On the efficiency of quantum lithography

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Abstract. Quantum lithography promises, in principle, unlimited feature resolution, independent of wavelength. However, in the literature, at least two different theoretical descriptions of quantum lithography exist. They differ in the extent to which they predict that the photons retain spatial correlation from generation to absorption, and although both predict the same feature size, they vastly differ in predicting how efficiently a quantum lithographic pattern can be exposed. Until recently, essentially all quantum lithography experiments have been performed in such a way that it is difficult to distinguish between the two theoretical explanations. However, last year an experiment was performed that gives different outcomes for the two theories. We comment on the experiment and show that the model that fits the data unfortunately indicates that the trade-off between resolution and efficiency in quantum lithography is very unfavourable.

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1. Introduction

For a long time it was believed that the resolution of optical systems was limited by the so-called Rayleigh criterion (see e.g. [1]), which roughly states that the resolution of the system is diffraction limited by half of the wavelength $\lambda$ of the used light. Hence, it was believed that only by going to shorter and shorter wavelengths could one improve the resolution substantially. However, going from the visible to ultraviolet light, x-rays or $\gamma$ radiation means using light with such a high photon energy that the radiation may damage or destroy the object one wants to expose or investigate. Furthermore, it is difficult to build optical systems, since ordinary lenses are no longer transparent in these wavelength regions, and even gases start to absorb the light quite substantially. Moreover, building mirrors is also challenging due to absorption and very small refractive index differences between different materials. This makes it complicated and expensive to build Rayleigh-limited optical systems for wavelengths shorter than about 200 nm.

There have been proposals for ‘classical’ systems that achieve higher resolution than the Rayleigh limit [2]–[4], but such lithographic schemes suffer from visibility problems and cannot simultaneously generate a sub-Rayleigh feature size and a sub-Rayleigh feature separation, at least not until new kinds of detectors are developed [5]–[7].

It therefore came as a promise when Boto et al showed that entanglement could be used to beat the diffraction limit given by the Rayleigh criterion [8]. Boto et al showed that the use of $N$ entangled photons would allow one to surpass the diffraction limit by a factor of $1/N$ compared with the Rayleigh criterion. Lithographic systems would, potentially, profit substantially from the idea of Boto et al by allowing the generation of features arbitrarily smaller than the Rayleigh limit [9]–[16]. In order to make these systems work in practice, however, there is still a plethora of problems to be solved. For example, one has to find a material that changes its behaviour only when absorbing $N$ photons [2, 17, 18], and one needs also to optimize the absorption rate of this material [19]. Unfortunately, there are limits to the extent to which this is feasible [20]. Additional difficulties involve the rather peculiar, multi-photon, entangled photon states that need to be generated [21, 22] and the fact that there are limits to the generation efficiency of these states while still surpassing the Rayleigh limit [23]. Alternatively, other ways of generating indistinguishable $N$-tuples of photons have to be developed [24]. There have been investigations
Figure 1. Setup for quantum imaging or lithography for \( N = 2 \). Each of the in total \( S + 1 \) coordinate points on the screen to the right consists of a detector with width \( b \) that detects two-photon absorption. The double slit on the left has a distance \( d \) between the slits and is at a distance \( R \) from the screen.

for the case of multiple frequencies or in the broadband limit [25, 26]. For some schemes even photon losses may become an issue [27]. Furthermore, one has to think about how to generate arbitrary patterns [10]–[12], [28]. Still, more than 10 years after the proposal of Boto et al, there has been no demonstration of quantum lithography coming close to the performance of conventional state-of-the-art lithography. Meanwhile, ‘conventional’ lithography is becoming even better, e.g. efforts have been taken to surpass the Rayleigh limit with these systems by using innovative methods, for example nonlinear effects [29].

In this paper, we address an open question that may make quantum lithography an even more remote contender for future lithographic applications. The question we pose is whether quantum lithography can work efficiently in practice. The main aspect is whether or not the time needed to irradiate a lithographic film would be short enough to make this method efficient.

We start this paper by describing in section 2 two different models for most of the experiments so far reported in the field of quantum lithography, where one of the explanations predicts a favourable time scaling whereas the other one does not. We then discuss in section 3 an experiment which would give different outcomes for the two explanations and which thereby could decide which of the explanations is correct (or at least applicable to this kind of experiment) and how the time for imaging or radiation scales with the number of photons \( N \) per state and the number of image ‘pixels’ \( S \). The outcome of an experiment recently performed by Peeters et al [30] allows us then to deduce which of the explanations seems to be applicable, and the result suggests that the time-scaling behaviour of such a quantum lithography scheme is, unfortunately, very unfavourable. In section 4 we relate our paper to other papers in the field and in section 5 we present our conclusions. The paper also has two appendices: one where we give a thorough mathematical analysis and derivation of our model and one where we look, in more detail, into the original proposal of Boto et al.

2. Different theoretical explanations

Imagine a setup as depicted in figure 1. A light beam of wavelength \( \lambda \) is incident on a double slit with a slit distance \( d \gg \lambda \). Due to diffraction at the slit, the light will produce an interference pattern on a screen at a distance \( R \gg d \) behind the slit. Assume that our screen now consists of
S + 1 detectors, each of the same width b. Below, we have assumed that S is even (representing an odd number of detectors). However, the conclusions we draw are equally valid if an even number of detectors is used. The detectors cover the whole length of the screen and b is small enough to clearly be able to resolve the interference pattern; that is, the distance between two intensity peaks in the detector plane is many times longer than b. A photon impinging on the double slit can pass through either the upper or the lower slit, i.e. the state at the slit will be

$$|\Psi_{1,sl}\rangle = \frac{1}{\sqrt{2}} \left( |1\rangle_{\text{upper}} \otimes |0\rangle_{\text{lower}} + |0\rangle_{\text{upper}} \otimes |1\rangle_{\text{lower}} \right).$$  \hspace{1cm} (1)

Since we have decided to remain ignorant as to which path the photon took through the double slit, but only detect the photon with the detectors at the screen, we will there have the state

$$|\Psi_{1,sc}\rangle = \frac{1}{N_1} \sum_{s=-S/2}^{S/2} \left[ e^{ikr(s)} \hat{a}_s^\dagger + e^{ik[r(s)+\Delta r(s)]} \hat{a}_s^\dagger \right] |0\rangle,$$

$$= \frac{1}{N_1} \sum_{s=-S/2}^{S/2} e^{ikr(s)} \left[ 1 + e^{ik\Delta r(s)} \right] \hat{a}_s^\dagger |0\rangle,$$  \hspace{1cm} (2)

where $N_1 = \sqrt{2(S+1)}$; the summation is over all detector modes s; $k$ is the wave vector of the incident light; $r(s) = \sqrt{R^2 + s^2b^2} \approx R \left( 1 + \frac{s^2b^2}{2R^2} \right)$; and $\Delta r(s) = d\sin \theta \approx d\theta = dbs/R$. (For the normalization in expression (2), we have for simplicity assumed that the pattern falling on the S + 1 detectors contains exactly an integer number of fringes.) The ensuing measurement probability at the detection plane is

$$P_{1,sc}(s) \propto \cos^2 \left( \frac{k dbs}{2R} \right).$$  \hspace{1cm} (3)

The distance between two peaks in the measured interference pattern is given by $\lambda_0 R/d$, where $\lambda_0$ is the used wavelength. In the equation above, we have neglected the envelope function induced by the exact shape of each slit. Instead we have assumed that the photons have equal probability amplitude to hit any detector. This is a simplification justified by the fact that it does not alter our main results or conclusion.

If we look at the quantum lithography scheme, the setup is slightly changed. We are no longer interested in single photons but in ‘bunches’ of N photons. For most of this paper, we restrict ourselves, for practical reasons, to $N = 2$ photons, but our whole argumentation will be valid for any number $N$. To show that one can obtain a higher resolution (i.e. a shorter distance between peaks) in this setup, one has to ensure that two indistinguishable photons pass through the same slit, i.e. that one has a two-photon, NOON state at the slits [8]

$$|\Psi_{2,sl}\rangle = \frac{1}{\sqrt{2}} \left( |2\rangle_{\text{upper}} \otimes |0\rangle_{\text{lower}} + |0\rangle_{\text{upper}} \otimes |2\rangle_{\text{lower}} \right).$$  \hspace{1cm} (4)

This can be implemented, for example, by placing a short type-I crystal in front of the double slit and pumping this crystal so that it produces pairs of degenerate photons that leave the crystal collinearly. One has to also ensure that the pump beam is broader than the distance between the slits and that the two slits are centred on the setup’s optical axis. Note that the state at the screen is isomorphic to a Bell state; that is, it has maximal transverse entanglement of its photon number. Also note that, since the slits can be viewed as the source illuminating the screen, an analysis of how this state is produced is not needed. Indeed, in appendix A it turns
out that even long crystals, narrow beams and a long distance between the crystal and slits can produce \(|\Psi_{2,0}\rangle\). It suffices to conclude that the state \(|\Psi_{2,0}\rangle\) is the optimal two-photon state for a demonstration of two-photon interference fringes, giving unity two-photon interference visibility and interference fringe doubling as compared to the classical Young’s experiment at the same wavelength of light.

The interesting question is: What happens to the state \(|\Psi_{2,0}\rangle\) after the photons have passed the double slit? One could use the concept of photonic de Broglie waves [31] to explain the effect of quantum imaging or lithography. Boto et al [8] analysed the probability for the two-photon state to arrive at a specific point on the screen depending on the intensity \(I\) of the incoming light as follows: ‘For two-photon absorption with entangled photon pairs, the absorption cross section scales as \(I\). . . . If the optical system is aligned properly, the probability of the first photon arriving in a small absorptive volume of space time is proportional to \(I\). However, the remaining \(N - 1\) photons are constrained to arrive at the same place at the same time, and so each of their arrival probabilities is a constant, independent of \(I\).’

Transferring this idea to the experiment in figure 1 would give the state

\[
|\Psi_{B,sc}\rangle = \frac{1}{N_B} \sum_{s=-S/2}^{S/2} e^{i2kdx(s)} \left[ 1 + e^{i2k\Delta r(s)} \right] (\hat{a}_{s}^\dagger)^2 |0\rangle,
\]

where \(N_B = \sqrt{4(S + 1)}\). At the screen, the distance between two peaks \(\lambda_0R/(2d)\) is only half as long as that in (3) and one could treat the two photons as a photonic de Broglie wave. The implication of this state is that the two photons can hit essentially any detector (except those at the interference anti-nodes) but that both photons will hit the same detector.

Steuernagel [32], however, opposed that description by stating that ‘... it is not true that the first arriving photon greatly constrains the arrival location of the following ones . . . Very few photons will be absorbed in one point since they typically arrive far apart’. In a mathematical description, his argumentation can be given the following description:

\[
|\Psi_{St,sc}\rangle = \frac{1}{N_{St}} \sum_{s=-S/2}^{S/2} \sum_{s=-S/2}^{S/2} \left( e^{ik[r_1(s) + r_2(t)]} \hat{a}_{s}^\dagger\hat{a}_{r}^\dagger + e^{ik[r_1(t) + r_2(s)]} \hat{a}_{s}^\dagger\hat{a}_{r}^\dagger \right) |0\rangle,
\]

where

\[
r_{1,2}(s) = R \left( 1 + \frac{s^2 b^2}{2R^2} \right) \mp \frac{db}{2R} s,
\]

is the distance from the upper and the lower slit, respectively, and \(N_{St} = \sqrt{2}(S + 1)\). Although one might argue that our derivation of (5) and (6) is somewhat short-handed or hand waving, a complete quantum mechanical analysis of the state from the pump beam to the screen leads to the same result. Such an analysis is given in appendix A. Steuernagel’s assertions have some support in a pioneering paper by Ghosh et al [33]. Ghosh et al show that photon pairs, generated by spontaneous emission downconversion, will give rise to sinusoidal intensity correlations at spatially separated points, and not only when two photons are detected at the same point.

Also Tsang opposed the assertions of Boto et al. He based his conclusion on a rather involved analysis, and found that the multi-photon absorption rate should actually be lower for spatially entangled states, such as NOON-states (of which the state \(|\Psi_{2,0}\rangle\) is an \(N = 2\) realization), than for classical states [34]. However, in the following, we will follow Steuernagel’s somewhat simpler model.
One difference between the descriptions is that the photons retain a spatial correlation ('stick together') in the description of Boto et al and hence will always arrive at the same position. In Steuernagel's model the photons will propagate independently of each other after the slit, and hence can arrive at different detectors. When looking at the case where both photons arrive at the same detector (i.e. $s = t$, see below) the models coincide when it comes to the spatial behaviour; thus, when only looking at two-photon, same-position events, one will see identical interference patterns, both compressed by a factor of $1/2$ as compared to the single-photon interference described in (2) and (3).

Despite predicting the same two-photon, same-position absorption pattern, the two descriptions in (5) and (6) will give quite different scaling behaviour for how efficient a two-photon exposure would be. In (5) both photons in every two-photon state passing through the slits arrive at the same detector, and going from 2 to $N$ photons in (4), the exposure efficiency will be invariant. (However, the absorption efficiency of any detector or material will in general be considerably lower, but this effect will be the same in both cases.) Changing the number of detectors $S$ will only linearly increase the time one has to expose to get a certain number of $N$-photon events per detector. Hence, the ‘exposure time’ of a quantum lithographic image will scale as $\tau_B(S, N) \propto S$.

Equation (6), however, assumes that the arrival of the first photon at a certain detector does not in any way determine where the other photon will arrive and therefore, in the two-photon case, the exposure time $\tau_{St}(S, 2) \propto S^2$. The photons in an $N$-photon state will also hit the detectors in a random fashion, but only if all photons arrive at the same detector will the detector fire. Hence, the exposure time scaling for an $N$-photon state will be $\tau_{St}(S, N) \propto S^N$. Hence, as pointed out by Steuernagel [32], the exposure time will increase exponentially with $S$, and for realistic values of the number of detectors (pixels) and the efficiency of $N$-photon absorbers, this will essentially rule out the practical use of quantum lithography. For example, for a $5 \times 5$ pixel image the exposure time will multiply 25 times going from a two-photon state to a three-photon state. For a more realistic image resolution of $100 \times 100$ pixels, the exposure time will increase by $10^4$ going from an $N$-photon state to an $(N + 1)$-photon state. Therefore, finding out which description is applicable is an important task to determine to what extent quantum lithography will be efficient.

3. A conclusive experiment

3.1. Description of the experiment

For the experiment in figure 1 both theoretical explanations predict the same interference pattern when looking at two-photon absorption. Also for most other experiments performed so far (see, for example, [9] for double slits and [35, 36] for gratings instead of a double slit), either description could be used to explain the obtained results since only same-position correlations were measured, and not the detection efficiency. To differentiate the theoretical models an experiment such as the one schematically depicted in figure 2 can be used. It is similar to the experiment in figure 1 in the sense that we ascertain that the state $|\Psi_{2, sl}\rangle$ is generated directly after the double slit. This can be achieved in the same manner as in the experiment described previously. However, instead of looking at same-position, two-photon correlation, we split the beam by an ordinary 50/50 beam splitter (BS) and detect the photons by two detectors that can have arbitrary positions relative to the respective optical axis. We look at only the case
Figure 2. Schematic experiment for measuring spatial correlations. See the text for more details.

Figure 3. Difference between the two theoretical explanations. In (a) the model of Boto et al is displayed. In (b) the model of Steuernagel is illustrated. The axes correspond to the positions $s$ and $t$ of the two detectors in figure 2, respectively. The distance between two peaks along the line $s = t$ is half as long as it would be for the state in (1). The lighter the colour, the higher the coincidence rate. The colour schemes in the two figures are only relative. One can see that for $s = t$ the two theoretical models give the same pattern.

where both detectors give a click, i.e. we are only looking at coincidences. We denote $s$ as the position of one of the detectors and $t$ as the position of the other detector. Subsequently, the coincidence counts as a function of $s$ and $t$ are recorded. For such an experiment the two theoretical explanations predict different results. The detection probability of the two photons corresponding to (5) is

$$P_{B,sc}(s, t) \propto \cos^2 \left( \frac{kds}{R} \right) \delta_{st},$$

where $\delta_{st}$ denotes the Kronecker delta. For (6), the detection probability becomes

$$P_{St,sc}(s, t) \propto \cos^2 \left[ \frac{kdb}{2R} (s + t) \right].$$

In figure 3 the detection coincidence rates are plotted as a function of $s$ (horizontal) and $t$ (vertical). (The lighter the colour, the higher the coincidence rate.) Figure 3(a) illustrates the prediction of the model of Boto et al (5) and (8), where the photons retain spatial correlation.
Figure 3(b) illustrates the case of Steuernagel’s description, corresponding to (6) and (9), where the photons can propagate independently after the slit and hence allow an interference pattern even for the case when the two detectors are at different positions. The case where \( s = t \), that is, along the figures’ lower left to upper right diagonals, illustrates the results of most of the experiments reported so far. For this case, one can see that both theoretical models predict identical correlation patterns.

3.2. Performing the experiment

An experiment similar to that described in figure 2 was recently performed by Peeters et al [30] in another context. The result of the measurement can be seen in figure 5(a) of their paper [30] and the figure is more or less identical to figure 3(b), indicating that the two photons through the double slit are diffracted independently, but the two-photon character of the state will still be manifested in the coincidence correlation.

Our understanding of the result is that since photons are non-interacting exchange particles, they will diffract and propagate independently in a linear medium. This means that in the absence of some additional ‘guiding’, the photons prepared in the state \( |\psi_{2,\text{sl}}\rangle \) will spread out according to normal diffraction theory, and it will be unlikely to detect the two photons at the same position far from the slit. This is also the case for more complicated states such as the inverse binomial states discussed in [11]. The entanglement will dictate the ensuing \( N \)-photon interference pattern, but each photon will in principle have its probability amplitude spread over the whole illuminated area, resulting in very low \( N \)-photon, single-detector, correlation. In appendix B, we show that this is also true for a slightly different setup discussed in [8]. In other words, Steuernagel seems to be correct in his analysis that both spatial and temporal concentration cannot be achieved at the same time at the required level for quantum lithography [37].

In principle, one could try to keep the spatial correlation by making the propagating media nonlinear and, for example, form spatial solitons [38]. However, to manage this on a few-photon level seems farfetched and not very realistic.

Finally, one may ask: Is our conclusion valid also for a quantum lithography involving more than two modes ([10]–[12], [28])? The answer is yes. Invoking more modes will allow the synthesis of more complicated patterns than a single-slit interference pattern superimposed by a sinusoidal modulation, which is what can be achieved by a double slit. By adding more modes one can achieve additional superpositions of modes with different transverse wave vectors. Thus, one can make more involved lithographic patterns. However, if two photons emerging from one of two slits diffract independently, which is what the results of Peeters et al indicate, then two (or more) photons will also diffract independently from any one of any number of slits or ‘holes’ in a lithographic mask. Therefore, our conclusions about the exposure efficiency are applicable to any lithographic mask when combined with propagation through a linear media.

4. Can spatial correlation be retained?

One might wonder how this paper relates to other papers like [39] or [40], where authors actually predict that \( N \)-photon absorption proportional to \( I \), where \( I \) is the normalized classical intensity of single-photon absorption, is theoretically possible. These two papers are also cited by Boto et al to motivate their argument that quantum lithography works with \( I \) instead of \( I^N \).
dependence. For the results of these paper to be valid, the authors postulate, without giving any specific mechanism, that the photons are somehow transported to exactly the same place; for example, the paper [39] states: ‘We shall assume that, by focusing or otherwise, the photons of a correlated pair are brought at the same time onto a three-level atom…’ This, however, is not the situation in a typical quantum lithography scheme, since the photons are diffracted independently at a double slit (or more generally, in some lithographic mask) as the present paper shows. In the proposal of Boto et al [8] and the experimental realization of D’Angelo et al [9], which does not use a double slit, the photons are spreading independently as an analysis can show (see appendix B). This independent spread could, in principle, also have been deduced from Ghosh et al’s paper [33]. Therefore, as long as light is propagating through linear media (such as air or vacuum), the results of [39] and [40] are not applicable in the existing quantum lithography schemes.

5. Conclusion

We showed that one can experimentally determine the scaling behaviour of the needed exposure time in quantum lithography as a function of the number of photons \( N \) and the number of detectors (i.e. pixels) \( S \). Such an experiment was recently reported, and it supports the view that photons, spatially entangled at a double slit to subsequently form a sub-wavelength pattern, will diffract independently. The photon state will still form the desired pattern (given that only two-, or in general, \( N \)-photon coincidences are recorded). However, the experiment unfortunately also indicates that by increasing the resolution through the number of entangled photons and by increasing the number of pixels, one has to wait an exponentially longer time to get the desired interference pattern. The explicit exposure time scaling is \( \tau_S S^N \). Therefore, quantum lithography involving a moderate number of pixels, unfortunately, seems to be impractical even for a rather small number of photons.

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Appendix A. A thorough analysis of the theoretical explanations

Only assuming cw-pumping one can write the down-converted two-photon state as

\[
|\tilde{\Psi}_{\text{crystal}}(k''_1, k''_2)\rangle \propto \tilde{E}_\rho(k''_1, k''_2) \tilde{\Xi} (k''_1, k''_2) a_{k''_1}^\dagger a_{k''_2}^\dagger |0, 0\rangle,
\]

where the tilde sign \( \tilde{\cdot} \) denotes Fourier transform and \( k''_1 (k''_2) \) is the transverse component of the wave vector of the signal (idler) photon. \( \tilde{E} \) is the momentum representation of the pump profile in the crystal centre plane and \( \tilde{\Xi} \) is the phase matching profile. Their exact shape can be specified later. Transferring the above function from momentum to position space gives

\[
\Psi_{\text{crystal}}(x''_1, x''_2) \propto \int \int dk''_1 dk''_2 \tilde{E}_\rho(k''_1, k''_2) \tilde{\Xi} (k''_1, k''_2)e^{-ik''_1 x''_1} e^{-ik''_2 x''_2},
\]

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where we have suppressed the Dirac notation with kets, as the information about subsequent interference lies in the probability amplitudes. The vector \( x''_{\ell} = (x''_{\ell,1}, y''_{\ell}) \), \( \ell \in \{1, 2\} \), is the transverse coordinates in the crystal centre plane and all the integrals here, and in the rest of the paper, are to be taken over the whole real axis.

In the paraxial approximation, i.e. when the angle between the signal and the idler is not too large, one can write

\[
\Psi_{\text{slit}}(x''_1, x''_2) \propto \int \int dx''_1 \, dx''_2 \, \Psi_{\text{cystal}}(x''_1, x''_2) \, h'(x''_1, x''_2) \, h'(x''_2, x''_2), \quad (A.3)
\]

where \( h' \) is the field propagator ([41] chapter 20) which can be specified explicitly later on as well. The propagator \( h' \) allows any paraxial system including lenses and other components, but is especially simple for free-space propagation. The transverse coordinates just at the slit plane are denoted as \( x'' \). After the double slit, we obtain

\[
\tilde{\Psi}_{\text{slit}}(k''_1, k''_2) \propto \int \int \, dx''_1 \, dx''_2 \, \tilde{\Psi}_{\text{slit}}(x''_1, x''_2) \, f(x''_1) \, f(x''_2) \, e^{ik''_1 \cdot x''_1} \, e^{ik''_2 \cdot x''_2}, \quad (A.4)
\]

where \( k''_\ell \) is the transverse wave vector of the signal (\( \ell = 1 \)) or idler (\( \ell = 2 \)), just after the slit, but in the slit plane, and where we have assumed that the thickness of the slit is negligible. With \( a \) being the width of a slit, \( \lambda < a < d \), and \( d \) being the distance between the two slits, one can define

\[
f(x''_\ell) = \begin{cases} 
1 & \text{for } (d/2) - a/2 < |x''_\ell| < (d/2) + a/2, \\
0 & \text{else},
\end{cases} \quad (A.5)
\]

i.e. the function is unity at the slit openings and 0 otherwise. (The slit function \( f \) of course also depends on the width \( a \), but we shall suppress that in the following as we shall be interested in relative rather than absolute values of fields and intensities.) We have oriented the slits so that they are oriented along the \( y'' \)-direction and have their slit modulation structure in the \( x'' \)-direction.

Finally, we transfer the system to position space again and let it propagate to a detection plane so that we obtain

\[
\Psi_{\text{detect}}(x_1, x_2) \propto \int \int \int \int dx'_1 \, dx'_2 \, dk'_1 \, dk'_2 \, \tilde{\Psi}_{\text{slit}}(k''_1, k''_2) \, h(x_1, x'_1) \, h(x_2, x'_2) \, e^{-ik'_1 \cdot x'_1} \, e^{-ik'_2 \cdot x'_2}, \quad (A.6)
\]

where \( x'_1, x'_2 \) are the transverse coordinates directly after the slit and \( x_1 \) and \( x_2 \) are the transverse coordinates of the detector positions for detecting the signal and the idler, respectively. Combining all the above formulae, one ends up with

\[
\Psi_{\text{detect}}(x_1, x_2) \propto \int \, dx'_1 \, \int \, dx'_2 \, \int \, dk'_1 \, \int \, dk'_2 \, \int \, dx''_1 \, \int \, dx''_2 \, \int \, dx''''_1 \, \int \, dx''''_2 \, \int \, dk''''_1 \times \int \, dk''''_2 \, \tilde{E}_p(k''''_1, k''''_2) \, \tilde{\Omega}(k''''_1, k''''_2) \, h(x_1, x'_1) \, h(x_2, x'_2) \, h'(x''_1, x''_2) \, h'(x''_2, x''_2) \times f(x''_1) \, f(x''_2) \, e^{-ik''''_1 \cdot x'_1} \, e^{-ik''''_2 \cdot x'_2} \, e^{ik''_1 \cdot x''_1} \, e^{ik''_2 \cdot x''_2} \, e^{-ik''_1 \cdot x''''_1} \, e^{-ik''_2 \cdot x''''_2}. \quad (A.7)
\]

Without loss of generality we can assume that the detectors are moved only perpendicular to the slits. It is then sufficient to evaluate all the integrals above only in the first \( x \) component of
the vectors. Thus it suffices to write scalars instead of vectors, so in the paraxial approximation the field propagator can be written as

\[ h'(x''_1, x'''_1) = \sqrt{\frac{i}{\lambda \Delta z}} e^{i \pi \Delta z (x''_1 - x'''_1)^2} \]  

(A.8)

and correspondingly for the other variables, where \( \Delta z = z'' - z''' \) is the distance of propagation. If \( \Delta z \) is small, then \( h'(x''_1, x'''_1) \) can be approximated by a \( \delta \)-function \( \delta(x''_1 - x'''_1) \).

Now we will make some assumptions. If the slits are very close to the crystal, we have \( \Delta z \to 0 \) and therefore \( h'(x''_1, x'''_1) \to \delta(x''_1 - x'''_1) \). Assuming that the slits are narrower than any spatial structure of the two-photon field, we can approximate

\[ f(x''_1) f(x'''_1) \propto \delta(x''_1 - d/2) \delta(x''_1 + d/2) \]

where we have assumed that the two slits are identical.

To proceed, the integration over \( k'_\ell \) can be simplified as

\[ \int dk'' \ e^{ik''x'} e^{-ik''x} \to \delta(x' - x') \]  

(A.10)

Combining all equations and carrying out the integrals leads to

\[ \Psi(x_1, x_2) \propto \int dk'' \int dk'''' \tilde{E}_p(k''_1, k''_2) \tilde{Z'}(k''_1, k''_2) \left[h(x_1, d/2)h(x_2, d/2) e^{-ik''d/2} e^{-ik''d/2} \right. \\
+ h(x_1, -d/2)h(x_2, -d/2) e^{-ik''d/2} e^{-ik''d/2} + h(x_1, -d/2)h(x_2, d/2) e^{ik''d/2} e^{-ik''d/2} \\
+ h(x_1, d/2)h(x_2, -d/2) e^{ik''d/2} e^{-ik''d/2} \right] \\
= c_{11}h(x_1, d/2)h(x_2, d/2) + c_{12}h(x_1, d/2)h(x_2, -d/2) + c_{21}h(x_1, -d/2)h(x_2, d/2) \\
+ c_{22}h(x_1, -d/2)h(x_2, -d/2), \]  

(A.11)

where for example,

\[ c_{11} = \int dk'' \int dk'''' \tilde{E}_p(k''_1, k''_2) \tilde{Z'}(k''_1, k''_2) e^{-ik''d/2} e^{-ik''d/2} \]

\[ = \int du \int dv E_p(u, v) \Xi (d/2 - u, d/2 - v) \]  

(A.12)

In the following, since we discuss interference that requires indistinguishability, we shall only consider the preparation of the symmetric state subspace at the slits. This requires symmetry of the functions \( E \) and \( \Xi \), implying that \( c_{11} = c_{22} \) and \( c_{12} = c_{21} \). Removing any global phase factor, and ignoring information about the absolute amplitudes, i.e. renormalizing \( c_{ij} \) to \( \tilde{c}_{ij} \), such that \( |\tilde{c}_{12}|^2 + |\tilde{c}_{11}|^2 = 1 \), we can hence parametrize the amplitudes \( \tilde{c}_{11} = \tilde{c}_{22} = e^{i\theta} \sin(\alpha/2), \tilde{c}_{12} = \tilde{c}_{21} = \cos(\alpha/2). \) Apart from determining the probability amplitudes, the form of the wave function at the screen is independent of how the pump beam looks and the particulars of the crystal, if one only imposes symmetry with respect to the two slits. Of course, the particulars of the pump beam and the crystal determine the absolute magnitudes of \( c_{11} \) and \( c_{12} \).

We want to stress that the requirement for symmetry \( (c_{11} = c_{22} \text{ and } c_{12} = c_{21}) \) is automatically fulfilled if \( \tilde{E} \) and \( \tilde{Z} \) are invariant under the simultaneously exchange of sign in both \( k''_1 \) and \( k''_2 \). The pump beam profile \( \tilde{E} \) is usually a Gaussian beam depending only
on the squared sum of the arguments, i.e. \((k_1'' + k_2'')^2\), which makes it invariant under sign exchange, whereas the phase matching profile \(\Xi\) under usual crystal orientation only depends on \((k_1'' - k_2'')^2\), making it invariant under sign exchange as well. However, it turns out that even under more advanced crystal orientations (the general case is given in equation (2) of [42]) the function \(\Xi\) stays invariant under sign exchange of their arguments, although the dependence of the argument becomes more complex. Thus our symmetry requirement is almost naturally fulfilled for any crystal system as long as the slits are symmetrical to the optical axis.

Above, we assumed that the crystal is sufficiently close to the slit so that \(\Delta z \to 0\). This assumption, however, can be dropped since any \(\Delta z\) would transfer \(\Psi_{\text{crystal}}(k_1'', k_2'')\) to some \(\Psi_{\text{slit}}(k_1'', k_2'')\), and from there on, the analysis will be identical, ending up with expression (A.11) in this case too. This expression is, in other words, very general. The assumptions we cannot drop are that the slits are symmetrically excited by the two-photon field and that the field does not vary appreciably over the width of the slits.

Inserting (A.8) into (A.11) and considering a NOON state \((\alpha = \pi)\) where both photons either take one slit or the other, substituting \(2\pi/\lambda = k\), and assuming a propagation distance \(R\) between the slits and the detector, we obtain

\[
\Psi_{\text{NOON}}(x_1, x_2) \propto e^{i(kd/2R)(x_1^2 + x_2^2)} \left[ e^{-i(kd/2R)(x_1 + x_2)} + e^{i(kd/2R)(x_1 + x_2)} \right].
\]

The detection probability at the detectors hence becomes

\[
P_{\text{NOON}}(x_1, x_2) \propto \cos^2 \left( \frac{kd}{2R} (x_1 + x_2) \right),
\]

or, if \(x_1 = x_2\), to \(P_{\text{NOON}}(x_1) \propto \cos^2 \left( \frac{kd}{R} x_1 \right)\). Equation (A.14) is of the same form as (9) (with \(bs = x_1\) and \(bt = x_2\)), but now we have arrived at it through a more thorough mathematical analysis.

If we instead consider a state where the photons ‘chose’ slit independently, so that \(c_{11} = c_{12} = c_{21} = c_{22}\), or \(\alpha = \pi/2\) and \(\varphi = 0\), one obtains the coincidence probability

\[
P(x_1, x_2) \propto \cos^2 \left( \frac{kd}{2R} x_1 \right) \cos^2 \left( \frac{kd}{2R} x_2 \right),
\]

or, if \(x_1 = x_2\), to \(P(x_1) \propto \cos^4 \left( \frac{kd}{2R} x_1 \right)\). Comparing these expressions shows that the NOON state leads to a doubling of the period, whereas allowing the photons to ‘chose’ slit independently does not.

Inserting (A.8) into (A.11) and varying the probability amplitude parameters \(\alpha\) and \(\varphi\), one obtains figure A.1. These pictures correlate extremely well with figure 5 in Peeters et al’s paper [30] and show that the theory above provides an excellent description of their experiment. (Or rather, that Peeters et al have to be applauded for their ability to generate the various two-photon, two-slit states they aim for.)

Note that these figures are not ‘ordinary’ interference pictures, as they represent the correlations between two ‘point-detectors’ along the same line, and not the intensity as a function of two-dimensional space. The idea of quantum lithography is illustrated in figure A.1(a), which describes NOON-state interference, (A.14), and should be compared to figure 3. The same pattern as for single-photon detection in (3) can be seen when one detector, say \(x_2\), is kept at a fixed position and only the other detector is moved. In that case one finds four and a half interference maxima along the \(x_1\)-axis. The two-photon absorption pattern is given by requiring that both detectors have the same coordinate, i.e. \(x_1 = x_2\). This corresponds
Figure A.1. Different plots of (A.11) where each axis represents one detector coordinate, i.e. $x_1$ and $x_2$. A lighter colour indicates a higher intensity. The parameters are (a) $\alpha = \pi$, $\varphi = 0$, (b) $\alpha = 3\pi/2$, $\varphi = 0$, (c) $\alpha = \pi/2$, $\varphi = \pi/2$, (d) $\alpha = 0$, $\varphi = 0$, (e) $\alpha = \pi/2$, $\varphi = 0$ and (f) $\alpha = \pi/4$, $\varphi = 0$. These calculated plots should be compared to the experimental results of Peeters et al, presented in figure 5 of their paper [30].

to the diagonals in the figures, and going from the lower left to the upper right corner along the diagonal in (a), one counts nine maxima, i.e. a doubling of the number of fringes as compared to the single-photon case. In contrast, one does not see this period doubling for $x_1 = x_2$ in (e), which corresponds to (A.16), a state without spatial entanglement at the slits. This state will be produced, for example, if the slits are located far from the crystal. In this case, the state after the slits becomes

$$
\Psi(x_1, x_2) = (|2, 0\rangle + \sqrt{2} |1, 1\rangle + |0, 2\rangle)/2 = \frac{(a^\dagger + b^\dagger)^2}{2\sqrt{2}}|0, 0\rangle,
$$

(A.16)

where $a^\dagger$ ($b^\dagger$) is the creation operator of a photon through slit 1 (slit 2). Hence, the transmission probabilities are those expected for two independent particles. As can be seen from (A.15), the resulting two-photon interference pattern is the product of two single-photon interference fringe patterns.

The peculiar pattern in figure A.1(c) is the result of a superposition $(|2, 0\rangle - i\sqrt{2} |1, 1\rangle + |0, 2\rangle)/2$ between a NOON state and the state $|1, 1\rangle$ in such a way that it cannot be written as the product of two independent photons. Figure A.1(d) describes the interference of the state $|1, 1\rangle$, and as expected, it does not show any same-position ($x_1 = x_2$), two-photon interference but does show different position two-photon correlation fringes. However, what we ask the reader to take
home from figure (A.1) is that Steuernagel’s description of quantum lithography describes the experiment of Peeters et al (correlations are predicted also where \( s \neq t \)), whereas the theory of Boto et al does not.

**Appendix B. Closer inspection of the original proposal of Boto et al**

We now assume the same situation as in the proposal of Boto et al [8] and in the experiment of D’Angelo et al [9], i.e. two Gaussian beams are meeting at a point \( z = 0 \), where the angle between the direction of propagation of a Gaussian beam and the \( z \)-axis is \( \pm \alpha \) and one of the beams is ‘above’ the \( z \)-axis, whereas the other one is ‘below’ so that the situation is symmetrical with respect to the \( z \)-axis. The two positions where the waists of the Gaussian beams are minimal are at a distance \( L \) from the point \( x = y = z = 0 \), where we put the \( x \)-axis such that the two positions with minimum waists are at \( x = \pm L \sin(\alpha) \), \( y = 0 \) and \( z = -L \cos(\alpha) \). The envelope function of a Gaussian beam at this minimum waist position can be written as

\[
\psi_G(r) = \psi_0 e^{-ikz} \exp \left( -\frac{r^2}{2q} \right),
\]

where \( \psi_0 \) is a normalization constant and

\[
\frac{1}{q} = \frac{1}{\rho} - \frac{i}{\pi w^2},
\]

where \( \rho \) is the beam’s phase radius of curvature, \( w \) the beam waist radius and \( r \) the radial distance from the beam’s optical axis. Moreover, at the minimum waist position we have \( \rho = \infty \).

The state for \( N \) photons in a Gaussian beam is then given by

\[
|\Psi_0(r_1, r_2, \ldots, r_N)\rangle = \psi_G(r_1)\psi_G(r_2) \cdots \psi_G(r_N) a_{r_1}^\dagger a_{r_2}^\dagger \cdots a_{r_N}^\dagger |0, 0, \ldots, 0\rangle.
\]

Here too, we shall suppress the Dirac notation in the following. The probability amplitude of the state after having propagated a distance \( L \) along its axis of propagation is given by

\[
\Psi_L(r'_1, r'_2, \ldots, r'_N) = e^{-iNkL} \left( \int \right)^N \psi_G(r_1) h(r'_1, r_1) h(r'_2, r_2) \cdots h(r'_N, r_N)
\]

\[
\times \psi_G(r'_1) \psi_G(r'_2) \cdots \psi_G(r'_N) d\mathbf{r}_1 d\mathbf{r}_2 \cdots d\mathbf{r}_N,
\]

where \( h(r', r) \) is the field propagator. For Gaussian beams propagating a distance \( L \) these integrals can be solved and one can write

\[
\Psi_L(r'_1, r'_2, \ldots, r'_N) = e^{-iNkL} \psi_G(r'_1) \psi_G(r'_2) \cdots \psi_G(r'_N),
\]

where [41]

\[
\psi_G(r') = \psi_0 \sqrt{\frac{1}{1 + L/q}} \exp \left( -i k \frac{r'^2}{2(q + L)} \right).
\]

Coming back to the case of the proposal and the experiment, the detector is placed at \( y = z = 0 \) with \( x \) as the free variable and not with \( r' \) as in our description above. We therefore have to change the coordinates in (B.6). To do that, however, only photons arriving at \( x = 0 \) will have travelled the distance \( L \), whereas photons arriving at any other position will have travelled the distance \( L \mp \Delta L(x) \), where the minus sign is for the case of the Gaussian beam coming from above the \( z \)-axis and the plus sign is for the case where the Gaussian beam comes from...
below the $z$-axis. One has the following geometrical relations:

$$
\sin(\alpha) = \Delta L/r', \quad \cos(\alpha) = x/r', \quad \tan(\alpha) = \Delta L/x,
$$

where $r'$ is the coordinate of $r'$ that is in the plane spanned by $x$ and $z$ (the other coordinate of $r'$ is parallel to $y$ and will not change by our coordinate transformation). From now on, we will for simplicity, but without loss of generality, set $y = 0$, since the detector is placed at $y = 0$ and the coordinate systems of both Gaussian beams have their $y$-axis oriented in the same direction. Equation (B.6) then transforms to

$$
\psi'_G(x) = \psi_0 \sqrt{\frac{1}{1 + [L \mp x \sin(\alpha)]/q}} \exp \left( -ik \frac{[x \cos(\alpha)]^2}{2[q + L \mp x \sin(\alpha)]} \right). \quad (B.8)
$$

Since $q$ in general is a complex number both the square root and the exponential function in (B.8) are complex numbers. To be able to calculate any interference effects between the two Gaussian beams, we will split the two functions explicitly into their real and imaginary parts.

Introducing the dimensionless parameter $eta = \lambda^2 L^2/(\pi^2 w^4)$ and looking only at the case when the minimum waist of the Gaussian beam is not too close to the detectors, i.e. $L \gg x \sin(\alpha)$, we can simplify the above expressions to

$$
\psi'_G(x) = \psi_0 (1 + \beta)^{-1/4} e^{i(\sqrt{\beta})} e^{-i(k \sqrt{\beta} x^2 \cos^2(\alpha)/2L(1+\beta^{-1}))} e^{-i\frac{k x^2 \cos^2(\alpha)}{2L(1+\beta^{-1})}}
$$

The second and the last exponential functions on the right-hand side are the important terms here. The second term will be the Gaussian envelope and the last term will depend on whether a photon originated from the upper or the lower path. Having changed the coordinates of $\psi'_G$, we have to be careful to also change the expression $e^{-i\sqrt{\beta}L}$ in (B.4) and (B.5), since $L$ there differs as well, depending on where the detector is placed. This leads to the replacement

$$
e^{-i\sqrt{\beta}L} \rightarrow e^{-ik \sum_{i=1}^N x_i} = e^{-i\sqrt{\beta}L} e^{-ik \sin(\alpha) \sum_{i=1}^N x_i}, \quad (B.10)
$$

where the $\mp$ sign is defined in the same manner as above.

So far all the calculations have considered a single Gaussian beam. In the proposal of Boto et al, and in the experiment by D’Angelo et al, we have a NOON state, i.e. a superposition of $N$ photons in an upper Gaussian beam and zero in a lower one and vice versa. For such a case, the probability amplitude $\Psi$ at the place of the detectors $(z = 0)$ considering the results above can, after some algebra, be written as

$$
\Psi_{\text{NOON}}(x_1, x_2, \ldots, x_N) = \sqrt{2} \left[ \frac{\psi_0}{(1 + \beta)^{1/4}} \right]^N e^{-i\sqrt{\beta}L} e^{i(\sqrt{\beta})} e^{-i(k \sqrt{\beta} \sum_{i=1}^N x_i^2/2L(1+\beta^{-1}))}
$$

$$
\times e^{-i(k \cos^2(\alpha) \sum_{i=1}^N x_i^2/2L(1+\beta^{-1}))} \cos \left( k \sin(\alpha) \sum_{i=1}^N x_i - \frac{k \cos^2(\alpha) \sin(\alpha) \sum_{i=1}^N x_i^3}{2L^2(1+\beta^{-1})} \right). \quad (B.11)
$$

If we only look at the case of detecting all the $N$ photons at the same place (i.e. $x_1 = x_2 = \ldots = x_n = x$), we obtain the detection probability

$$
P_{\text{NOON}}(x, \ldots, x) = 2 \left[ \frac{\psi_0}{(1 + \beta)^{1/4}} \right]^{2N} e^{-i(k \cos^2(\alpha) \sqrt{\beta} N x^2/L(1+\beta^{-1}))}
$$

$$
\times \cos^2 \left[ k \sin(\alpha) N x - \frac{k \cos^2(\alpha) \sin(\alpha) N x^3}{2L^2(1+\beta^{-1})} \right]. \quad (B.12)
$$

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To see any interference effects we have to demand that the cosine function oscillates faster than the exponential envelope function vanishes. This is fulfilled either for

\[ k \sin(\alpha)Nx > \frac{k \cos^2(\alpha)\sqrt{\beta}Nx^2}{L(1+\beta^{-1})} \]  
(B.13)

or for

\[ \frac{k \cos^2(\alpha)\sin(\alpha)Nx^3}{2L^2(1+\beta^{-1})} > \frac{k \cos^2(\alpha)\sqrt{\beta}Nx^2}{L(1+\beta^{-1})}. \]  
(B.14)

The above relations are fulfilled if we have \( x < \frac{(1+\beta^{-1})\sin(\alpha)\pi w^2}{\lambda} \) or \( x > 2\pi w^2/[\lambda(\sin(\alpha))] \), where we have used the definition of \( \beta \). For \( \alpha > \pi/4 \), one of the two conditions is always fulfilled. To get a feel for these terms, we can rewrite the cosine function in (B.12) as

\[ \cos^2 \left[ 2\pi \sin(\alpha)Nx \frac{x}{\lambda} - \frac{2\pi k \cos^2(\alpha)\sin(\alpha)N}{2L^2(1+\beta^{-1})}\left(\frac{x}{\lambda}\right)^3 \right] \]  
(B.15)

and look at a typical example where we set \( \alpha = 30^\circ, \lambda = 1 \mu m, L = 10 \text{ cm}, w = 1 \text{ mm} \) and \( x = 0.1 \text{ mm} \). For these parameters the pre-factor of an \( x/\lambda \)-term in the cosine function argument is 14 orders of magnitude higher than the pre-factor of the \( (x/\lambda)^3 \) term, so that the latter can safely be neglected in this case.

Irrespective of which condition is fulfilled, the distance between two interference fringes scales with \( 1/N \) when going from 1 to \( N \) photons and the width of the envelope function scales with \( 1/\sqrt{N} \). This is the behaviour predicted by Boto et al. However, if we look again at (B.11) for the case \( N = 2 \) and with the assumption that \( L \) is sufficiently large, so that the \( x^3 \)-term in the cosine function can be neglected, we obtain

\[ P(x_1, x_2) = 2\psi_0^4 \frac{\psi_1^4}{1+\beta} e^{-ik\cos^2(\alpha)\sqrt{\beta}(x_1^2+x_2^2)}/L(1+\beta^{-1}) \cos^2 [k \sin(\alpha)(x_1+x_2)]. \]  
(B.16)

This function produces the same kind of pattern as in Steuernagel’s model and shows that our somewhat simple description, (6), not only describes the case of the double-slit experiment, but describes also the case of a NOON state shared between two Gaussian beams, such as the setup of D’Angelo et al. In the derivation of (B.16), we have assumed nothing but Gaussian beams and that the original state is in a NOON state distributed over the two arms.

We will now redo the previous analysis, but now assuming that all the photons are always produced at the same spatial coordinate. Instead of (B.3) we use

\[ \psi_0(r_1, r_2, \ldots, r_N) = \psi_0' e^{-ikzN} e \left( -\frac{k r_1^2}{2q} \right) \delta(r_1-r_2)\delta(r_1-r_3)\cdots\delta(r_1-r_N) \]  
(B.17)

as our input state probability amplitude, where \( \psi_0' \) is the new normalization factor. This state is at the origin of the theory of Boto et al. The implication is that the first photon can essentially be found anywhere within the beam’s Gaussian envelope, but the rest of the photons will subsequently be found at the same position as the first.

In the same manner as before, we can calculate the state after a distance \( L \) with the help of the propagators. This leads to

\[ \psi_L(r_1', r_2', \ldots, r_N') = \psi_0' e^{-i\hbar k L} \sqrt{\frac{1}{N+L/q}} \left[ \frac{1}{\lambda L} \right]^{N-1} \exp \left( -\frac{k}{2L} \left[ \sum_{i=1}^{N} (r_i')^2 / (N+L/q) - \sum_{i=1}^{N} r_i'^2 \right] \right). \]  
(B.18)
which, looking at the case where all the photons are detected at the same place, i.e. \( r_1 = r_2 = \cdots = r_N := (x, 0) \), and assuming the NOON state, leads to

\[
\Psi_{\text{NOON}}(x) = \sqrt{2} \psi_0' e^{-iNkL} \left( \frac{1}{N^2 + \beta} \right)^{1/4} \sqrt{\frac{N-1}{\lambda L}} e^{(6/2) \arctan(\sqrt{\beta}/N)} e^{i(kN^2 \cos^2(\alpha)/2L(\beta+N^2))} \\
\times e^{-(kN^2 \sqrt{\beta} x^2 \cos^2(\alpha)/2L(\beta+N^2))} \cos \left[ k (\sin \alpha) N x \right],
\]

where we did the same replacements as in (B.7) and (B.10), setting \( y = 0 \) as before, and ignoring terms of higher order than \( x^2 \) as before.

This gives the \( N \)-photon detection probability

\[
P_{\text{NOON}}(x) = 2\psi_0'^2 \sqrt{\frac{1}{N^2 + \beta}} \left( \frac{1}{\lambda L} \right)^{N-1} e^{-(kN^2 \sqrt{\beta} x^2 \cos^2(\alpha)/L(\beta+N^2))} \cos^2 \left[ k (\sin \alpha) N x \right].
\]

This expression should be compared with (B.12), where the \( x^3 \) term is neglected. One sees immediately that both equations lead to the same distance between nearby peaks, given by the cosine term. (This distance halving is what has been mainly been looked for in experiments so far.) Writing the equivalent of (B.16) when starting with (B.17) and (B.18) leads to

\[
P(x_1, x_2) = 2\psi_0'^2 \sqrt{\frac{1}{4 + \beta/\lambda L}} e^{-(k \cos^2(\alpha) \sqrt{x_1 + x_2^2}/L(\beta+4))} \cos^2 \left[ k (\sin \alpha) (x_1 + x_2) \right].
\]

Thus, in neither (B.16) nor (B.21) are the photons constrained to arrive at the same place as assumed by Boto et al.

A difference between (B.12) and (B.20) lies in the exponential function. If one chooses \( L \) and \( w \) carefully so that one has \( \beta \gg 1 \), the argument of the exponential function is proportional to \( \beta^{1/2} \propto w^{-2} \) in (B.12), but proportional to \( \beta^{-1/2} \propto w^2 \) in (B.20). Since the argument of the exponential function determines the number of fringes that one can see, one could, by changing \( w \), determine which of the states one has produced, since in one case one would see more fringes, but in the other case fewer fringes. This may be of interest for further experimental work in this field.

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