Optimal preparation of quantum states on an atom chip device

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Atom chips provide compact and robust platforms towards practical quantum technologies. A quick and faithful preparation of arbitrary input states for these systems is crucial but represents a very challenging experimental task. This is especially difficult when the dynamical evolution is noisy and unavoidable setup imperfections have to be considered. Here, we experimentally prepare with very high small errors different internal states of a Rubidium Bose-Einstein condensate realized on an atom chip. As a possible application of our scheme, we apply it to improve the sensitivity of an atomic interferometer.

Microscopic magnetic traps, as atom chips\textsuperscript{1,2}, represent an important advance in the field of degenerate quantum gases towards their realization of practical quantum technological devices. Indeed, applications outside the laboratory depend on the compactness and robustness of these setups. Atom chips have already enabled, for instance, the demonstration of miniaturized interferometers\textsuperscript{3,4}, or the creation of squeezed atomic states\textsuperscript{5}. Very recently, they have been also applied to the realization of atom interferometers based on non-classical motional states\textsuperscript{6}. Furthermore, atom chips can be integrated with nanostructures\textsuperscript{7}, or photonic components\textsuperscript{8}, to implement new quantum information processing tools, or used as novel platforms for quantum simulations\textsuperscript{9} and controllable coherent dynamics\textsuperscript{10}.

As in most quantum technological applications, the ability to quickly and faithfully prepare arbitrary input states is of utmost importance. In particular, it is essential to speedup the initialization protocol of these quantum devices in order to reduce the effects of the inevitably present noise and decoherence sources. In this context, quantum optimal control provides powerful tools to tackle this problem by finding the optimal way to transform the system from an experimentally readily prepared initial condition to a desired state with high fidelity from which further quantum manipulations can be performed\textsuperscript{11,13}. It has been successfully implemented in several physical systems, ranging from cold atoms\textsuperscript{14} to molecules\textsuperscript{15}. Recently optimal control algorithms have been developed and successfully applied to many-body quantum systems\textsuperscript{16-21}.

In addition, optimal control allows one to speed up the state preparation process up to the ultimate bound imposed by quantum mechanics, the so-called quantum speed limit (QSL)\textsuperscript{22,25}. Indeed, the change of a state into a different one cannot occur faster than a time scale that is inversely proportional to the associated energy scale. This is especially relevant for quantum systems like Bose-Einstein condensates (BECs) where it is crucial to perform some desired (coherent) task before dephasing processes unavoidably occur\textsuperscript{26}.

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In this work we present experimental results obtained on a Rubidium ($^{87}$Rb) BEC, trapped on an atom chip, evolving in a five-level Hilbert space given by the five spin orientations of the $F = 2$ hyperfine ground states — see Fig. 1. We use optimal control theory to design frequency modulated radio-frequency (RF) pulses that, starting from a given initial state fixed by the BEC evaporation procedure, prepare different arbitrary coherent superpositions of states with high precision. We also show how the latter scales with the preparation time, with a highly satisfactory agreement between theory and experiment. We apply, finally, this technique to improve the sensitivity of an atom interferometer.

\textbf{Model:} In the presence of a magnetic field the degeneracy between the spin orientations of an atomic hyperfine state is lifted. The phenomenon is described by the Breit-Rabi formula which determines the energies of all different sublevels. For the five fold $F = 2$ hyperfine ground state of $^{87}$Rb in the presence of a constant magnetic bias field arbitrarily set to 6.179 G\textsuperscript{27}, the atomic part of the Hamiltonian is $H_0 = \hbar \text{ diag}(\omega_1, \omega_2, \omega_3, \omega_4, \omega_5) = \hbar [2\omega_1, 2\omega_2, 2\omega_3, 2\omega_4, 2\omega_5]$, where $\omega_i$ are the atomic transition frequencies.
we assume the same dephasing rate for all five sublevels, i.e. by \( \gamma \) the phase of each sublevel \( f \) is determined by the bias field, we also add in our model a Lindblad super-driving noise, mainly originated in our experiment by the presence of magnetic noise. In all our experiments, we set \( \Omega = 2 \pi \times 8657 \) kHz in order to keep the Rabi frequency \( \tilde{\Omega} \) additionally the optimal control protocols on a BEC of \( ^{87} \text{Rb} \) atoms produced in a magnetic micro-trap realized on an atom chip. The trap has a longitudinal frequency of 46 Hz while the radial trapping frequency is 950 Hz. The atoms are evaporated to quantum degeneracy by ramping down the frequency of a radio frequency (RF) field. The BEC has typically \( 8 \cdot 10^4 \) atoms, at a critical temperature of 0.5 \( \mu \text{K} \) and is 300 \( \mu \text{m} \) from the chip surface. The magnetic fields for the micro-trap are provided by a \( z \)-shaped wire on the atom chip and an external pair of Helmholtz coils. The RF fields for evaporation and manipulation of the Zeeman states are produced by two additional conductors also integrated on the atom chip. Finally, the homogeneous bias field that lifts the degeneracy between the spin orientations is created by two further external Helmholtz coils [10,30].

All manipulations are performed 0.7 ms after turning off the magnetic trap to guarantee bias field homogeneity and strongly reduce the effects of atomic collisions that can produce decoherence. In this way the most relevant source of noise on the evolution turns out to be the instability of the environmental magnetic field, which has a spectrum distribution made of several peaks on a decreasing envelope in the low frequency range. To record the number of atoms in each of the \( m_F \) states of the \( F = 2 \) hyperfine state we apply a Stern-Gerlach method. After the state manipulation has been performed, in addition to the homogeneous bias field, an inhomogeneous magnetic field is applied along the quantization axis for 10 ms. This causes the different \( m_F \) states to spatially separate. After a time of 23 ms of expansion a standard absorption imaging sequence is performed. The atomic populations in each \( m_F \) state is normalized to the total number of observed atoms.

\textbf{Results:} Let us define the theoretical state preparation error \( \epsilon_F \) and the experimentally measured one \( \tilde{\epsilon}_F \). We then apply ten times the same preparation pulse and the reported values of \( \epsilon_F \) are the corresponding means and standard deviations. We take into account the presence of magnetic noise in

\[
H_1 = \hbar f(t) \begin{pmatrix}
0 & \Omega & 0 & 0 & 0 \\
0 & 0 & \sqrt{3/2} \Omega & 0 & 0 \\
0 & \sqrt{3/2} \Omega & 0 & \sqrt{3/2} \Omega & 0 \\
0 & 0 & \sqrt{3/2} \Omega & 0 & \Omega \\
0 & 0 & 0 & 0 & \Omega \\
\end{pmatrix},
\]

(1)

where \( f(t) = \frac{d}{dt} \omega(t) \) describes the energy differences between the five \( F = 2 \) levels within the RWA, \( \omega(t) \) is the time dependent frequency of the coupling field that will be modulated in order to implement our optimal control scheme. The Rabi frequency \( \Omega \) is proportional to the RF field intensity and, in all our experiments, we set \( \Omega = 2 \pi \times 60 \) kHz. Hence, the driving field Hamiltonian is \( H_1 + V(t) \).

In order to include the unavoidable presence of dephasing noise, mainly originated in our experiment by the presence of environmental magnetic noise superimposed on the bias field, we also add in our model a Lindblad super-operator term \( \mathcal{L} \), acting on the density matrix \( \rho \) as \( \mathcal{L}(\rho) = \sum_{j=1}^{5} \gamma_j \{ -\{ j \rho | j \} + 2 \{ j \rho j \} \} \), which randomizes the phase of each sublevel \( j \) with a rate \( \gamma_j \). For simplicity, we assume the same dephasing rate for all five sublevels, i.e. \( \gamma_j = \gamma \). The time evolution of the density matrix is then given by

\[
\frac{d}{dt} \rho(t) = -\frac{i}{\hbar} [H_0 + H_1 + V(t), \rho(t)] + \mathcal{L}(\rho(t)).
\]

(2)

\textbf{Optimal control:} Here, we use a recently introduced versatile and efficient optimization algorithm (CRAB) [16,17] for the time dependence \( f(t) = f_0 (1 + \sum_{k=-n_f}^{n_f} A_k \exp(i \nu_k t)) \) with \( \nu_k = 2\pi k / T \), \( n_f \) being the number of frequencies, \( A_k \) complex numbers, and \( T \) the time duration of the modulation. The optimization is reflected in the modulation of the field’s frequency \( \omega(t) \), which we constrain within the range \( \omega(t) \in 2\pi [4000, 4700] \) kHz in order to keep the Rabi frequency \( \tilde{\Omega} \) constant (see Supplementary Information for further details). The optimization of the control parameters \( A_k \) in \( f(t) \) is done by means of the Subplex variant of the Nelder-Mead algorithm [29]. The error function to be minimized is \( \epsilon = \sum_i |\rho_i - b_i|/2 \), with \( \rho_i \) being the final population (after time \( T \), i.e. at the end of the pulse) of the sublevel \( i \), while \( b_i \) are the corresponding target populations we want to achieve. The initial state \( \rho_0 \) is given by all population in the energy level \( m_F = 2 \), i.e. \( \rho_{11}(t = 0) = 1 \). In the left half of Tab. \ref{tab:results} we show the target populations of different target states (\( A, \ldots, L \)). Note that there are no constraints on the coherence terms (i.e., the off-diagonal elements \( \rho_{ij} \)), which are however relevant in the context of quantum state preparation. Even if state tomography would have been in principle possible [28], it would have been hard to perform it for all cases studied here. For this reason, in the experiment we measure only the populations of the five levels, while in the numerical simulations we take into account the full density matrix during the optimization. However, we have strong experimental evidence that the coherence terms (i.e. relative phases) are also prepared according to our theoretical predictions, because: 1) as shown in the fine details of the inset of Fig. 3 there is a very good agreement between theory and experiment in the population dynamical evolution, 2) when we prepare eigenstates of the Hamiltonian (see Fig. 2 and in the Supplementary Information Fig. 6), they do not evolve anymore, as expected if the Hamiltonian is kept fixed, 3) finally, we consider a multi-state interferometer and our theoretical results are reproduced very well by the experimental data. These three tests would have failed if all the relative phases had not been prepared as in the target quantum state that is considered in the numerical optimization protocol.
In the absence of dephasing noise, the theoretically predicted magnetic field in a state for every initial state is a measure of distance (Uhlmann fidelity $F$) agreement between theory and experiment. We include also the population of the levels $m_F = 2$ (blue), $m_F = -2$ (red), and the sum of the remaining three levels (black) during the application of optimal pulse of length $100 \mu s$.

As noted above, these experimental results only show that we are able to prepare different population distributions in the $F = 2$ manifold. To directly verify our ability to prepare superposition of levels, we have prepared the ground state and highest excited state of the Hamiltonian $H_0 + H_1 + V(t)$ in which $f(t) = \overline{f} = 2\pi 4323 \text{ kHz}$, and observed the absence of any evolution of the population distribution of these states after applying the preparation pulse, i.e. for $t \geq T$, while keeping $f(t) = \overline{f}$ ($T = 20 \mu s$ is the pulse length). The results are reported in Fig. 2 where we plot the experimental points and expected theoretical behaviour. The theoretical prediction without any free parameter is shown as a continuous line. Some typical examples of optimal pulses are shown in the Supplementary Information.

We then proceed to analyse the time limits for the preparation of the states, in other words we are interested in keeping the pulse length as short as possible while still suppressing errors in the state preparation. For this analysis we choose the target state $A$, fixing different pulse times and minimizing the error function $\epsilon$ under the same constraint on magnetic field and Rabi frequency. In Fig. 3 we report the results obtained. The theoretical noiseless results saturate very close to vanishing errors for a pulse length of $90 \mu s$. For the experimental results we used the same procedure outlined above to determine the error value. Experimental results are perfectly matched to the theoretical noiseless value at short pulse times but deviate for longer pulses, due to the increasing effect of the low frequency experimental noise on the magnetic field.

\begin{figure}[h]
\centering
\includegraphics[width=\textwidth]{fig2.png}
\caption{Theoretical (lines) and experimental (dots) time evolution of $\epsilon$ for the preparation (grey area) of a ground state and the highest excited state of the Hamiltonian $H_0 + H_1 + V(t)$ with $V(t)$ being the harmonic potential that oscillates at $f(t) = \overline{f} = 2\pi 4323 \text{ kHz}$ for every $t \geq T = 20 \mu s$ (white area).}
\end{figure}

\begin{figure}[h]
\centering
\includegraphics[width=\textwidth]{fig3.png}
\caption{State error preparation $\epsilon$ as a function of pulse length $T$, obtained theoretically (black dots), in the absence of dephasing noise, and experimentally (red dots), to prepare the state $A$ by optimal control. The effect of dephasing noise in the range $2\pi [20, 200] \text{ Hz}$ is also considered (shaded region). The results achievable with constant $f(t) = \overline{f} = 2\pi 4323 \text{ kHz}$ are also shown (dashed line). Inset: Theoretical (lines) and experimental (dots) dynamical evolution of the population of the levels $m_F = 2$ (blue), $m_F = -2$ (red), and the sum of the remaining three levels (black) during the application of optimal pulse of length $100 \mu s$.}
\end{figure}

\begin{table}
\centering
\begin{tabular}{|c|c|c|c|c|c|c|}
\hline
Target State & $\rho_{11}$ & $\rho_{22}$ & $\rho_{33}$ & $\rho_{44}$ & $\rho_{55}$ & $\epsilon_T$ & $\epsilon_E$ & $\mathcal{F}$ \\
\hline
A & 1/2 & 0 & 0 & 0 & 1/2 & 0.04(3) & 0.07(1) & 0.71 \\
B & 1/2 & 0 & 0 & 1/2 & 0 & 0.04(2) & 0.02(1) & 0.67 \\
C & 0 & 1/2 & 0 & 1/2 & 0 & 0.04(3) & 0.04(1) & 0.11 \\
D & 1/2 & 1/2 & 0 & 0 & 0 & 0.03(2) & 0.02(1) & 0.71 \\
E & 0 & 1/5 & 1/3 & 1/3 & 0 & 0.04(2) & 0.03(1) & 0.02 \\
F & 1/5 & 1/5 & 1/5 & 1/5 & 1/5 & 0.02(1) & 0.03(1) & 0.45 \\
G & 0 & 1 & 0 & 0 & 0 & 0.05(4) & 0.04(1) & 0.15 \\
H & 0 & 0 & 0 & 1 & 0 & 0.04(3) & 0.03(1) & 0.07 \\
I & 0 & 0 & 1 & 0 & 0 & 0.07(3) & 0.07(1) & 0.15 \\
\hline
\end{tabular}
\caption{Target populations $\rho_{ii}$ of different target states (A, ..., I), theoretical ($\epsilon_T$) and experimental ($\epsilon_E$) error functions in their preparation, and fidelity $\mathcal{F}$ between the initial and the target state. The error bars for $\epsilon_T$ are obtained considering the maximal deviation when the noise rate $\gamma$ is varied in the range $\gamma \in [20, 200] \text{ kHz}$ and the magnetic field in a 1 mG range around the ideal value $B = 6.179 G$. In the absence of dephasing noise, the theoretically predicted $\epsilon_T$ are always smaller than 0.02.}
\end{table}
If we include in the simulation the magnetic field and dephasing noise term fluctuations to correct all the theoretical predictions, we obtain a good agreement with the experimental observation. The precise matching between theory and experiment is also shown in the inset of Fig. 4 where we report the experimentally measured values of the atomic populations in the five levels during the application of the 100 µs-pulse. No fitting procedure was used to get the theoretical results (shown as continuous lines).

**Application to interferometry:** We have implemented an atomic Ramsey interferometer scheme using some of the state preparation pulses listed above. In a conventional Ramsey interferometer the application of a π/2 pulse creates a balanced superposition of two atomic states separated by an energy difference ΔE. After the pulse, the atoms are let free to evolve for a time τ, during which the states acquire a phase difference ΔE/hτ. Then, the application of a second π/2 pulse maps this phase difference onto the atomic populations, which then oscillate as cos(ΔE/hτ). Even if the smallest signal that an interferometer can ultimately resolve is limited by the signal to noise ratio, typically its sensitivity is determined by the maximum slope of the interferometric fringe. The best achievable sensitivity of a conventional Ramsey interferometer is found to improve linearly with ΔE.

The presence of five levels in the F = 2 manifold complicates the analysis of the interferometer albeit offering a possibility to realize a multi-state interferometer with increased sensitivity [30]. Whereas optimal control would allow to maximize the performance of an interferometer [6], for our system this would require an extensive analysis of the content of information of the interferometric output that goes beyond the focus of the present work. Nonetheless, as an application of our protocol, we restrict ourselves to interferometric configurations where, for simplicity, only two levels play a role during the free evolution. To this aim, we use the state preparation pulses A, B, C, D (all of them populating two initial states only), both as input and output beam splitters.

For each preparation pulse we record the interferometric pattern, i.e. the behaviour of the relative populations in the F = 2 manifold after applying the two pulses and a variable time delay τ between them. Among the five components of the interferometer’s signal we select the one with the highest fringe amplitude A and we calculate the point-to-point slope selecting its maximum value S. Since the fringe amplitude can be different for each preparation pulse, we take into account its effect on the sensitivity by plotting in Fig. 4 the values S/A with respect to the energy difference ΔE between the levels populated via the state preparation pulse. The results confirm, within the errors, the expected linear dependency between the two quantities. Let us point out that none of these interferometers can be realized without optimal control since no one of the states above (A, B, C, D) can be reached with constant pulses. It must be noted that, in our experiment, we have not taken into consideration the fringe amplitude which could be manipulated by a suitable choice of optimization parameters. However, this simple application still allows us to achieve a four-fold improvement of the interferometer sensitivity, without changing neither the applied magnetic field nor the time interval between the two pulses.

**Conclusions:** The preparation of specific quantum states is a crucial step to control atom-chip devices, especially in the context of quantum information and communication technologies. However, this is highly non-trivial to perform even theoretically, for instance for complex quantum systems, and especially from the experimental side since one has to deal with imperfections and the presence of noise due to the external environment, limiting the observation time scale. Over the last years, the theory of optimal control has allowed to address this issue in an efficient way, also when working with fragile quantum system and non-unitary (noisy) evolutions. Here, we have applied optimal control tools to prepare a family of different internal states of a Rb Bose-Einstein condensate produced in an atom chip-based micro-trap. In particular, we theoretically and experimentally find that these states can be prepared with a very small error (ε ~ 0.03), with very short (control) pulses successfully overcoming the unavoidable presence of dephasing noise. Hence, we also test these protocols to implement improved interferometric configurations that would have been not accessible without the preparation of some input quantum states by optimal control. Therefore, these results might pave the way for new schemes towards, among others, better control of quantum dynamics, preparation of squeezed states, time-inversion of the system dynamics, realization of quantum gates, and so more powerful quantum information protocols.
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SUPPLEMENTARY INFORMATION

As introduced in the main text, we use RF induced evaporation on magnetically trapped $^{87}$Rb atoms to reach the BEC. The magnetic trap is provided by the chip structures and, before its loading, we perform optical pumping in the $m_F=2$ magnetic sublevel. This condensation scheme imply that the initial state of the condensate is the $|F=2,m_F=2\rangle$ state, so as initial condition every state preparation pulse uses a system in that pure state. After the BEC creation we let expand the atomic cloud in the homogeneous magnetic field for 0.7 ms and then, using a programmable function generator, we apply to a wire on the chip the signal $V(t) = V_0 \sin[\omega(t)]$, which build up the driving field. In the frequency range [4000, 4700] Hz in which we constrain $\omega(t)$ the radiation efficiency of the used wire can be considered constant, so the amplitude $V_0$ of the signal can be determined with high accuracy by performing a series of atomic population oscillation to have a Rabi frequency $\Omega = 2\pi \times 60 \text{ kHz}$.

To calibrate the homogeneous magnetic field applied by the Helmholtz coils, we close a multi-state atom interferometer with constant amplitude pulses as in [30] from which we deduce the right current to apply to the coils to produce the value $B_{1h} = 6.179 \text{ G}$. To compensate for very slow variations of the magnetic field we repeat the calibration procedure before the sequence of ten measurements of a given state preparation pulse.

From a statistical analysis of the interferometer output we also derive information on the residual noise present in the field. In particular, by the reduction of the fringe contrast we deduce the dephasing rate interval $\gamma \in 2\pi \times [20, 200]$, while by the variations of the fringe phase we can estimate the amount of shot-to-shot offset field fluctuations, to 1 mG. Introducing these parameters in the theoretical model it is possible to derive also a maximal theoretical error by taking the worst and the best output of the model evolution. Those errors are reported in table I.

To observe the evolution during the optimization pulse, we splice it in subpulses, measuring the output at the end of each of them. Figure 5 shows an example of experimental points and theoretical prediction, reported in solid line, relative to the five population $m_F = +2, \ldots, -2$ driven from the initial state to the ground state of the Hamiltonian $H_0 + H_1 + \overline{V}(t)$ with $\overline{V}(t)$ being the harmonic potential that oscillates at $f(t) = \overline{f} = 2\pi \times 4323 \text{ kHz}$. The optimized preparation pulse is $20\mu s$ long and, after the preparation, we keep the frequency of the potential at the constant target value. The stability of the populations shows that the system is really in the target state, with the right coherence between the $|F,m_F\rangle$ components. Finally we show in Fig. 6 the optimal pulses for the preparation of the ground and excited states as in Fig. 5. Similar pulses (with $n_F = 7$) have been obtained as a result of our optimization scheme to prepare the other states discussed in the main text.
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