s outcomes into different partitions. Often, partitioning class measurements possess some advantage versus direct measurements, because they can exploit correlations between different partitions. The simplest partitioning class measurement involves discarding all data points which fail to meet some criterion assessed by the second measurement. This is postselection. While others have pointed out that instead of discarding outcomes, they should be weighted and optimally analyzed, our assertion is just that exploiting such correlations may have significant advantages, and that while the actual throwing out of data per se can never increase the amount of information available, there may be regimes where some (perhaps most) of the advantage survives even this procrustean approach.

Once the “measurement design” has been specified, the space of available estimators is infinite. As the variance associated with different estimators can vary wildly, the “choice of estimator” is highly nontrivial. Conveniently, a technique called maximum likelihood estimation [27] can be used to identify the estimator with minimum variance in the large data set limit. Furthermore, there is a mathematical theorem [43] that bounds the minimum variance of an unbiased estimator to be larger than the inverse Fisher information. This allows us to forgo discussion of estimators entirely, skipping directly to the calculation of the Fisher information, which is defined as the inverse variance of the optimal estimator. Find the Fisher information for a given measurement scenario, invert it, and one will have found the minimum variance of any unbiased estimator.

### III. Fisher Information in a Two Measurement Outcome Example

It is instructive to consider a pedagogical example that will clarify the meaning of the Fisher information, the importance of the choice of estimator, and the significance of the covariance matrix to this task. If we return to the scenario in the previous section, but imagine collecting only two data points, we will have a data set \( \{s_1, s_2\} \) with assumed knowledge of the covariances and variances. As before, our goal is to estimate the mean \( d \). The covariance matrix, which is a \( 2 \times 2 \) matrix, is

\[
C(x) = \begin{pmatrix} \text{Var}(s_1) & \text{Cov}(s_1,s_2) \\ \text{Cov}(s_2,s_1) & \text{Var}(s_2) \end{pmatrix},
\]

where \( \text{Var}(s_1), \text{Var}(s_2) \geq 0 \), and \( \text{Cov}(s_1,s_2) = \text{Cov}(s_2,s_1) \) are all real numbers. The Cauchy-Schwartz inequality \( \text{Cov}(s_1,s_2)^2 < \text{Var}(s_1) \text{Var}(s_2) \) guarantees both eigenvalues be non-negative which also guarantees \( C \) to be positive semidefinite. The \( \text{Cov}(s_1,s_2) \) itself can be positive or negative, representing either positive or negative correlation between \( s_1 \) and \( s_2 \). Our goal is to estimate the expectation value of our two data points while minimizing the variance of our estimate. Were we to use the first data point alone to estimate the mean, the variance would be \( \text{Var}(s_1) \); similarly, use of the second data point alone gives \( \text{Var}(s_2) \). Instead, we can use a combination of the two data points as our estimator, \( s_{\text{est}} = \alpha s_1 + \beta s_2 \), where \( \alpha \) and \( \beta \) are constant weighting factors. To keep this estimator unbiased we require that \( \alpha + \beta = 1 \). The variance of our estimator \( s_{\text{est}} \) is given by

\[
\text{Var}(s_{\text{est}}) = \alpha^2 \text{Var}(s_1) + \beta^2 \text{Var}(s_2) + 2\alpha\beta \text{Cov}(s_1,s_2).
\]

We minimize the variance with respect to \( \alpha \), keeping \( \text{Var}(s_1), \text{Var}(s_2), \text{Cov}(s_1,s_2) \) fixed. Doing so gives a minimum variance of

\[
\text{Var}(s_{\text{est}})_{\text{min}} = \frac{\text{Var}(s_1)\text{Var}(s_2) - \text{Cov}(s_1,s_2)^2}{\text{Var}(s_1) + \text{Var}(s_2) - 2\text{Cov}(s_1,s_2)},
\]

which is smaller than or equal to either \( \text{Var}(s_1) \) or \( \text{Var}(s_2) \) for any allowed values of \( \text{Var}(s_1), \text{Var}(s_2), \text{Cov}(s_1,s_2) \). We note that unlike uncorrelated random variables, the inverse variance is not additive. If instead we took equal weighting of the two data points, \( \alpha = \beta = 1/2 \), this would give

\[
\text{Var}(s_{\text{est}})_{\text{equal}} = \frac{\text{Cov}(s_1,s_2)}{2} + [\text{Var}(s_1) + \text{Var}(s_2)]/4.
\]

If the two outcomes are perfectly negatively correlated \([\text{Cov}(s_1,s_2) \rightarrow -\sqrt{\text{Var}(s_1)\text{Var}(s_2)}]\), then both the optimal and the equal weighting estimators have zero variance. This can be understood as resulting from anticorrelated fluctuations canceling each other out, for example if \( s_1 = d + x_1 \) and \( s_2 = d - x_1 \), a straightforward averaging of \( s_1, s_2 \) will result in the perfect cancellation of the noise \( (x_1) \). If the two outcomes are perfectly positively correlated, but the variances are not equal \([\text{Var}(s_1) \neq \text{Var}(s_2)]\), the optimal variance vanishes, whereas the equal weighting variance limits to that of using just a single outcome. This can be understood by considering the
case where $s_1 = d + 2x_1$, and $s_2 = d + x_1$. If our estimator is $s_{\text{est}} = 2s_2 - s_1$, we can eliminate the noise just like in the anticorrelated case. When the variances are equal and the correlations are positive, the optimal variance does not vanish because the noise cannot be canceled without making the estimator biased. In this case, the equal-weighting estimator turns out to be optimal. We recall that the Fisher information is equal to the inverse of the variance of the optimal estimator in the large data limit. The Fisher information is defined for smooth distributions as

$$I = -\int ds_1 \ldots ds_n P(s_1, s_2, \ldots, s_n|d) \frac{\partial^2}{\partial d^2} \ln P,$$

(7)

where $P = P(s_1, s_2, \ldots, s_n|d)$ is the probability distribution of $\{s_1, \ldots, s_n\}$ given a fixed value of $d$. In our model, it is taken as a multidimensional Gaussian distribution with mean $d\mathbf{1}$ and covariance matrix $C_{ij}$. From Ref. [27] the Fisher information about the mean for Gaussian correlated noise is given by

$$I = \sum_{i,j}^N [C^{-1}]_{i,j},$$

(8)

and the Fisher information for $N = 2$ is simply the minimized variance we found previously,

$$I = \frac{\text{Var}(s_1) + \text{Var}(s_2) - 2\text{Cov}(s_1, s_2)}{\text{Var}(s_1)\text{Var}(s_2) - \text{Cov}(s_1, s_2)^2} = \text{Var}(s)^{-1}. \quad (9)$$

It is convenient to introduce two parameters, an asymmetry parameter, $x = \text{Var}(s_1)/\text{Var}(s_2) \in [0, \infty)$, and a relative correlation parameter, $r = \text{Cov}(s_1, s_2)/\sqrt{\text{Var}(s_1)\text{Var}(s_2)} \in [-1, 1]$. Dividing the inverse Fisher information by $\sqrt{\text{Var}(s_1)\text{Var}(s_2)}$ gives a function that depends only on $r$ and $x$. We plot the inverse Fisher information in this case in Fig. 2, noting the asymmetry in both $r$ and $x$ which we will now explore.

We note that (9) indicates that negative values of $\text{Cov}(s_1, s_2)$ (or $r$) typically have higher Fisher information than positive values; that is, anticorrelation is more informative than correlation. In Figs. 3–5, we plot the variance of $s_{\text{est}}$ versus $\alpha$ for different values of $x$ and $r$. The minimum value of the estimator corresponds to the inverse Fisher information for that choice of covariance matrix. These figures highlight how the information that can be extracted from a probability distribution depends in a complicated way on the parameters of that distribution even in the simplest (two-dimensional) cases. Importantly, they show that while anticorrelations typically increase the available information, positive correlations can in certain circumstances also boost the amount of information that is available.

In what follows, we will show that due to the properties of the Covariance matrix the Fisher information can be expressed simply in terms of the eigenvalues and eigenvectors of $C$. This will pave the way for an exactly solvable noise model, which we will use to address the questions raised in Sec. I by comparing the Fisher information of the various measurement strategies. Finally, we will present a numerical investigation where the conclusions reached with the exactly solvable noise model are shown to hold in experimentally realistic scenarios.
obtain orthogonal decomposition of it as follows: positive, semidefinite matrix. Therefore, we can make an is a diagonal matrix with eigenvalues between the outcomes (symmetric case (14)). We define a vector \( \mathbf{v} \) with components, \( v_k = \sum_i O_{i,k} \). This then gives

\[
\mathcal{I} = \sum_k \sigma_k^{-2} v_k^2. \tag{13}
\]

Next we note that the \( O \) matrix is orthogonal, and therefore \( O^T = I \), the identity matrix. Summing over both indices we obtain

\[
\sum_{i,j,k} O_{i,k} O_{i,j}^T = \sum_k v_k^2 = \sum_{i,j} \delta_{i,k} = N. \tag{14}
\]

Since the sum of the squares of the \( v \) vector components must be \( N \) by orthogonality, we define a weight vector \( \mathbf{w} \), whose elements are \( w_i = v_i^2 / N \), so that the sum of the weights is 1. With this definition, the Fisher information is

\[
\mathcal{I} = N \sum_k \sigma_k^{-2} w_k. \tag{15}
\]

Thus, the Fisher information is given by the weighted average of the eigenvalues of the covariance matrix, with the weights related to the eigenvectors of covariance matrix.

According to the Cramér-Rao inequality (CRI), the variance of any unbiased estimator \( d_{\text{est}} \) must be greater than the inverse Fisher information [43],

\[
\text{Var}[d_{\text{est}}] \geq \mathcal{I}^{-1}. \tag{16}
\]

Let us see if this is true in the case of the simple estimator \( d_{\text{est}} = (1/N) \sum_j s_j \). The variance is given by

\[
\text{Var}[d_{\text{est}}] = (1/N^2) \sum_{i,j} C_{i,j}. \tag{17}
\]

We can make a similar analysis as above, \( C = ODO^T \), to find

\[
\text{Var}[\hat{d}] = (1/N^2) \sum_{i,j,k} \sigma_i^2 O_{i,k} O_{j,k}. \tag{18}
\]

We can rewrite this as

\[
\text{Var}[d_{\text{est}}] = (1/N) \sum_k \sigma_k^2 w_k. \tag{19}
\]

The CRI can be restated in this case as the inequality

\[
\sum_i w_i \sigma_i^2 \sum_j w_j \sigma_j^{-2} \geq 1. \tag{20}
\]

This relation can be proved directly with the Cauchy-Schwarz inequality. We define a vector \( \mathbf{u}_1 \) of dimension \( N \) with elements \( u_{1,i} = \sqrt{w_i} \sigma_i \), and another vector \( \mathbf{u}_2 \) of the same dimension, with elements \( u_{2,i} = \sqrt{w_i} \sigma_i^{-1} \). The Cauchy-Schwarz inequality,

\[
|\mathbf{u}_1 \cdot \mathbf{u}_2| \leq ||\mathbf{u}_1|| \cdot ||\mathbf{u}_2||, \tag{21}
\]

applied to these vectors gives \( \sum_i w_i \sigma_i / \sigma_i = 1 \) for the left-hand side of (21), since \( w_i \) are weighting factors, and \( \sum_i w_i \sigma_i^2 \sum_j w_j \sigma_j^{-1} \) for the right-hand side of (21). If a square root of a quantity is greater than 1, that quantity is greater than 1 as well, establishing the desired relation (20).

We note that a sufficient condition on \( C \) to make the simple estimator efficient is that \( C \)’s rows (or columns) all sum to the same number. This is equivalent to the statement that \( C \) has an eigenvector \( \mathbf{e}_1 = (1,1, \ldots ,1)^T \). By construction, all other eigenvectors are orthogonal to this one, so the vector \( \mathbf{v} \) will have components 0, except for the first entry. Therefore, the weighting factors are given by \( w_j = (1,0,0, \ldots ,0) \). This gives the variance (19) \( \text{Var}[d_{\text{est}}] = \sigma_1^2/N \), which saturates the CRI, as seen from Eq. (15). In this case, \( \sigma_1^2 \) is just the sum of any row or column of \( C \). As we will see in Sec. VII this result will facilitate calculating the Fisher information for our simple correlated noise model.

V. INCLUDING THE WEAK-VALUE AMPLIFICATION EFFECT

Let us now consider the weak-value case. In a weak-value-type metrology experiment, a system is weakly coupled to a meter via an interaction whose coupling strength we would like to determine. We are interested in the case where the weak value gets large, which occurs when the initial and final states are nearly orthogonal. Experimental observations
of weak-value amplification must, therefore, involve choosing the initial and final state to be orthogonal. When this occurs, and when the coupling parameter $d$ is small, one may still get a large pointer shift, which in turn suggests the technique might be useful as a way to measure the value of otherwise small coupling constants $d$.

The interaction term can be written as the product of some system observable $A$ and an operator on the pointer that we call $P$. The interaction Hamiltonian is therefore

$$H_I = d(t) \mathbf{P} \cdot \mathbf{A},$$

where $d(t)$ is a function with compact support near the time of measurement. Usually, the interaction is approximated as happening very quickly, for example: $d(t) = d \delta(t)$, where $d$ is the interaction strength. The pointer operator is termed $\mathbf{P}$, because the effect of $H_I$ is to generate translations in the conjugate variable ($x \ d A$) consequently interpreted as a pointer position [44]. For the measurement to be considered “weak” the pointer shift must be much less than the uncertainty, $(\sigma)$, in the position of the pointer, $d (A) \leq \sigma$. In the standard weak-value approach, the system is pre- and postselected to be in state $|i\rangle$ and $|f\rangle$. For an $M$-dimensional system, there are $M$ simultaneously possible outcomes for a given projective measurement, with the probability of a given outcome $|f\rangle$ given by $\gamma = |\langle f | i \rangle|^2$. While it is possible that in general there may be information in the probability of the selection [28,45–47], in the usual approach, at la Aharonov, the probability of selection is independent of the parameter of interest, and all of the available information is in the meter deflections [1,28]. If the average deflection in the absence of a second selective measurement is given by $d (|i|A|j\rangle = d(A)$, then in the weak limit, the Fisher information about $d$ is multiplied by a factor of $\langle A \rangle^2$ [see Eqs. (7) and (8)]. In contrast, in the presence of the second selective measurement, the deflection is given by $A_w d$; where

$$A_w \equiv \frac{\langle f | A | i \rangle}{\langle f | i \rangle}.$$  

is defined as the weak value. We note that the weak value can be imaginary and that it is not bounded by the eigenspectrum of the operator $A$. If the probability of postselection is made very small, the weak value can become very large, hence the term amplification. In this work, we focus on real weak values; see Refs. [20,27,48] for a discussion of imaginary weak-value amplification.

If we make a weak-value amplification-type experiment, with postselection of probability $\gamma$, the resulting data set $[s_j]$ contains on average only $\gamma N$ data points (where we recall that $N$ was the number of data points in the nonpostselected measurement). Whereas before we would rescale the meter deflection by the expectation value $A$ in order to isolate the parameter of interest ($d$), we now have to account for the amplification effect. Our signal is boosted from $d (\langle A \rangle)$ to $d A_w$ and the correlation matrix changes to $C^\prime$, and is now a smaller approximately $\gamma N \times \gamma N$ matrix. The Fisher information is given by [27]

$$\mathcal{I}_{w} = A_w^2 \sum_{i,j} (C^\prime)^{-1}_{i,j}.$$  

We can now treat this case like we did in the previous section. We have a new covariance matrix $C'$ with which we can make a similar decomposition, $C' = O' D' O'^T$. The dimension is reduced by a factor of $\approx \gamma$. We make exactly the same treatment as before, calling $\sigma_{j}^2$ the eigenvalues of $C'$, and $w_j'$ the new weights.

The Fisher information is now given by

$$\mathcal{I}_{w} = A_w^2 (\gamma N) \sum_{k} \sigma_{k}^{-2} w_k'^2,$$  

where the weights $w_k'$ are normalized. In order to account for the effect of the weak-value amplification, we will later give a detailed model for the precise form the weak values take on. For the moment, we estimate $A_w = (\langle A \rangle)^2/\gamma$, as is true in many weak-value amplification experiments. If that is so, we have

$$\mathcal{I}_{w} = N (\langle A \rangle^2) \sum_{k} \sigma_{k}^{-2} w_k'^2.$$  

We note that the amplification factor has canceled the factor of $\gamma$ which arose due to the reduced size of the new covariance matrix. Comparing this relation to (15), accounting for the multiplication of the Fisher information with $\langle A \rangle^2$, we see that both scale as $N$, and the main change to the Fisher information is how the correlations are affected by the postselection. If the randomly postselected events have the same type of correlations as the nonpostselected case, then the Fisher information is comparable—it is still a weighted average of (a smaller number of) inverse eigenvalues. If the correlations are reduced because retained measurements are further separated in time reducing temporal correlations, for example, then the Fisher information could be larger and WVA would possess an advantage over the direct measurement approach. We will soon see that this is indeed the case.

VI. OPTIMAL PARTITIONING MEASUREMENT

We next consider improving on the weak-value amplification scheme by incorporating the discarded measurement results in our estimation strategy. This involves optimally implementing a partitioning measurement so that all output channels and resulting correlations are used. We will refer to this as an optimal partitioning class measurement (OPM), which we can compare to the simpler WVA case.

For an $M$-dimensional system, there are $M$ possible outcomes of the second projective measurement on the system, $|f_1\rangle, |f_2\rangle, \ldots, |f_M\rangle$. For each of those possibilities, there is a weak value, so $A_{w,f_j}, j \equiv 1, \ldots, M$. The distribution of events is assumed to be a multivariable Gaussian distribution $P(|s_j||d)$, with mean $\mu = \tilde{A}_w d$ and covariance matrix $C$. Here, we define a vector of weak values, $\tilde{A}_w$, associated with each outcome with elements $A_{w,f_j}$, where $j = 1, \ldots, M$. We will now focus on the $M = 2$ case.

Given $N$ measurement outcomes, the selection tags $\gamma N$ of the outcomes with one postselection associated with the final state $|f_1\rangle$, and the remaining $(1 - \gamma)N$ outcomes with the postselection associated with the final state $|f_2\rangle$. We reorder the outcomes and label them $i = 1, \ldots, \gamma N; \gamma N + 1, \ldots, N$. This will not typically be the temporal ordering. However, the first selection is associated with the weak-value $A_w$,
and the second with $A_w^\dagger$. We write the covariance matrix in $2 \times 2$ block form (this can be generalized for multiple partitionings in the higher dimensional case). The probability distribution is a multivariable Gaussian distribution. It has mean $\mu = (A_w d_1 y N, A_w^\dagger d_1 (1 - y) N)'$, where $1_y$ is a vector of $y$ 1s. The covariance matrix in block form is

$$C = \begin{pmatrix} C_{11} & C_{12} \\ C_{21} & C_{22} \end{pmatrix},$$

where $C_{11}$ (of dimension $y N \times y N$), and $C_{22}$ [of dimension $(1 - y) N \times (1 - y) N$] are the covariance matrices associated with the two selections, and $C_{12} = C_{21}^T$ is the correlation matrix between the selected outcomes, which has dimension $y N \times (1 - y) N$ or $(1 - y) N \times y N$.

The Fisher information for such a situation is given by [43]

$$I = \partial d^T \cdot C^{-1} \cdot \partial d$$

$$= I_1 + I_2 + I_3$$

$$= A_w^2 \sum_{ij} (C^{-1})_{11}^i j + (A_w^\dagger)^2 \sum_{ij} (C^{-1})_{22}^i j$$

$$+ A_w A_w^\dagger \left( \sum_{ij} (C^{-1})_{11}^i j + \sum_{i' j'} (C^{-1})_{12}^i j' \right),$$

where the sums run over the appropriate ranges. It is important to note that $C^{-1}$ is, where $k, l = 1, 2$ refers to the block-matrix form of the inverse of the entire $C$ matrix, not the inverses of each of the subblocks. Just as in the simple example of the $2 \times 2$ covariance matrix, the total information exceeds using either selection state alone, however, there may be cases where all the information is in one of the selection states, and the other may be discarded without any loss in variance. We will see such an example in the next section.

It is also of interest to find the optimal estimators in this case, using maximum likelihood methods. Each event is filtered according to the selection state, and is associated with a mean of that weak value times $d$. The covariance matrix is assumed to have the same values as before (with a mean meter shift of $A_w d$ or $A_w^\dagger d$) in the weak measurement limit, but the indices are relabeled to put the matrix into block form, associated with each weak value.

We can find the optimal estimator by solving for the value of $d$ that maximizes the log likelihood, $\ln P(\{s_i\}|d) = \{-\bar{s} - \bar{A}_w d\}^T \cdot C^{-1} \cdot (-\bar{s} - \bar{A}_w d) + \text{const}$ with respect to the parameter $d$ [27]. The maximum likelihood estimator is given by

$$d_{\text{est}} = \frac{\bar{A}_w^T \cdot C^{-1} \cdot \bar{s}}{\bar{A}_w^T \cdot C^{-1} \cdot \bar{A}_w},$$

where $\bar{s}$ is a vector of outcomes.

For the special case of two selection states, with weak values $A_w$ and $A_w^\dagger$, so $\bar{A}_w = (A_w 1_y N, A_w^\dagger 1_{1 - y} N)$, we assume that each selection has outcomes $1, \ldots, y N$, and $y N + 1, \ldots, N$. The inverse covariance matrix takes the form of a $2 \times 2$ block matrix. We further break the vector of outcomes in two, corresponding to each selection label, $\bar{s} = (\bar{s}_1, \bar{s}_2)$. To give an explicit expression, we express the estimator as a ratio, $d_{\text{est}} = N/D$, in which the numerator $N$ is given by

$$N = (A_w 1_y N, A_w^\dagger 1_{1 - y} N) \left( \begin{array}{cc} (C^{-1})_{11} & (C^{-1})_{12} \\ (C^{-1})_{21} & (C^{-1})_{22} \end{array} \right) \left( \begin{array}{c} \bar{s}_1 \\ \bar{s}_2 \end{array} \right),$$

$$= A_w \sum_{i=1}^{\gamma N} \sum_{j=1}^{\gamma N} ((C^{-1})_{11})_{ij} s_{1, j} + \sum_{j=\gamma N + 1}^N ((C^{-1})_{12})_{ij} s_{2, j} + A_w^\dagger \sum_{i=1}^{\gamma N} \sum_{j=\gamma N + 1}^N ((C^{-1})_{21})_{ij} s_{1, j} + \sum_{j=\gamma N + 1}^N ((C^{-1})_{22})_{ij} s_{2, j}.$$

and the denominator $D$ is the Fisher information, (30).

VII. EXACTLY SOLVABLE MODEL

A. Direct case

We now consider a simplified model for correlated noise, in which the covariance matrix corresponds to a combination of uncorrelated noise in time (a) and perfectly correlated noise (c). The $N \times N$ covariance matrix $C$ is then a sum of a diagonal matrix and a matrix of identical elements,

$$C_{ij} = a \delta_{ij} + c,$$

where the constant term in Eq. (34) reflects a shift common to all elements of the data set (e.g., a systematic error), which is drawn from a zero-mean Gaussian distribution with a variance of $c$.

Equation (34) has a simple eigensystem, because the characteristic equation for the eigenvalues, $\det(C - \sigma^2 I) = 0$, may be solved by substituting $\sigma^2 = c a^2 + a$, and noting that $\sigma^2$ solves for the eigenvalues of the matrix of all 1s. The latter matrix has one eigenvalue of $N$, and the rest 0. The first normalized eigenvector is $e_1 = (1, 1, \ldots, 1)/\sqrt{N}$, corresponding to the $N$ eigenvalue, and the other $N - 1$ eigenvectors corresponding to the 0 eigenvalue are constructed orthogonal to $e_1$.

This result indicates that the eigenvalues of $C$ are

$$\sigma^2_j = (Ne + a, a, a, \ldots, a),$$

with the same eigenvectors mentioned above. The positive semidefinite condition requires that $c \geq -a/N$. We can now apply the procedure outlined above by writing the orthogonal matrix formed from the orthonormal eigenvectors as

$$O = (e_1, e_2, \ldots, e_N).$$

The vector $\bar{s}_k = \sum_j O_{1,k} e_j$ may be computed for this model by using the orthonormality conditions of the vectors $e_j$. This is because the sum runs down the column of each unit
basis vector. Using the properties
\[ \sum_{i} (e_i)_j = \sqrt{N}, \quad e_i \cdot e_j = \sum_{i} (e_j)_i = 0, \]
for every unit vector with \( j \neq 1 \), we find
\[ v = (\sqrt{N}, 0, 0, \ldots, 0), \quad w = (1, 0, 0, \ldots, 0), \]
for this model. Noting that Eq. (14) is correct, we find that the
weight, the result is
\[ \text{measurements into a retained and a discarded partition.} \]
and where the Fisher information is
\[ I_{ac} = \langle A \rangle^2 \frac{N}{c + a}. \]
We see that the effect of the correlations is to contribute to the
coefficient matrix as previously shown. As
\[ c \rightarrow c^2 \gamma N \]
but with the effective change of
\[ c \rightarrow \gamma c. \]
That is, postselection reduces the size of the correlation.
However, the resulting advantage depends on the kind of
noise: If \( c < 0 \), the postselection increases the variance,
whereas if \( c > 0 \), the postselection reduces the variance of
the optimal estimator. The latter is in accordance with the
findings of Ref. [41] for this model. We note that, as before,
there is no difference between the optimal estimator and the
equal-weighting estimator.

We note that while \( \gamma \) can be made arbitrarily small, the
Fisher information is bounded by the necessity of sampling
some high-information content events, selected from the
covariance matrix elements \( a + c \), giving the bound
\[ I_{wv} \leq \langle A \rangle^2 \frac{N}{a + c}. \]
This result represents an enormous suppression of the detri-
mental effects of noise accompanied by a significant reduction
in the size of the data set. The WVA technique is able to
recover the performance of the conventional method in the
uncorrelated noise limit by simply selecting a small enough
\( \gamma \). If for some experimental reason, \( \gamma \) is bounded to be larger
than some minimum, \( \gamma_{\text{min}} \) gives a practical limit to the noise
reduction (44).

C. Using the other selection: Optimal partitioning measurement

We now apply the results of Sec. VI to our model to see how
much the Fisher information may be improved by the optimal
partitioning method, which involves optimally weighting both
the “postselected” and the “postselection rejected” partitions
of the data set in order to estimate the parameter of interest.
Comparing this approach to the WVA strategy will tell us how
much information was discarded by the postselection step in
WVA. The Fisher information may be explicitly evaluated in
our exactly solvable model. This is because the exact inverse
of matrix (34) is given by
\[ C^{-1}_{ij} = \frac{(a + c)\delta_{ij} - c}{a^2 + Nac}, \]
as can be checked by direct calculation, \( C C^{-1} = I \). It is
straightforward to see that summing over both indices in (46)
returns the Fisher information (39). The Fisher information
with weak-value amplification, Eq. (30) in this special case,
has the exact form
\[ I = a N [\gamma A_w^2 + (1 - \gamma) (A_w^L)^2] + \gamma (1 - \gamma) N^2 (A_w - A_w^L)^2 \]
\[ a^2 + Nac. \]

In order to compare the different methods, we adopt the
standard model of the weak value, taken from [20]:
\[ A_w = -\cot(\phi/2), \quad A_w^L = \tan(\phi/2), \quad \gamma = \sin^2(\phi/2), \]
where \( \phi/2 \) is the overlap angle between \( |f \rangle \) and \( |i \rangle \), the pre-
and postselected states. This model leads to the simplification
of the three terms defined in (30) as
\[ I_1 = \frac{a N \cos^2(\phi/2) + c N^2 \cos^4(\phi/2)}{a^2 + Nac}. \]
FIG. 6. Different contributions to the Fisher information in Eq. (30) vs \( \phi \), in units of \( N/a \), the uncorrelated Fisher information. Term 1 \((I_1)\) is proportional to \( A_w^+ \), term 2 \((I_2)\) is proportional to \( (A_w^-)^2 \), and term three \((I_3)\) is proportional to \( A_w A_w^+ \). We take \( N = 100 \), \( c/a = 0.5 \).

\[
I_1 = \frac{a N \sin^2(\phi/2) + c N^2 \sin^4(\phi/2)}{a^2 + N ac},
\]

\[
I_2 = \frac{2c N^2 \sin^2(\phi/2) \cos^2(\phi/2)}{a^2 + N ac}.
\]

The sum of these three terms is

\[
I = \frac{N}{a}.
\]

Remarkably, the total Fisher information is now independent of the value of \( \phi \), and is larger than the uncorrelated Fisher information. Assuming \( c > 0 \), we recall that just using the one selection state had the effect of reducing \( c \to \gamma c \). We see here that adding in the other selection state and their correlation allows us to eliminate the effect of \( c \) entirely. In Fig. 6, we show the various contributions to the combined Fisher information using both output selections. The Fisher information (30) comes from the three terms from each subblock. We note that if it is the case that \( c < 0 \), then it is clearly advantageous to use the direct measurement scheme. In Table I we tabulate the Fisher information (FI) of the three approaches we consider in white- and slow-noise limits.

Summarizing the chart above, we have found that all three approaches are identical in terms of measurement performance in the uncorrelated noise limit. Whichever is experimentally easier to implement, therefore, has the advantage. In the correlated noise limit, WVA has a clear advantage over the direct approach, confirming the work of Feizpour, Xing, and Steinberg [41]. What Feizpour et al. did not consider, however, is that unlike in the white-noise limit, once there are correlations between measurements, retaining all partitions and the correlations between them becomes advantageous. In fact, in the same regime where WVA has a real quantitative advantage over the conventional approach, it turns out to be slightly inferior to the optimal partitioning method.

D. Limiting optimal estimators

We see from the above analysis that with the OPM approach, the selection angle \( \phi \) simply changes how the information is distributed in the various outcomes. It is instructive to focus on the form of the optimal estimators (32) in the two extreme limits for our exactly solvable model: \( \gamma < 1 \), the weak-value amplification limit, and \( \gamma = 1/2 \), the balanced limit, in order to understand how the optimal partitioning approach is able to completely eliminate \( c \).

In the very unbalanced limit, the term proportional to \( A_w \) dominates the Fisher information. The weak values are given approximately by \( A_w \approx -2/\phi \), \( A_w^+ \approx \phi/2 \), and the asymmetry is given by \( \gamma \approx \phi^2/4 \). We find the optimal estimator to be

\[
d_{\text{est}} \approx d_{\text{est}}^{uw} = \frac{A_w \gamma}{a + N \gamma} \left( \gamma N \sum_{j=1}^{N/2} s_{1,j} + \gamma N \sum_{j=N/2+1}^{N} s_{2,j} \right),
\]

where the weak-value estimator \( d_{\text{est}}^{uw} \) is defined as

\[
d_{\text{est}}^{uw} = \frac{A_w \gamma}{N} \sum_{j=1}^{N/2} s_{1,j}.
\]

The total (correlated) optimal estimator is just the weak-value estimator, plus another term involving the sum of all the data, whose average is approximately 0, and its prefactor vanishes as \( c \to 0 \). The additional term is able to account for the (known) correlations in the system and make a further suppression of the variance, at the cost of having to process all collected data.

We now turn to the balanced case, where \( \phi = \pi/4 \), \( \gamma = 1/2 \), and the weak values are \( A_w = -1 \), \( A_w^+ = 1 \). In this case, the weighting prefactors in front of the collected data cancel out, and we find the estimator

\[
d_{\text{est}}^b = \frac{1}{N} \left( \sum_{j=1}^{N/2} s_{2,j} - \sum_{j=N/2+1}^{N} s_{1,j} \right),
\]

Table I. In this table, the Fisher Information is summarized in the uncorrelated and correlated noise limits for the three different strategies we consider: the direct approach, the weak value approach, and the optimal partitioning approach. From the table, \( N \) is the number of measurements carried out, \( \gamma N \) is the number of retained measurements in the WVA approach, \( a \) is the variance of the white noise, and \( c \) is the variance of the unknown systematic. From the figure we can see OPM removes all effect of \( c \), while WVA reduces its effect by a factor of \( N \), recovering exactly the variance achievable in the absence of correlations.

<table>
<thead>
<tr>
<th>Approach</th>
<th>Uncorrelated noise</th>
<th>Correlated noise</th>
</tr>
</thead>
<tbody>
<tr>
<td>FI direct</td>
<td>( \frac{N}{a+c} )</td>
<td>( \frac{N}{a+Ne} )</td>
</tr>
<tr>
<td>FI WVA</td>
<td>( \frac{N}{a} )</td>
<td>( \frac{N}{a+N} )</td>
</tr>
<tr>
<td>FI OPM</td>
<td>( \frac{N}{a+c} )</td>
<td>( \frac{N}{a} )</td>
</tr>
</tbody>
</table>
used in experimental laboratories to eliminate correlated noise. We discuss this in more detail in the next section.

VIII. BACKGROUND SUBTRACTION

Background subtraction typically involves partitioning measurement outcomes into two types: measurements of signal plus background noise, and measurements of just background noise. By comparing the two partitions, unknown offsets or errors in calibration can be corrected. Background subtraction can also be used to suppress temporally correlated noise by sampling the noise background at least once per correlation time and subtracting off the slowly evolving offset. Unfortunately, optimizing how often to sample the background noise is usually a hard problem. This is because usually both white and slow noise are present, and oversampling the background (sampling more than once per correlation time) ceases to reduce either. Optimizing the amount of background subtraction (as opposed to normal measurements) is only possible with knowledge of the correlation time of the noise. This makes it preferable, when possible (as it is in our example), to alternate the sign of the signal instead of chopping it on and off. In this variant of background subtraction, measurement outcomes are partitioned into two types: measurements of a background noise plus signal, and measurements of background noise minus signal. By subtracting the two partitions and dividing by two (just like in Sec. III), we can eliminate any slow noise present while also averaging down the white noise.

While we have considered partitioning measurements using postselection, it is possible to achieve the same result by alternating the preparation. It is worth noting that the additional information gained by this more complex use of the partitioning measurement is very marginal compared to the “average” number of correlated measurements) to represent the correlation time of the noise. While we have considered two limiting cases of this kind is quite challenging, hence the interest in a generalized noise model. In the “white-noise limit,” \( \eta = 0 \) and measurement outcomes are uncorrelated. In the opposite limit, \( \eta \to \infty \) and the covariance matrix reduces to the directly solvable one from Sec. VII,

\[
C_{i,j} = a \delta_{i,j} + c.
\]

We refer to this as the “slow-noise limit.” This limit represents taking the correlation time of the noise to infinity resulting in a scenario where estimation error is increased by some unknown offset \( c \). We have previously treated the two limits of our noise model for all three measurement approaches considered (direct, WVA, and OPM). We expect the effects of the correlations to grow as the correlation time \( \eta \) increases, smoothly connecting our two limits for all the measurement approaches considered (direct, WVA, and OPM). We refer to this as the “slow-noise limit.”

IX. NUMERICAL INVESTIGATION

We will now return to our original (and more experimentally relevant) noise model. This noise model represents a familiar scenario in a laboratory. Measurements are made sequentially in time and are compiled into a data set \( \{s_i\} \). As before, a measurement outcome \( s_i \) can be decomposed into the parameter of interest \( d \) and a Gaussian distributed, zero-mean, random variable \( x_i \), such that

\[
s_i = d + x_i,
\]

where \( i = 1, 2, \ldots, N \). We take the covariance matrix to be

\[
C_{i,j} = a \delta_{i,j} + ce^{-|i-j|/\Delta t/\tau}.
\]

Here, \( \Delta t \) sets the time between subsequent measurements, \( \tau \) represents the correlation time of the noise, and the ratio of \( a \) to \( c \) sets the relative amount of white noise to slow noise (except in the limit where \( \tau \to 0 \) where all noise is white).

For simplicity we will use the unitless quantity \( \eta = \tau / \Delta t \) (the “average” number of correlated measurements) to represent the correlation time of the noise.

This noise model captures an experimentally common scenario, wherein there is a white-noise floor (represented by \( a \)) and some correlated noise arising due to experimental imperfections (represented by \( c \)). Overcoming technical noise of this kind is quite challenging, hence the interest in a technique that is robust against it.

We have already considered two limiting cases of this generalized noise model. In the “white-noise limit,” \( \eta \to 0 \) and measurement outcomes are uncorrelated. In the opposite limit, \( \eta \to \infty \) and the covariance matrix reduces to the directly solvable one from Sec. VII,

\[
C_{i,j} = a \delta_{i,j} + c.
\]

We refer to this as the “slow-noise limit.” This limit represents taking the correlation time of the noise to infinity resulting in a scenario where estimation error is increased by some unknown offset \( c \). We have previously treated the two limits of our noise model for all three measurement approaches considered (direct, WVA, and OPM). We expect the effects of the correlations to grow as the correlation time \( \eta \) increases, smoothly connecting our two limits for all the measurement approaches considered. Finding the Fisher information in this intermediate regime, however, is challenging due to the complex structure of the covariance matrix. In order to find the Fisher information, we generate large covariance matrices with the appropriate structure and invert and sum them while varying the correlation time. This allows us to smoothly connect the two limits. In Fig. 7, the Fisher information for
both the WVA method, the direct method, and background subtraction (equivalent to OPM) is plotted as a function of the average number of correlated measurements $\eta$. In each scenario, the Fisher information is found through construction of the covariance matrix that would be produced by the measurement design and then inverted and summed. For comparison, the inverse variance of the mean associated with the weak-weighting estimator is also found analytically and plotted in solid lines. In the two limits ($\eta \rightarrow 0$ and $\eta \rightarrow \infty$) the two estimators are equivalent. In the intermediate regime, the equal-weighting estimator is no longer strictly optimal. It deviates only very slightly from the Fisher information as a result of the finite size of the covariance matrices we consider (due to the boundaries of our data set certain measurements are correlated to more measurements than others). In Fig. 7, as expected, as $\eta \rightarrow 1$, the Fisher information in the direct method degrades swiftly, while the Fisher information of the weak-value method is unaffected. Only after $\eta$ approaches $1/\gamma$ ($\gamma$ was arbitrarily chosen to be 0.5%) does the performance of the weak-value method begin to diminish. If $\gamma$ is fixed, then in the limit of $\eta \rightarrow \infty$ the minimum variance becomes $a/N + \gamma c$. Even in the intermediate regime, with weak correlations, postselection suppresses the detrimental effects of the correlations by increasing the time between retained measurement outcomes.

These results confirm the work of Feizpour, Xing, and Steinberg who argued that WVA-based measurement estimation strategies afforded an advantage in a correlated noise limit over direct methods. They also agree with Ferrie and Combe's observation [24] that WVA does not always achieve globally optimal performance. We see that the OPM–background subtraction technique slightly outperforms WVA as soon as the correlation time becomes comparable to the time between measurements. This difference can be substantial if the minimum probability of postselection is large; however, if it is on the order of $1/\gamma$, it is quite small. In Fig. 7 the relative size of $a$ and $c$ that represent the white and slow portions of the noise was chosen in order to visually differentiate the different approaches. In typical situations, $c$ is much less than $a$, for otherwise an experimentalist would be easily able to detect it in preparation for the experiment. If, however, the value of $c$ is on the order of $a/N$, then an experimentalist would have to spend more time characterizing the noise then performing the experiment—an unlikely scenario in all but the most heroic metrology experiments. A realistic value for $c$ is therefore on the order of $a/N$, and we see that in this scenario the additional advantage afforded by optimally using all the data is

$$\Delta I = \frac{N}{a} - \frac{N}{a + c} \approx \frac{1}{a + a/N}. \tag{59}$$

This cost in precision limits to $1/a$ as $N \rightarrow \infty$ and is equal to the information which would be gained by a single additional measurement in this scenario. These numerical results confirm that the conclusions reached using our simplified model for correlated noise hold generically in more realistic laboratory situations with noise environments with time decaying correlations.

X. CONCLUSIONS

We have investigated how the introduction of a second partitioning measurement affects parameter estimation. Using a realistic model for additive, Gaussian, time-correlated noise we found that, for all the cases considered, introducing a second partitioning measurement affords an advantage in the Fisher information over a direct measurement method (no second partitioning measurement) once the correlation time of the noise becomes longer than the measurement rate. Furthermore, we have found that if one or more output channels from the partitioning measurement are filtered (corresponding to a WVA-like postselection), the majority of this advantage is retained. Thus, WVA can help dramatically suppress the detrimental effects of slow noise, recovering the performance achieved in the white-noise limit by effectively decreasing the correlations between retained measurement outcomes. The informational cost of the postselection step is studied by comparing the Fisher information of the WVA approach with an optimal partitioning measurement scheme, which utilizes all the data and correlations between partitions. The cost of discarding all but one partition is found to be related to the ratio of the magnitudes of the slow and white noise. We argue that in a typical laboratory situation, this cost will be comparable to the information gained by performing a single extra measurement and is, therefore, negligible. An analysis of the estimators used by OPM leads us to the realization that in the balanced case, OPM is equivalent to probabilistic background subtraction. This insight provides a unified framework for understanding when partitioning-class measurements, and specifically weak-value amplification, can be useful. In conclusion, in experimental settings with time-correlated noise, if background subtraction or other optimal partitioning measurement methods are technically challenging to implement or if a substantially reduced data set is desirable, WVA vastly outperforms conventional measurement approaches and is near optimal.

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