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Dissipative preparation of W states in trapped ion systems

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Abstract

We present protocols for dissipative entanglement of three trapped-ion qubits and discuss a scheme that uses sympathetic cooling as the dissipation mechanism. This scheme relies on tailored destructive interference to generate any one of six entangled W states in a three-ion qubit space. Using a beryllium–magnesium ion crystal as an example system, we theoretically investigate the protocol's performance and the effects of likely error sources, including thermal secular motion of the ion crystal, calibration imperfections, and spontaneous photon scattering. We estimate that a fidelity of $\sim 98\%$ may be achieved in typical trapped ion experiments with ~ 1 ms interaction time. These protocols avoid timescale hierarchies for faster preparation of entangled states.

1. Introduction

Dissipation arises in quantum systems through interaction with the environment and presents a challenge for applications in quantum simulation, computation, communication, and metrology. However, controlled dissipation can also be introduced and leveraged to manipulate quantum systems. Familiar examples in atomic physics include optical pumping and laser cooling. These techniques allow removal of entropy and approximate preparation of a desired pure state from an uncontrolled and unknown initial state. This cannot be accomplished by unitary operations. Recently, attention has focused on using dissipation for quantum information processing [1–3], and in particular for generating entanglement. While not inherently superior to unitary entanglement-generation strategies, dissipative schemes are less sensitive to certain error mechanisms. Moreover, they allow resource states to be created and stabilized in the presence of noise so that they are available on demand. Dissipative protocols for generation and stabilization of entangled and other nonclassical states have been demonstrated in a number of systems, including macroscopic atomic ensembles [4], trapped ions [5–7], and superconducting qubits [8–10]. Numerous proposals describe additional schemes to generate entanglement [11–17], perform error correction [18, 19], and initialize quantum simulators [20]. Broadly, the full scope over which engineered dissipation may be applied for quantum information processing is not yet clear, and practical protocols that can accomplish new tasks expand the frontier.

A first set of experiments has demonstrated preparation of steady-state entanglement using couplings that are applied continuously [6, 8]. An important ingredient in these schemes is timescale hierarchies, for example the application of a strong dressing drive at rate G alongside other interactions with characteristic rates g_i . The dressing drive creates resonances that are resolved by the other drives in the limit $g_i \ll G$, and this hierarchy of timescales $g_i^{-1} \gg G^{-1}$ protects the target state. However, the steady-state entanglement fidelity of this kind of scheme only asymptotically approaches unity as the relative strength $r = G/\max\{g_i\}$ of the dressing drive increases. Even more importantly, timescale hierarchies limit the speed of entanglement generation, because the other interactions g_i that populate the target state must be driven slowly compared to experimentally achievable rates for G . This puts practical limits on the speed of state preparation and the achievable fidelities in the presence of various error sources, and these limits are worse

than the fundamental limit determined by r . Recent proposals for generation of bipartite entanglement of trapped ions [21] and superconducting qubits [22] avoid imposing timescale hierarchies. As a result, they are predicted to generate entanglement more efficiently with the same experimental resources and achieve unity steady-state fidelities for finite rates in the absence of errors. In this work, we describe how couplings available in trapped-ion systems can extend these concepts for efficient dissipation engineering to the preparation of tripartite entanglement.

We present concrete schemes for the dissipative production of three-qubit W states. For single-qubit basis states $|\uparrow\rangle$ and $|\downarrow\rangle$, W states are N -qubit entangled states of the general form $(e^{i\phi_1}|\uparrow\downarrow\downarrow\dots\rangle + e^{i\phi_2}|\downarrow\uparrow\downarrow\dots\rangle + \dots + e^{i\phi_N}|\dots\downarrow\uparrow\rangle) / \sqrt{N}$, where each phase ϕ_i is defined modulo 2π and, as usual, the state is defined up to a global phase. W states are notable in part because their entanglement is particularly robust to particle loss⁴ [23]. Unitary trapped-ion W-state generation was first demonstrated in 2005 using $^{40}\text{Ca}^+$ ions [24], and individual addressing of the ions was crucial to that demonstration. A related proposal using rapid adiabatic passage and targeted addressing of a portion of an ion chain demonstrated that this approach could be used to create generalized W states [25]. It has also been proposed that W states can be generated using globally addressed composite pulses on the motional sideband [26]. Another experiment using Hilbert-space engineering relied on global couplings and hence did not require individual qubit addressing [27]; however, it imposed a hierarchy of timescales akin to those in early dissipative state-preparation schemes. Dissipative preparation of W states was considered in reference [28], which discusses quasi-local couplings that drive a multipartite quantum system into a steady state. The dissipative protocols we present here require neither timescale hierarchies nor individual addressing. Instead, they exploit symmetries to enable the efficient generation of any one of six target W states in the three-qubit Hilbert space (section 1.1, equations (3)–(8)).

We structure our presentation as follows: in section 1.1 we present a useful notation for the W states of a three-qubit system. This notation naturally describes the couplings that our schemes employ. In section 1.2 we review one possible application for these W states, the noiseless-subsystem encoding for a logical qubit. In section 2 we describe the experimental platform in which we envision implementing these schemes. We discuss a paradigmatic system, a mixed-species, five-ion linear crystal composed of three qubit ions and two auxiliary ions of a different species for sympathetic cooling, and review the relevant features of the interactions our schemes use: resolved-sideband couplings, Mølmer–Sørensen (MS) spin–spin interactions, and sympathetic cooling. Section 3 describes dissipative protocols for W-state generation. In section 3.1 we consider a natural attempt to extend the dissipative singlet-generation scheme presented by Horn *et al* [21], and discuss the failure of this scheme to generalize efficiently to more qubits. Section 3.2 presents a scheme for dissipative generation of W states of three ions that is more promising for experimental implementation, in which dissipation is incorporated through sympathetic cooling of the collective motion. We discuss expected error mechanisms, which indicate from simulations that dissipative W-state generation in a typical trapped-ion experiment can be achieved with $\sim 98\%$ fidelity and ~ 1 ms interaction time. Section 4 presents a concluding discussion.

1.1. W state notation

To span the eight-dimensional Hilbert space for three qubits, we choose orthonormal basis states $|W_{n_\uparrow s}\rangle$ parameterized by $n_\uparrow$, the number of qubits that are in the $|\uparrow\rangle$ state, and the chirality eigenvalue $s = 0, +1$, or -1 under a chirality operator [29]:

$$\chi = \frac{1}{2\sqrt{3}} \sum_{\alpha\beta\gamma} \epsilon_{\alpha\beta\gamma} \sigma_\alpha^{(1)} \sigma_\beta^{(2)} \sigma_\gamma^{(3)}. \quad (1)$$

Here the indices α, β , and γ run over x, y , and z , $\sigma_\alpha^{(j)}$ is a Pauli matrix acting on qubit j , and $\epsilon_{\alpha\beta\gamma}$ is the Levi-Civita symbol, which contributes a factor of 0 or ± 1 depending on the permutation $\alpha\beta\gamma$. Explicitly, the basis states $|W_{n_\uparrow s}\rangle$ are:

$$|W_{00}\rangle = |\downarrow\downarrow\downarrow\rangle, \quad (2)$$

$$\sqrt{3}|W_{10}\rangle = |\uparrow\downarrow\downarrow\rangle + |\downarrow\uparrow\downarrow\rangle + |\downarrow\downarrow\uparrow\rangle, \quad (3)$$

$$\sqrt{3}|W_{1+}\rangle = w^*|\uparrow\downarrow\downarrow\rangle + |\downarrow\uparrow\downarrow\rangle + w|\downarrow\downarrow\uparrow\rangle, \quad (4)$$

$$\sqrt{3}|W_{1-}\rangle = w|\uparrow\downarrow\downarrow\rangle + |\downarrow\uparrow\downarrow\rangle + w^*|\downarrow\downarrow\uparrow\rangle, \quad (5)$$

⁴ A straightforward example is the effect of measurement of a single qubit on a W state, in which case an $(N-1)$ -qubit W state is preserved with probability $(N-1)/N$. This compares favorably with the same measurement on a generalized Greenberger–Horne–Zeilinger state, when entanglement is always destroyed.

$$\sqrt{3}|W_{20}\rangle = |\downarrow\uparrow\uparrow\rangle + |\uparrow\downarrow\uparrow\rangle + |\uparrow\uparrow\downarrow\rangle, \quad (6)$$

$$\sqrt{3}|W_{2+}\rangle = w^*|\downarrow\uparrow\uparrow\rangle + |\uparrow\downarrow\uparrow\rangle + w|\uparrow\uparrow\downarrow\rangle, \quad (7)$$

$$\sqrt{3}|W_{2-}\rangle = w|\downarrow\uparrow\uparrow\rangle + |\uparrow\downarrow\uparrow\rangle + w^*|\uparrow\uparrow\downarrow\rangle, \quad (8)$$

$$|W_{30}\rangle = |\uparrow\uparrow\uparrow\rangle, \quad (9)$$

where $w = \exp(2\pi i/3)$ and phases are defined modulo 2π . The six W states are the states with $n_\uparrow = 1$ or 2 (equations (3)–(8)), and we refer to the four of these states $|W_{n_\uparrow\pm}\rangle$ with chirality eigenvalue $s = \pm 1$ as chiral W states, in contrast with the achiral states for which $s = 0$.

1.2. Noiseless-subsystem qubit encoding

Here we review one possible application for three-ion W states. In order to implement a practical quantum computer or memory, it is necessary to employ either active error correction or encodings for logical qubits that are inherently robust to noise, for which several strategies have been proposed and realized [30–36]. Chiral W states $|W_{n_\uparrow\pm}\rangle$ can be used in a noiseless-subsystem (NSS) qubit encoding [29, 37]. This encoding generalizes a two-qubit decoherence-free subspace (DFS) encoding, which is robust against global σ_z dephasing, to a three-qubit encoding that is insensitive to any global Pauli operator. Explicitly, the noise operator against which the encoding is protected is:

$$\left(\vec{\sigma}_1 \cdot \vec{n}(t)\right) \otimes \mathbb{1}_2 \otimes \mathbb{1}_3 + \mathbb{1}_1 \otimes \left(\vec{\sigma}_2 \cdot \vec{n}(t)\right) \otimes \mathbb{1}_3 + \mathbb{1}_1 \otimes \mathbb{1}_2 \otimes \left(\vec{\sigma}_3 \cdot \vec{n}(t)\right), \quad (10)$$

where $\vec{\sigma}_j$ is the vector of Pauli matrices and $\mathbb{1}_j$ is the identity operator, each acting on qubit j , and the slowly varying $\vec{n}(t)$ describes the magnitude and direction of the global noise as a function of time. Examples of NSS encodings include an exchange-only qubit [35, 38, 39] and the related encoding that we discuss below, which has been realized and explored in nuclear magnetic-resonance systems [40, 41].

In a subsystem encoding, a qubit is encoded in a tensor factor of a Hilbert (sub-)space. A noiseless subsystem encoding is realized when the qubit degree of freedom does not couple to the global environment; only the remaining ‘gauge’ degrees of freedom do. In the NSS encoding that we consider, a qubit is encoded in the total-spin- $\frac{1}{2}$ subspace that is spanned by the chiral W states. States within this subspace can be factored into joint eigenstates of a gauge S_z degree of freedom with eigenvalues $\pm \frac{\hbar}{2}$ (related to the label $n_\uparrow$) and a chiral qubit degree of freedom with eigenvalues ± 1 . Global noise of the form described by equation (10) couples only to the former, so that the qubit is unaffected if the two degrees of freedom are separable.

Concretely, we associate the logical state $|\downarrow\rangle_L$ with chirality $s = +1$ and $|\uparrow\rangle_L$ with chirality $s = -1$. With this choice, the pair $|W_{1+}\rangle, |W_{1-}\rangle$ and the pair $|W_{2+}\rangle, |W_{2-}\rangle$ each constitute a DFS qubit that is protected from global σ_z noise, and global σ_x and σ_y noise map between these qubits. Arbitrary logical states are protected against any global noise as defined in equation (10) when they are encoded in terms of gauge coefficients G_i (which do not affect the logical state) as:

$$\alpha_L|\downarrow\rangle_L + \beta_L|\uparrow\rangle_L \rightarrow G_1(\alpha_L|W_{1+}\rangle + \beta_L|W_{1-}\rangle) + G_2(\alpha_L|W_{2+}\rangle + \beta_L|W_{2-}\rangle), \quad (11)$$

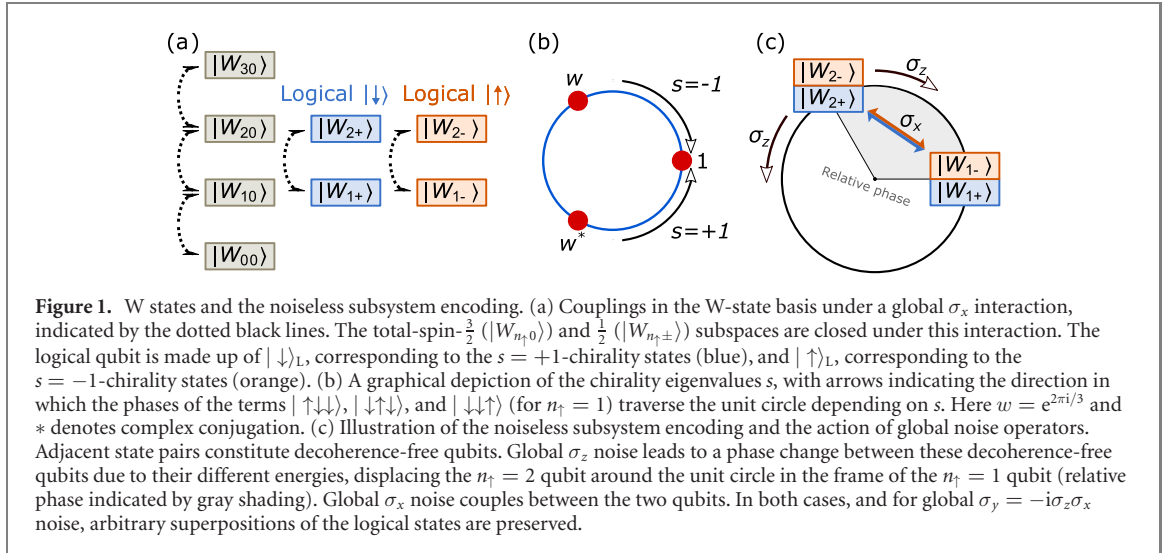
which can be factored into the gauge and qubit degrees of freedom as:

$$\alpha_L|\downarrow\rangle_L + \beta_L|\uparrow\rangle_L \rightarrow (G_1|n_\uparrow = 1\rangle + G_2|n_\uparrow = 2\rangle) \otimes (\alpha_L|s = +1\rangle + \beta_L|s = -1\rangle). \quad (12)$$

In the above, $|\alpha_L|^2 + |\beta_L|^2 = 1$ and $|G_1|^2 + |G_2|^2 = 1$. Global noise only changes the coefficients G_i . The separability represented in equation (12) is crucial. For example, $(|\downarrow\rangle_L + |\uparrow\rangle_L)/\sqrt{2}$ is not a valid encoding of $(|\downarrow\rangle_L + |\uparrow\rangle_L)/\sqrt{2}$ because the qubit and gauge degrees of freedom are entangled, and the logical state is not preserved under global σ_z noise. The encoded logical qubit can be manipulated using suitable operators [29], with the chirality operator χ (equation (1)) acting as the single-qubit logical Z operator. This NSS encoding is illustrated schematically in figure 1.

Implicit in equation (10) is a choice about the orientation of each qubit’s Bloch sphere. The NSS can protect against either spatially uniform noise or a noise field that has spatially uniform intensity but a phase that varies with position, such as a laser field. In this case the orientation of each qubit’s Bloch sphere can be fixed by stipulating the direction of the noise field at a given time so that equation (10) applies.

The NSS encoding improves upon the DFS encoding by adding immunity to collective σ_x and σ_y rotations to the DFS’s immunity to collective dephasing. It is useful to consider under what circumstances this might be particularly beneficial. One application that would use the full power of the NSS’s robustness against general uniform noise would be to preserve quantum information during application of global



control pulses (the ‘noise’ field) that affect all physical qubits in a quantum register, including those making up the logical NSS qubits. For example, the NSS encoding can protect against over- and under-rotation errors in global pulses used for dynamical decoupling [42], which themselves can reduce decoherence due to uniform or spatially varying fluctuations in, e.g. a quantization field. In this way, the NSS encoding can be used as part of a strategy for mitigating nonuniform noise. The NSS encoding may also be used to realize a reference-frame-free qubit for quantum communication [43].

2. Experimental ingredients

The dissipative protocols we present make use of standard techniques for ion-trap experiments: resolved sideband transitions, effective spin–spin couplings driven by MS-type interactions, and sympathetic cooling with an auxiliary species. In this section we introduce an example system that could be used, a linear crystal of beryllium-9 ($^9\text{Be}^+$) and magnesium-25 ($^{25}\text{Mg}^+$) ions, and then review the relevant features of these techniques.

2.1. Mixed-species ion crystal

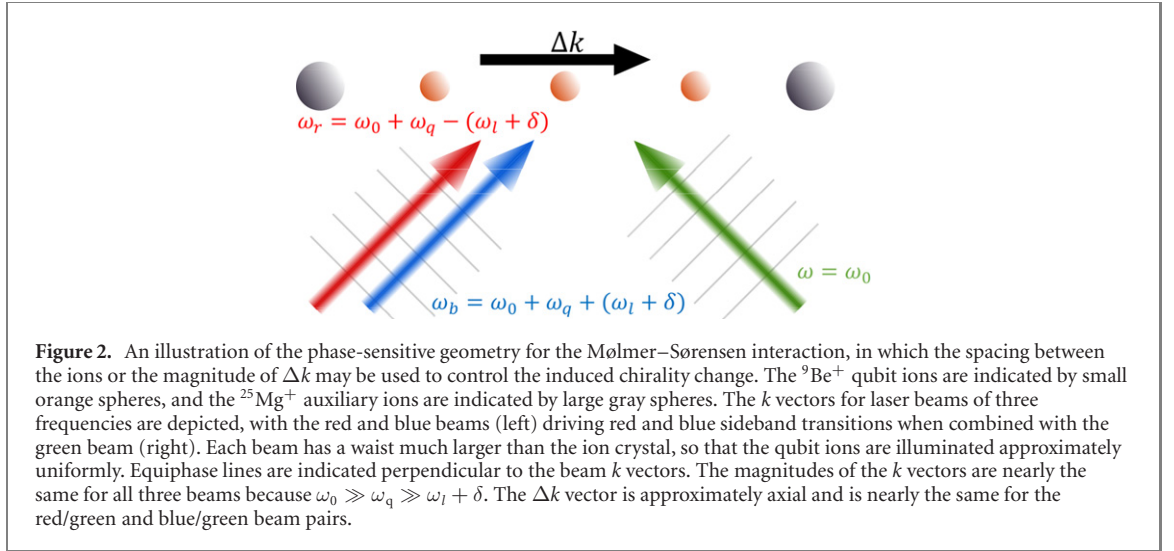
We consider a system in which three qubit ions and two auxiliary ions in the same trapping potential form a one-dimensional Coulomb crystal along the axis of weakest confinement [44]. The Coulomb interaction leads to 15 quantized normal modes of collective motion. We assume that the ten modes of radial motion along the dimensions of tighter confinement factor from the five axial modes, and that the interactions we discuss below couple only to the axial degrees of freedom. For concreteness, we discuss an ion crystal of the form $^{25}\text{Mg}^+ - ^9\text{Be}^+ - ^9\text{Be}^+ - ^9\text{Be}^+ - ^{25}\text{Mg}^+$ (‘MBBBM’), but these schemes could be realized with a variety of qubit and auxiliary species. Confining this crystal in a parity-symmetric axial potential provides symmetric crystal geometry and axial mode eigenvectors that are either symmetric or antisymmetric under parity [45]. The l^{th} eigenvector can be described by entries $z_j^{(l)}$ that specify the participation of ion j in the mode’s zero-point motion in units of length.

2.2. Resolved sideband transitions

When ions of different species with optical transitions of different wavelengths are used for qubits and for cooling, it is typically possible to address the two species independently. Here we consider driving stimulated-Raman sideband transitions in a ground-state hyperfine qubit such as $^9\text{Be}^+$. In the Lamb–Dicke approximation and the interaction picture for the qubit and the vibrational levels, and neglecting rapidly oscillating terms, a blue-sideband interaction resonant with mode l of the ion crystal driven by addressing only the qubit ions can be described by the Hamiltonian [46, 47]:

$$H_{\text{sb}} = a^\dagger \sum_{j=j_{\text{qubit}}} (\pm)_j \frac{|\Omega_j|}{2} \sigma_+^{(j)} e^{i(\Delta k \cdot Z_{0j} + \Delta\phi)} + \text{H.c.}, \quad (13)$$

and for a red-sideband interaction the creation operator $a^\dagger$ for mode l is replaced by the annihilation operator a . The notation $j = j_{\text{qubit}}$ indicates that the sum runs only over the qubit ions. The qubit raising



operator is $\sigma_+^{(j)} = |\uparrow\rangle_j \langle \downarrow|_j$, and $|\Omega_j|$ is the magnitude of the effective Rabi frequency of the interaction for the j^{th} ion. For a ground-state hyperfine qubit, the Rabi frequency Ω_j is proportional to the electric field strengths of the two laser beams driving a stimulated Raman transition and to $\eta_j^{(l)} = \Delta k \cdot |z_j^{(l)}|$, the Lamb-Dicke parameter quantifying the j^{th} ion's participation in the interaction. Here Δk is the axial projection of the difference wavevector between the Raman beams. In addition to a global phase coming from the phase difference $\Delta\phi$ between the laser beams, there is a j -dependent phase arising from the ions' relative positions in the interference pattern between the beams. This phase is represented by $e^{i\phi_j} = (\pm)_j e^{i\Delta k \cdot Z_{0,j}}$, where $Z_{0,j}$ is the equilibrium axial coordinate of ion j and $(\pm)_j = \text{Sign}(z_j^{(l)})$. We can control these phases by adjusting the ions' relative positions along the trap axis or by adjusting the magnitude of Δk . For particular values of these phases, the sideband interaction can induce definite changes Δs in the chirality eigenvalue s . We associate Δs with the coupling:

$$|W_{00}\rangle \leftrightarrow |W_{1(\Delta s)}\rangle \leftrightarrow |W_{2(-\Delta s)}\rangle \leftrightarrow |W_{30}\rangle, \quad (14)$$

and then, for example, $\Delta s = +1$ for $\phi_j = 2\pi/3 \cdot j$ for the MBBBM crystal.

Any non-uniformity of $|\Omega_j|$ across the qubit ions will introduce errors in the schemes we describe below. Important sources of non-uniformity are different participation amplitudes $|z_j^{(l)}|$ in the mode that is driven by the sideband interaction and laser beam intensity differences across the ions. The latter effect can be mitigated by illuminating the ions with beams that have approximately uniform intensity over the extent of the ion crystal, and we address the issue of mode participation in section 2.4.

The Hamiltonian discussed above is also applicable to optical qubits with minor modifications: sideband transitions on optical qubits are driven by a single beam whose axial wavenumber k replaces Δk throughout, and whose phase ϕ replaces the phase difference $\Delta\phi$. With these changes, equation (13) and the discussion of the MS interaction in the following section apply to optical qubits, and our protocols may be used to entangle these qubits as well.

2.3. Mølmer-Sørensen effective spin-spin interaction

In the MS interaction [48–52], an effective spin-spin coupling is driven by red and blue sidebands with opposite detunings $\pm\delta$ from resonance with mode l . Assuming that the Rabi frequency magnitudes $|\Omega_j|$ are uniformly equal to Ω for the qubit ions (and zero for the auxiliary ions), the near-resonant terms in the Hamiltonian are:

$$H_{\text{MS}} = \frac{\Omega}{2} \sum_{j=\text{qubit}} (\pm)_j \sigma_+^{(j)} \left[a^\dagger e^{i(\Delta\phi_b + \Delta k_b \cdot Z_{0,j} + \delta t)} + a e^{i(\Delta\phi_r + \Delta k_r \cdot Z_{0,j} - \delta t)} \right] + \text{H.c.} \quad (15)$$

Here Δk_b , Δk_r , $\Delta\phi_b$, and $\Delta\phi_r$ (or k_b , k_r , ϕ_b , and ϕ_r for an optical qubit) represent the corresponding quantities from equation (13) that describe the beams that drive each sideband. For interaction times $T = 2\pi m/\delta$ for integer m , the MS interaction realizes the propagator for a spin-spin Hamiltonian on the qubit ions, returning the motion to its initial state.

The exact propagator that is obtained depends both on the ratio Ω/δ and on the experimental geometry, as follows. We consider an implementation of the ‘phase sensitive’ geometry for driving stimulated Raman

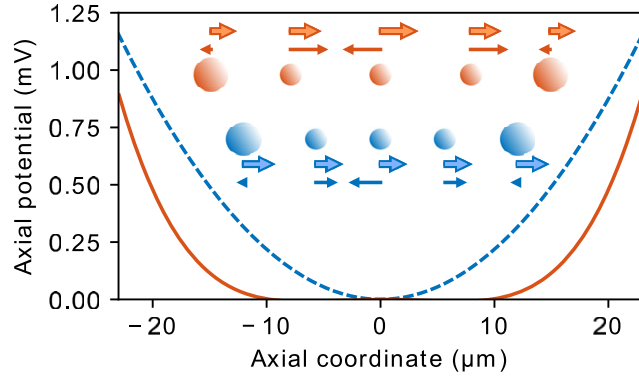


Figure 3. Two possibilities for the static axial trapping potential for the MBBBM crystal described in the text. A harmonic potential is shown in dashed blue, and a potential with significant but practical anharmonicity is shown in solid orange. The confinement is chosen so that the frequency of in-phase oscillation is 0.81 MHz for both crystals. The ion crystals at equilibrium for the anharmonic (top, orange) and harmonic (bottom, blue) potentials are shown, with the beryllium qubit ions corresponding to the smaller spheres, along with relative participation vectors for the lowest-frequency (in-phase) mode of motion (thicker, light-filled arrows) and the highest-frequency mode of motion (dark arrows) for each crystal. The zero-point amplitudes of the in-phase mode in the harmonic potential are $z_{\text{harm}}^{(1)} = (9.56, 7.89, 7.62, 7.89, 9.56)$ nm, and for the highest-frequency mode of the anharmonic potential, with 2.56 MHz frequency, they are $z_{\text{anhar}}^{(5)} = (-1.38, 8.35, -8.35, 8.35, -1.38)$ nm. Because the latter has reduced center-of-mass motion, it is expected to have a heating rate due to excitation by global fields that is 5.4% that of hypothetical in-phase motion at the same frequency (see appendix A).

transitions (see reference [47], the dependence of the propagator on the spacing between the ions would be removed for the ‘phase insensitive’ geometry), which we illustrate in figure 2. In this configuration, copropagating beams at frequencies $\omega_o + \omega_q \pm (\omega_l + \delta)$ intersect on the ions with an orthogonal beam of frequency ω_o , such that $\Delta \vec{k}$ is axial for each orthogonal pair. Here ω_o is an optical frequency detuned from a transition between the ground-state qubit manifold and an electronic excited state, ω_q is the qubit frequency, and ω_l is the frequency of the l^{th} motional mode, upon which a and $a^\dagger$ act. The phases and axial wavenumbers for the beam pairs are approximately the same: $\Delta \phi_r \simeq \Delta \phi_b = \Delta \phi$ and $\Delta k_r \simeq \Delta k_b$.

In this configuration, population transfer between basis states is maximized for $\delta = 2\Omega$ and interaction time $T = 2\pi/\delta$. The MS interaction is driven by simultaneous application of two detuned sidebands, and the chiralities of these sidebands determine the chirality of the couplings under the MS gate. For red- and blue-sidebands driving chirality change $-\Delta s$ as described above, the MS interaction couples $|W_{00}\rangle \leftrightarrow |W_{2(\Delta s)}\rangle$. In matrix notation, the action of the gate on the state vector $(|W_{00}\rangle, |W_{1(-\Delta s)}\rangle, |W_{2(\Delta s)}\rangle, |W_{30}\rangle)^T$ is:

$$U = e^{i\pi/8} \begin{pmatrix} \frac{1}{2} & 0 & -\frac{\sqrt{3}}{2}u^* & 0 \\ 0 & -\frac{1}{2} & 0 & -\frac{\sqrt{3}}{2}u^* \\ -\frac{\sqrt{3}}{2}u & 0 & -\frac{1}{2} & 0 \\ 0 & -\frac{\sqrt{3}}{2}u & 0 & \frac{1}{2} \end{pmatrix}, \quad (16)$$

where $u = e^{2i\Delta\phi}$. The subspace spanned by the four other basis states is left invariant. For $\Delta s = 0$, the chiral W states are left invariant. For $\Delta s = \pm 1$, two chiral and two achiral W states are left invariant. A maximum of 75% population transfer can be achieved between the pairs of the remaining states separated by $\Delta n_\uparrow = \pm 2$.

2.4. Achieving uniform $|\Omega_j|$ using anharmonic potentials

As discussed in section 2.2 and assumed in section 2.3, it is important that $|\Omega_j|$ is uniform across the qubit ions. The lowest-frequency, in-phase axial mode for the MBBBM crystal has nearly uniform participation. However, this mode is especially susceptible to heating by homogeneous electric field noise [53] because of the in-phase motion and the inverse scaling of the heating rate with frequency [54]. Therefore, it may be better to use one of the higher-frequency modes in which some of the ions oscillate out of phase. Uniform participation of the qubit ions in one of these modes can be achieved by introducing some controlled anharmonicity to the trapping potential, as we illustrate in figure 3. The anharmonicity assumed here can be realized in existing ion traps, e.g. the one described in reference [55] with $\sim 150 \mu\text{m}$ ion–electrode distance. We discuss this issue in greater depth for a variety of ion species and configurations in appendix A.

2.5. Sympathetic cooling

In section 3.2 we use sympathetic [56, 57] resolved-sideband cooling [58–60] as a dissipation mechanism that allows one-way population flow into a target W state. Resolved-sideband cooling on auxiliary ions can be used to ground-state cool any crystal mode in which they participate without disturbing the qubit states, provided that the radiation used to manipulate the auxiliary species is far-detuned from the qubit-species transitions.

3. Dissipative preparation of W states

In this section we present protocols for dissipative preparation of W states. First, we consider attempts to extend the rapid, symmetry-based dissipative singlet-generation scheme presented by Horn *et al* in reference [21] to more ions. This scheme incorporates dissipation through spontaneous decay from an electronic excited state. Direct extension is not promising because the rate at which the target state can be populated decreases exponentially with the number of ions. We then present a more promising approach in which dissipation is provided by sympathetic cooling, and we provide an analysis of likely error sources.

The simulation results we present are obtained by numerical simulation of the quantum-mechanical master equation in Lindblad form:

$$\partial_t \rho = -i[H, \rho] + \mathcal{L}_D \rho, \quad (17)$$

where ρ is the system's density matrix, H is the Hamiltonian (of the form prescribed by equations (13) and (15) for a given interaction), and $\mathcal{L}_D$ is the Lindblad dissipator:

$$\mathcal{L}_D \rho = \sum_k \left[L_k \rho L_k^\dagger - \frac{1}{2} \{ \rho, L_k^\dagger L_k \} \right], \quad (18)$$

where L_k are the jump operators for the system that represent deliberate controlled dissipation and also undesirable dissipation mechanisms such as heating of the crystal. Our simulations are carried out with QuTiP's master equation solver [61].

3.1. A natural extension of a scheme for dissipative singlet preparation

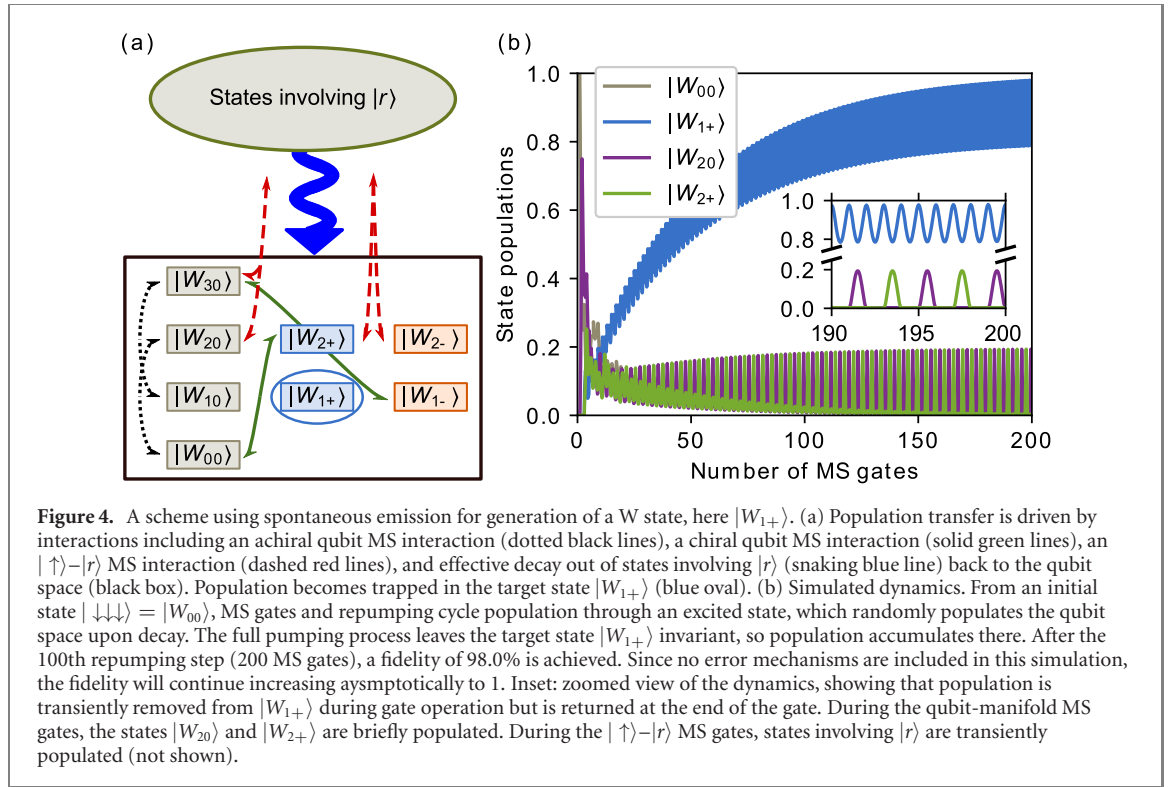
The scheme presented by Horn *et al* [21] generates a singlet state $|S\rangle = (|\uparrow\downarrow\rangle - |\downarrow\uparrow\rangle) / \sqrt{2}$ using engineered decay through an electronic excited state. We briefly review the basic principle: a third stable ground-state level $|r\rangle$ is pumped to an electronic excited state, which then decays back to $|\uparrow\rangle$, $|\downarrow\rangle$, or $|r\rangle$, randomly populating the nine-dimensional two-qutrit Hilbert space. By implementing global carrier and motion-adding ($|n\rangle \rightarrow |n+1\rangle$) interactions that act on the qubits as $|\downarrow\downarrow\rangle \leftrightarrow (|\downarrow\uparrow\rangle + |\uparrow\downarrow\rangle) / \sqrt{2} \leftrightarrow |\uparrow\uparrow\rangle$, and a motion-subtracting coupling from $|\uparrow\rangle$ to $|r\rangle$, population can be cycled through the excited state until it becomes trapped in $|S\rangle \otimes |n=0\rangle$, which is the sole dark state of the full system dynamics. A specific set of levels must be chosen for this scheme so that the excited state can decay only to $|\uparrow\rangle$, $|\downarrow\rangle$, or $|r\rangle$. For the concrete case of the $^9\text{Be}^+ \ ^2S_{1/2}$ manifold, one possible choice is to use $|F=2, m_F=2\rangle = |\downarrow\rangle$, $|F=1, m_F=1\rangle = |\uparrow\rangle$, and $|F=2, m_F=1\rangle = |r\rangle$, with a repumper laser driving the transition to the excited state $|r\rangle \rightarrow |^2P_{1/2}, F=2, m_F=2\rangle$.

Naively applying the same interactions on a three-ion system results in continuous population cycling that leads to a steady-state distribution of population in a mixed state. So far we have not identified a promising three-qubit protocol that uses spontaneous emission as the dissipation mechanism. However, we briefly present one example scheme; this scheme's flaw is that it is prohibitively slow.

We replace both the global qubit-manifold drive and the drive from $|\uparrow\rangle$ to the auxiliary state $|r\rangle$ of the singlet scheme with MS interactions. For the drive to $|r\rangle$, an MS interaction enforces pair-wise transitions and prevents one-up states $|W_{1s}\rangle$ from coupling to states that include $|r\rangle$. Within the qubit manifold, an MS interaction achieves for the three-qubit Hilbert space what a global interaction achieves for the two-qubit Hilbert space: it leaves the target state invariant, but makes connections between other states. Unfortunately, the MS interaction leaves too many states invariant. We can work around this problem by iterating over application of an achiral qubit MS gate, driving population from $|\uparrow\rangle$ to $|r\rangle$ with an MS gate, and repumping, and then applying a *chiral* qubit MS gate and carrying out the repumping process again. The chirality change $\Delta s = \pm 1$ of the chiral MS gate is chosen to leave the desired chiral W state invariant. The alternating chirality of the MS gates can be achieved by periodic adjustment of the spacing of the ion crystal or the magnitude of Δk .

In figure 4, we present an illustration of the concept and a simulation of the dynamics for a chiral qubit MS gate with $\Delta s = +1$. In the simulation, we use jump operators to model the repumping as:

$$L_q^{(j)} = \sqrt{b_q \Gamma} |q\rangle_j \langle r|_j, \quad (19)$$



where $q = \uparrow, \downarrow$, or r indexes over the states, j indexes over the ions, and Γ is the rate of effective decay out of $|r\rangle$. This rate is related to the lifetime of the excited state $|e\rangle$ and to the Rabi frequency and detuning of the $|r\rangle \leftrightarrow |e\rangle$ coupling [62]. It can be determined experimentally through measurement of depletion of $|r\rangle$ by the repumper laser. The parameters $b_{\uparrow} = 5/12$, $b_{\downarrow} = 1/3$, and $b_r = 1/4$ approximately describe the branching of the decay out of $|e\rangle$ [6]. Since the duration of the MS gates is expected to be the speed bottleneck for this scheme, the simulations implement the repumping step for long enough that $|r\rangle$ is effectively fully depleted, and the repumping dynamics are not shown explicitly. As shown in figure 4(b), the dissipative dynamics lead to trapping of population in the state $|W_{1+}\rangle$ that is invariant under these combined dynamics.

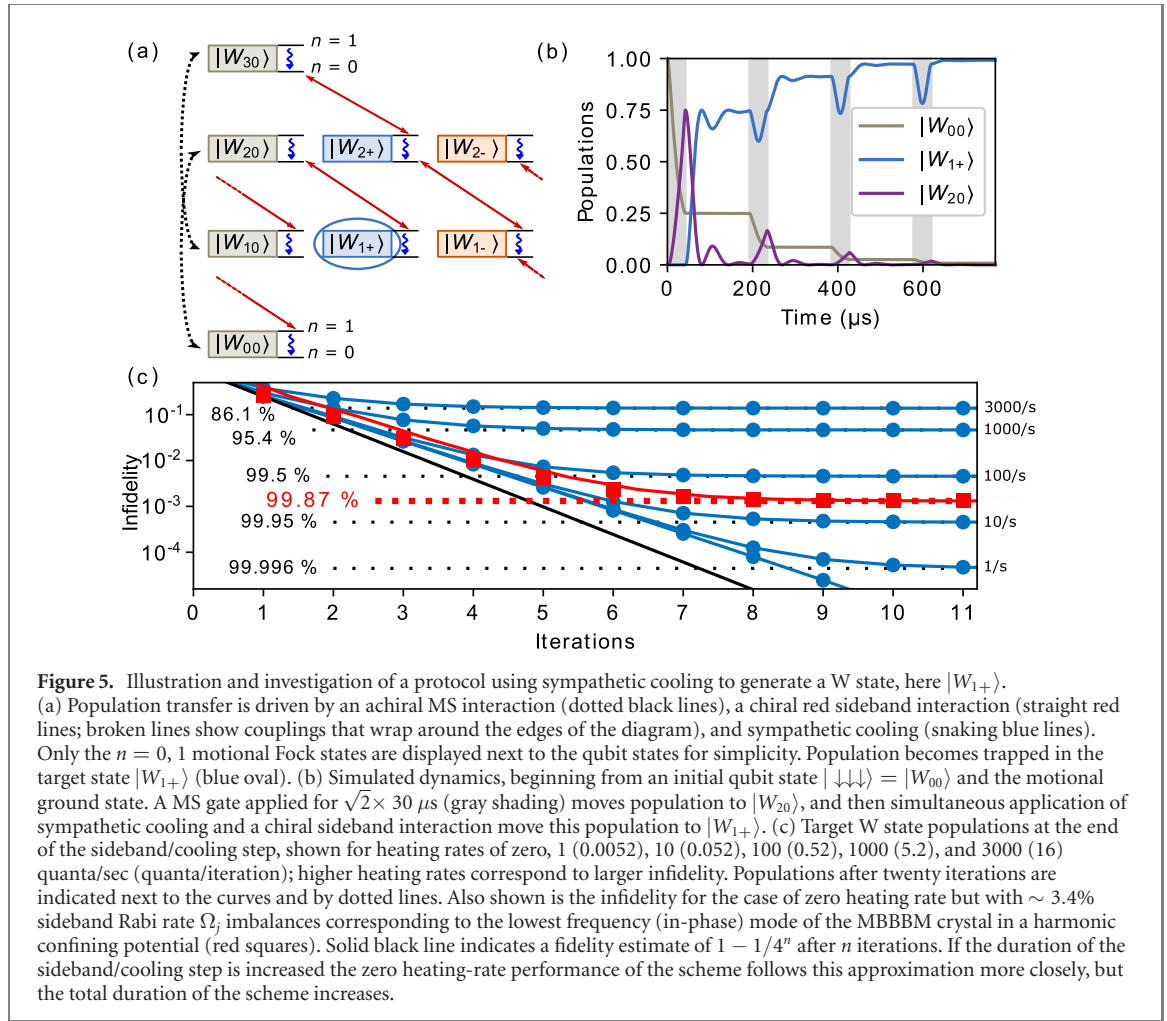
Figure 4 shows that hundreds of MS gates are needed to approach 100% fidelity, even in the absence of errors and imperfections. This illustrates a fundamental limitation of the extension of the singlet scheme to larger numbers of qubits: as the dimension of the qubit space increases exponentially, the rate at which the target state is populated through random decay out of the electronic excited state decreases correspondingly. Therefore, protocols that rely on randomly populating the target state scale inherently poorly.

3.2. Dissipative preparation of chiral W states using sympathetic cooling

Now we present a more promising protocol that incorporates dissipation through sympathetic cooling, which can convert unitary, periodic dynamics to one-way population flow. Unfortunately, it is not possible to generate a W state by simply evolving out of the state $|\downarrow\downarrow\downarrow\rangle$ with a coupling $|\downarrow\rangle \leftrightarrow |\uparrow\rangle$. Although a state $|W_{1s}\rangle$ can be populated under this coupling, the overall state remains separable as population flows to states with increasing $n_{\uparrow}$. This flow can be truncated by preparing a Fock state $|n = 1\rangle$ and then operating on a red sideband $|\downarrow\downarrow\downarrow, n = 1\rangle \leftrightarrow |W_{1s}, n = 0\rangle$ such that a W state is created under unitary evolution [24], but implementing this strategy is not straightforward without individual addressing (or targeted subset addressing [25]). Promising conceptually distinct unitary approaches exist, such as the composite-pulse approach proposed in reference [26], and the best choice of strategy for W state generation will likely depend on system-dependent considerations and capabilities.

The protocol we present gets around the challenges discussed above by iterating over two steps: first, some fraction of the population less than 100% is transferred to the target W state through operations that also induce a motional transition $n \rightarrow n + 1$ on a collective mode. Second, sympathetic cooling removes phonons from this mode. As this mode approaches its ground state, these steps can be repeated without removing population that has already reached the target state, so that population can accumulate there over several iterations.

From an initial state $|W_{00}\rangle \otimes |n = 0\rangle$, the basic steps proceed as follows: an achiral MS gate transfers 75% of the population to $|W_{20}\rangle$. A $\Delta s = \mp 1$ red sideband interaction then moves this population to the



state $|W_{1\pm}\rangle \otimes |n = 1\rangle$. When the ion crystal is sympathetically cooled, the acquired phonon is removed. In practice, the cooling and sideband interaction can be applied simultaneously. Repeated iteration of these steps pumps population to the target state $|W_{1\pm}\rangle \otimes |n = 0\rangle$, from which it cannot escape: the chiral W states are invariant under the achiral MS gate, the $\Delta s = \mp 1$ sideband does not couple $|W_{00}\rangle \leftrightarrow |W_{1\pm}\rangle$, and $|W_{1\pm}\rangle \otimes |n = 0\rangle$ is prevented from evolving to $|W_{20}\rangle$ under the ∓ 1 -chirality red sideband because it is in the ground state of the motion.

Figure 5 presents an illustration of the concept and a numerical investigation of the dynamics for chirality change $\Delta s = +1$. As can be seen from figure 5(a), there is a path to the target state $|W_{1+}\rangle$ from any of the eight basis states. In fact, the protocol will move all population in the qubit space to the target state, as we verify by simulating the dynamics beginning from each of 64 states in a complete basis for the space of three-qubit density matrices [63]. Moreover, it is straightforward to adjust the interactions to generate any of the six W states defined in equations (3)–(8).

The simulation shown in figure 5(b) iterates between application of an achiral MS pulse and simultaneous application of the chiral sideband and sympathetic cooling. To set the timescale of the dynamics, for concreteness we assume that a sideband transition on both the qubit and the auxiliary⁵ ions can be driven in $30 \mu\text{s}$. Therefore, the time for the MS gate is $\sqrt{2} \times 30 \mu\text{s}$ in the phase-sensitive geometry, as half the optical power in one of the Raman beams, and all the power in the other, contributes to driving each $\pm\delta$ -detuned MS sideband and the sideband Rabi rate scales with the field amplitude [46]. With the same sideband transition rate and two cooling ions, sympathetic cooling can extract phonons at a maximum rate of approximately $\kappa = 2/30 \mu\text{s}$. The jump operator that implements sympathetic cooling at this rate is $L = \sqrt{\kappa}a$. Interactions of alternating chirality can be implemented by periodically changing the spacing of the ion crystal or the magnitude of Δk .

⁵ For the specific case of the MBBM crystal described in section 2.1, the significantly smaller relative participation of the magnesium ions in the driven mode can be compensated for by using Raman beams with smaller detuning for sympathetic cooling, because a higher photon scattering rate can be tolerated on the auxiliary species.

Table 1. W-state fidelity after twenty iterations as a function of phenomenological heating rate Γ applied to model a given GSC limit $\bar{n}_{\text{GSC}} = \Gamma/\kappa$, where $\kappa = 2/(30 \mu\text{s})$ is the cooling rate. This heating is applied only during the sympathetic cooling/sideband step.

$\bar{n}_{\text{GSC}}$	Phenomenological heating rate Γ (s^{-1})	Fidelity (%)
0.001	67	99.8
0.003	200	99.4
0.01	667	98.0
0.03	2000	94.4
0.1	6667	84.0
0.3	20 000	65.2

If the population transfer from $|W_{20}\rangle$ to $|W_{1+}\rangle$ is complete, the fidelity after n iterations is $1 - \frac{1}{4^n}$. In practice, this depends on the details of the sideband/cooling step. Our simulations apply the sideband/cooling interaction for $150 \mu\text{s}$, which achieves 99% one-way population transfer from $|W_{20}\rangle \otimes |n=0\rangle$ to $|W_{1+}\rangle \otimes |n=0\rangle$.

3.2.1. Effect of motional occupation on preparation fidelity

This scheme relies on occupation of the motional ground state to prevent population from leaking out of the target state during the simultaneous sideband/cooling step. Therefore, heating of the driven mode will reduce the fidelity, and this is expected to be the limiting error mechanism if the heating is significant. On the other hand, since sympathetic cooling is built in, the final fidelity is asymptotically insensitive to the initial motional state. In figure 5(c) we present numerical calculations of the fidelity for several heating rates. As described previously, this error mechanism can be mitigated by using a mode that has anti-parallel participation vectors for different ions, which reduces its sensitivity to global fields.

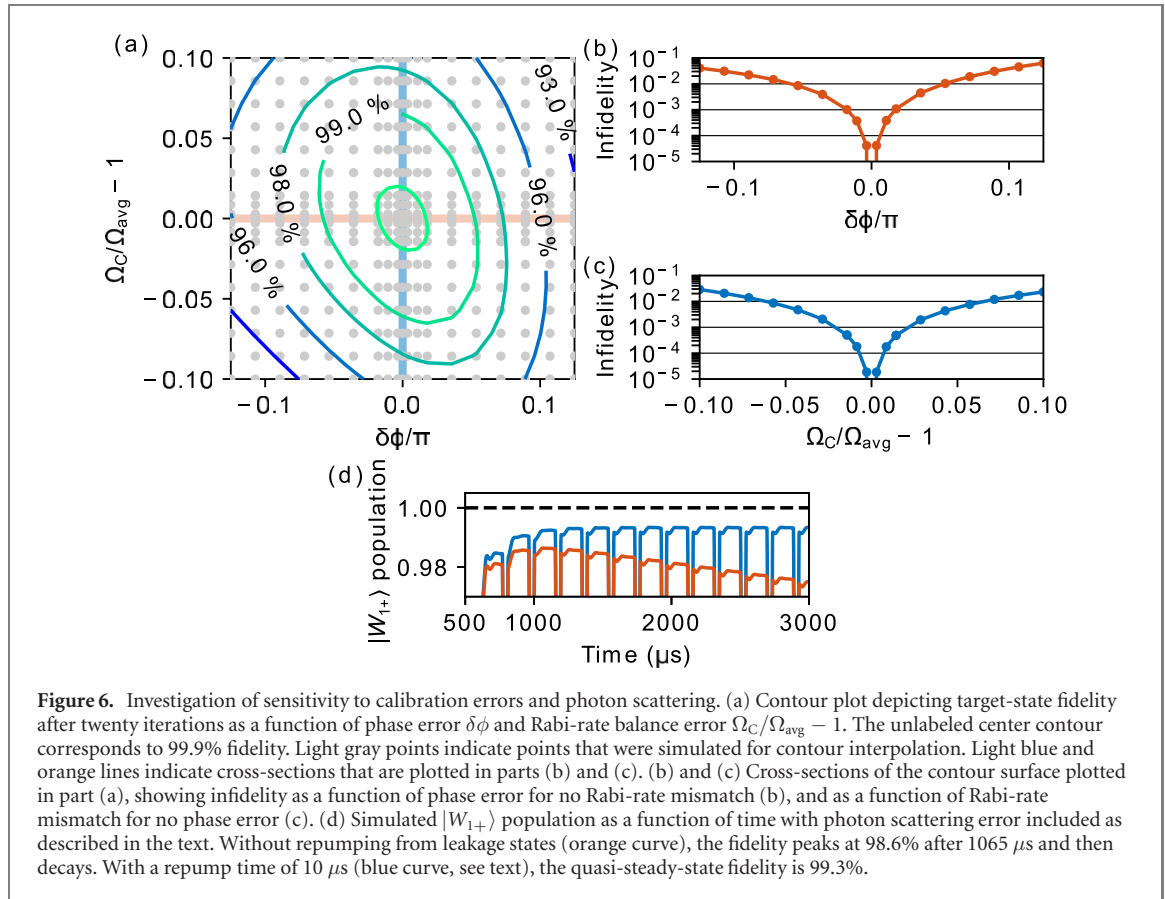
Heating of the motion with rate Γ is implemented in the master equation with two collapse operators $\sqrt{\Gamma}a$ and $\sqrt{\Gamma}a^\dagger$ [46, 51]. When this is combined with cooling at rate κ , implemented with collapse operator $\sqrt{\kappa}a$, the steady-state mean phonon number is $\bar{n} = \Gamma/\kappa$. However, when Γ is determined experimentally by measuring the change in phonon population after a variable probe delay [64], this ratio may predict a limit $\bar{n}_{\text{GSC}}$ for ground-state cooling (GSC) that is far below what is experimentally achieved. This is because the sideband cooling process itself includes sources of heating, such as off-resonant excitation of motion-adding transitions and recoil heating during scattering of repump photons. To address this, in table 1 we present additional calculations of the fidelity in the presence of heating applied only during the sympathetic cooling step to phenomenologically model a given GSC limit. These results show that the scheme benefits from excellent GSC, but does not require unreasonably low $\bar{n}_{\text{GSC}}$.

3.2.2. Sensitivity to calibration errors

We investigate the protocol's sensitivity to calibration errors and present the results in figure 6. We consider failure to properly set the ion spacing, which results in interactions that do not neatly change the chirality according to $\Delta s = 0$ or ± 1 , and failure to achieve uniformity of the sideband Rabi frequencies $|\Omega_j|$ across the ions. For simplicity we focus on the case of symmetrical errors. This covers several important failure mechanisms, for example imperfect adjustment of the depth of the potential well or of the strength of an even-order anharmonic term. We parameterize non-uniformity of the Rabi frequencies by the ratio $\Omega_C/\Omega_{\text{avg}}$ between the Rabi rate on the center qubit ion and the Rabi rate averaged over the qubit ions, with the rates for the left and right ions being the same. We parameterize ion-spacing errors by the deviation $\delta\phi$ of the phase difference between the ions of the Raman interaction. For example, for the $\Delta s = -1$ sideband, the phases across the ions on the $a^\dagger\sigma_-$ term in the Hamiltonian, which is relevant for generation of $|W_{1+}\rangle$, are $-2\pi/3 - \delta\phi$, 0 , and $2\pi/3 + \delta\phi$. Since the MS gate can be made insensitive to this phase error by using the phase-insensitive geometry [47], and because in the phase-sensitive geometry the MS gate requires a different ion spacing than for the sideband-transition step, we simulate the phase error only on the sideband-transition step. As can be seen in figure 6, the fidelity is sensitive to these errors, but not prohibitively so.

3.2.3. Choice of qubit states

Because this scheme does not use closed population cycling through an electronic excited state, there is more freedom in the choice for the pair of states that constitutes the qubit than for schemes that rely on spontaneous emission. A natural choice is a pair of states connected by a first-order magnetic-field-insensitive transition. This choice of a 'clock' transition has benefits such as reduced sensitivity to fluctuating magnetic field gradients, which otherwise reduce the scheme's fidelity.



As a concrete example, we envision using a qubit consisting of $|F = 2, m_F = 0\rangle = |\downarrow\rangle$ and $|F = 1, m_F = 1\rangle = |\uparrow\rangle$ within the ${}^9\text{Be}^+ {}^2S_{1/2}$ manifold, which is first-order field-insensitive at an applied magnetic field of 11.96 mT [34].

3.2.4. Infidelity due to spontaneous photon scattering

Spontaneous photon emission after excitation of a qubit ion to an electronic excited state is an important source of error in trapped-ion systems that use stimulated Raman transitions [65]. To numerically investigate this effect, we simulate the case of ${}^9\text{Be}^+$ qubit ions addressed by laser beams red-detuned by 1 THz from the ${}^2S_{1/2} \leftrightarrow {}^2P_{1/2}$ transition, with beam power such that first-sideband π times are 30 μs , as above. We assume an experimental geometry in which one Raman beam is aligned with the quantizing magnetic field at 45° to the trap axis, and the other lies perpendicular to the first beam in the plane defined by the magnetic field and the trap axis, so that the wavevector difference between the two beams is axial. We choose $\hat{\sigma}_-$ polarization and lower frequency ω_- for the beam parallel to the magnetic field and $\hat{\pi}$ polarization and higher frequency $\omega_\pi > \omega_-$ for the beam perpendicular to it, so that $\omega_\pi - \omega_- = (E_\uparrow - E_\downarrow)/\hbar$ and resonant stimulated Raman transitions are driven between the $|\uparrow\rangle$ and $|\downarrow\rangle$ qubit states with energies $E_\uparrow$ and $E_\downarrow$, respectively. These polarization and frequency choices maximize the Raman Rabi rate and minimize the total scattering rate out of the $|\uparrow\rangle$ and $|\downarrow\rangle$ states at a given detuning from the $S_{1/2} \leftrightarrow P_{1/2}$ transition.

The model includes Raman scattering that leads to population loss from the qubit space by populating ground-state levels other than $|\uparrow\rangle$ and $|\downarrow\rangle$, as well as Raman scattering that leads to transitions between $|\uparrow\rangle$ and $|\downarrow\rangle$. Independently, these two errors lead to decay of the fidelity over time and to a reduction in the steady-state fidelity, respectively, since the latter can be corrected by the dissipative dynamics. The model also includes Rayleigh scattering, which induces decoherence between $|\uparrow\rangle$ and $|\downarrow\rangle$ and reduces the steady-state fidelity to the extent that the scattering amplitudes differ for the two qubit states. Finally, we include the recoil associated with these scattering events, whereby they induce heating of the mode used in the scheme. We use a Lamb–Dicke parameter of $\eta = 0.24$ for these calculations, corresponding to the highest-frequency mode in the anharmonic potential shown in figure 3, which has uniform participation of the qubit ions. The calculations are conducted for an applied quantization field of 11.96 mT, as discussed above.

The results of the simulation are shown in figure 6(d). In this simulation a maximum fidelity of 98.6% is reached after 1065 μs of operation, and then the fidelity decays due to loss of population from the qubit space. Since this scheme uses iterated dissipative dynamics to populate the target state, population leakage can be corrected by using repumper lasers that drive population back to e.g. $|\downarrow\rangle$. In the limit of instantaneous repumping, population leakage error is then converted into decay from $|\uparrow\rangle$ to $|\downarrow\rangle$ and dephasing of the $|\downarrow\rangle$ state. For a reasonable repumping time of $\tau_{\text{rep}} = 10 \mu\text{s}$ from each leakage state (where the corresponding collapse operator is scaled by $\sqrt{1/\tau_{\text{rep}}}$), we find a quasi-steady-state fidelity of 99.3%. The model includes heating due to recoil from scattering on repump transitions, but it neglects off-resonant driving of other transitions by the repump beams.

Error due to spontaneous photon scattering is qubit-species dependent [65]. In general, independent of this choice, the rate of spontaneous Raman scattering can be decreased while maintaining the rate of the desired stimulated transitions by increasing the laser power and detuning. The Rayleigh scattering rate cannot be reduced by the same approach, but the decoherence effect of Rayleigh scattering is typically quite small. For example, under the conditions discussed above we calculate the rate of the decoherence-inducing differential Rayleigh scattering to be $\sim 2 \times 10^{-4}$ times the total leakage rate. On the other hand, the total Rayleigh scattering rate is ~ 174 times the total leakage rate, but this effect leads only to heating through recoil, not to state changes or decoherence. Finally, it would be possible to entirely eliminate photon scattering error if all sideband transitions were driven by a different mechanism, such as a combination of microwave radiation and magnetic-field gradients [66–69], but this would come with its own technical challenges.

3.2.5. Other error mechanisms

We have discussed these error mechanisms to present an overview of what must be considered for successful implementation of this protocol. Ultimately, the errors that contribute to an experimental demonstration will be system-dependent, and factors that we have not quantitatively analyzed above will also contribute. For example, off-resonant coupling of the driving fields to other motional modes may be an important error mechanism, depending on the relationship between the Rabi frequency of the sideband transitions and the frequency separation between the modes. For the MBBBM crystal considered as our example system, the separation between the highest frequency mode and the nearest neighbor is ~ 220 kHz, and the separation from the closest second-order sideband is ~ 194 kHz. These transitions are far-detuned relative to the 17 kHz Rabi frequency assumed for the desired sideband transitions. To further reduce the effect of off-resonant coupling, the strength of the axial confinement may be increased, which increases these frequency spacings and also further suppresses the second-sideband transitions by decreasing the Lamb–Dicke parameter, or the Rabi frequencies of the sideband transitions may be reduced.

Another possible error source is thermal populations of the other modes of the ion crystal. A primary effect is fluctuations of the Rabi rates of the desired interactions due to Debye–Waller factors coming from $\bar{n}_l > 0$ on spectator modes l [46], which can affect the fidelity of the MS gates. This effect can be mitigated by fully re-cooling each axial mode of the crystal before each MS gate, and this also presents an interesting opportunity for system-dependent optimization: for example, if the heating and final temperature of the lowest frequency (in-phase) mode of axial motion during the MS gate is the limiting error mechanism, each full MS gate may be broken into a series of smaller phase-space loops, in between which this mode may be ground-state cooled.

4. Discussion

Using the couplings available in trapped-ion systems, we have designed dissipative protocols for the preparation of W states of three ions. The most promising protocol combines unitary couplings in the form of sideband interactions and MS gates with dissipation through sympathetic cooling, and it avoids the need for addressing of individual ions. The resulting effective dissipation is engineered through cooling of motional excitations, i.e. of an external degree of freedom. Despite being dissipative, this interaction allows correlations between internal degrees of freedom, the qubits, to survive. Damping of harmonic oscillator auxiliary subsystems [6, 12, 19] or coupled systems [5] has previously allowed for the implementation of dissipative schemes, and this approach holds promise for future protocols extending the range of dissipative quantum computation. Moreover, the application of the MS gate for dissipative preparation of W states shows again that time-dependent interactions can be useful to aid the generation of quasi-steady states, as previous work has demonstrated [5, 6].

The unitary entangling gates that these approaches use can themselves generate maximally entangled states, either alone or when combined with global Pauli rotations, depending on the number of ions participating in the interaction [49]. However, the entangled states that are directly accessible this way are

generalized Greenberger–Horne–Zeiling states of the form $|\text{GHZ}\rangle_N = (|\downarrow_1 \dots \downarrow_N\rangle + |\uparrow_1 \dots \uparrow_N\rangle) / \sqrt{N}$, which fall into a fundamentally different class of entangled states from the W states that we have considered: using local operations and classical communication, it is impossible to obtain a three-qubit W state from a $|\text{GHZ}\rangle_3$ state, even with a low chance of success [23]. Therefore, one way to conceptualize the dissipative protocols presented here is that they facilitate conversion between entanglement classes.

There is at least one challenge that must be addressed to allow scaling of these protocols. In order to pump population into a target state, engineered dissipative dynamics must either decouple from the target state entirely or periodically remove population from the target state and then return it, as is the case for the MS gates used in our schemes. Naive extension of the schemes we have presented fails at this point: in general, for odd $N > 3$, achiral MS gates do not leave chiral W states invariant. For N qubits, $\sigma_x^{(N-1)}$ interactions could prepare $|W_{(n-1)0}\rangle$ from $|W_{00}\rangle$ while leaving the target state invariant, thereby playing the same role as the MS gate in the protocol presented in section 3.2. However, it is not clear how to achieve an $(N - 1)$ -ion interaction in practice. There may be creative solutions to this problem, and investigation could be an avenue for future work.

The protocols we have presented are part of a new generation of strategies for dissipative entanglement preparation that eschew timescale hierarchies in favor of symmetry-based dissipation engineering. As a consequence, the fidelities predicted by numerical simulations are comparable to those achievable with unitary schemes [27]. Future efforts may explore these concepts beyond state preparation, such as in new protocols for autonomous quantum error correction and quantum sensing [19], as well as for quantum simulation [20, 70].

Acknowledgments

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Data availability statement

The data that support the findings of this study are available upon reasonable request from the authors.

Appendix A. Achieving uniform Rabi rates using anharmonicity

As discussed in the main text, the fidelity achieved by the protocols we present is maximized when the Raman beam intensity is equal on the qubit ions and the qubit ions have uniform participation in the driven mode. Ideally, this mode also has only weak coupling to electric field noise that is homogeneous over the length of the ion chain (in order to reduce heating) and is separated in frequency from other modes to avoid off-resonant coupling. For a linear ion crystal these requirements can be satisfied by adjusting the axial trapping potential, which affects the ion equilibrium positions and mode vectors. Specifically, we consider electrostatic trapping voltages of the form:

$$V(x) = \frac{1}{2}k_2x^2 + \frac{1}{4}k_4x^4, \quad (\text{A1})$$

which represents the simplest symmetrical deviation from a harmonic potential. By adjusting the terms k_2 and k_4 , modes with the desired relative ion participation can be engineered.

With three qubit ions, the smallest symmetrical mixed-species crystals have two auxiliary ions and always have a qubit ion in the center. Modes with even parity (e.g. $(\leftarrow, \leftarrow, 0, \rightarrow, \rightarrow)$, schematically) do not couple to uniform electric fields and consequently have very low heating rates, but they also have no participation from the center ion. In a crystal with ions of non-uniform masses, the odd-parity modes typically have some center-of-mass motion and therefore couple to homogeneous electric fields. The heating rate $\dot{n}_l$ of a given mode l with frequency ω_l due to excitation by a uniform external field at the same

Table 2. Sample ion chain configurations for various species combinations along with k_2 parameter that achieves uniform coupling amplitudes for the qubit ion (bold). All k_4 parameters used are of amplitude $0.0156 \mu\text{V } \mu\text{m}^{-4}$ (an ‘experimental maximum’ that we enforce, although it is very likely that many systems can achieve higher amplitudes), with the exponent $(-)$ on the corresponding k_2 parameter when the value of k_4 used was negative. $C^{(l)}/C_0$ quantifies the coupling to spatially uniform electric field noise compared to the coupling for the lowest-frequency mode of the same crystal in a harmonic potential. Unless otherwise stated, crystals in configurations listed above have a >130 kHz separation between nearest frequency modes, minimizing the effects of off-resonant coupling.

Ion chain	Normal-mode amplitudes $z_j^{(l)}$ (nm)	Freq. (MHz)	k_2 ($\mu\text{V } \mu\text{m}^{-2}$)	$C^{(l)}/C_0$
Mg–Be–Be–Be–Mg ^a	–1.38, 8.35 , – 8.35 , 8.35 , –1.38	2.56	–0.729	0.0542
Be–Mg–Mg–Mg–Be	–4.17, 2.66 , – 2.66 , 2.66 , –4.17	6.00	25.0 ⁽⁻⁾	0.153
Ca–Be–Be–Be–Ca ^a	–0.823, 8.58 , – 8.58 , 8.58 , –0.823	2.48	–0.894	0.123
Ca–Ca–Ca–Ca–Ca ^b	–2.74, 4.74 , – 4.74 , 4.74 , –2.74	1.42	0.436	0.001
Ca–Sr–Sr–Sr–Ca	–3.53, 2.70 , – 2.70 , 2.70 , –3.53	1.73	6.12	0.093
Ba–Yb–Yb–Yb–Ba ^c	–2.32, 3.13 , – 3.13 , 3.13 , –2.32	0.778	1.05	0.011
Ba–Yb–Yb–Yb–Ba	3.45, 3.51 , 3.51 , 3.51 , 3.45	0.527	19.2 ⁽⁻⁾	1.00

^aCrystals with lighter ions in the center may be more difficult to arrange, depending on the system’s capability to rearrange ions.

^bHere we use ‘Ca’ to mean $^{43}\text{Ca}^+$. All other ‘Ca’ in this table refer to $^{40}\text{Ca}^+$.

^cIn this configuration, the next highest frequency mode is at 0.737 MHz, so the scheme may suffer from significant off-resonant coupling. If the experimental setup is capable of higher k_4 and k_2 values, the frequencies, as well as the frequency difference, can be increased. For this ion chain we list the in-phase mode in the next row, which does not suffer from this issue but has a higher heating rate.

frequency is given by (generalized from reference [57]):

$$\dot{n}_l = \frac{q^2 S_E(\omega_l)}{4\hbar\omega_l} \left(\sum_{j=1}^N \frac{v_j^{(l)}}{\sqrt{m_j}} \right)^2, \quad (\text{A2})$$

$$= \frac{q^2 S_E(\omega_l)}{4\hbar\omega_l} C^{(l)}, \quad (\text{A3})$$

where q is the fundamental charge, $S_E(\omega_l)$ is the power spectral density of spatially uniform electric-field noise, $v^{(l)}$ is the (root-mass-weighted) normalized eigenvector for mode l , and m_j are the ion masses. We define $C^{(l)}$ as a frequency-independent measure of the mode’s coupling to global fields. Typically, a given mode’s heating rate may be decreased by increasing the mode’s frequency.

By considering the axial trapping potential and the Coulomb force, the equilibrium positions of the ions and the normal mode vectors $v^{(l)}$ can be numerically determined [44, 71]. In order to investigate the general practicality of the schemes we described in the main text, we performed this analysis for several mixed-species ion crystal configurations.

We considered four mixed-species ion combinations: $^9\text{Be}^+ / ^{25}\text{Mg}^+$, $^9\text{Be}^+ / ^{40}\text{Ca}^+$, $^{40}\text{Ca}^+ / ^{88}\text{Sr}^+$, and $^{138}\text{Ba}^+ / ^{171}\text{Yb}^+$. We also consider an equal-mass, mixed-state configuration of $^{43}\text{Ca}^+$ where ions of the same species but in a different state act as coolant ions [72]. We explored this range of masses and mass ratios to illustrate the effect mass has on the feasibility of achieving ideal engineered modes for this experiment. Engineering modes with equal ion participation where the frequencies of all five axial modes are well-separated (>100 s of kHz away) is particularly difficult for the $^{138}\text{Ba}^+ / ^{171}\text{Yb}^+$ combination due to the inherently low mode frequencies of these heavier ions. For all the ions we consider, the value of k_2 varies from 30 to $60 \mu\text{V } \mu\text{m}^{-2}$ for typical single-ion harmonic trapping potentials [73–75].

The values of k_4 that can be practically realized depend on geometric factors such as the size of the electrodes and distance from the ion, as well as the maximum electrode voltage that the system can achieve [76]. We used a maximum anharmonicity of $k_4 \approx 0.0156 \mu\text{V } \mu\text{m}^{-4}$ because this should be achievable in traps such as those described in references [55, 71]. In general, ion traps with smaller ion-to-electrode distance and higher voltages should be capable of higher k_4 .

Most of the ion chains we considered are symmetric combinations in the order $m_L m_H m_H m_H m_L$, where m_L (m_H) denotes the lighter (heavier) ion of the two. We also investigated crystals of the form $m_L m_H m_L m_H m_L$, and we found that in general the modes were not suitable when tuned to have equal ion participation in each species: using the highest frequency mode, the difference between the highest two frequencies was consistently less than 4%; the middle frequency mode has low m_H ion participation; and the lowest frequency mode has high heating rate. With the exception of the MBMMB crystal presented in the main text (which is preferred over BMMMB because $^9\text{Be}^+$ has a lower spontaneous emission error than $^{25}\text{Mg}^+$ for a given laser intensity, so it serves as a better qubit), we do not consider crystals of the form $m_H m_L m_L m_L m_H$ when a good set of parameters can be found for other configurations, because these may be more difficult to arrange in practice.

Table 2 summarizes the results of the analysis. We present the normal mode participation vectors and frequencies for optimized values of k_2 and k_4 , along with the mode's coupling to homogeneous external fields as defined in equation (A2). For the participation vector, we list $z_j^{(l)} = \sqrt{\hbar/2m_j\omega_l}v_j^{(l)}$, the ground state wavefunction root-mean square size for ion j in mode l [77]. The Lamb–Dicke parameter $\eta_j^{(l)} = \Delta k \cdot |z_j^{(l)}|$ quantifies ion j 's coupling in mode l to a pair of Raman beams with axial wavenumber difference Δk . The heating rate parameter $C^{(l)}$ is heavily dominated by the ion mass, so we normalize it by the C_0 , the value of $C^{(l)}$ for the in-phase mode of the same crystal in a harmonic potential (which is the closest to a 'true' center-of-mass mode for a mixed-species ion crystal in a harmonic potential, and thus has the highest heating rate).

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References

- [1] Poyatos J F, Cirac J I and Zoller P 1996 Quantum reservoir engineering with laser cooled trapped ions *Phys. Rev. Lett.* **77** 4728–31
- [2] Kraus B, Büchler H P, Diehl S, Kantian A, Micheli A and Zoller P 2008 Preparation of entangled states by quantum Markov processes *Phys. Rev. A* **78** 042307
- [3] Verstraete F, Wolf M M and Ignacio Cirac J 2009 Quantum computation and quantum-state engineering driven by dissipation *Nat. Phys.* **5** 633–6
- [4] Krauter H, Muschik C A, Jensen K, Wasilewski W, Petersen J M, Ignacio Cirac J and Polzik E S 2011 Entanglement generated by dissipation and steady state entanglement of two macroscopic objects *Phys. Rev. Lett.* **107** 080503
- [5] Barreiro J T et al 2011 An open-system quantum simulator with trapped ions *Nature* **470** 486–91
- [6] Lin Y, Gaebler J P, Reiter F, Tan T R, Bowler R, Sørensen A S, Leibfried D and Wineland D J 2013 Dissipative production of a maximally entangled steady state of two quantum bits *Nature* **504** 415–8
- [7] Kienzler D, Lo H-Y, Keitch B, de Clercq L, Leupold F, Lindenfesler F, Marinelli M, Negnevitsky V and Home J P 2015 Quantum harmonic oscillator state synthesis by reservoir engineering *Science* **347** 53–6
- [8] Shankar S et al 2013 Autonomously stabilized entanglement between two superconducting quantum bits *Nature* **504** 419–22
- [9] Kimchi-Schwartz M E, Martin L, Flurin E, Aron C, Kulkarni M, Tureci H E and Siddiqi I 2016 Stabilizing entanglement via symmetry-selective bath engineering in superconducting qubits *Phys. Rev. Lett.* **116** 240503
- [10] Liu Y, Shankar S, Ofek N, Hatridge M, Narla A, Sliwa K M, Frunzio L, Schoelkopf R J and Devoret M H 2016 Comparing and combining measurement-based and driven-dissipative entanglement stabilization *Phys. Rev. X* **6** 011022
- [11] Plenio M B, Huelga S F, Beige A and Knight P L 1999 Cavity-loss-induced generation of entangled atoms *Phys. Rev. A* **59** 2468–75
- [12] Kastoryano M J, Reiter F and Sørensen A S 2011 Dissipative preparation of entanglement in optical cavities *Phys. Rev. Lett.* **106** 090502
- [13] Carr A W and Saffman M 2013 Preparation of entangled and antiferromagnetic states by dissipative Rydberg pumping *Phys. Rev. Lett.* **111** 033607
- [14] Bhaktavatsala Rao D D and Mølmer K 2013 Dark entangled steady states of interacting Rydberg atoms *Phys. Rev. Lett.* **111** 033606
- [15] Morigi G, Eschner J, Cormick C, Lin Y, Leibfried D and Wineland D J 2015 Dissipative quantum control of a spin chain *Phys. Rev. Lett.* **115** 200502
- [16] Reiter F, Reeb D and Sørensen A S 2016 Scalable dissipative preparation of many-body entanglement *Phys. Rev. Lett.* **117** 040501
- [17] Shao X Q, Wu J H, Yi X X and Long G-I 2017 Dissipative preparation of steady Greenberger–Horne–Zeilinger states for Rydberg atoms with quantum Zeno dynamics *Phys. Rev. A* **96** 062315
- [18] Pastawski F, Clemente L and Cirac J I 2011 Quantum memories based on engineered dissipation *Phys. Rev. A* **83** 012304
- [19] Reiter F, Sørensen A S, Zoller P and Muschik C A 2017 Dissipative quantum error correction and application to quantum sensing with trapped ions *Nat. Commun.* **8** 1822
- [20] Raghunandan M, Wolf F, Ospelkaus C, Schmidt P O and Weimer H 2020 Initialization of quantum simulators by sympathetic cooling *Sci. Adv.* **6** eaaw9268
- [21] Horn K P, Reiter F, Lin Y, Leibfried D and Koch C P 2018 Quantum optimal control of the dissipative production of a maximally entangled state *New J. Phys.* **20** 123010
- [22] Doucet E, Reiter F, Ranzani L and Kamal A 2020 High fidelity dissipation engineering using parametric interactions *Phys. Rev. Res.* **2** 023370
- [23] Dür W, Vidal G and Cirac J I 2000 Three qubits can be entangled in two inequivalent ways *Phys. Rev. A* **62** 062314
- [24] Häffner H et al 2005 Scalable multiparticle entanglement of trapped ions *Nature* **438** 643–6
- [25] Linington I E and Vitanov N V 2008 Decoherence-free preparation of Dicke states of trapped ions by collective stimulated Raman adiabatic passage *Phys. Rev. A* **77** 062327
- [26] Ivanov S S, Vitanov N V and Korolkova N V 2013 Creation of arbitrary Dicke and NOON states of trapped-ion qubits by global addressing with composite pulses *New J. Phys.* **15** 023039
- [27] Lin Y et al 2016 Preparation of entangled states through Hilbert space engineering *Phys. Rev. Lett.* **117** 140502
- [28] Ticozzi F and Viola L 2014 Steady-state entanglement by engineered quasi-local Markovian dissipation: Hamiltonian-assisted and conditional stabilization *Quantum Inf. Comput.* **14** 265–94
- [29] Viola L, Knill E and Laflamme R 2001 Constructing qubits in physical systems *J. Phys. A: Math. Gen.* **34** 7067–79
- [30] Zanardi P and Rasetti M 1997 Noiseless quantum codes *Phys. Rev. Lett.* **79** 3306–9
- [31] Lidar D A, Chuang I L and Whaley K B 1998 Decoherence-free subspaces for quantum computation *Phys. Rev. Lett.* **81** 2594–7
- [32] Duan L-M and Guo G-C 1998 Reducing decoherence in quantum-computer memory with all quantum bits coupling to the same environment *Phys. Rev. A* **57** 737–41

- [33] Kielpinski D, Meyer V, Rowe M A, Sackett C A, Itano W M, Monroe C and Wineland D J 2001 A decoherence-free quantum memory using trapped ions *Science* **291** 1013–5
- [34] Langer C *et al* 2005 Long-lived qubit memory using atomic ions *Phys. Rev. Lett.* **95** 060502
- [35] Andrews R W *et al* 2019 Quantifying error and leakage in an encoded Si/SiGe triple-dot qubit *Nat. Nanotechnol.* **14** 747–50
- [36] Gyenis A, Mundada P S, Paolo A D, Hazard T M, You X, Schuster D I, Koch J, Blais A and Houck A A 2021 Experimental realization of a protected superconducting circuit derived from the $0-\pi$ qubit *PRX Quantum* **2** 010339
- [37] Knill E, Laflamme R and Viola L 2000 Theory of quantum error correction for general noise *Phys. Rev. Lett.* **84** 2525–8
- [38] DiVincenzo D P, Bacon D, Kempe J, Burkard G and Whaley K B 2000 Universal quantum computation with the exchange interaction *Nature* **408** 339–42
- [39] Russ M and Burkard G 2017 Three-electron spin qubits *J. Phys.: Condens. Matter* **29** 393001
- [40] Viola L, Fortunato E M, Pravia M A, Knill E, Laflamme R and Cory D G 2001 Experimental realization of noiseless subsystems for quantum information processing *Science* **293** 2059–63
- [41] Fortunato E M, Viola L, Pravia M A, Knill E, Laflamme R, Havel T F and Cory D G 2003 Exploring noiseless subsystems via nuclear magnetic resonance *Phys. Rev. A* **67** 062303
- [42] Viola L, Knill E and Lloyd S 1999 Dynamical decoupling of open quantum systems *Phys. Rev. Lett.* **82** 2417–21
- [43] Bartlett S D, Rudolph T and Spekkens R W 2003 Classical and quantum communication without a shared reference frame *Phys. Rev. Lett.* **91** 027901
- [44] James D F V 1998 Quantum dynamics of cold trapped ions with application to quantum computation *Appl. Phys. B* **66** 181–90
- [45] Morigi G and Walther H 2001 Two-species Coulomb chains for quantum information *Eur. Phys. J. D* **13** 261–9
- [46] Wineland D J, Monroe C, Itano W M, Leibfried D, King B E and Meekhof D M 1998 Experimental issues in coherent quantum-state manipulation of trapped atomic ions *J. Res. Natl Inst. Stand. Technol.* **103** 259–328
- [47] Lee P J, Brickman K-A, Deslauriers L, Haljan P C, Duan L-M and Monroe C 2005 Phase control of trapped ion quantum gates *J. Opt. B: Quantum Semiclass. Opt.* **7** S371–83
- [48] Sørensen A and Mølmer K 1999 Quantum computation with ions in thermal motion *Phys. Rev. Lett.* **82** 1971–4
- [49] Mølmer K and Sørensen A 1999 Multiparticle entanglement of hot trapped ions *Phys. Rev. Lett.* **82** 1835–8
- [50] Solano E, de Matos Filho R L and Zagury N 1999 Deterministic Bell states and measurement of the motional state of two trapped ions *Phys. Rev. A* **59** R2539–43
- [51] Sorensen A and Molmer K 2000 Entanglement and quantum computation with ions in thermal motion *Phys. Rev. A* **62** 022311
- [52] Milburn G J, Schneider S and James D F V 2000 Ion trap quantum computing with warm ions *Fortschr. Phys.* **48** 801–10
- [53] King B E, Wood C S, Myatt C J, Turchette Q A, Leibfried D, Itano W M, Monroe C and Wineland D J 1998 Cooling the collective motion of trapped ions to initialize a quantum register *Phys. Rev. Lett.* **81** 1525–8
- [54] Brownnutt M, Kumph M, Rabl P and Blatt R 2015 Ion-trap measurements of electric-field noise near surfaces *Rev. Mod. Phys.* **87** 1419–82
- [55] Bradford Blakestad R 2010 Transport of trapped-ion qubits within a scalable quantum processor *Doctoral Thesis* University of Colorado Boulder
- [56] Larson D J, Bergquist J C, Bollinger J J, Itano W M and Wineland D J 1986 Sympathetic cooling of trapped ions: a laser-cooled two-species nonneutral ion plasma *Phys. Rev. Lett.* **57** 70–3
- [57] Kielpinski D, King B E, Myatt C J, Sackett C A, Turchette Q A, Itano W M, Monroe C, Wineland D J and Zurek W H 2000 Sympathetic cooling of trapped ions for quantum logic *Phys. Rev. A* **61** 032310
- [58] Eschner J, Morigi G, Schmidt-Kaler F and Blatt R 2003 Laser cooling of trapped ions *J. Opt. Soc. Am. B* **20** 1003–15
- [59] Monroe C, Meekhof D M, King B E, Jefferts S R, Itano W M, Wineland D J and Gould P 1995 Resolved-sideband Raman cooling of a bound atom to the 3D zero-point energy *Phys. Rev. Lett.* **75** 4011–4
- [60] Chen J S, Brewer S M, Chou C W, Wineland D J, Leibbrandt D R and Hume D B 2017 Sympathetic ground state cooling and time-dilation shifts in an $^{27}\text{Al}^+$ optical clock *Phys. Rev. Lett.* **118** 1–5
- [61] Johansson J R, Nation P D and Nori F 2013 QuTiP 2: a Python framework for the dynamics of open quantum systems *Comput. Phys. Commun.* **184** 1234–40
- [62] Reiter F and Sørensen A S 2012 Effective operator formalism for open quantum systems *Phys. Rev. A* **85** 032111
- [63] Chuang I L and Nielsen M A 1997 Prescription for experimental determination of the dynamics of a quantum black box *J. Mod. Opt.* **44** 2455–67
- [64] Turchette Q A *et al* 2000 Heating of trapped ions from the quantum ground state *Phys. Rev. A* **61** 063418
- [65] Ozeri R *et al* 2007 Errors in trapped-ion quantum gates due to spontaneous photon scattering *Phys. Rev. A* **75** 042329
- [66] Mintert F and Wunderlich C 2001 Ion-trap quantum logic using long-wavelength radiation *Phys. Rev. Lett.* **87** 257904
- [67] Johanning M, Braun A, Timoney N, Elman V, Neuhauser W and Wunderlich C 2009 Individual addressing of trapped ions and coupling of motional and spin states using RF radiation *Phys. Rev. Lett.* **102** 073004
- [68] Warring U, Ospelkaus C, Colombe Y, Jördens R, Leibfried D and Wineland D J 2013 Individual-ion addressing with microwave field gradients *Phys. Rev. Lett.* **110** 173002
- [69] Srinivas R, Burd S C, Sutherland R T, Wilson A C, Wineland D J, Leibfried D, Allcock D T C and Slichter D H 2019 Trapped-ion spin-motion coupling with microwaves and a near-motional oscillating magnetic field gradient *Phys. Rev. Lett.* **122** 163201
- [70] Reiter F, Lange F, Jain S, Grau M, Home J P and Lenarčič Z 2019 Engineering generalized Gibbs ensembles with trapped ions (arXiv:1910.01593)
- [71] Brown K R, Ospelkaus C, Colombe Y, Wilson A C, Leibfried D and Wineland D J 2011 Coupled quantized mechanical oscillators *Nature* **471** 196–9
- [72] Moore I, Metzner J, Quinn A, Wineland D and Allcock D 2020 DAMOP presentation meetings.aps.org/Meeting/DAMOP20/Session/D03.1
- [73] Wright J A 2015 Mixed species ion chains for scalable quantum computation *Doctoral Thesis* University of Washington
- [74] Bruzewicz C D, McConnell R, Stuart J, Sage J M and Chiaverini J 2019 Dual-species, multi-qubit logic primitives for Ca^+/Sr^+ trapped-ion crystals *npj Quantum Inf.* **5** 102
- [75] Negnevitsky V 2018 Feedback-stabilised quantum states in a mixed-species ion system *Doctoral Thesis* ETH Zurich
- [76] Home J P and Steane A M 2006 Electrode configurations for fast separation of trapped ions *Quantum Inf. Comput.* **6** 289–325
- [77] Home J P, Hanneke D, Jost J D, Leibfried D and Wineland D J 2011 Normal modes of trapped ions in the presence of anharmonic trap potentials *New J. Phys.* **13** 073026