

Observation of narrow Autler-Townes components in the resonant response of a dense atomic gas

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We have experimentally studied the reflection of a weak probe beam from a dense atomic potassium vapor in the presence of a strong laser field tuned to the atomic resonance transition. We have observed an Autler-Townes doublet under hitherto unexplored conditions, namely that the Rabi frequency induced by the strong laser field is much smaller than the self-broadened width of the resonance transition of the unexcited vapor. We attribute our observation to a reduction of the atomic decoherence by the strong drive field. We present a theoretical model of nonlinear processes in a dense atomic gas to explain the observed results.

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Spectral line broadening is a universal phenomenon and a multitude of techniques has been devised to reduce or eliminate it altogether, enabling the study of spectral features that would, otherwise, remain hidden [1, 2, 3, 4]. In particular this applies to atomic or molecular gases where the collision rate can be made sufficiently small by, for instance, rarefying the vapor. However, when the interest lies with the (strongly) interacting gas, line broadening becomes the essence, and the study of the width of the fundamental resonance transitions in high-density atomic vapors is well-documented [5, 6].

Although, in the general case, the broadening of the resonance line has many contributions [7], for dense homogeneous gases, the major contribution is caused by the resonant dipole-dipole interaction between ground- and excited-state atoms of the same species, and is known as self- or resonance broadening [5, 6]. The interaction is long range, so that already at reasonably modest densities of order 10^{17} cm⁻³ one leaves the binary-collision regime and multi-perturber effects may come into play [6, 8]. It has, for instance, been shown that the Zeeman effect is modified at atomic densities of this order [9]. We note that multi-perturber effects were observed in high pressure buffer gas [10, 11, 12] when the duration of collisions cannot be neglected.

In a high-density vapor the atom responds not just to the externally applied EM field but also to the field re-radiated by other atoms in the vapor, i.e. it responds to the *local* field. As first shown by Lorentz, this causes the position of the resonance to be shifted [13]. During the last decade this local-field shift, named after Lorentz, has been studied by both frequency [14, 15, 16, 17, 18, 19] and time-domain [20] techniques. It is proportional to the atomic density, and of the same order of magnitude as the line width [21].

It has been predicted that the Lorentz local-field shift depends on the degree of (incoherent) excitation of the vapor [21, 22], a prediction that stimulated the study of a variety of nonlinear optical phenomena in dense atomic vapors, such as piezophotonic switching and lasing without inversion [23]. The experimental demonstration of the excitation dependence of the Lorentz shift [17, 19] brought to light that the self-broadened width itself is also excitation dependent, an effect that could be explained in terms of a quasistatic (multi-perturber, quasi-molecular) collision model [6, 17]. Recently, it was shown by using a time-domain technique, that non-Markovian collision dynamics can be observed in these high density

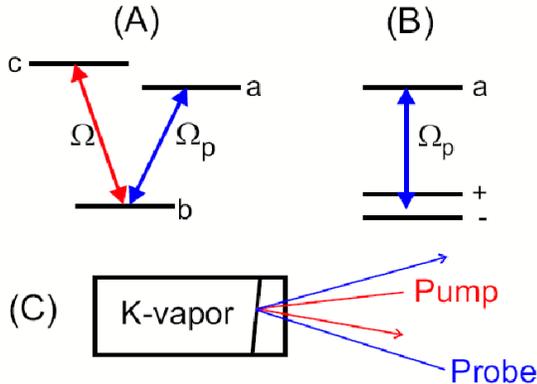


FIG. 1: (Color online) Energy-level scheme for potassium. The bare atomic states are shown in (A), where level a corresponds to the $4p \ ^2P_{1/2}$ excited state, level b to the $4s \ ^2S_{1/2}$ ground state, and level c to the $4p \ ^2P_{3/2}$ excited state. The strong drive field, with Rabi frequency Ω , couples states b and c , while the probe laser beam, with Rabi frequency Ω_p , is scanned over the $b - a$ transition. Frame (B) shows the probe transition in the dressed basis, the states being written as $|+\rangle$ and $|-\rangle$, and (C) provides a sketch of the experimental setup.

resonantly-broadened vapors, indicating that the duration of a collision can no longer be neglected [24].

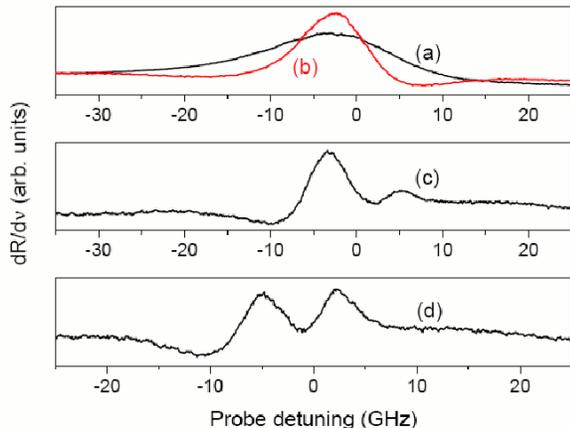


FIG. 2: (Color online) Frequency-modulated reflectivity spectra of a potassium vapor at a density of $4.9 \times 10^{17} \text{ cm}^{-3}$, as measured on the D_1 transition. Curves (a) and (b) display the FM reflection spectrum of the vapor when the pump laser is switched off, and when it operates at a power of $\approx 0.5 \text{ W}$ and is tuned far (130 GHz) above the D_2 transition. When the detuning of the pump laser is reduced to $\approx 3 \text{ GHz}$ the spectral line in the FM spectrum becomes an Autler-Townes doublet (curve (c)). Curve (d) shows the Autler-Townes components at zero pump detuning.

The experimental study of resonance broadening remains a topical issue in view of theoretical claims that, because of the long-range nature of the interaction, many-body effects play a role at all densities [6]. Here we show that at a higher number density of $4.9 \times 10^{17} \text{ cm}^{-3}$ the atomic ensemble can be made to evolve coherently as if it is a two-level system with intensity dependent line width. We deduce this behavior from our measurements of the Autler-Townes splitting [25] of the potassium atomic ground state. Observation of the Autler-Townes splitting of optical transition in dilute gases were previously reported in [26]. Essential to our observation is the effect of collision-induced line narrowing [16, 17, 19].

At the number density of our experiment, the self-broadened line width of both resonance lines of potassium is considerably larger than either their hyperfine structure or their Doppler width. Under these conditions the level structure of potassium is well-represented by a three-level scheme, shown in Fig. 1. A strong driving field (Rabi frequency Ω) couples levels b and c , while a weak probe field, with Rabi frequency Ω_p is scanned over the $a \leftrightarrow b$ transition.

The experiment is performed in reflection (see Fig. 1C), using an experimental arrangement similar to that of Ref. [19]. The strong coupling field is produced by a CW Ti-

sapphire laser tuned close to the D_2 -line ($\lambda = 766 \text{ nm}$) of atomic potassium. The laser beam, having a maximum power of $\approx 0.5 \text{ W}$, has its focus (the area of the beam is $5 \times 10^{-4} \text{ cm}^2$) at the window of a vapor cell that is constructed entirely out of sapphire. A weak probe beam from an external cavity diode laser is focussed to overlap with the pump beam at the entrance window of the cell. The probe laser is scanned over the D_1 -line at $\lambda = 770 \text{ nm}$, at a power level of $\approx 10^{-4} \text{ W}$, far below the onset of saturation (saturation parameter [3] is < 0.1). Both the pump and probe beams are linearly polarized and are incident on the sapphire cell at near-normal incidence. While being scanned, the probe laser is frequency modulated over a range of 100 MHz at a frequency of 400 Hz. The reflected probe beam is captured by a photodiode whose signal output is fed to a lock-in amplifier. Our experiment runs at an atomic density of $4.9 \times 10^{17} \text{ cm}^{-3}$, yielding a self-broadened linewidth $\Gamma_{\text{self}}/2\pi$ of 28.4 GHz for the D_1 line [9].

Figure 2 shows the measured FM-modulated spectra. The zero on the frequency axis refers to the center frequency of the absorption spectrum of a low density ($N \simeq 10^{11} \text{ cm}^{-3}$) potassium vapor in a reference glass cell. Curve (a) displays the reflection spectrum when the pump beam is switched off, i.e. it shows the FM linear reflectivity of the sapphire/high-density vapor interface. When the pump laser is switched on, operates at an output level of 0.5 W, and is tuned 130 GHz above the D_2 line, the FM spectrum of curve (b) is obtained. At this large detuning the pump-laser beam penetrates a considerable distance into the vapor cell. Due to radiation trapping the decay of the excited-state population is slowed down and, as a result of radiative transport, the vapor near the sapphire window becomes homogeneously and incoherently excited. In line with earlier observations ([19]) one notices that the FM spectrum of the strongly excited vapor (curve (b)) is narrower than its unexcited counterpart (curve (a)), and that their centers do not coincide. However, both spectra are described by the same spectral function, with adjusted values for the Lorentz local-field shift and self-broadened width [16, 17, 19].

The spectral width Γ_{self} of the resonance line can be estimated from the interval $\Delta\omega_{\text{mm}}$ between the maximum and minimum of the reflection spectrum or, equivalently, from the zeros of the FM spectrum. For the D_1 transition one has $\Gamma_{\text{self}} = 0.87\Delta\omega_{\text{mm}}$ [9, 19]. By this method, we determine the line width for the unexcited vapor as 31 GHz (Fig. 2a), in reasonable agreement with the value $\Gamma_{\text{self}}/2\pi = 28.4 \text{ GHz}$, calculated using the known self-broadening coefficient and our estimate for the atomic density. Similarly, we estimate the line width for the incoherently excited vapor as just 11 GHz (Fig. 2b). Hereby we show that the excitation dependence of the self-broadened line width is a valid concept also at the vapor density of the present experiment.

In order to observe coherent effects we tune the pump

laser close to the center of the atomic line. Curves (c) and (d) show the experimental FM reflectivity spectra for a pump detuning of 3 and 0 GHz, respectively. In both cases one observes that the single resonance of curve (b) is split into an Autler-Townes doublet [25, 26]. The asymmetry of the doublet in curve (c) reflects the fact that, for this curve, the pump laser is tuned slightly above the D_2 resonance transition, while the near-perfect symmetry of curve (d) shows that the pump laser is, in that case, tuned on resonance. Note that, as a result of the various line shifts, there is no easy method, beyond looking at the symmetry of the Autler-Townes doublet, to establish whether the pump laser is tuned exactly to resonance or not.

To obtain a good estimate for the width and separation of the the components of the Autler-Townes doublet we have made a non-linear fit of the experimental data of curve (d) in Fig. 2, yielding a width of 8.5 GHz and a splitting of 7.6 GHz.

It is well known that the separation between the two components of the Autler-Townes doublet equals the Rabi frequency induced by the pump laser [3, 27]. For the conditions of the present experiment where hyperfine and Zeeman substructure can be ignored, the Rabi frequency Ω for the potassium D_2 line can be expressed as

$$\Omega/2\pi = 8 \times 10^7 \sqrt{I \left[\frac{\text{W}}{\text{cm}^2} \right]} \text{ Hz}, \quad (1)$$

where $I = cE^2/(8\pi)$ is the pump laser intensity (in units of W/cm^2). Here the Rabi frequency is written as

$$\Omega = E\varphi_x/\hbar, \quad (2)$$

with E the optical electric field and φ_x the projection of the atomic dipole moment $\vec{\varphi}$ along that field. The value of $\varphi_x = \varphi/3$ has been obtained from experimental data for the spontaneous lifetime [28]. Using Eq. (1) we calculate the Autler-Townes splitting to be 8 GHz at a pump power of 0.5 W, in satisfying agreement with the experimental value (7.6 GHz).

We have recorded the Autler-Townes doublet for several values of the detuning Δ and intensity of the pump laser. When $|\Delta|$ increases one of the components of the doublet becomes stronger while the other fades away. When $|\Delta/2\pi| > 6$ GHz the weak component can no longer be measured, and only a single resonance is observed. For this case, where the pump laser is detuned from exact resonance, the Autler-Townes splitting is given by the generalized Rabi frequency [27]

$$\tilde{\Omega} = \sqrt{\Omega^2 + \Delta^2}. \quad (3)$$

The measured values of the Autler-Townes splitting as a function of the generalized Rabi frequency $\tilde{\Omega}$ are presented in Fig. 3. The three points for $\tilde{\Omega}/2\pi \leq 8$ GHz have been measured at zero detuning and pump laser powers

equal to 0.3, 0.4 and 0.5 W, respectively. The points at $\tilde{\Omega}/2\pi \geq 8.5$ GHz were measured by using a pump power equal to 0.5 W and detunings of $\Delta/2\pi = \pm 3$ GHz and ± 6 GHz, respectively. Also shown is a linear fit through the data; its slope equals 0.92, in quite good agreement with the expected value, equal to 1 (the error of slope is 2% and the difference between theoretical and experimental curves is about 6% due to systematic uncertainties).

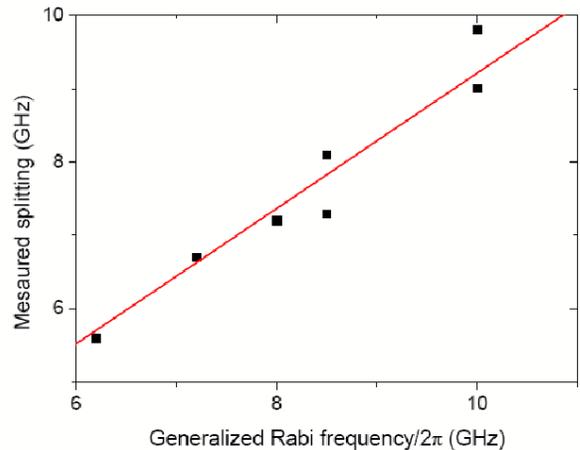


FIG. 3: The measured Autler-Townes splitting versus generalized Rabi frequency. The solid line shows the result of a linear fit.

To explain our results we use the optical Bloch equations for a V-type three-level atom, suitably modified for the fact that the system under study is a dense rather than a dilute vapor. From these equations we evaluate the susceptibility of the probe at frequency ω_p as [29]

$$\chi(\omega_p) = i \frac{3\lambda^2\gamma}{8\pi} \frac{(n_a - n_b) + \frac{\Omega^2}{\Gamma_{cb}\Gamma_{ca}}(n_c - n_b)}{\Gamma_{ab} + \frac{\Omega^2}{\Gamma_{ca}}}. \quad (4)$$

Here the level populations n_a , n_b and n_c depend on the Rabi frequency of the driving field Ω , γ is the radiative rate on the $a \rightarrow b$ transition, and the generalized coherence decay rates are $\Gamma_{cb} = \tilde{\gamma}_{cb} + i\delta_{cb}$, $\Gamma_{ca} = \tilde{\gamma}_{ca} + i\delta_{ca}$, $\Gamma_{ab} = \tilde{\gamma}_{ab} + i\delta$. The frequency of the pump is denoted as $\bar{\omega}$, and its detuning is given by $\delta_{cb} = \omega_{cb} - \bar{\omega} + \Delta\omega_L + \Delta\omega_c$, with $\Delta\omega_L$ the Lorentz local field shift, and $\Delta\omega_c$ the collisional frequency shift, which is independent of the level of excitation [14, 15, 19]. The detuning of the probe from resonance is $\delta = \omega_{ab} - \omega_p + \Delta\omega_L + \Delta\omega_c$, with ω_p the frequency of the probe laser. For the two photon detuning we write $\delta_{ca} = \omega_{ca} - \bar{\omega} + \omega_p + \Delta\omega_L + \Delta\omega_c$. For $\tilde{\gamma}_{cb}$, $\tilde{\gamma}_{ca}$ and $\tilde{\gamma}_{ab}$ we write $\tilde{\gamma}_{ab} = \gamma/2 + \Gamma_{\text{self}} \simeq \Gamma_{\text{self}}$, $\tilde{\gamma}_{cb} = \gamma/2 + \Gamma_{\text{self}} \simeq \Gamma_{\text{self}}$, and $\tilde{\gamma}_{ca} = \gamma + \Gamma_{\text{self}} \simeq \Gamma_{\text{self}}$, since the self-broadened line width is much larger than the natural width of the transitions.

For the self-broadened width we write [17, 19]

$$\Gamma_{\text{self}} = kn_b, \quad (5)$$

with k the self-broadening coefficient [5]. The Lorentz local-field shift is given by [21, 22]

$$\Delta\omega_L = \frac{k}{3}(n_b - n_a). \quad (6)$$

The reflectivity of the coherently driven dense vapor can now be evaluated using

$$R(\omega) = \left| \frac{n(\omega) - n_0}{n(\omega) + n_0} \right|^2, \quad (7)$$

with $n(\omega) = \sqrt{1 + 4\pi\chi(\omega)}$ and n_0 the refractive indices of the vapor and of the window material of the vapor cell, respectively. Using Eq. (7) and appropriate values for the experimental parameters we calculate the FM reflection spectra, as shown in Fig. 4, for a pump detuning of zero and 3 GHz, respectively, and a pump Rabi frequency of 12 GHz. Fully in line with the experimental result our theoretical model predicts a well-resolved Autler-Townes doublet under circumstances that the Rabi frequency is much smaller than the unsaturated self-broadened line width.

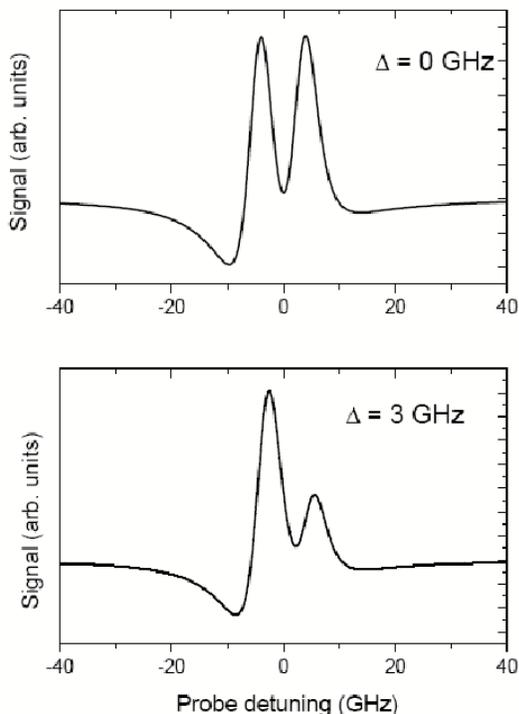


FIG. 4: Results of the theoretical model for the FM-reflection spectrum of a vapor of coherently driven V-type three-level atoms at high density, using Eqns. (4) and (7).

It is worth to mention here that, even though the experiments and theoretical analysis have been done for conditions corresponding to hot dense gases, it will be interesting to study discussed multy-perturber effects in dense cold atomic and molecular gases, and possible influence on coherent backscattering in cold atoms [30].

In conclusion, we have observed the Autler-Townes splitting of the atomic ground state in the reflection spectrum of a dense potassium vapor in a regime where the separation of the two Autler-Townes components, as given by the Rabi frequency, is much smaller than the nominal self-broadened width of the transition. We attribute our observation to the fact that the strong drive field, in addition to inducing coherent evolution of the atomic levels, also causes a dramatic narrowing of the spectral line. Taking into account the excitation dependence of both the local-field shift and the resonance line width we obtain good agreement between our experimental results and an effective three-level model based on modified Bloch equations.

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