The logical density matrix elements take the form

$$\rho_{ij;kl}(t) = \text{Tr}_{\mathcal{E}}(\rho_{\mathcal{E}}(0)X_{kl}^{\dagger}X_{ij}), \tag{31}$$

with Tr_E indicating the trace over the environment degrees of freedom and $\rho_E(0)$ the initial state of the environment.

To keep the discussion concise we assume that the photon pulses are very sharp. We then can replace the squares of the wavefunctions by δ distributions and get

$$X_{ij} = c_{ij}e^{-\mathrm{i}i(1-j)\phi_2}R_1(v_1t + \delta z_i^{(1)}, t)R_2(d + v_2t + \delta z_j^{(2)}, t).$$
(32)

The density matrix elements then become

$$\rho_{ij;kl}(t) = c_{ij}c_{kl}^* e^{-i(i(1-j)-k(1-l))\phi_2} N_{ij;kl}(t), \qquad (33)$$

with the Hermitian decoherence matrix

$$N_{ij;kl}(t) \equiv \left\langle R_2^{\dagger}(d + v_2 t + \delta z_l^{(2)}, t) R_1^{\dagger}(v_1 t + \delta z_k^{(1)}, t) \right. \times \left. \left. \left. \left. \left(x_1 t + \delta z_i^{(1)}, t \right) R_2(d + v_2 t + \delta z_j^{(2)}, t) \right. \right\rangle \right).$$

$$(34)$$

In App. C3) we show that it can be written as

$$N = \begin{pmatrix} 1 & c_2 & c_1 e^{i(\phi_1 - \phi_2)} & N_{00;11} \\ c_2^* & 1 & N_{01;10} & c_1 \\ c_1^* e^{-i(\phi_1 - \phi_2)} & N_{01;10}^* & 1 & c_2 \\ N_{00;11}^* & c_1^* & c_2^* & 1 \end{pmatrix},$$
(35)

$$c_n \equiv \left\langle R_n^{\dagger}(v_n t + \delta z, t) R_1(v_n t, t) \right\rangle. \tag{36}$$

Hence the influence of decoherence in our model can be described by four complex parameters $N_{00;11}$, $N_{01;10}$, and c_n . This is a consequence of the approximation of very short pulses. For extended pulses additional decoherence contributions are anticipated.

We again calculate the concurrence of density matrix (33) for the initial product state $c_{ij} = \frac{1}{2}$ for all i, j. The specific values of the parameters $N_{00;11}$, $N_{01;10}$, and c_n depend on the detailed decoherence model that one employs so that the predictions may vary substantially depending on the assumptions behind the model. We present one particular decoherence model in Appendix D. However, all decoherence models must be consistent with all eigenvalues of the density matrix (22) taking values between 0 and 1.

To simplify the discussion we consider the special case $\phi_2 = \pi$ for the oft-used assumption that the two interaction potentials $V_i(z)$ are equal and independent of the group velocities [33, 34] so that $\mathcal{V}_1 = \mathcal{V}_2$. The phase factor that appears in the decoherence matrix (35) is then $\phi_1 - \phi_2 = \pi \Delta v/v_2 > 0$. For a given value of $\phi_1 - \phi_2$ we can derive a "minimal decoherence model" by finding those values of the parameters $N_{00;11}$, $N_{01;10}$, and c_n

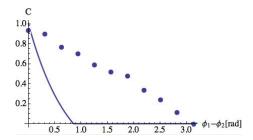


FIG. 3: Concurrence as a function of the XPM phase shift for minimal decoherence (dots) and the decoherence model presented in Appendix D (solid line).

which maximize the concurrence C for a consistent density matrix. Fig. 3 shows the concurrence for this minimal decoherence model as a function of $\phi_1 - \phi_2$. The dots are found through a random search for the parameters $N_{00:11}$, $N_{01:10}$, and c_n of the minimal decoherence model. Each dot is based on a sample of typically 10⁷ random events. The error in each value is estimated to be about 5%. The concurrence decreases with $\phi_1 - \phi_2$ and becomes zero for $\phi_1 - \phi_2 = \pi$. In this case decoherence completely destroys the entangling capacity of the XPM interaction. The numerically determined density matrix then corresponds to a nearly equal mixture of two highly entangled states. Realistic decoherence models would typically predict a less than optimal performance. For instance, the blue line in Fig. 3 shows C for the decoherence model of App. D under the fairly optimistic assumption $\Delta vt = 4\sqrt{\pi}w$, where w is the width of a Gaussian potential $V_1(z) = V_2(z) = V(0) \exp(-z^2/w^2)$. For more realistic values $\Delta vt \gg w$ [47] the concurrence would be non-zero only in a very narrow range around $\phi_1 = \phi_2$.

Our phenomenological model therefore suggests that a CPG may not be achievable with XPM unless ϕ_1 is very close to ϕ_2 . In most proposals for XPM based on double EIT this could be achieved by a suitable preparation of the atomic gas, although it may require fine tuning of parameters like magnetic fields or pump field intensities. A better way to achieve $\phi_1 = \phi_2$ would be to find a system in which this is guaranteed through microscopic symmetries.

B. Quantum non-demolition measurement of photon numbers

The scheme for a QND measurement of the photon number discussed in Sec. III A can easily be extended to the non-unitary case by calculating the measurement signal $\langle \mathcal{E}_1(z,t) \rangle$ using solution (27) instead of (5). The only change is that the measurement result (15) is multiplied by $\langle R_1(z,t) \rangle$. This factor depends strongly on the specific decoherence model but generally will lead to a decrease in the contrast of the phase measurement.

To give a rough estimate we consider the decoherence