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Fast generations of tree-type three-OPENdimensional entanglement via Lewis-Riesenfeld invariants and transitionless quantum driving

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Recently, a novel three-dimensional entangled state called tree-type entanglement, which is likely to have applications for improving quantum communication security, was prepared via adiabatic passage by Song *et al.* **Here we propose two schemes for fast generating tree-type three-dimensional entanglement among three spatially separated atoms via shortcuts to adiabatic passage. With the help of quantum Zeno dynamics, two kinds of different but equivalent methods, Lewis-Riesenfeld invariants and transitionless quantum driving, are applied to construct shortcuts to adiabatic passage. The comparisons between the two methods are discussed. The strict numerical simulations show that the tree-type three-dimensional entangled states can be fast prepared with quite high fidelities and the two schemes are both robust against the variations in the parameters, atomic spontaneous emissions and the cavity-fiber photon leakages.**

Entanglement plays a crucial role in quantum information processing¹⁻⁴. Some of typical entangled states, such as Bell state⁵, Greenberger-Horne-Zeilinger (GHZ) state^{[6](#page-10-2)} and W state^{[7](#page-10-3)}, have attracted great attention in the last decades. However, all of these states are entangled states that are defined in Hilbert spaces with two dimensions. Recently, high-dimensional entanglement has attracted more and more attention due to their superior security than qubit systems, especially in the aspect of quantum key distribution. Besides, it has been demonstrated that violations of local realism by two entangled high-dimensional systems are stronger than that by two-dimensional systems⁸. Thus, much interest has been focused on the generation of high-dimensional entanglement in theory via various techniques including quantum Zeno dynamics (QZD[\)9–11,](#page-10-5) stimulated Raman adiabatic passage $(STIRAP)^{12,13}$ $(STIRAP)^{12,13}$ $(STIRAP)^{12,13}$ $(STIRAP)^{12,13}$ $(STIRAP)^{12,13}$, and dissipative dynamics^{[14](#page-10-8),[15](#page-10-9)}. Also, experimental generations of high-dimensional entanglement have been already achieved^{[16](#page-11-0),[17](#page-11-1)}

With no doubt, a lot of remarkable achievements have been made with regard to high-dimensional entangled states. However, most of these high-dimensional entangled states are two-body but few multi-body. For a dozen years, some attention has been paid to multi-body high-dimensional entangled states such as singlet state¹⁸, and lots of schemes have been proposed for generations of singlet state¹⁹⁻²². A short time before, a novel three-body three-dimensional entangled state called tree-type entanglement was prepared via adiabatic passage by Song *et al.*[23.](#page-11-4) In the ref. [23,](#page-11-4) the tree-type three-dimensional entanglement was prepared among one single atom and two BECs and the authors predicted that the tree-type three-dimensional entanglement is likely to have great applications in improving quantum communication security.

Among the techniques mentioned above for generations of high-dimensional entanglement, there are two techniques widely used for their robustness against decoherence in certain conditions. One is STIRAP^{[12](#page-10-6),[13](#page-10-7),[21,](#page-11-5)23}, and the other is QZ[D9–11](#page-10-5)[,19](#page-11-3)[,20](#page-11-6)[,22](#page-11-7). STIRAP is widely used in time-dependent interacting fields and robust against the atomic spontaneous emission and variations in the experimental parameters, but a relatively long interaction time is usually required. QZD is usually robust against photon leakages and does not need a long interaction time. However QZD is sensitive to the atomic spontaneous emissions and variations in the experimental parameters. Thus some of researchers introduce detuning between the atomic transitions to restrain the influence of atomic spontaneous emissions^{[10](#page-10-10)}, but the interaction time significantly increases unavoidably. Therefore, in order to solve the problem of long interaction time, researchers have paid more attention to "shortcut to adiabatic passage"

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which employs a set of techniques to speed up a slow quantum adiabatic process^{24–35}. Further more, the multiple Schrödinger pictures and dynamics are introduced to implement physically feasible transitionless quantum driving (TQD) for two-^{[36,](#page-11-9)37} and three-level^{[38](#page-11-11)} quantum systems, and it offers an effective way for experimental realizations of robust quantum information processing. Recently, the transitionless superadiabatic protocols were experimentally implemented, which make the system follow the instantaneous adiabatic ground state nearly perfectly[39.](#page-11-12) The nonadiabatic holonomic quantum computation, which possesses the advantage of robustness against decoherence and enjoys a short operation time at the same time, was realized in experiment⁴⁰. Assisted quantum adiabatic passage was experimentally implemented in a single spin⁴¹. Up to the present, Lewis-Riesenfeld invariants (LRI) and TQD for shortcuts to adiabatic passage, were used widely to speed up the operations of multi-particle systems, such as fast multi-particle quantum state transfer^{[42](#page-11-15),[43](#page-11-16)}, fast generations of entanglement⁴⁴⁻⁴⁶ and fast constructions of quantum gates $47-51$.

In this paper, we propose two schemes for fast generating of tree-type three-dimensional entanglement among three spatially separated atoms via LRI and TQD, respectively. Based on LRI and TQD we construct effective shortcuts to adiabatic passage for fast generating tree-type three-dimensional entanglement among three atoms trapped respectively in three spatially separated cavities connected by two fibers. We will give the interesting comparisons between the LRI method and the TQD method. The generations of tree-type three-dimensional entanglement in our schemes are implemented within a short time, and the strict numerical simulations demonstrate that our schemes are both robust against the decoherence caused by the atomic spontaneous emissions, photon leakages and the variations in the parameters.

Preliminary Theory

Lewis-Riesenfeld invariants. Here we give a brief description about Lewis-Riesenfeld invariants theory⁵². A quantum system is governed by a time-dependent Hamiltonian *H*(*t*), and the corresponding time-dependent Hermitian invariant *I*(*t*) satisfies

$$
i\hbar \frac{\partial I(t)}{\partial t} = [H(t), I(t)]. \tag{1}
$$

The solution of the time-dependent Schrödinger equation $i\hbar\partial_t |\Psi(t)\rangle = H(t) |\Psi(t)\rangle$ can be expressed by a superposition of invariant *I*(*t*) dynamical modes $|\Phi_n(t)\rangle$

$$
|\Psi(t)\rangle = \sum_{n} C_n e^{i\alpha_n} |\Phi_n(t)\rangle,
$$
\n(2)

where *Cn* is the time-independent amplitude, *αn* is the time-dependent Lewis-Riesenfeld phase, and |Φ*n*(*t*)〉 is one of the orthogonal eigenvectors of the invariant $I(t)$ satisfying $I(t)|\Phi_n(t)\rangle = \lambda_n |\Phi_n(t)\rangle$ with a real eigenvalue λ_n . The Lewis-Riesenfeld phases are defined as

$$
\alpha_n(t) = \frac{1}{\hbar} \int_0^t dt' \left\langle \Phi_n(t') \left| i\hbar \frac{\partial}{\partial t'} - H(t') \right| \Phi_n(t') \right\rangle.
$$
 (3)

Transitionless quantum driving. Suppose a system is dominated by a time-dependent Hamiltonian $H_0(t)$ with instantaneous eigenstates $|\phi_n(t)\rangle$ and eigenvalues $E_n(t)$,

$$
H_0(t) |\phi_n(t)\rangle = E_n(t) |\phi_n(t)\rangle.
$$
 (4)

When a slow change satisfying the adiabatic condition happens, the state of the system governed by $H_0(t)$ can be expressed as^{[26](#page-11-20),[53](#page-11-21)}

$$
\begin{array}{rcl}\n\langle \psi(t) \rangle & = & e^{i\xi_n(t)} \langle \phi_n(t) \rangle, \\
\xi_n(t) & = & -\frac{1}{\hbar} \int_0^t dt' E_n(t') + i \int_0^t dt' \langle \phi_n(t') | \partial_{t'} \phi_n(t') \rangle.\n\end{array} \tag{5}
$$

Because the instantaneous eigenstate $|\phi_n(t)\rangle$ do not meet the Schrödinger equation, there may be transitions between the eigenstates of $H_0(t)$ with a finite probability during the whole evolution process even under the adiabatic condition. In order to construct the Hamiltonian *H*(*t*) that exactly drives the instantaneous eigenstates $|\phi_n(t)\rangle$ s, i.e., there are no transitions between different eigenstates during the whole evolution process, the simplest choice of the Hamiltonian *H*(*t*) can be written as

$$
H(t) = i\hbar \sum_{n} |\partial_{t} \phi_{n}\rangle \langle \phi_{n}|. \tag{6}
$$

Therefore, as long as *H*(*t*) is constructed, the system will evolves with no transitions between different eigenstates.

Quantum Zeno dynamics. Assume that a quantum system's dynamics evolution is governed by the Hamiltonian

$$
H_K = H_{\text{obs}} + KH_{\text{meas}},\tag{7}
$$

Figure 1. (**a**) The schematic setup for generating tree-type three-dimensional entanglement; (**b**) the level configurations and relevant transitions. (**c**) The APF design of the schematic setup for TQD to fast generate treetype three-dimensional entanglement.

where H_{obs} can be viewed as the Hamiltonian of the quantum system investigated and H_{meas} as an additional interaction Hamiltonian performing the measurement. *K* is a coupling constant, and in the strong coupling limit $K \rightarrow \infty$, the whole system is governed by the Hamiltonian⁵⁴

$$
H(t) = \lim_{K \to \infty} \sum_{n} (K \lambda_n P_n + P_n H_{\text{obs}} P_n),
$$
\n(8)

where P_n is one of the eigenprojections of H_{meas} with eigenvalues $\lambda_n (H_{\text{meas}} = \sum_n \lambda_n P_n)$. Interestingly, it is easy to deduce that the system state will remain in the same Zeno subspace as that of its initial state. In particular, if the system is initially in the dark state $|\Psi_d\rangle$ of $H_{\rm meas}$, i.e., $H_{\rm meas}|\Psi_d\rangle$ = 0, the the Hamiltonian (8) reduces to

$$
H(t) = \sum_{n} P_n H_{\text{obs}} P_n. \tag{9}
$$

Description of the physical model

The schematic setup for generating the tree-type three-dimensional entanglement is shown in [Fig. 1\(a\).](#page-2-0) Three atoms are trapped respectively in three spatially separated optical cavities which are connected by two fibers. Under the short fiber limit $(l\nu)/(2\pi c) \leq 1$, only the resonant modes of the fibers interact with the cavity modes⁵⁵, where *l* is the length of the fiber and *v* is the decay rate of the cavity field into a continuum of fiber modes. The atomic level configurations and relevant transitions are shown in [Fig. 1\(b\).](#page-2-0) As shown in [Fig. 1\(b\)](#page-2-0), the five-level atom1 and atom3 are both *M*-type with two excited states $|e_L\rangle$ and $|e_R\rangle$ and three ground states $|g_L\rangle$, $|g_0\rangle$ and $|g_R\rangle$. The four-level atom2 is tripod-type with one excited state $|e_0\rangle$ and three ground states $|g_L\rangle$, $|g_0\rangle$ and $|g_R\rangle$. The atomic transition $|e_{L(R)}\rangle_i \leftrightarrow |g_0\rangle_i$ (*j* = 1, 3) is resonantly coupled to the left-circularly (right-circularly) polarized mode of *j*th cavity with corresponding coupling constant $g_{j,l(R)}$, and $|e_0\rangle_2 \leftrightarrow |g_{L(R)}\rangle_2$ is resonantly coupled to the left-circularly (right-circularly) polarized mode of cavity2 with corresponding coupling constant $g_{2,L(R)}$). The tran- $\text{stions} |e_{L(R)}\rangle_j \leftrightarrow |g_{L(R)}\rangle_j \text{ and } |e_0\rangle_2 \leftrightarrow |g_0\rangle_2 \text{ are resonantly driven by classical laser fields with the time-dependent$ Rabi frequencies $\Omega_j(t)$ and $\Omega_2(t)$, respectively. Then the whole system can be dominated by the interaction Hamiltonian $(h=1)$:

$$
H_{\text{total}} = H_{al} + H_{acf},
$$

\n
$$
H_{al} = \sum_{i=L,R} \sum_{j=1,3} \Omega_j(t) |e_i\rangle_j \langle g_i| + \Omega_2(t) |e_0\rangle_2 \langle g_0| + \text{H. c. },
$$

\n
$$
H_{af} = \sum_{i=L,R} \sum_{j=1,3} \sum_{j,1} g_{j,i} a_{j,i} |e_i\rangle_j \langle g_0| + g_{2,i} a_{2,i} |e_0\rangle_2 \langle g_i| + \nu_{1,i} (a_{1,i} + a_{2,i}) b_{1,i}^{\dagger} + \nu_{2,i} (a_{2,i} + a_{3,i}) b_{2,i}^{\dagger}] + \text{H. c. },
$$
\n(10)

where H_{total} is the total Hamiltonian of the whole system, H_{al} (H_{act}) is the interaction between the atoms and the classical laser fields (the cavity-fiber system), $v_{1(2),L(R)}$ is the coupling strength between the modes of the cavity1,2 (cavity2,3) and the modes of the fiber1(2), $a_{k,L(R)}$ ($k = 1, 2, 3$) is the annihilation operator of left-circularly (right-circularly) polarized mode of *k*th cavity, and $b_{1(2),L(R)}^{\dagger}$ is the creation operator of fiber1(2) left-circularly (right-circularly) polarized mode. For simplicity, we assume $g_{k,L(R)}$ and $v_{1(2),L(R)}$ are real, $g_{k,L(R)} = g$, and $v_{1(2),L(R)} = v$.

Suppose that the total system is initially in the state $|\phi_1\rangle = |g_0\rangle_1 |g_0\rangle_2 |g_0\rangle_3 |00000\rangle_{cf_1 c_2 f_2 c_3}$ denoting *k*th atom in the state $|g_0\rangle_k$ and all of three cavities and two fibers in the vacuum state. Thus dominated by the total Hamiltonian in Eq. (10), the whole system evolves in the Hilbert space spanned by

$$
\begin{aligned}\n|\phi_1\rangle &= |g_0\rangle_1 |g_0\rangle_2 |g_0\rangle_3 |00000\rangle_{cf_1cf_2f_2c_3}, \quad |\phi_2\rangle = |g_0\rangle_1 |e_0\rangle_2 |g_0\rangle_3 |00000\rangle_{cf_1cf_2f_2c_3}, \\
|\phi_3\rangle &= |g_0\rangle_1 |g_L\rangle_2 |g_0\rangle_3 |00L00\rangle_{cf_1cf_2f_2c_3}, \quad |\phi_4\rangle = |g_0\rangle_1 |g_R\rangle_2 |g_0\rangle_3 |00R00\rangle_{cf_1cf_2f_2c_3}, \\
|\phi_5\rangle &= |g_0\rangle_1 |g_L\rangle_2 |g_0\rangle_3 |0L000\rangle_{cf_1cf_2f_2c_3}, \quad |\phi_6\rangle = |g_0\rangle_1 |g_L\rangle_2 |g_0\rangle_3 |000L0\rangle_{cf_1cf_2f_2c_3}, \\
|\phi_7\rangle &= |g_0\rangle_1 |g_R\rangle_2 |g_0\rangle_3 |0R000\rangle_{cf_1cf_2f_2c_3}, \quad |\phi_8\rangle = |g_0\rangle_1 |g_R\rangle_2 |g_0\rangle_3 |000R0\rangle_{cf_1cf_2f_2c_3}, \\
|\phi_9\rangle &= |g_0\rangle_1 |g_L\rangle_2 |g_0\rangle_3 |L0000\rangle_{cf_1cf_2f_2c_3}, \quad |\phi_{10}\rangle = |g_0\rangle_1 |g_L\rangle_2 |g_0\rangle_3 |0000L\rangle_{cf_1cf_2f_2c_3}, \\
|\phi_{11}\rangle &= |g_0\rangle_1 |g_R\rangle_2 |g_0\rangle_3 |00000\rangle_{cf_1cf_2f_2c_3}, \quad |\phi_{12}\rangle = |g_0\rangle_1 |g_R\rangle_2 |g_0\rangle_3 |00000\rangle_{cf_1cf_2f_2c_3}, \\
|\phi_{13}\rangle &= |e_L\rangle_1 |g_L\rangle_2 |g_0\rangle_3 |00000\rangle_{cf_1cf_2f_2c_3}, \quad |\phi_{14}\rangle = |g_0\rangle_1 |g_R\rangle_2 |e_R\rangle_3 |00000\rangle_{cf_1cf_2f_2c_3}, \\
$$

Obviously, the system is initially in the dark state of H_{ac} , i.e., H_{ac} , ϕ_1 = 0. Therefore, under the Zeno limit condition $\Omega_k(t) \ll g$, $\nu(k=1, 2, 3)$, the whole system can approximatively evolve in an invariant Zeno subspace consisting of dark states corresponding to the zero eigenvalue of *Hacf*:

$$
H_P = \{ |\phi_1\rangle, |\Psi_D\rangle, |\phi_{17}\rangle, |\phi_{18}\rangle, |\phi_{19}\rangle, |\phi_{20}\rangle \},
$$
\n(12)

corresponding to the projections

$$
P^{\alpha} = |\alpha\rangle\langle\alpha|, \quad (|\alpha\rangle \in H_p). \tag{13}
$$

Here,

$$
\begin{split} |\Psi_D\rangle \; &= \; \frac{1}{2\sqrt{2\nu^2 + g^2}} [2\nu|\phi_2\rangle - g(|\phi_5\rangle + |\phi_6\rangle + |\phi_7\rangle + |\phi_8\rangle) \\ &+ \nu(|\phi_{13}\rangle + |\phi_{14}\rangle + |\phi_{15}\rangle + |\phi_{20}\rangle)]. \end{split} \tag{14}
$$

Therefore, the system Hamiltonian can be rewritten as the following form based on Eq. (9):

$$
H_{\text{total}} \simeq \sum_{\alpha} P^{\alpha} H_{al} P^{\alpha}
$$

=
$$
\frac{\nu}{\sqrt{2\nu^{2} + g^{2}}} \Big[\frac{1}{2} \Omega_{1}(t) (\vert \phi_{17} \rangle + \vert \phi_{18} \rangle) \langle \Psi_{D} \vert
$$

+
$$
\Omega_{2}(t) \vert \phi_{1} \rangle \langle \Psi_{D} \vert + \frac{1}{2} \Omega_{3}(t) (\vert \phi_{19} \rangle + \vert \phi_{20} \rangle) \langle \Psi_{D} \vert \Big] + \text{H.c.}
$$
 (15)

Here setting $v = g$ and $\Omega_3(t) = \Omega_1(t)$, we can obtain an effective Hamiltonian of the system

$$
H_0(t) = \frac{1}{\sqrt{3}} (\Omega_2(t) |\Psi_1\rangle + \Omega_1(t) |\Psi_2\rangle) \langle \Psi_D| + \text{H.c.},
$$
\n(16)

in which $|\Psi_1\rangle = |\phi_1\rangle$, and $|\Psi_2\rangle = \frac{1}{2}(|\phi_{17}\rangle + |\phi_{18}\rangle + |\phi_{19}\rangle + |\phi_{20}\rangle$). The instantaneous eigenstates of *H*₀(*t*) corresponding respectively to the eigenvalues $\lambda_0 = 0$ and $\lambda_+ = \pm \Omega(t)/\sqrt{3}$ are

$$
|n_0(t)\rangle = \begin{pmatrix} -\cos\theta(t) \\ 0 \\ \sin\theta(t) \end{pmatrix}, \quad |n_{\pm}(t)\rangle = \frac{1}{\sqrt{2}} \begin{pmatrix} \sin\theta(t) \\ \pm 1 \\ \cos\theta(t) \end{pmatrix}, \tag{17}
$$

where $\Omega(t) = \sqrt{\Omega_1(t)^2 + \Omega_2^2(t)}$ and $\tan \theta(t) = \Omega_2(t)/\Omega_1(t)$.

Two methods used to generate tree-type three-dimensional entanglement

The method of Lewis-Riesenfeld invariants. In order to construct a shortcut by the LRI method for fast generating of tree-type three-dimensional entanglement, we are supposed to chase down the Hermitian invariant operator *I*(*t*) satisfying *ih* $\partial_t I(t) = [H_0(t), I(t)]$. Because of SU(2) dynamical symmetry of $H_0(t)$ in Eq. (16), *I*(*t*) can be easily given by^{27,[56](#page-11-25)}

$$
I(t) = \frac{1}{\sqrt{3}} \chi \begin{pmatrix} 0 & \cos \nu \sin \beta & -i \sin \nu \\ \cos \nu \sin \beta & 0 & \cos \nu \cos \beta \\ i \sin \nu & \cos \nu \cos \beta & 0 \end{pmatrix},
$$
(18)

where χ is an arbitrary constant in the unit of frequency keeping *I*(*t*) in the unit of energy, and ν and β are time-dependent auxiliary parameters satisfying the equations

$$
\dot{\nu} = \frac{1}{\sqrt{3}} (\Omega_2(t) \cos \beta - \Omega_1(t) \sin \beta),
$$

\n
$$
\dot{\beta} = \frac{1}{\sqrt{3}} \tan \nu (\Omega_1(t) \cos \beta + \Omega_2(t) \sin \beta).
$$
 (19)

So $\Omega_1(t)$ and $\Omega_2(t)$ can be easily deduced as follows:

$$
\Omega_1(t) = \sqrt{3} (\beta \cot \nu \cos \beta - \nu \sin \beta), \n\Omega_2(t) = \sqrt{3} (\beta \cot \nu \sin \beta + \nu \cos \beta).
$$
\n(20)

The solution of Shrödinger equation $i\hbar\partial_t |\Psi(t)\rangle = H_0(t) |\Psi(t)\rangle$ with respect to the instantaneous eigenstates of *I*(*t*) can be written as $|\Psi(t)\rangle = \sum_{n=0,\pm} C_n e^{i\alpha_n} |\phi_n(t)\rangle$, where $\alpha_n(t)$ is the Lewis-Riesenfeld phase in Eq. (3), $C_n = \langle \phi_n(0) | \Psi_1 \rangle$, and $| \phi_n(t) \rangle$ is the eigenstate of the invariant *I*(*t*) as

$$
|\phi_0(t)\rangle = \begin{pmatrix} \cos \nu & \cos \beta \\ -i & \sin \nu \\ -\cos \nu & \sin \beta \end{pmatrix}, \quad |\phi_{\pm}(t)\rangle = \frac{1}{\sqrt{2}} \begin{pmatrix} \sin \nu & \cos \beta \pm i & \sin \beta \\ i & \cos \nu \\ -\sin \nu & \sin \beta \pm i & \cos \beta \end{pmatrix}.
$$
 (21)

Then we consider a series of boundary conditions satisfying $[H_0(0), I(0)] = 0$ and $[H_0(t_f), I(t_f)] = 0$ to give Į

$$
\lim_{t \to 0} \frac{\Omega_2(t)}{\Omega_1(t)} = 0, \quad \lim_{t \to t_f} \frac{\Omega_2(t)}{\Omega_1(t)} = 2,
$$
\n(22)

where t_f is the operation time. The Eq. (22) is the guarantee for the system to evolve along $|n_0(t)\rangle$ in Eq. (17) so that we obtain the target state $|\Psi_{LRI}\rangle = \frac{1}{\sqrt{5}} |\Psi_1\rangle - \frac{2}{\sqrt{5}} |\Psi_2\rangle$. Therefore, to avoid infinite Rabi frequencies, we can choose the boundary conditions for ν and β as follows:

$$
\nu(0) = \varepsilon, \quad \nu(0) = 0, \quad \nu(t_f) = \varepsilon, \quad \nu(t_f) = 0,
$$

\n
$$
\beta(0) = 0, \quad \beta(t_f) = \arctan 2,
$$
\n(23)

where ε is a time-independent small value. Then the parameters can be easily set as

$$
\nu(t) = \varepsilon, \quad \beta(t) = \frac{\arctan 2t}{t_f}, \tag{24}
$$

and thus we can deduce

$$
\Omega_1(t) = \frac{\sqrt{3} \arctan 2}{t_f} \cot \varepsilon \cos \frac{\arctan 2t}{t_f},
$$

$$
\Omega_2(t) = \frac{\sqrt{3} \arctan 2}{t_f} \cot \varepsilon \sin \frac{\arctan 2t}{t_f}.
$$
 (25)

Based on the parameters above, we can determine the value of ε by calculating the fidelity

$$
F = |\langle \Psi_{LRl} | \Psi(t_f) \rangle|^2 = \left[\cos^2 \varepsilon + \sin^2 \varepsilon \cos \left(\frac{\arctan 2}{\sin \varepsilon} \right) \right]^2, \tag{26}
$$

 $\langle \text{where}|\Psi(t_f)\rangle = \sum_{n=0,\pm}C_ne^{i\alpha_n(t_f)}|\phi_n(t_f)\rangle$ with the Lewis-Riesenfeld phases

$$
\alpha_0(t_f) = 0, \quad \alpha_{\pm}(t_f) = \pm \frac{\arctan 2}{\sin \varepsilon},
$$
\n(27)

 $\text{and }|\Psi_{LRI}\rangle=\frac{1}{\sqrt{5}}(|\phi_1\rangle-|\phi_{17}\rangle-|\phi_{18}\rangle-|\phi_{19}\rangle-|\phi_{20}\rangle)$ is the tree-type three-dimensional entanglement generated by the LRI method. Therefore, for the appropriate Rabi frequencies and the fidelity $F=1$, we can choose

$$
\frac{\arctan 2}{\sin \varepsilon} = 2\pi, \quad \text{i.e.} \quad \varepsilon = \arcsin\left(\frac{\arctan 2}{2\pi}\right) = 0.177. \tag{28}
$$

Thus, the transformation $|\Psi_1\rangle \to |\Psi_{LR}r\rangle$ is achieved and we have constructed a shortcut by the LRI method to speed up the generation of the tree-type three-dimensional entanglement.

The method of transitionless quantum driving. Because the instantaneous eigenstates of $H_0(t)$ do not meet the Schrödinger equation, there may be transitions between the eigenstates in Eq. (17). Therefore, we need to construct the TQD Hamiltonian $H(t)$ that exactly drives the instantaneous eigenstates with no transitions between different eigenstates. Based on Eq. (6), we learn the simplest Hamiltonian *H*(*t*) is derived in the form

$$
H(t) = i \sum_{k=0,\pm} |\partial_t n_k(t)\rangle \langle n_k(t)|.
$$
\n(29)

Substituting Eq. (17) into Eq. (29), we obtain

$$
H(t) = i\dot{\theta}|\Psi_1\rangle\langle\Psi_2| + \text{H.c.},\tag{30}
$$

in which $\hat{\theta}(t) = (\Omega_2(t)\Omega_1(t) - \Omega_2(t)\Omega_1(t))/\Omega(t)^2$. The Hamiltonian (30) is constructed by a direct coupling of $|\Psi_1\rangle$ and $|\Psi_2\rangle$, and it is too hard to be implemented in experiment for such a complex system we use^{[24](#page-11-8),[35](#page-11-26)}. However, fortunately inspired by the references^{[44,](#page-11-17)46}, we find an alternative physically feasible (APF) Hamiltonian whose effect is equivalent to $H(t)$. The APF design is shown in [Fig. 1\(c\).](#page-2-0) Comparing [Fig. 1\(b,c\),](#page-2-0) we change all of the resonant atomic transitions into non-resonant atomic transitions with detuning Δ .

The interaction Hamiltonian of the non-resonant system reads

$$
H'_{total} = H'_{al} + H'_{acf} + H_e,
$$

\n
$$
H'_{al} = \sum_{i=L,R} \sum_{j=1,3} \Omega'_{j}(t) |e_{i}\rangle_{j} \langle g_{i}| + \Omega'_{2}(t) |e_{0}\rangle_{2} \langle g_{0}| + \text{H.c.},
$$

\n
$$
H'_{acf} = \sum_{i=L,R} \sum_{j=1,3} \sum_{j, i} a_{j,i} a_{j,i} |e_{i}\rangle_{j} \langle g_{0}| + g_{2,i} a_{2,i} |e_{0}\rangle_{2} \langle g_{i}| + \nu_{1,i} (a_{1,i} + a_{2,i}) b_{1,i}^{\dagger} + \nu_{2,i} (a_{2,i} + a_{3,i}) b_{2,i}^{\dagger}| + \text{H.c.},
$$

\n
$$
H_{e} = \sum_{i=L,R} \sum_{j=1,3} \Delta |e_{i}\rangle_{j} \langle e_{i}| + \Delta |e_{0}\rangle_{2} \langle e_{0}|.
$$
\n(31)

Then similar to the approximation for the Hamiltonian from Eqs (10–16), an effective Hamiltonian for the present non-resonant system can be obtained

$$
H_0'(t) = \frac{1}{\sqrt{3}} \left[|\Psi_D\rangle (\Omega_2'(t) \langle \Psi_1| + \Omega_1'(t) \langle \Psi_2|) + \text{H.c.} \right] + \Delta |\Psi_D\rangle \langle \Psi_D|.
$$
\n(32)

Under the limit condition $\Delta \gg Ω_1'(t)/\sqrt{3}$, $Ω_2'(t)/\sqrt{3}$, by adiabatically eliminating the state $|\Psi_D\rangle$, the effective Hamiltonian $H'_0(t)$ becomes

$$
H_{\text{eff}}(t) = \frac{1}{3\Delta} \left[\left| \Omega_2'(t) \right|^2 \left| \Psi_1 \right\rangle \left\langle \Psi_1 \right| + \left| \Omega_1'(t) \right|^2 \left| \Psi_2 \right\rangle \left\langle \Psi_2 \right| + \left(\Omega_1'^*(t) \Omega_2'(t) \left| \Psi_2 \right\rangle \left\langle \Psi_1 \right| + \text{H.c.} \right) \right]. \tag{33}
$$

The first two terms can be removed by setting $\Omega'_{1}(t) = i\Omega'_{2}(t)$ and the final effective Hamiltonian becomes

$$
H'_{eff}(t) = i\frac{\Omega_2'(t)^2}{3\Delta} |\Psi_1\rangle \langle \Psi_2| + \text{H. c.}
$$
\n(34)

This effective Hamiltonian is equivalent to the Hamiltonian *H*(*t*) in Eq. (30) if we set $\dot{\theta}(t) = \Omega_2'(t)^2/3\Delta$, i.e.,

$$
\Omega_2'(t) = \sqrt{\frac{3\Delta(\dot{\Omega}_2(t)\Omega_1(t) - \Omega_2(t)\dot{\Omega}_1(t))}{\Omega(t)^2}},\tag{35}
$$

which is the correlation between the Rabi frequencies of the TQD method and the Rabi frequencies of STIRAP. By setting the Rabi frequencies of STIRAP to satisfy the same boundary conditions as Eq. (22), we can achieve the

Figure 2. The fidelity for the LRI method versus (a) t_f/g^{-1} with $\varepsilon = 0.177$ and (b) ε with $t_f = 80/g$, respectively; (**c**) the three dimensional image of the fidelity for the LRI method versus t_f/g^{-1} and ε .

transformation $\ket{\Psi_\psi}\to\ket{\Psi_{TQD}}$ to implement the fast generation of tree-type three-dimensional entanglement, where $|\Psi_{TQD}\rangle = \frac{1}{\sqrt{5}} (|\phi_1\rangle + |\phi_{17}\rangle + |\phi_{18}\rangle + |\phi_{19}\rangle + |\phi_{20}\rangle)$ is the tree-type three-dimensional entanglement generated by the TQD method.

Numerical simulations and comparisons between LRI and TQD

In the following, we will give the numerical simulations in three subsections to discuss respectively the selections of parameters of the two methods, the feasibility of generating tree-type three-dimensional entanglement and the robustness of our schemes. Also the comparisons between the LRI method and the TQD method will be included in every subsection.

Selections of parameters. Firstly, to determine the parameters of the LRI method, we plot the fidelity $F = |\langle \Psi_{LRl} | \Phi(t_f) \rangle|^2$ versus the operation time t_f and ε in [Fig. 2](#page-6-0), where $|\Phi(t_f)\rangle$ is the state at the time $t = t_f$ of the whole system governed by the total Hamiltonian *H*_{total} in Eq. (10). In [Fig. 2\(a\),](#page-6-0) we plot the relation between the fidelity and the operation time t_f with ε = 0.177 which is determined in Eq. (28). And we can see that in a very short operation time t_f = 80/*g* the fidelity is already almost unity: $F(t_f=80/g)$ = 0.996. From [Fig. 2\(b\)](#page-6-0), we can find that under $t_f = 80/g$ when $\varepsilon = 0.177$ the fidelity is highest. Thus we can choose $t_f = 80/g$ and $\varepsilon = 0.177$ as the parameters of the LRI method in the following discussion. Furthermore, in order to consider the joint effects of *t*_f and ε on the fidelity we plot the three dimensional image of the fidelity versus *t_i*/g^{−1} and ε in [Fig. 2\(c\)](#page-6-0). From the three dimensional image, it is clear that the effects of t_f and ε on the fidelity are not dependent on each other.

Next we determine the parameters of the TQD method. In order to satisfy the boundary conditions in Eq. (22), the Rabi frequencies $\Omega_1(t)$ and $\Omega_2(t)$ in the original Hamiltonian H_{total} are chosen as^{[46](#page-11-27)}

$$
\Omega_1(t) = \frac{1}{\sqrt{5}} \Omega_0 e^{-(t-t_f/2-\tau)^2/T^2} + \Omega_0 e^{-(t-t_f/2+\tau)^2/T^2},
$$
\n
$$
\Omega_2(t) = \frac{2}{\sqrt{5}} \Omega_0 e^{-(t-t_f/2-\tau)^2/T^2},
$$
\n(36)

where Ω_0 is the pulses' amplitude, t_f is the operation time, and τ and T are related parameters. The time-dependent $\Omega_1(t)$ and $\Omega_2(t)$ are shown in [Fig. 3.](#page-7-0)

Based on the correlation in Eq. (35), the Rabi frequencies of the TQD method can be figured out. As an illustration, we plot the fidelity $F=|\langle\Psi_{TQD}|\Phi(t_f)\rangle|^2$ versus the detuning Δ and t_f in [Fig. 4](#page-7-1), where $|\Phi(t_f)\rangle$ is the state at the time $t = t_f$ of the whole system governed by the total Hamiltonian H'_{total} in Eq. (31). To compare with each other effectively, we choose the same operation time $t_f = 80/g$ in the TQD method as that in the LRI method. From [Fig. 4\(a\),](#page-7-1) we can find that under $t_f=80/g$ when $\Delta=6g$ the fidelity is highest. Besides, we can see that the fidelity is almost unity: $F(t_f = 80/g) = 0.996$ at the point $t_f = 80/g$ from [Fig. 4\(b\).](#page-7-1) Thus we choose $t_f = 80/g$ and $\Delta = 6g$ as the parameters of the TQD method in the following discussion. Similar to the LRI method, in order to consider the joint effects of t_f and Δ on the fidelity we plot the three dimensional image of the fidelity versus t_f/g^{-1} and Δ/g in [Fig. 4\(c\).](#page-7-1) However, from [Fig. 4\(c\),](#page-7-1) we are not able to judge whether the effects of t_f and Δ on the fidelity are dependent or not dependent on each other. We will make a detailed discussion about the joint effects of t_f and Δ on the fidelity later in the last subsection.

Next, a brief discussion is very necessary to show our schemes are fast. Firstly, it is quite apparent that the operation time $t_f = 80/g$ is far shorter than $t = 3000/g$ which is the generation time of the tree-type three-dimensional entanglement generated in the ref. [23](#page-11-4). Further more, for a more effective discussion, we consider the adiabatic condition $|\dot{\theta}(t)| \ll \Delta E^{57}$, where $\Delta E = \Omega(t)/\sqrt{3}$ is the minimum separation between the eigenvalues of the Hamiltonian (16). For convenience, we just consider the LRI method for the shortcut to adiabatic passage. With a simple calculation, for the shortcut to adiabatic passage, we can obtain the constants $\vert \dot{\theta}_{LRI} \vert = 0.014$ and $\Delta E_{LRI} = 0.077$ by using Eq. (25). However, for ref. [23](#page-11-4), we obtain time-dependent but $\frac{dE}{dt}$ single-peak $|\dot{\theta}(t)|$ and ΔE , and the amplitudes are $|\dot{\theta}|_{max} = 0.003$ and $\Delta E_{max} = 0.067$, respectively. Obviously,

according to the information above, the adiabatic condition $|θ(t)| ≤ ∆E$ is met in the ref. [23](#page-11-4) but needs not to be met in our schemes, which means that the slow process of the generation of the tree-type three-dimensional entanglement is sped up.

Discussion of feasibility. In this subsection, we will give the numerical simulations for discussing the feasibility of our two schemes. For the LRI method, we plot the time-dependent Rabi frequencies $\Omega_1(t)$ and $\Omega_2(t)$ which are described by Eq. (25) in [Fig. 5\(a\).](#page-8-0) The evolutions of populations $P_{1(17\sim20)}^{LRI}$ of states $|\phi_{1(17\sim20)}\rangle$ governed by H_{total} are shown in [Fig. 5\(b\).](#page-8-0) For the TQD method, we plot the time-dependent Rabi frequency $\Omega'_2(t)$ $(|\Omega'_1(t)|)$ which is described by Eq. (35) in [Fig. 5\(c\)](#page-8-0). The evolutions of populations $P_{1(17\sim20)}^{TQD}$ of states $|\phi_{1(17\sim20)}\rangle$ governed by H'_{total} are shown in Fig. 5(d). In addition, in Fig. 6(a) we plot the atomic excited populations $P_a^{LRT} = \sum_{m=2,13\sim 16} P_m^{LRI}$ (red solid line) and $P_a^{TQD} = \sum_{m=2,13\sim 16} P_m^{TQD}$ (blue solid line) corresponding to the LRI method and the TQD method respectively, and the cavity-fiber excited populations $P_{cf}^{LRI} = \sum_{n=3 \sim 12} P_n^{LRI}$ (red dashed line) and $P_{cf}^{TOP} = \sum_{n=3 \sim 12} P_{n}^{TOP}$ (blue dashed line) corresponding to the cases of the LRI method and the TQD method respectively. In [Fig. 6\(b\)](#page-8-1), we plot the fidelities of the tree-type three-dimensional entanglement generated by the LRI method (red dashed line) and the TQD method (blue solid line), respectively.

Here we first consider [Fig. 6\(b\)](#page-8-1). From [Fig. 6\(b\)](#page-8-1), we know that both of the two lines of fidelity illustrated based on the two methods reach a very high point close to unity at the time *t*= 80/*g*, and thus both of our schemes are feasible. [Figure 6\(b\)](#page-8-1) also shows that the fidelity of the TQD method can reach a very high value within a shorter time than that of the LRI method. Next we consider the populations of the target states in [Fig. 5\(b,d\).](#page-8-0) We can see that a near perfect result we expect is obtained in [Fig. 5\(b\)](#page-8-0), but a little bit imperfect result appears in [Fig. 5\(d\)](#page-8-0) in which there exists a small gap between two lines of the populations of $|\phi_1\rangle$ and $|\phi_{17-20}\rangle$. Therefore, for the transformation of populations, the LRI method is a bit better than the TQD method. However, when compare the pulse types in [Fig. 5\(a,c\)](#page-8-0), we find that the TQD method is more feasible than the LRI method. Because the pulses in LRI method are short-time truncations of two harmonic pulses and the truncations of the two harmonic pulses in a short time are too hard to be achieved. But the pulses in TQD method are almost complete Gaussian pulses which are relatively easy to be achieved. Moreover, the populations of atomic and cavity-fiber excited states for two methods are shown in Fig. $6(a)$, and all of the populations of excited states are near zero at the time $t = 80/g$.

Figure 5. For the LRI method, (**a**) the time dependence of the Rabi frequencies $\Omega_1(t)$ (blue solid line) and $\Omega_2(t)$ (red dashed line); (**b**) the populations of states $|\phi_1\rangle$ and $|\phi_{17-20}\rangle$ governed by H_{total} . For the TQD method, (**c**) the time dependence of the Rabi frequency $\Omega'_2(t)$ ($\Omega'_1(t)$); (**d**) the populations of states $|\phi_1\rangle$ and $|\phi_{17-20}\rangle$ governed by $H'_{\rm total}$. The parameters used here are $t_f{=}80/g, \varepsilon{=}0.177, \Delta{=}6g, \tau{=}0.14t_f$ and $T{=}0.19t_f$

Figure 6. (**a**) The time dependence of atomic excited populations (red solid line) and cavity-fiber excited populations (red dashed line) in the LRI method, and the time dependence of atomic excited populations (blue solid line) and cavity-fiber excited populations (blue dashed line) in the TQD method; (**b**) the fidelities of treetype three-dimensional entanglement for the LRI method (red dashed line) and the TQD method (blue solid line). The parameters used here are same as in [Fig. 5](#page-8-0).

So we can deduce that whichever method is employed, the state of the whole system almost populates in tree-type three-dimensional entanglement.

It is worth explaining the gap between the two lines of the populations of $|\phi_1\rangle$ and $|\phi_{17\sim20}\rangle$ in [Fig. 5\(d\)](#page-8-0). For the TQD method, there are two limit conditions $\Omega'_k(t) \ll g$, $v(k=1, 2, 3)$ and $\Delta \gg g$, *v* applied to prepare tree-type three-dimensional entanglement. However, as shown in [Fig. 5\(c\),](#page-8-0) the amplitude of $\Omega'_2(t)$ ($|\Omega'_1(t)|$) is 0.8 *g* which does not strictly meet the limit condition $\Omega'_k(t) \ll g$, *v*. And also the detuning $\Delta = 6g$ does not strictly meet the limit condition $\Delta \gg g$, *v*. In fact, these two limit conditions are difficult to be coordinated. With no assignments of t_f and Δ , we calculate the amplitude of $\Omega'_2(t)$ with the parameters $\tau = 0.14t_f$ and $T = 0.19t_f$ to obtain

$$
\Omega_0' \approx \Omega_2'(t)|_{t=\frac{1}{2}t_f} = 2.9 \sqrt{\frac{\Delta}{t_f}},\tag{37}
$$

which is not dependent on the amplitude Ω_0 of $\Omega_{1,2}(t)$ in Eq. (36) but only proportional to $\sqrt{\Delta/t_f}$. Thus, Ω_0' roughly equals to $\sqrt{\Delta/3}$ if the operation time is chosen as $t_f=80/g$. Nevertheless, the ratio 1/3 is not small enough to satisfy both two limit conditions $\Omega'_k(t) \ll g$, ν and $\Delta \gg g$, ν , i.e., the condition $\Omega'_k(t) \ll g$, ν will not be satisfied if the limit condition ∆ *g*, *v* is satisfied and vice versa. Therefore, there exists a gap between the two lines of the populations of $|\phi_1\rangle$ and $|\phi_{17\sim20}\rangle$ in [Fig. 5\(d\).](#page-8-0) In addition, Eq. (37) reveals that $\Omega'_2(t)$'s amplitude $\Omega'_0 \propto \sqrt{\Delta t_f}$. It is known that the fidelity of the TQD method is strongly dependent on Ω'_0 , so we can deduce that the fidelity of the TQD method is strongly dependent on the value of Δ/t_f . As a result, in [Fig. 4\(c\),](#page-7-1) the effects of *t_f* and Δ on the fidelity are dependent on each other.

Based on the discussion above, for fast generating tree-type three-dimensional entanglement, both the LRI method and the TQD method are feasible. Besides, the two methods have their own advantages and disadvantages and we can choose a certain method depending on the conditions in experiment.

Figure 7. (**a**) The fidelity of the LRI method versus *δtf* /*tf* and *δε*/*ε*; (**b**) the fidelity of the TQD method versus *δt_f*/ t_f and *δ*Δ/Δ. The parameters used here are same as in [Fig. 5](#page-8-0).

Discussion of robustness. In the above discussion, the whole system are perfect and considered as absolutely isolated from the environment. Therefore, it is necessary to give the discussions of robustness of our schemes against the variations in the parameters and decoherence induced by the atomic spontaneous emissions and photon leakages of the cavity-fiber system. For discussing the effects of the variations in the parameters, we plot the fidelity of the LRI method versus the variations in t_f and ε in [Fig. 7\(a\)](#page-9-0) and the fidelity of the TQD method versus the variations in t_f and Δ in [Fig. 7\(b\).](#page-9-0) Here we define $\delta x = x' - x$ as the deviation of *x*, in which *x* denotes the ideal value and x' denotes the actual value. In [Fig. 7\(a\)](#page-9-0), the fidelity decreases with the increase of $|\delta \varepsilon|$ as described in [Fig. 2\(b\).](#page-6-0) From Eq. (25), we know that the Rabi frequencies decrease with the increase of the operation time t_f . According to the limit condition $\Omega_k(t) \ll g$, *v* we use, the values of the Rabi frequencies are the smaller the better, so the operation time t_f is the longer the better as described in [Fig. 2\(a\).](#page-6-0) Therefore, the fidelity of the LRI method increases with the increase of δt_f in [Fig. 7\(a\).](#page-9-0) In [Fig. 7\(b\)](#page-9-0), we can clearly see that the effects of *t*_{*j*} and Δ on the fidelity of the TQD method are dependent on each other and even the fidelity of the TQD method is apparently dependent on the value of Δ/t_f as mentioned in the last subsection. Significantly, we notice that the fidelities of the two methods are both over 0.98 even when $|\delta x/x| = 0.1$ ($x = t_f$, ε , Δ). Therefore, both of our schemes are robust against the variations in the parameters.

We can also see that the smallest fidelity of the TQD method in [Fig. 7\(b\)](#page-9-0) is slightly higher than the smallest fidelity of the LRI method in [Fig. 7\(a\).](#page-9-0) This fact can easily be found by comparing Figs $2(c)$ and $4(c)$, in which there is a greater advisable range of the parameters in the TQD method than in the LRI method for preparing tree-type three-dimensional entanglement with a high fidelity.

Next taking the decoherence induced by the atomic spontaneous emissions and photon leakages of the cavity-fiber system into account, the whole system is dominated by the master equation

$$
\rho(t) = -i[H, \rho(t)]
$$
\n
$$
- \sum_{i=L,R} \sum_{j=1,3} \frac{\gamma_i^j}{2} (\sigma_{e_je_i}^j \rho - 2\sigma_{g_je_i}^j \rho \sigma_{e_je_i}^j + \rho \sigma_{e_je_i}^j)
$$
\n
$$
- \sum_{i=L,R} \sum_{j=1,3} \frac{\gamma_i^j}{2} (\sigma_{e_je_i}^j \rho - 2\sigma_{g_0e_i}^j \rho \sigma_{e_je_0}^j + \rho \sigma_{e_je_i}^j)
$$
\n
$$
- \sum_{i=L,0,R} \frac{\gamma_i^2}{2} (\sigma_{e_0e_0}^2 \rho - 2\sigma_{g_je_0}^2 \rho \sigma_{e_0g_i}^2 + \rho \sigma_{e_0e_0}^2)
$$
\n
$$
- \sum_{i=L,R} \sum_{j=1,2} \frac{\kappa_{j,i}^f}{2} (b_{j,i}^{\dagger} b_{j,i} \rho - 2b_{j,i} \rho b_{j,i}^{\dagger} + \rho b_{j,i}^{\dagger} b_{j,i})
$$
\n
$$
- \sum_{i=L,R} \sum_{j=1,2,3} \frac{\kappa_{j,i}^c}{2} (a_{j,i}^{\dagger} a_{j,i} \rho - 2a_{j,i} \rho a_{j,i}^{\dagger} + \rho a_{j,i}^{\dagger} a_{j,i}), \qquad (38)
$$

where *H* is the total Hamiltonian H_{total} of the LRI method in Eq. (10) or H'_{total} of the TQD method in Eq. (31). γ_i^j ($i = L$, 0, R, $j = 1, 2, 3$) is the spontaneous emission rate of *j*th atom from the excited state $|e_i\rangle_j$ to the ground state $|g_i\rangle_j$; $\kappa^f_{j,i}$ and $\kappa^c_{j,i}$ denote the photon leakage rates from the fiber modes and the cavity modes, respectively; $\sigma_{mn}^j = |m\rangle_j \langle n|(m, n = e_i, g_i)$ is Pauli operators. For simplicity, we assume $\gamma_i^j = \gamma$ and $\kappa_{j,i}^f = \kappa_{j,i}^c = \kappa$.

Based on the master equation, we plot the fidelities of the LRI method and the TQD method versus *κ*/*g* and *γ*/*g* in [Fig. 8](#page-10-11). As we can see from the decrease of the fidelity of the LRI method with the increases of *κ*/*g* and *γ*/*g* in [Fig. 8\(a\)](#page-10-11), we learn that the influence of atomic spontaneous emissions on the fidelity is greater than that of photon leakages of the cavity-fiber system. However, in [Fig. 8\(a\)](#page-10-11) the influence of cavity-fiber photon leakages on the fidelity of the TQD method plays a full role, but that of atomic spontaneous emissions is little. As a cross reference, we can get some inspiration from [Fig. 6\(a\).](#page-8-1) In [Fig. 6\(a\),](#page-8-1) the highest value of the cavity-fiber excited populations (blue dashed line) of the TQD method is over 0.2 during the evolution process but that of the atomic excited populations (blue solid line) of the TQD method is near zero which caused by the detuning Δ . The highest values

of the atomic excited populations (red solid line) and the cavity-fiber excited populations (red dashed line) of the LRI method are slightly higher and slightly lower than 0.05, respectively. Therefore, it is no doubt that the results of [Figs 6](#page-8-1) and [8](#page-10-11) are corresponding to each other. Finally, it is necessary to emphasize that the fidelities of the LRI method and the TQD method are near 0.94 and over 0.955 respectively, even when $\kappa = \gamma = 0.02 g$. Therefore, our two schemes of the LRI method and the TQD method both are robust against the decoherence induced by the atomic spontaneous emissions and photon leakages of the cavity-fiber system.

Experimental feasibility and conclusion

Now we show the experimental feasibility of our schemes. As mentioned in the ref. [23](#page-11-4), ⁸⁷Rb can be used in our schemes to construct the required atomic level configurations. For ⁸⁷Rb, $|F = 1$, $m_F = -1$), $|F = 1$, $m_F = 0$) and $|F = 1$, $m_F = 1$ of 5*S*_{1/2} can be used as the ground states $|g_L\rangle$, $|g_0\rangle$ and $|g_R\rangle$ respectively, and $|F = 1$, $m_F = -1$), $F = 1$, $m_F = 0$ and $F = 1$, $m_F = 1$ of $5P_{3/2}$ can be used as the excited states $|e_L\rangle$, $|e_0\rangle$ and $|e_R\rangle$ respectively. Based on a set of cavity QED parameters *g*=2*π*×750MHz, *γ*=2*π*×3.5MHz and *κ*=2*π*×2.62MHz, which are achieved in recent experiments^{58–60}, we can obtain the very high fidelities $F_{LRI}=$ 0.984 and $F_{TQD}=$ 0.990 corresponding to the LRI method and the TQD method respectively, which show our schemes for generating tree-type three-dimensional entangled states both are feasible in the experiment.

In conclusion, we have proposed two schemes to speed up the generations of the tree-type three-dimensional entanglement via Lewis-Riesenfeld invariants and transitionless quantum driving. The two different tree-type three-dimensional entangled states are prepared among three atoms trapped respectively in three spatially separated optical cavities connected by two fibers. The strict numerical simulations show that the LRI method and the TQD method both are feasible and robust against the variations in the parameters, atomic spontaneous emissions and photon leakages of the cavity-fiber system. Besides, comparing the two methods, we know they both have their own advantages and disadvantages. So we can choose different methods depending on different conditions in experiment. In short, both of our schemes are fast, robust and feasible. We hope that the tree-type three-dimensional entanglement would contribute to the improvement of quantum communication security and our work would be useful for the experimental realization of quantum information processing in the near future.

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Author Contributions

J.-L.W. designed the scheme under the guidance of X.J. and S.Z. All authors contributed to the interpretation of the work and the writing of the manuscript. All authors participated in the writing and revising of the text.

Additional Information

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