Multiphoton simulator for non-Abelian anyons

Gavin K Brennen

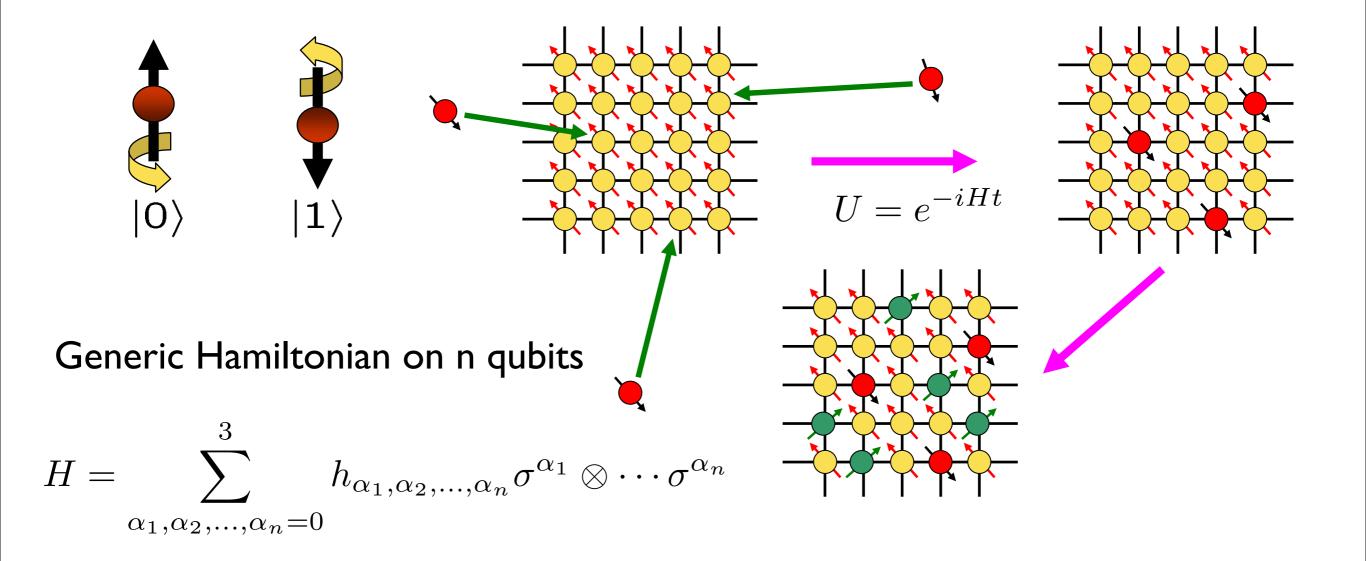


QSciTech, Macquarie University, Sydney

Joint work with: Dominic Berry (IQC, Waterloo) Miguel Aguado (MPQ, Garching) Alexei Gilchrist (Macquarie, Sydney)

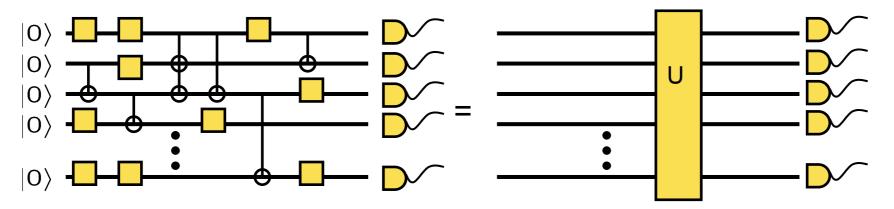
arXiv:0906.4578

- ``We should try to find out what kinds of quantum mechanical systems are mutually intersimulatable, and try to find a specific class, or character of that class which will simulate everything'' Feynman 1982
- Many physical systems have an isomorphic state space, so if we can control the state of one then that simulates another
 - one qubit=polarization space of a photon= spin space of an electron= occupation space of a single photon in two orthogonal modes, etc.

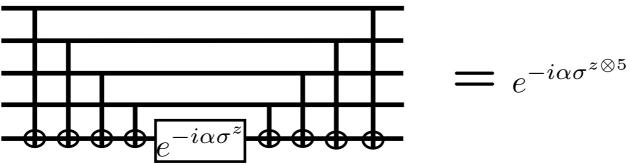


Digital Simulators

• A fault tolerant quantum computer can simulate any dynamics over finite dimensional systems



- Problem: This may not be efficient. General unitaries on n qubits systems require 2^n gates (for qudits d^{2n})
- Some dynamics are efficient

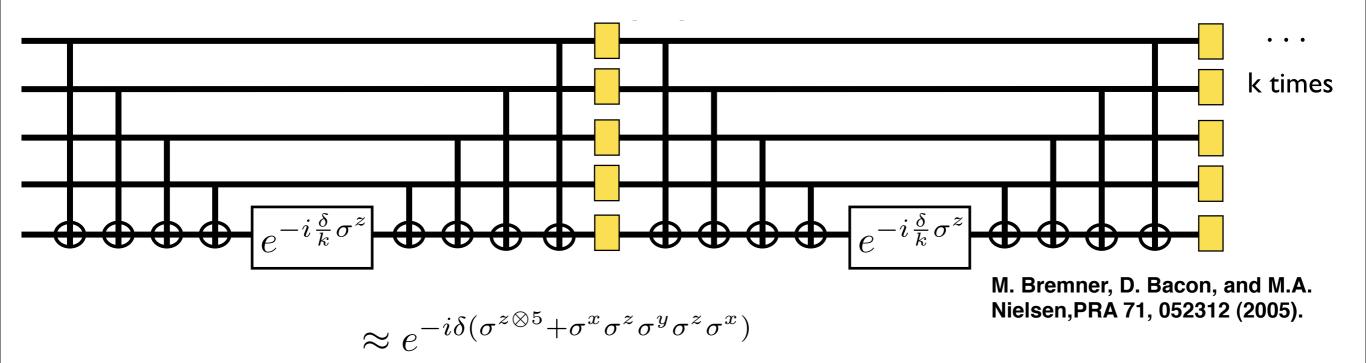


- Is there a constructive approximate method? Yes

Stroboscopic techniques

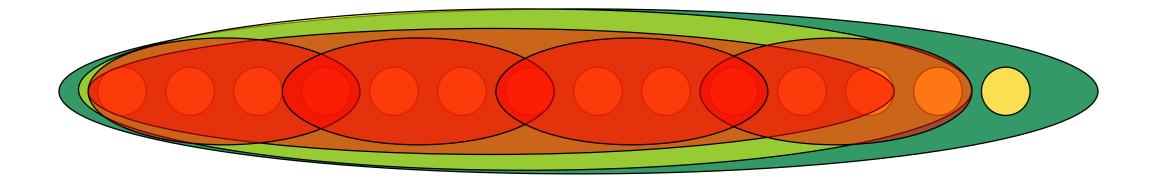
- Control Hamiltonian $H(t) = H_{ent} + \sum f_m(t)H_m$
 - System governed by TDSE $i\hbar \dot{U} = \overset{m}{H}(t)U(t)$
- Using properties of commutators (Trotterization) we can digitize a simulation.
 - Advantage: Digitized circuits can be made fault tolerant

$$e^{-i\Delta H_a}e^{i\Delta H_b}e^{i\Delta H_a}e^{-i\Delta H_b} = e^{-i(i[H_a, H_b])\Delta^2} + O(\Delta^3)$$
$$e^{-i\Delta(H_a + H_b)} = [e^{-i\Delta H_a/k}e^{-i\Delta H_b/k}]^k + O(\Delta^2/k)$$



k-local Hamiltonians

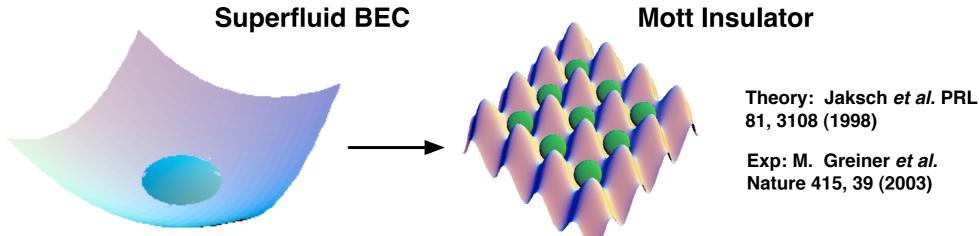
 Bounded Hamiltonians that can be written as a sum of terms involving tensor products of no more than k terms



- It is frequently the case that k is small for physically relevant Hamiltonian (often k=2)
- Complexity of stroboscopic circuit for k-local Hamiltonians is $O(n^k)$

Analogue simulators

- Engineer an "always on" interaction that mimics a physics you want to simulate
- Ex: Superfluid to Mott insulator phase transition



Analogue simulation of Bose-Hubbard interaction

Pros/Cons

- Analogue simulators:
 - Pros: A more direct method, much simpler control
 - Cons: Not fault tolerant, architectural limitations, physical limitations on locality and strength of interactions in the simulating system
- Digital simulators:
 - Pros: Universality, can be made fault tolerant, no architectural constaints in principle
 - Cons: Complicated control pulse sequences, evidence* that fault tolerant simulation has complexity that scales like $1/\epsilon$ for a tolerated error ϵ

*K. Brown, R.J. Clark, I.L. Chuang, PRL 97, 050504 (2006).

A quantum simulation algorithm

- Problem: Compute the energy gap ΔE between the ground and first excited state of a Hamiltonian H
- Algorithm:
 - Map the Hilbert space of the system to be simulated to n qubits
 - Prepare QC in the ground state of a local Hamiltonian H_0
 - e.g. $|\psi(0)\rangle = |+_x\rangle^{\otimes n}$ $H_0 = -\sum \sigma_j^x$
 - Evolve, using stroboscopic circuits, according to the following adiabatic Hamiltonian $\lim \partial H(s) \lim$

$$H(s) = (1-s)H_0 + sH \qquad s = t/T \quad T \gg \frac{||\overline{\Delta s}||}{(\Delta E)^2}$$

PRL 89, 057904 (2002).

- Failure to remain adiabatic results in a final state which has some small admixture of ground and excited states of H $|\psi\rangle_i = c_a |\lambda_a\rangle + c_e |\lambda_e\rangle$
- Evolve by simulated H for time t_i
- Measure some operator that couples ground and excited states $\langle O(t_i) \rangle$
- Repeat for a polynomial number of times steps (polynomial number of steps enough to resolve ΔE). Compute Fourier transform of $\{\langle O(t_i) \rangle\}$

Caveats

- To estimate of the gap with error # of digits precision needed is $\log(1/\epsilon)$
 - A good quantum algorithm would have a complexity $O(poly \log(1/\epsilon))$
- Error in trotter approximation of evolution due to a sum of Hamiltonians scales like $O(t^2/k)$. Higher order Trotter expansion has error $O(t^{m+1}/k^m)$ using gates $O(2^m)$
- Total time to implement algorithm is independent of k. If control gates are perfect, then precision improves with larger k.
- If not, then each subcircuit must be replaced with a fault tolerant version.
- But time to implement $e^{-iHt/k}$ is independent of k (for standard control Hamiltonians)! Hence for a fixed error the total time for the simulation is $O(1/\epsilon)$.

K. Brown, R.J. Clark, I.L. Chuang, PRL 97, 050504 (2006).

- Workarounds:
 - Find better gate libraries (maybe global pulses with correlated errors) combined with noiseless subsystems
 - Hard wire the Hamiltonian into a system so no Trotterization is needed!

What to simulate?

- Hubbard models
 - 2D Fermi-Hubbard Model

$$H = -\sum_{\langle \mathbf{r}, \mathbf{r}' \rangle, \sigma} t_{\mathbf{r}, \mathbf{r}'} \left(c_{\mathbf{r}, \sigma}^{\dagger} c_{\mathbf{r}', \sigma} + H.c. \right) + U \sum_{\mathbf{r}} n_{\mathbf{r}, \uparrow} n_{\mathbf{r}, \downarrow}$$

- Observables: Energy gap, correlation functions
- Proposed model for high Tc superconductivity. A q. simulat
- 2D and 3D lattice gauge theory
 - Ring exchange model: Emergent U(I) gauge theory

$$H_{\rm RE} = K \sum_{\Box} (b_1^{\dagger} b_2 b_3^{\dagger} b_4 + b_1 b_2^{\dagger} b_3 b_4^{\dagger} - n_1 n_3 - n_2 n_4)$$

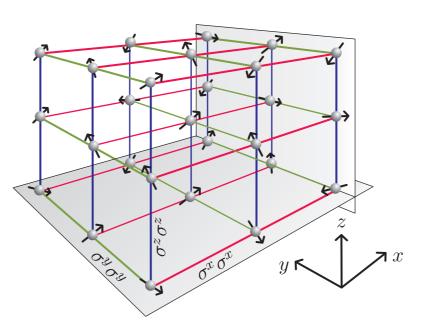
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A. Auerbach, 1994

 $\{c_{\mathbf{r},\sigma}, c^{\dagger}_{\mathbf{r}',\sigma'}\} = \delta_{\mathbf{r},\mathbf{r}'}\delta_{\sigma,\sigma'}$

H.P. Buchler, et al. PRL 95, 040402 (2005)

• Spin lattice models



- Study quantum phase transitions (simulate spin liquids, valence bond solids, etc)
- Resource states for measurement based quantum computation, e.g. AKLT model
- Emergent physics and topological order
 - Loop gas models, string net models, discrete gauge models (this talk)
 - General idea: prepare a highly entangled state of a spin network that is the vacuum state of a physical theory.

A spin lattice model* of discrete gauge theory

- Pick a finite group ${\cal G}$
 - spins on each edge: local basis $\{|g\rangle; g \in G\}$
- Local gauge transformation

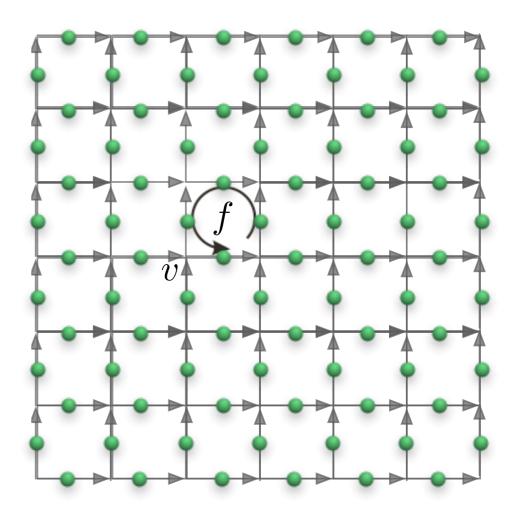
$$T_g(v) = \prod_{e_j \in [v,*]} L_g(e_j) \prod_{e_j \in [*,v]} R_{g^{-1}}(e_j)$$

• Want ground states local gauge invariant

$$H_{\text{TO}} = -\sum_{v} A(v) - \sum_{f} B(f)$$

$$A(v) \left| \begin{array}{c} g_{3} \\ g_{3} \\ g_{4} \\ g_{4} \end{array} \right\rangle = \frac{1}{|\overline{G}|} \sum_{h \in G} \left| \begin{array}{c} g_{3}h \\ g_{3}h \\ g_{4}h \\ g_{4}h \end{array} \right\rangle$$

$$B(f) \left| \begin{array}{c} g_{3} \\ g_{3} \\ g_{4} \\ g_{4}$$



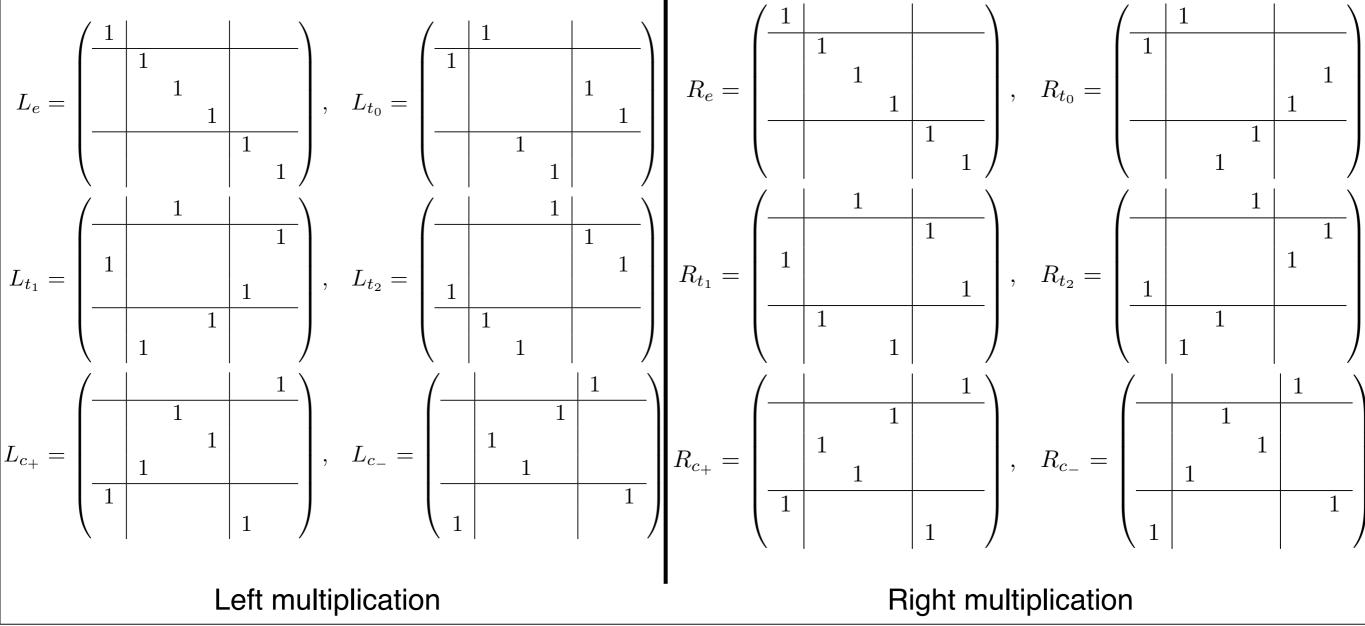
Realizes the quantum double ${\cal D}({\cal G})$

^{*}A. Yu. Kitaev, Annals of Physics **303**, 2 (2003)

- Gauge transformations $T_g(v) = \prod_{e_j \in [v,*]} L_g(e_j) \prod_{e_j \in [*,v]} R_{g^{-1}}(e_j)$
- Simplest non-Abelian group $S_3 = \{e, c_+, c_-, t_0, t_1, t_2\}$

cyclic perms transpositions

• Regular rep:



Sunday, 5 July 2009

• Simplification using semi-direct product structure of group $S_3 \cong \mathbb{Z}_3 \rtimes_{\phi} \mathbb{Z}_2$

$$L_{e} = \mathbf{1}_{3} \otimes \mathbf{1}_{2}, \quad L_{t_{0}} = F(1,2) \otimes \sigma^{x}, \quad L_{t_{1}} = F(0,2) \otimes \sigma^{x}, \quad L_{t_{2}} = F(0,1) \otimes \sigma^{x},$$
$$L_{c_{+}} = X^{-1} \otimes \mathbf{1}_{2}, \quad L_{c_{-}} = X \otimes \mathbf{1}_{2}, \quad R_{e} = \mathbf{1}_{3} \otimes \mathbf{1}_{2}, \quad R_{t_{0}} = \mathbf{1}_{3} \otimes \sigma^{x},$$
$$R_{t_{1}} = X^{-1} \otimes \sigma^{-} + X \otimes \sigma^{+}, \quad R_{t_{2}} = X^{-1} \otimes \sigma^{+} + X \otimes \sigma^{-},$$
$$R_{c_{+}} = X \otimes |0\rangle \langle 0| + X^{-1} \otimes |1\rangle \langle 1|, \quad R_{c_{-}} = X^{-1} \otimes |0\rangle \langle 0| + X \otimes |1\rangle \langle 1|,$$

$$F(i, j) = (|i\rangle\langle j| + |j\rangle\langle i|) \oplus 1$$

- Suggests a qutrit/qubit encoding of spins
- Efficient quantum circuit exists for preparing vacuum state of model + manipulation of anyonic excitations^{*}
 - works with or without a background Hamiltonian present

*M. Aguado, GKB, F. Verstraete, J.I. Cirac, Phys. Rev. Lett. **101**, 260501 (2008), GKB, M. Aguado, J.I. Cirac, New J. Phys. **11** 053009 (2009).

Particle spectrum of $D(S_3)$

• Labels Γ	$\mathbf{I}_{R(N_{[\alpha]})}^{[\alpha]} \bigstar$	conjugacy class> magnetic charge
	$\mathbf{I} (\mathbf{I} [\alpha])$	irrep of centralizer of conjugacy class> electric charge
particle type		quantum dimension $= [\alpha] R $
Vacuum	$\Pi^{[e]}_{R_1^+}$	1
Pure magnetic charge	$\Pi^{[c]}_{\beta_0}$	$\Pi_{\gamma_0}^{[t]} \tag{2,3}$
Pure electric charge	$\Pi^{[e]}_{R_1^-}$	$\Pi_{R_2}^{[e]} \tag{1,2}$
Dyonic combination	$\Pi^{[c]}_{\beta_1}$	$ \Pi_{\beta_2}^{[c]} \Pi_{\gamma_1}^{[t]} \tag{2,2,3} $

• Particles with quantum dimension >1 are non-Abelian anyons

Braid relations

- All excitations come in particle/anti-particle pairs
 - Magnetic flux pair $|a,a^{-1}\rangle$ \Box ----
 - Electric charge pairs transform under the irrep R and conjugate R*.

$$M^{R}\rangle = \frac{1}{\sqrt{|R|}} \sum M^{R}_{\mu,\nu} |\mu\rangle_{R} \otimes |\nu\rangle_{R^{*}} \qquad \sum_{\mu,\nu} |M^{R}_{\mu,\nu}|^{2} = |R| \qquad \diamondsuit \cdots \blacklozenge$$

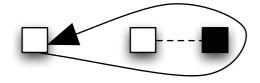
Interchanging two fluxes

$$\mathcal{R} |a\rangle |b\rangle = \sigma |a\rangle |aba^{-1}\rangle = |aba^{-1}\rangle |a\rangle$$

Braiding two fluxes

$$\mathbb{Q}^{2}|a\rangle|b\rangle = |abab^{-1}a^{-1}\rangle|abbb^{-1}a^{-1}\rangle = |(ab)a(ab)^{-1}\rangle|aba^{-1}\rangle$$

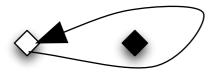
• Braiding one flux around flux pair



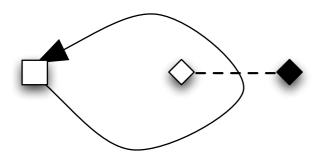
$$\mathcal{R}_{1,2}^{2} \otimes \mathcal{R}_{1,3}^{2} |b\rangle |a,a^{-1}\rangle = |b\rangle |bab^{-1}, ba^{-1}b^{-1}\rangle$$

- action is trivial if pair prepared in *chargeless* state $|0_{[\ell]}\rangle = \frac{1}{\sqrt{|[\ell]|}} \sum_{\ell \in [\ell]} |\ell, \ell^{-1}\rangle$

• Electric charges moving past each other have no effect

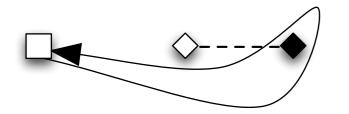


• Braiding a flux around one charge in a pair



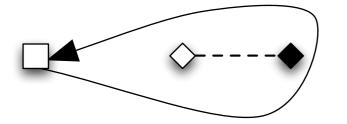
$$\mathcal{R}_{1,2}^{2}|h\rangle|M^{R}\rangle=|h\rangle|R(h)M^{R}\rangle$$

• Braiding a flux around the anti-charge in a pair



$$\mathcal{R}_{1,3}^{2}|h\rangle|M^{R}\rangle=|h\rangle|M^{R}R(h^{-1})\rangle$$

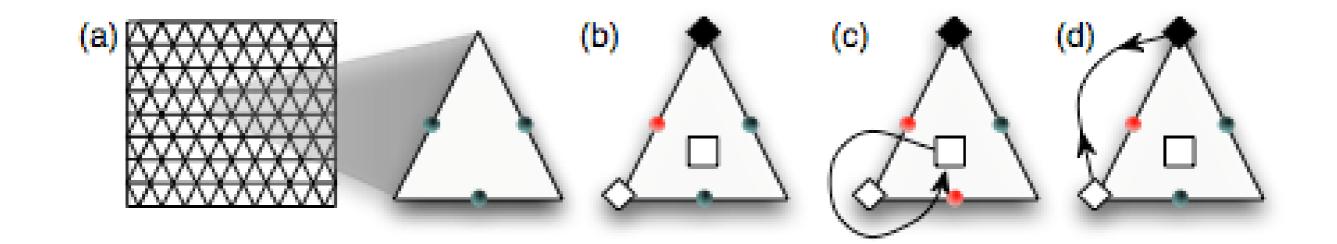
• Braiding a flux around the pair acts like conjugation



$$\mathcal{R}_{1,2}^2 \otimes \mathcal{R}_{1,3}^2 |h\rangle |M^R\rangle = |h\rangle |R(h)M^R R(h^{-1})\rangle$$

• For each irrep R, a unique *fluxless* state invariant under conjugation $|1_{|R|}\rangle$

Sketch of proposed interferometry



(a) A spin lattice model of $D(S_3)$, we simulate a single plaquette

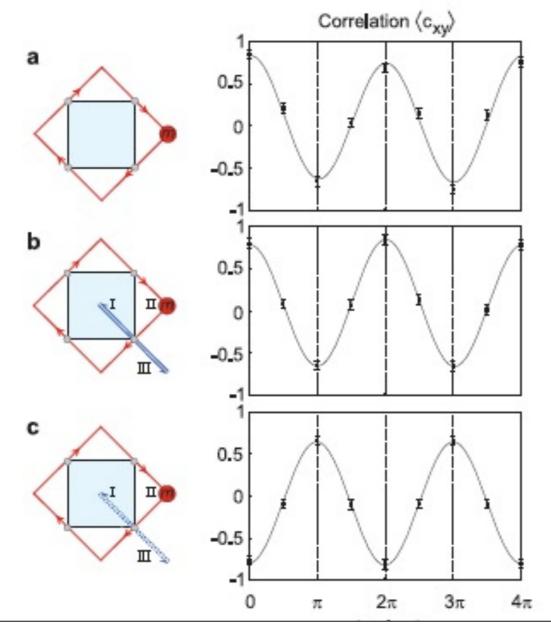
- (b)Acting on one spin (red) produces a anyonic electric charge pair (diamonds) and also can produce a flux (square)
- (c) Braiding the flux around one charge by acting on (red) spins
- (d)Fusion of the electric charge pair. Incomplete fusion to vacuum is signature of non-Abelian anyonic statistics

Experimental Simulation of Abelian anyons

• Algorithmic simulation of $D(\mathbb{Z}_2)$ (toric code) with entangled photons

J.K. Pachos, W. Wieczorek, C. Schmid, N. Kiesel, R. Pohlner, H. Weinfurter, arXiv:0710.0895 New J Phys (in press); Chao-Yang Lu, Wei-Bo Gao, Otfried Gühne, Xiao-Qi Zhou, Zeng-Bing Chen, Jian-Wei Pan, PRL **102**, 030502 (2009)

Background Hamiltonian is zero!



Ground state of

$$|\xi\rangle = \prod_{s} \frac{1}{\sqrt{2}} (\mathbbm{1} + \sigma_{s,1}^x \sigma_{s,2}^x \sigma_{s,3}^x \sigma_{s,4}^x) |00...0\rangle$$

Prepare GS on one plaquette $|\xi\rangle = (|0000\rangle + |1111\rangle)/\sqrt{2}$

Create flux pair braid around plaquette, annihilate

$$\sigma_1^x \sigma_2^x \sigma_3^x \sigma_4^x |\xi\rangle = |\xi\rangle$$

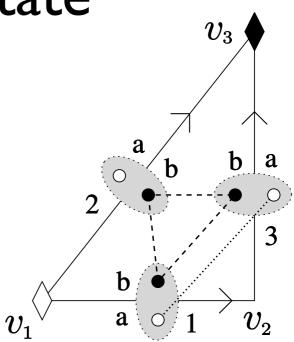
Create charge pair, braid flux pair around charge, annihilate

$$\sigma_3^z [\sigma_1^x \sigma_2^x \sigma_3^x \sigma_4^x] \sigma_3^z |\xi\rangle = -[\sigma_1^x \sigma_2^x \sigma_3^x \sigma_4^x] |\xi\rangle = -|\xi\rangle$$

Measure interference of two processes $e^{-i\frac{\pi}{4}\sigma_1^z} [\sigma_1^x \sigma_2^x \sigma_3^x \sigma_4^x] e^{i\frac{\pi}{4}\sigma_1^z} |\xi\rangle = (|0000\rangle - |1111\rangle)/\sqrt{2}$

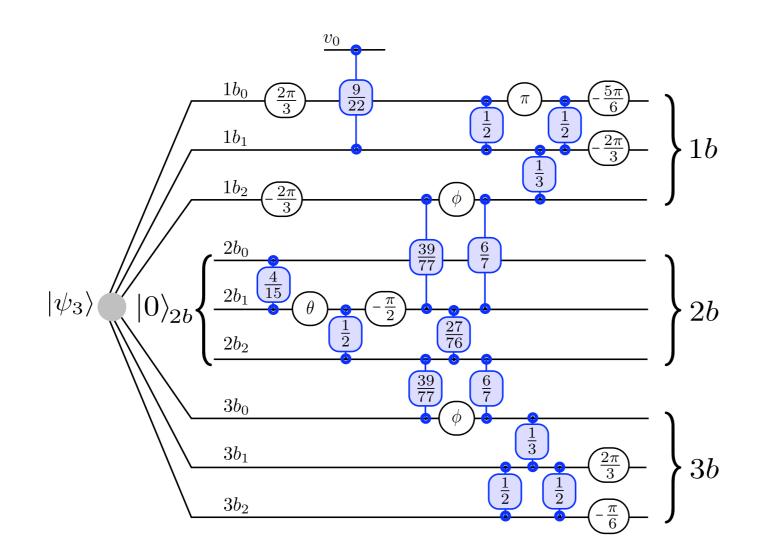
Building the initial state

- Resources needed:
 - 3 type-I SPDC crystals
 - 6 photons, 15 modes
 - 14 beam splitters + 11 phase shifters



- Creating 2 qutrit entanglement
 - type-I SPDC: strong pulse in, entangled photon in same polarization out $|SPDC\rangle = \sqrt{1-\lambda^2} \sum_{n=0}^{\infty} \lambda^n |nn\rangle$ - three crystals $|SPDC\rangle^{\otimes 3} = (1-\lambda^2)^{3/2} \sum_{n_1,n_2,n_3=0}^{\infty} \lambda^{n_1+n_2+n_3} |n_1n_1, n_2n_2, n_3n_3\rangle$
 - probability for one photon per triple above and below is $\lambda^2(1-\lambda^2)^3$.

• Input $|\psi_3\rangle_{1b,3b} = (|0\rangle_{1b}|0\rangle_{3b} + |1\rangle_{1b}|1\rangle_{3b} + |2\rangle_{1b}|2\rangle_{3b})/\sqrt{3}$



- Post-select on one photon per triple mode 1,2,3. Success probability 9/55
- **Output** $\{2|0\rangle_{2b}(|0\rangle_{1b}|0\rangle_{3b}+|1\rangle_{1b}|1\rangle_{3b}+|2\rangle_{1b}|2\rangle_{3b})-|2\rangle_{2b}(|0\rangle_{1b}|1\rangle_{3b}+|1\rangle_{1b}|2\rangle_{3b}+|2\rangle_{1b}|0\rangle_{3b})-|2\rangle_{2b}(|0\rangle_{1b}|1\rangle_{3b}+|1\rangle_{1b}|2\rangle_{3b}+|2\rangle_{1b}|1\rangle_{3b})\}/(3\sqrt{2}),$

Measuring fusion data (a simple method)

- Braid flux around one charge at vertex v $T_h(v)|\mathbf{1}_{R_2};(v,v')\rangle = |R_2(h);(v,v')\rangle$
 - Measure projector onto irrep for fluxless charge pair

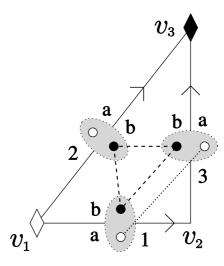
$$\langle R(h)|W_{R'}|R(h)\rangle = \sum_{a,b,d,e=1}^{|R|} \sum_{c=1}^{|R'|} Q_{acd,bce}^{[RR'*R^*]} R_{ab}^*(h) R_{de}(h)$$

• Projectors onto vacuum fusion channel for three irreps

$$Q_{ace,bcf}^{[R^{(1)}R^{(2)}R^{(3)}]} = \frac{1}{|G|} \sum_{g} R_{ab}^{(1)}(g) R_{cd}^{(2)}(g) R_{ef}^{(3)}(g)$$

- Just local (non-entangling) operations + measurement on photons in 2a and 2b
 - signature of non-Abelian statistics

$$\langle R_2(h); (v, v') | W_{R_2}(e) | R_2(h); (v, v') \rangle = \begin{cases} 1 & h = e \\ -\frac{1}{2} & h = c_{\pm} \\ 0 & h = t_j \forall j \end{cases}$$



Summary

- Good quantum simulation requires good quantum control
- Can simulate emergent physics
 - Build highly entangled spin networks corresponding to vacuum states of models with exotic excitations
 - Can manipulate these excitations and measure their properties
- Photonic spin networks look promising in the near term for demonstrating prototype models
 - Integrated photonics* with waveguides etched into glass is a good platform

*A. Politi, M. J. Cryan, J. G. Rarity, S. Yu, and J. L. O'Brien, Science **320**, 646 (2008).

 Larger simulations become inefficient due to bad scaling of probability to create many spin entangled states. There are work arounds but other systems such as Josephson junctions or trapped atoms/molecules in optical lattices may be better