Coherent Quantum Engineering of Free-Space Laser Cooling

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We perform a quantitative analysis of the cooling dynamics of three-level atomic systems interacting with two distinct lasers. Employing sparse-matrix techniques, we find numerical solutions to the fully quantized master equation in steady state. Our method allows straightforward determination of laser-cooling temperatures without the ambiguity often accompanied by semiclassical calculations, and more quickly than non-sparse techniques. Our calculations allow us to develop an understanding of the regimes of cooling, as well as a qualitative picture of the mechanism, related to the phenomenon of electromagnetically induced transparency. Effects of the induced asymmetric Fano-type lineshapes affect the detunings required for optimum cooling, as well as the predicted minimum temperatures which can be lower than the Doppler limit for either transition.

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Laser cooling has led to tremendous achievements, enabling extremely precise measurements (see, for example, Ref. 1), the development of highly accurate atomic clocks 2, and the production and manipulation of quantum-degenerate gases 3. The atomic internal structure greatly restricts the species for which laser cooling can be applied successfully, and determines whether ultracold temperatures (on the order of a few µK) can be achieved. The basic Doppler cooling mechanism 4 5, based on momentum transfer from a near-resonant light beam, has a minimum temperature (typically a few hundred µK) proportional to the scattering rate (linewidth) of the atomic transition, and given by the balance between friction (cooling) and diffusion (heating) in the atom-laser interaction. Lower temperatures can be achieved either by exploiting the multilevel hyperfine structure (Sisyphus 6 7, Raman 8, VSCPT 9), by modifying the atomic scattering (spontaneous emission) for example in a near-resonant cavity 10, or, more recently, by using weakly allowed, narrow two-level optical transitions in a second-stage for atoms that have been pre-cooled by other means 11 12. Here, we explore cooling by dissipative radiation-pressure forces 13 in three-level systems under two-color excitation. Similar schemes have been studied previously for trapped ions 14 15. In the electromagnetically-induced-transparency (EIT) regime, e.g., with a weak ”probe”, and a strong ”dressing” laser, an effective two-level system is engineered that allows sideband cooling of ions to the ground state of the trapping potential 16 17.

This Letter presents a general method that uses coherent quantum interference to tailor an atom's internal energy structure that makes it more amenable to simple Doppler cooling in free space. To exploit the effects of quantum interference, at least three internal atomic levels and two lasers with distinct frequencies must be utilized. There are three basic types of three-level systems — the Λ-, V-, and Ξ-systems — each classified according to the ordering of the bare quantum states in energy, and the possible decay pathways. The Λ-configuration is commonly used for studying of EIT 18 and related phenomena. Here we focus on a Ξ-type system, also known as a cascade configuration. We note, however, that our general conclusions can be applied to any type of three-level system, and can be extended to systems with more than three levels. In particular, we will focus our discussion on a cascade system suggested by Magno et al. 17 as a simpler second-stage cooling scheme for alkaline-earth atoms, which has recently been demonstrated for magnesium 19. The level structure and internal parameters for this system are depicted in Fig. 1. The cooling scheme is more effective for Λ-type three-level systems having metastable lower states, but also works for ladder systems with an upper state narrower than the intermediate one 17. This cooling technique seems well suited for alkaline-earths, which are good candidates for the next-generation optical atomic clocks, studies of ultracold collisions, optical Feshbach resonances 20, and achievement of quantum degeneracy 21. Other elements with similar structure include Zn, Cd, and Hg. For lighter alkaline-earth atoms, such as Ca and Mg, for which techniques other than Doppler cooling are impossible or too difficult to be used 17, it will facilitate trapping in optical potentials and lattices. Our treatment solves the fully quantum-mechanical equations of motion for the three-level atom in the field of both lasers. The dynamics of this system are determined exactly using numerical methods, giving quantitative predicted laser-cooling temperatures. The results of these calculations then provide us with a basis for a more intuitive interpretation of the cooling mechanism. Along these lines, we develop a dressed-state picture of the system dynamics that makes qualitative predictions that agree well with the numerical calculations.

Fig. 1 shows the internal atomic states in order of increasing energy as |0⟩, |1⟩, and |2⟩. The transition en-
energy of the lower transition, $|0\rangle \rightarrow |1\rangle$, and of the upper transition, $|1\rangle \rightarrow |2\rangle$, are $h\omega_0^{(1)}$ and $h\omega_0^{(2)}$, respectively. We include two lasers, of frequency $\omega_1$ and $\omega_2$, and define their detunings from the appropriate atomic transitions as $\delta_i = \omega_i - \omega_0^{(i)}$, for $i = 1, 2$. The intensities of lasers 1 and 2 are characterized by their Rabi frequencies $\Omega_1 = -\langle 0|d(1)|E_1(x)\rangle$ and $\Omega_2 = -\langle 1|d(2)|E_2(x)\rangle$, respectively, where $d$ is the electric-dipole operator of the atom and $E_i$ is the electric-field amplitude for the $i$th laser. The spontaneous-emission linewidths of states $|1\rangle$ and $|2\rangle$ are $\Gamma_1$ and $\Gamma_2$, respectively (88 MHz and 2.2 MHz, respectively, for $^{24}\text{Mg}$). The time evolution for the atom moving in the laser field, with mass $m$ and center-of-mass momentum operator $p$, is described by the master equation,

$$\dot{\rho}(t) = i\hbar^{-1}[H, \rho] + \mathcal{L}[\rho],$$

where $\rho$ is the reduced density matrix of the atom system, the vacuum photon field degrees of freedom having been traced over, $H = p^2/2m + h\omega_0^{(1)}|1\rangle\langle 1| + h\omega_0^{(2)}|2\rangle\langle 2| + V_{laser}^{(1)}(x) + V_{laser}^{(2)}(x)$ is the Hamiltonian of the atom-laser system, and the superoperator Liouvillian $\mathcal{L}$ describes effects due to coupling of the atom to the vacuum photon field, resulting in spontaneous emission.

Eq. (1) treats all the atomic degrees of freedom (internal and external) quantum mechanically, so its solutions generally provide an accurate description of the atom’s dynamics. We thus avoid much of the ambiguity and difficulty associated with semiclassical approximations of the system. For reasonable temperatures, however, the number of numerical basis states required to directly solve the problem, even in one dimension, is impractical for most computers. Here we resolve this problem without resorting to Monte Carlo methods by noting that the matrix for the linear system equivalent to Eq. (1) is very sparse — i.e., only a small fraction of its elements are nonzero. The particular structure shared by Liouvillian operators $\mathcal{L}$ describing relaxation processes simplify the numerics. Starting from the microscopic properties of atomic operators comprising this so-called Lindblad form of $\mathcal{L}$, we construct a matrix with the zero elements eliminated. The steady-state solution of Eq. (1) is then found using standard sparse-matrix inversion or propagation techniques. The result is an exact direct solution of a fully-quantized master equation.

The steady-state density matrix has been calculated this way in one dimension. The average kinetic energy is then $\langle p^2/2m \rangle = \text{Tr}(\rho p^2/2m)$, which we equate with $k_BT/2$, where $T$ is the temperature and $k_B$ is Boltzmann’s constant. The large parameter space of the three-level atom two-laser problem has been explored, e.g., varying the detunings, $\delta_1$ and $\delta_2$, as well as the strengths $\Omega_1$ and $\Omega_2$ of the two lasers. However, we find that the laser strengths are better characterized in terms of how strongly they dress the atom. The relevant non-resonant saturation parameters, $s_1$ and $s_2$, for the respective transitions, are

$$s_i = \frac{1}{2\delta_i^2 + (\Gamma_i/2)^2}.$$

The parameter space has three distinct regimes. In the first, with $s_1 \gg 1$ and $s_2$ arbitrary, only heating occurs, as expected from Doppler-cooling theory since the lower transition is driven strongly. In the second, with $s_1 \ll 1$ and $s_2 \ll 1$, cooling occurs only to the Doppler limit for the lower transition $T_D^{(1D)} = h\Gamma_1/2k_B$, and only in the range near $\delta_1 = -\Gamma_1/2$. In this case, laser 2 has no effect, as this amounts to simple two-level Doppler cooling on the lower transition with laser 1. In the final regime, when $s_1 \ll 1$ and $s_2 \gg 1$, cooling occurs down to substantially below $T_D^{(1D)}$. Figure (2) illustrates cooling in this regime for the partial case of $^{24}\text{Mg}$ mentioned above, with the temperature normalized to $T_D^{(1D)}$, and with the bare two-photon resonance ($\delta_1 + \delta_2 = 0$) denoted by the dashed line. The steady-state temperature is plotted as a function of $\delta_1$ and $\delta_2$ for $s_1(\delta_1) = 0.001$ and $s_2(\delta_2) = 1$. Note that the saturation parameters are being held fixed.
as the detunings are varied, so that the Rabi frequencies are being continuously adjusted. We see the lowest temperatures, on the order of $10^{-2} T_D(1D)$, in the quadrant with $\delta_1 > 0$ and $\delta_2 < 0$, as well as less extreme cooling in other regions. Observe that the lowest temperatures are obtained for frequencies detuned to the blue of the two-photon resonance. This seems counterintuitive, since a red detuning is usually required in order to have a net decrease of atomic kinetic energy in a photon-scattering event. Qualitative understanding of the cooling mechanism emerges from analysis of the simpler Hamiltonian,

$$H = \frac{\hbar}{2} \begin{pmatrix} 0 & \Omega_1 & 0 \\ \Omega_1 & 0 & i\Gamma_1 \\ 0 & -2\delta_1 - i\Gamma_1 & 0 \end{pmatrix} \begin{pmatrix} 0 \\ -2\delta_1 - i\Gamma_1 \\ -2(\delta_1 + \delta_2) - i\Gamma_2 \end{pmatrix}. \quad (3)$$

Its complex eigenvalues have real dressed energies, and imaginary parts giving dressed-state linewidths — that is, a measure of the coupling of the dressed states to the photon vacuum. These dressed energies and widths are plotted on the right side of Fig. 1 as functions of $\delta_2$, for the same parameters used in Fig. 2. The cancellation of one of the widths can be viewed as an EIT effect: since laser 1 is perturbative while laser 2 strongly dresses the upper transition, the new eigenstates of the system, denoted $|+\rangle$ and $|\rangle\rangle$, are well approximated as linear combinations of the bare states $|1\rangle$ and $|2\rangle$. These states have the modified energies and widths shown in Fig. 4. The linewidth modifications can be viewed as a Fano interference [23], in which the dressing laser transitions caused by the probe laser enable multiple coherent pathways among the bare states. Constructive or destructive interference respectively increases or decreases the atomic linewidth.

The cooling mechanism is thus qualitatively explained as ordinary two-level Doppler cooling. But instead of using a transition between two bare states of an atom, the transition occurs between a (mostly) unmodified ground state, and a dressed excited state, with a shifted energy and a new linewidth that can be narrower than the bare linewidth of the lower transition. As the probe laser is scanned, the detuning relative to the dressed energy levels is varied, but since these levels are shifted from their bare energies, resonance occurs for different detunings than are encountered in the bare system. In fact, the shifts of the eigen-energies in Fig. 4 from the bare energies explain the apparent observation of blue two-photon cooling in Fig. 2. In the dressed system, the bare two-photon resonance is no longer meaningful, and the cooling region is in fact to the red of a dressed resonance.

An additional caveat applies when mapping this system onto Doppler cooling theory: the lineshapes are not Lorentzian, but are in fact asymmetric Fano lineshapes for the dressed system, as shown in the upper part of Fig. 4 as a function of $\delta_1$ for a fixed value of $\delta_2 = -\Gamma_1/2$. This fact changes the optimum-detuning condition into a new one: maximum cooling for a given transition occurs when the probe laser is detuned from the dressed excited

FIG. 2: (Color online) Steady-state laser-cooling temperatures for a $^{24}\text{Mg}$ three-level cascade system, as a function of the two detunings $\delta_1$ and $\delta_2$, obtained from exact numerical solutions of Eq. (1). The detuning-dependent saturation parameter for the lower transition (perturbative probe laser) is $s_1(\delta_1) = 0.001$ for both plots, and for the upper transition (dressing pump laser) is $s_2(\delta_2) = 1$ and 5 in the upper and lower plot, respectively. The temperature is normalized to the one-dimensional Doppler limit for the lower transition $T_D(1D) = 7\hbar\Gamma_1/40k_B$ [22], which is the optimum temperature expected for cooling with just one laser. The dashed line indicates the location of the bare two-photon resonance. This example illustrates the main parameter regime where sub-Doppler cooling occurs. Note that, in order to hold the saturation parameters fixed as the detuning is varied, the Rabi frequencies $\Omega_i$, for $i = 1, 2$, are continuously adjusted.
where \( v \) is the atomic velocity. Since the detuning of the laser and the resonant atomic velocity are linearly related, the derivative of the absorption spectrum with respect to \( \delta_1 \) also yields a maximum in the cooling force. This is evident in normal Doppler cooling because the optimum detuning occurs when \( \delta = -\Gamma/2 \), which is the inflection point of the Lorentzian. In general then, for asymmetric lineshapes, the optimum detuning does not obey such a simple relation, but depends on the degree of asymmetry.

From this complete picture of the cooling mechanism as a weak probe applied to a dressed three-level system, the minimum expected temperatures can now be determined, allowing for the detuning modification due to asymmetric lineshapes. The lower part of Fig. 3 is a plot of the ratio of the maximum slope of a Lorentzian lineshape with width \( \Gamma_1 \) to the slope of the asymmetric lineshape, as a function of \( \delta_1 \) with \( \delta_2 = -\Gamma_1/2 \). For comparison, fully quantum numerical results are indicated by data points. This ratio provides an indication of the expected cooling for the dressed system relative to the Doppler limit for the lower transition. Note that the expected temperature, due to the asymmetric lineshape, is predicted to be lower than the upper-transition Doppler limit, indicated by the horizontal dotted line in the lower plot of Fig. 3. This prediction is supported by the numerical data.

In conclusion, coherent engineering of an atomic three-level system can optimize the effectiveness of two-level Doppler cooling. By strongly driving a particular transition between two excited internal states of the atom, dressed energy eigenstates are created with modified linewidths which, due to mixing, can vary anywhere in the range between the smallest and the largest of the two bare linewidths. Smaller linewidths lead to lower temperatures, but the additional effect of asymmetries in the linewidths of the dressed states can lead to even lower temperatures, below the Doppler limit of the upper transition. The ability to tailor the degree of cooling lends this technique additional utility. A dressing scheme can be suited to the characteristics of a particular atom, and real-time adjustment of the cooling properties can allow narrowing of the velocity-capture range as an atomic gas is cooled. Utilizing such coherent effects should lead to relatively simple schemes for cooling far below the typical Doppler limit.

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\[ \alpha = -\frac{d}{dv}f(v), \]  

where \( f(v) \) is the absorption rate.

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