The emission characteristics in the fluorescence of two laser-driven dipole-dipole-interacting three-level atoms is investigated. When the light from both atoms is detected separately, a correlation of the emission processes is observed in dependence of the dipole-dipole interaction. This opens the possibility to investigate the dipole-dipole interaction through the emission behavior. We present Monte-Carlo simulations which are in good agreement with the analytic solutions.

PACS numbers: 42.50.Fx, 42.50.Lc, 42.50.Ar

The problem of the influence of a neighboring atom on the atomic emission behavior [1–4] has been investigated since a very long time. Especially the problem of the correlation of the emission of two neighboring atoms has found quite some interest [5–7]. This is a complex issue and the answer is very much dependent on various system parameters such as the strength of the pumping field, the life times and the wavelengths of the different transitions involved. An early theoretical calculation concluded that the dipole-dipole interaction between adjacent atoms is irrelevant for quantum jumps within a 3-level-system under the usual experimental conditions where the Rabi frequency of the pump field is large compared to all other rates in the problem [8,9]. Those results have been confirmed by quantum Monte-Carlo calculations [10–12], [13]. In this paper we reconsider the problem of dipole-dipole interaction and investigate especially the correlations in the emissions of two neighboring atoms which are observed individually.

The arrangement is shown schematically in Fig. 1. We consider 2 identical nearby atoms in a trap, each with levels |1⟩, |2⟩ and |3⟩. One of the atoms (say atom 1) is initially prepared, e.g. by pulsed excitation [14], in a metastable state |2⟩. A resonant cw pump interacts with both atoms. The initial state of atom 1 is such that it can start interacting with the pump either due to decay to the state |3⟩ by emission of a photon or via excitation to the state |1⟩ as a result of the dipole-dipole interaction. In the second case atom 2 goes to the state |2⟩. In our analysis we assume that the level spacing |1⟩ ↔ |2⟩ is close enough, so that the dipole-dipole interaction is effective only on this transition. The transitions |1⟩ ↔ |3⟩ and |2⟩ ↔ |3⟩ are supposed to lie in the optical domain, where the distance between two atoms is assumed to be much larger than the wavelengths of the corresponding transitions. For example in In⁺ the wavelength on the |1⟩ ↔ |2⟩ corresponds to 9.3µm which could be two to three times the distance between two trapped ions [15]. We further assume that the fluorescence from each atom can be resolved individually by two distinct detectors 1 and 2. In what follows let us work in the limit

$$\Omega_R > \gamma_{13}^{(1)} \gg \gamma_{12}, \gamma_{23}$$

In our prepared system, detector 1 will not detect any fluorescence while atom 2 is fluorescing on the |1⟩ ↔ |3⟩ transition until the dipole-dipole interaction on the transition |1⟩ ↔ |2⟩ brings atom 1 towards the cycling transition |1⟩ ↔ |3⟩ and atom 2 towards the metastable state |2⟩. In this case detector 1 is switched on and at the same time detector 2 is switched off. Detector 2 will remain off until either the atom 2 makes a transition from the state |2⟩ to |3⟩ by spontaneous emission or again to the state |1⟩ by the dipole-dipole interaction. If atom 2 goes to the state |3⟩, both the detectors will be on. If on the other hand atom 2 goes to the state |1⟩ by dipole-dipole interaction, then again detector 1 will switch off. Under the above inequality the dipole-dipole interaction is more probable and hence the chance for both detectors 1 and 2 to switch on at the same time is rare. We thus conclude that the fluorescence records of the detectors 1 and 2 will be complementary. This complementary record of...
fluorescence is a clear signature of the dipole-dipole interaction between the two atoms. The key ingredients in the above argument are (a) the inequality (1), (b) preparation of one of the atoms in the metastable state [2], (c) capability to resolve the fluorescence from individual atoms which are separated by a distance much bigger than the wavelength of the strong transition. In what follows we will present results from the quantum Monte-Carlo simulation of the fluorescence from the two 3-level-atoms. These simulations validate the above ideas.

The two identical atoms are assumed to be at fixed positions $r_i$, and the field outside the laser-beam in the vacuum state. The dipole matrix elements of the atoms are defined as $d_{ij} := c(i|\sigma|j)$ and the atomic operators for the $i$-th atom as $\sigma^{nm}_i := |n\rangle_i \langle m|$. The interaction hamiltonian for two dipole-dipole interacting three-level-systems reads [16]

$$H = \hbar \sum_{n,m} \sum_{k} (g_{ks} \sigma^{nm}_k + \text{h.c.})$$

with

$$\sigma^{nm}_k := \sum_i \sigma^{nm}_i e^{ik \cdot r_i},$$

and

$$g_{ks} := -i \sqrt{\frac{\omega}{2\hbar c V}} \left( \mathbf{d} \cdot \mathbf{e}_{ks} \right).$$

For the time steps $\Delta t$ used in the simulations we need the time-development of the system under the condition that no photon is emitted in the time between $t = 0$ and $t = \Delta t$. This is described by a non-hermitian ‘conditional hamiltonian’, including the atom laser interaction, which is found to be [10–12,17]

$$H_{\text{cond}} = \frac{\hbar}{2} \left[ \sum_{n,m=1}^{3} \sum_{i,j=1}^{2} \gamma^{nm}_{ij} \sigma^{nm}_i \sigma^{nm}_j ight. + \left. \sum_{i=1}^{2} i \Omega_R \left( \sigma^{13}_i + \sigma^{31}_i \right) \right].$$

Here $2\gamma^{nm}_{ij}$ is the Einstein $A$-coefficient for the transition $|n\rangle \leftrightarrow |m\rangle$ for $i = j$ and $2\Omega_R$ is the Rabi frequency. For $i \neq j$, $\gamma^{nm}_{ij}$ is the complex parameter which describes the strength of the dipole-dipole interaction, where we have made the rotating-wave-approximation in the derivation of (5). In the Markov approximation they can be calculated analytically [16]:

$$\gamma^{nm}_{12} = \frac{3}{2} \gamma^{nm}_{11} e^{ik \cdot r} \left[ \frac{1}{ik^{nm}} (1 - \cos^2 \theta_{nm}) + \left( \frac{1}{ik^{nm}} \right)^2 + \left( \frac{1}{ik^{nm}} \right)^3 \right] (1 - 3 \cos^2 \theta_{nm})$$

Here $r = |r_1 - r_2|$ is the distance between the two atoms and $\theta_{nm}$ the angle between $\mathbf{d}_{nm}$ and $r_1 - r_2$. For the further calculations we define $\gamma^{nm} = \gamma^{nm}_{11} = \gamma^{nm}_{22} + \gamma^{nm}_{dd}$.

Their spatial dependence is shown in figure 2 and 3. Obviously $\gamma^{nm}_{dd}$ and $\Omega^{nm}_{dd}$ tend to zero for $r \to \infty$, but for $r \to 0 \Omega^{nm}_{dd}$ diverges. This means that $\Omega^{nm}(r)$ can have important consequences for the evolution of the system for $r < \lambda_{nm}$. By use of the Monte-Carlo simulations we indeed show, that this is the case for the quantum-jump-behavior of the system.

The evolution of the state vector between two consecutive emissions is given by [10–12,17]

$$|\Psi(t + \Delta t)\rangle = \left( 1 - \frac{i}{\hbar} H_{\text{cond}} \Delta t \right) |\Psi(t)\rangle$$

where we assume that $\Delta t$ is much larger than the inverse optical frequencies but much smaller than all the atomic decay times.

To show that the dipole-dipole interaction can have important consequences for the quantum-jump-behavior of the system we detect the fluorescence light of the two atoms separately with the detectors 1 and 2, which are modeled by the operators...
the detection rates, which are clearly distinguishable. As
and gets dark. Finally we have three possible changes of
cycle on the atom \( |2\rangle \) gets excited to state \( |1\rangle \) and shows fluorescence, i.e. detector 1 clicks, while atom 2 is dark. There are two possible decays which change the fluorescence behavior of the system (a) atom 1 falls to \( |2\rangle \) and both detectors do not detect any further photons (b) atom 2 falls to \( |3\rangle \) and both atoms show fluorescence at the same time. There is also a probability for a third event which is the most interesting in this context. As mentioned above with every time step where we have a non zero probability amplitude of the system being in state \( |1, 2\rangle \) we get a non zero probability amplitude for being in state \( |2, 1\rangle \), due to the dipole-dipole interaction. If this transition occurs, atom 2 gets excited to state \( |1\rangle \) from where it starts to cycle on the \( |1\rangle \leftrightarrow |3\rangle \) transition and to emit fluorescence photons, while atom 1 ends up in the metastable state \( |2\rangle \) and gets dark. Finally we have three possible changes of the detection rates, which are clearly distinguishable. As

\[ D_i = |3\rangle_i \langle 1| \]  \hspace{1cm} (8)

where the probability \( P_i \) for a measurement with the detector \( i \) is proportional to the excitation probability of the \( i \)-th atom. We find:

\[ P_i = 2\gamma^{13}\langle D_1 \rangle \]  \hspace{1cm} (9)

After each measurement we have to continue with the reduced state \( |\Psi\rangle_{\text{red}} = \frac{1}{N} D_i |\Psi\rangle \), with \( N \) being a normalization constant.

Next we focus on the dipole-dipole interaction on the transition \( |1\rangle \leftrightarrow |2\rangle \), where the \( r \)-dependent factor \( \frac{1}{r^3} \) is largest. In the time evolution of the system which is given by equation (7) the interaction parameter \( \gamma^{12}_{12} \) appears with the operators

\[ \sigma_1^{12}\sigma_2^{21} + \sigma_2^{12}\sigma_1^{21}. \]  \hspace{1cm} (10)

If the probability amplitude for the system being in state \( |1, 2\rangle = |1\rangle_1 \otimes |2\rangle_2 \) is not zero for the initial state the probability amplitude of being in state \( |2, 1\rangle \) is not zero one time step further. Consequently the dipole-dipole interaction drives the system from a non-entangled state to an entangled state [18]. This is a non-classical effect, which we can detect with our setup. The relevant time scale of this process is \( \frac{1}{\gamma^{12}_{12}} \) as is clear from equation (5).

To see the effect, on atom has to be in state \( |2\rangle \), which can be prepared by an appropriate laser pulse, while the other one must have a non zero probability of being in state \( |1\rangle \). What is achieved by the driving laser. The initial state for the simulation is therefore \( |1, 2\rangle \). It is evident, that every ‘click’ of one of our detectors destroys the entanglement and one starts again with a separable state. There are no problems with decoherence effects, as no coherence needs to be reserved longer than the lifetime of the upper state \( |1\rangle \).

If we compare the time scales and probabilities in the regime (1) it is clear that atom 1 shows fluorescence, i.e. detector 1 clicks, while atom 2 is dark. The emissions of atom 1 are plotted with positive, those of atom 2 with negative sign, so the plot shows an event where atom 1 gets dark and atom 2 bright at the same time.

The flip in the detection of fluorescence (Fig. 4) is a clear signature of the dipole-dipole interaction of the system and shows how the quantum jump statistic can be modified by cooperativity. Note that the effect of dipole-dipole interaction is not observable if we do not distinguish between the fluorescence of the two atoms [13]. To see how this behavior is connected to the interaction parameter \( \gamma^{12}_{12} \) we count the flips per unit time for different atom distances of interest (from \( r = 0.1\gamma^{12}_{12} \) to \( r = 3\gamma^{12}_{12} \)) and find, up to a scaling factor, very good functional agreement with \( |\gamma^{12}_{12}|^2 \) (Fig. 5, 6).
FIG. 5. Monte-Carlo data of Flips per time compared to $|\gamma_{12}|^2$ for $\frac{\gamma_{12}}{\gamma_{13}} = 2 \cdot 10^{-2}$ with $c_s = 2$ and $\Omega_R = 8$.

FIG. 6. Monte-Carlo data of Flips per time compared to $|\gamma_{12}|^2$ for $\frac{\gamma_{12}}{\gamma_{13}} = 5 \cdot 10^{-3}$ with $c_s = 2$ and $\Omega_R = 8$.

If we further analyze our data for different ratios of the decay rates $\gamma_{12}$ and $\gamma_{13}$ we find that the remaining scaling factor $c_s$ is nearly constant over reasonable magnitudes.

If we further analyze our data for different ratios of the decay rates $\gamma_{12}$ and $\gamma_{13}$ we find that the remaining scaling factor $c_s$ is nearly constant over reasonable magnitudes.

The result is shown in Fig. 7. This shows that indeed the flip rate is a consequence of the dipole-dipole interaction and that other parameters than $\gamma_{12}$ do not intervene.

In conclusion, we have outlined one solution to the problem of detecting the dipole-dipole interaction in the phenomena of quantum jumps. For that purpose we analyzed a system of two dipole-dipole interacting 3-level systems and proposed, in contrast to earlier calculations [8,9,17], a detection scheme, where we distinguish between the two atoms in order to monitor the consequences of this interaction. Within this scheme there are no problems related to decoherence, as no coherences need to be preserved over timescales longer than the shortest decay time of the system. Furthermore the dependence of the flip rate on the spatial variation of the dipole-dipole interaction has been analyzed.