

Interacting anyons in topological quantum liquids: The golden chain

Adrian Feiguin,¹ Simon Trebst,¹ Andreas W. W. Ludwig,^{2,3} Matthias Troyer,⁴
Alexei Kitaev,^{1,5} Zhenghan Wang,¹ and Michael H. Freedman¹

¹*Microsoft Research, Station Q, University of California, Santa Barbara, CA 93106*

²*Physics Department, University of California, Santa Barbara, CA 93106*

³*Kavli Institute for Theoretical Physics, University of California, Santa Barbara, CA 93106*

⁴*Theoretische Physik, Eidgenössische Technische Hochschule Zürich, 8093 Zürich, Switzerland*

⁵*Institute for Quantum Information, California Institute of Technology, Pasadena, CA 91125*

(Dated: November 26, 2024)

We discuss generalizations of quantum spin Hamiltonians using anyonic degrees of freedom. The simplest model for interacting anyons energetically favors neighboring anyons to fuse into the trivial ('identity') channel, similar to the quantum Heisenberg model favoring neighboring spins to form spin singlets. Numerical simulations of a chain of Fibonacci anyons show that the model is critical with a dynamical critical exponent $z = 1$, and described by a two-dimensional (2D) conformal field theory with central charge $c = 7/10$. An exact mapping of the anyonic chain onto the 2D tricritical Ising model is given using the restricted-solid-on-solid (RSOS) representation of the Temperley-Lieb algebra. The gaplessness of the chain is shown to have topological origin.

PACS numbers: 05.30.Pr, 73.43.Lp, 03.65.Vf

Introduction Non-Abelian anyons are exotic particles expected to exist in certain fractional quantum Hall (FQH) states [1, 2]. A set of several anyons supports very robust collective states that are degenerate to exponential precision; such states can potentially be used as quantum memory and for quantum computation [3]. However, this degeneracy can be lifted by a short-range interaction if the anyons are very close to each other. As a first step towards understanding interacting anyons, we describe a simple, exactly solvable model that is an anyonic analogue of the quantum Heisenberg chain.

We start by considering the well-known Moore-Read state [1], a candidate state, exhibiting non-Abelian statistics, for the topological nature of FQH liquids at filling fraction $\nu = 5/2$. It has two important types of excitations: quasiholes with electric charge $e/4$ and neutral fermions. Quasiholes may be trapped by an impurity potential while the fermions can still tunnel between them [4]. For a one-dimensional (1D) array of trapped quasiholes, the Hamiltonian can be described in terms of free Majorana fermions on a lattice, which is in turn equivalent to the 1D transverse field Ising model at the quantum phase transition point. The more interesting model discussed here is based on so-called 'Fibonacci anyons', which represent the non-Abelian part of the quasiparticle statistics in the $k = 3, Z_k$ -parafermion state [2], an effective theory for FQH liquids at filling fraction $\nu = 12/5$ [5]. Even without parameter fine-tuning, these 1D anyonic arrays will be shown to exhibit gapless excitations due to topological symmetry.

Model Our model describes pairwise interactions within an array of L anyons, for instance along a chain as shown in Fig. 1a). In the Fibonacci theory there are only two types of particles: the Fibonacci anyon, denoted by τ , and the trivial particle denoted by 1 with a fusion rule $\tau \times \tau = 1 + \tau$. We refer to the label 1 or τ as the topological charge. When two neighboring anyons interact, indicated in the figure by the ellipses, they can either fuse in the trivial channel, annihilating each other, or in the nontrivial one, becoming a single τ -anyon

[6]. We define our model by assigning an energy gain if they fuse along the trivial channel. This is an anyonic analogue of the spin-1/2 quantum Heisenberg antiferromagnet, which assigns an energy gain to two neighboring spin-1/2 fusing into a spin-0 singlet as compared to a spin-1 triplet.

To define the Hilbert space of τ -anyons we consider the tree-like fusion diagram in Fig. 1b). The basis corresponds to all admissible labelings $|x_1, x_2, \dots\rangle$ of the links, with $x_i = 1$ or τ . Each label represents the combined topological charge of the particles left to a given point. Not all possible values (x_1, x_2, \dots) represent allowed basis states due to the fusion rules: a 1 must always be preceded and followed by a τ , since the fusion of a 1 and a τ always gives a τ . This reduces the dimension of the Hilbert space of the open chain (with τ -labels at the boundary) to the Fibonacci sequence $\dim_L = F_{L+1}$, and for the periodic chain $\dim_L = F_{L-1} + F_{L+1}$. For large L it is well-known that these numbers grow at a rate $\dim_L \propto \varphi^L$, where $\varphi = (1 + \sqrt{5})/2$ is the golden ratio. This Hilbert space has no natural decomposition in the form of a tensor product of single-site states, in contrast to $SU(2)$ quantum spin chains.

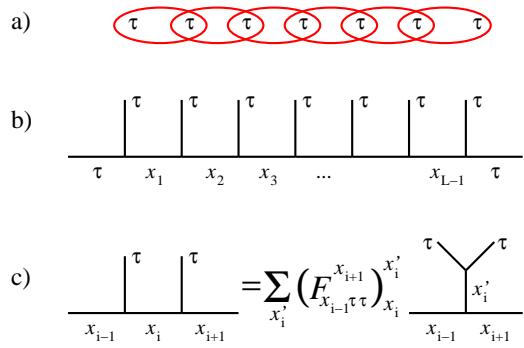


FIG. 1: a) Illustration of the Fibonacci chain with L τ -anyons.
b) The fusion path. c) Definition of the F -matrix.

In order to generate a local Hamiltonian $\mathbf{H} = \sum_i \mathbf{H}_i$ assigning an energy to the fusion of two neighboring τ -anyons we use the so-called F -matrix to transform the local basis as shown in Fig. 1c). In the transformed basis the state x'_i corresponds to the fusion of the two anyons. The Hamiltonian is then defined by assigning an energy $E_\tau = 0$ for $x'_i = \tau$, and $E_1 = -1$ for $x'_i = 1$. The resulting local terms \mathbf{H}_i contain three-body interactions in the link basis,

$$\mathbf{H}_i|x_{i-1}x_ix_{i+1}\rangle = \sum_{x'_i=1,\tau} (\mathbf{H}_i)_{x_i}^{x'_i} |x_{i-1}x'_{i+1}\rangle \quad (1)$$

with $(\mathbf{H}_i)_{x_i}^{x'_i} := -(F_{x_{i-1}\tau\tau}^{x_{i+1}})_{x_i}^1 (F_{x_{i-1}\tau\tau}^{x_{i+1}})_{x'_i}^1$.

It is diagonal in the subspace $\{|x_{i-1}x_ix_{i+1}\rangle\} = \{|1\tau 1\rangle, |1\tau\tau\rangle, |\tau\tau 1\rangle\}$, $\mathbf{H}_i = \text{diag}\{-1, 0, 0\}$, where the F -matrix is a number due to the constraints arising from the fusion rules. For the case $x_{i-1} = x_{i+1} = \tau$, the F -matrix and the corresponding Hamiltonian are the following 2×2 -matrices ($x_i, x'_i \in \{1, \tau\}$)

$$\mathbf{F}_{\tau\tau\tau}^\tau = \begin{pmatrix} \varphi^{-1} & \varphi^{-1/2} \\ \varphi^{-1/2} & -\varphi^{-1} \end{pmatrix}, \quad \mathbf{H}_i = -\begin{pmatrix} \varphi^{-2} & \varphi^{-3/2} \\ \varphi^{-3/2} & \varphi^{-1} \end{pmatrix}. \quad (2)$$

Looking at the matrix form of the Hamiltonian, it can be written in terms of standard Pauli matrices:

$$\mathbf{H}_i = (n_{i-1} + n_{i+1} - 1) - n_{i-1}n_{i+1} \left(\varphi^{-3/2} \sigma_i^x + \varphi^{-3} n_i + 1 + \varphi^{-2} \right),$$

where the sum runs over the links of the chain. In this expression, the operators n_i count the τ -particle occupation on link i , $n_i = \frac{1}{2}(1 - \sigma_i^z) = 0, 1$, and the Hamiltonian \mathbf{H} acts on the constrained Hilbert space defined above.

Entropy scaling and central charge We simulated this model numerically, and calculated the finite-size excitation gap by exact diagonalization for chains of up to $L = 32$ sites. A finite-size analysis shows that the gap vanishes linearly in $1/L$, indicative of a critical model with dynamical critical exponent $z = 1$ described by a conformal field theory. The central charge c of a CFT can be calculated from the finite-size scaling of the entanglement entropy. For two subsystems with equal size $L/2$ on systems with periodic (PBC) and open boundary conditions (OBC) the entanglement entropy scales as [7]

$$S_{\text{PBC}}(L) \propto \frac{c}{3} \log L \quad \text{and} \quad S_{\text{OBC}}(L) \propto \frac{c}{6} \log L. \quad (3)$$

The density matrix renormalization group method (DMRG) [8, 9] provides a natural framework for calculating these quantities. Fits of our numerical results according to Eqs. (3) shown in Fig. 2, give central charge estimates of $c_{\text{PBC}} = 0.701 \pm 0.001$ and $c_{\text{OBC}} = 0.70 \pm 0.01$ respectively. Since possible (unitary) CFTs in the vicinity of these estimates have central charges[19] $1/2, 7/10$ or $4/5$ we can unambiguously conclude that our results are consistent only with central charge $c = 7/10$.

Mapping and exact solution We now proceed to derive these results exactly. By construction the local contribution to the Hamiltonian $\frac{1}{\varphi} \mathbf{X}_i = -\mathbf{H}_i$ is a projector onto the trivial particle. One can then verify that the operators \mathbf{X}_i form a representation of the Temperley-Lieb algebra [10]

$$\begin{aligned} (\mathbf{X}_i)^2 &= d \mathbf{X}_i, \\ \mathbf{X}_i \mathbf{X}_{i \pm 1} \mathbf{X}_i &= \mathbf{X}_i, \\ [\mathbf{X}_i, \mathbf{X}_j] &= 0 \text{ for } |i - j| \geq 2, \end{aligned} \quad (4)$$

where the ‘d-isotopy’ parameter equals the golden ratio, $d = \varphi$. This representation can be seen to be identical to the standard Temperley-Lieb algebra representation associated with $SU(2)_k$ at level $k = 3$. For arbitrary k , the latter contains $k+1$ anyon species labelled by $j = 0, 1/2, 1, \dots, k/2$, satisfying the fusion rules of $SU(2)_k$ [11]. The operators \mathbf{e}_i defined by

$$\begin{aligned} \mathbf{e}[i]|j_{i-1}j_i j_{i+1}\rangle &= \sum_{j'_i} \left(\mathbf{e}[i]_{j_{i-1}}^{j_{i+1}} \right)_{j_i}^{j'_i} |j_{i-1}j'_i j_{i+1}\rangle \\ \text{and } \left(\mathbf{e}[i]_{j_{i-1}}^{j_{i+1}} \right)_{j_i}^{j'_i} &= \delta_{j_{i-1}, j_{i+1}} \sqrt{\frac{S_{j_i}^0 S_{j'_i}^0}{S_{j_{i-1}}^0 S_{j_{i+1}}^0}} \end{aligned} \quad (5)$$

are known [12] to form a representation of the Temperley-Lieb algebra (4) for any value of k , where $|j_i - j_{i+1}| = 1/2$ and $S_j^{j'} := \sqrt{\frac{2}{(k+2)} \sin[\pi \frac{(2j+1)(2j'+1)}{k+2}]} [13]$.

Our model of interacting Fibonacci anyons can be cast into this form at $k = 3$ by first mapping $x_i = 1 \rightarrow j_i = 0$, and $x_i = \tau \rightarrow j_i = 1$, and then applying the $SU(2)_3$ fusion rule $3/2 \times j = 3/2 - j$ to the even-numbered sites. This maps any admissible labeling $|\vec{x}\rangle := |x_1, x_2, \dots\rangle$ uniquely into $|\vec{j}\rangle :=$

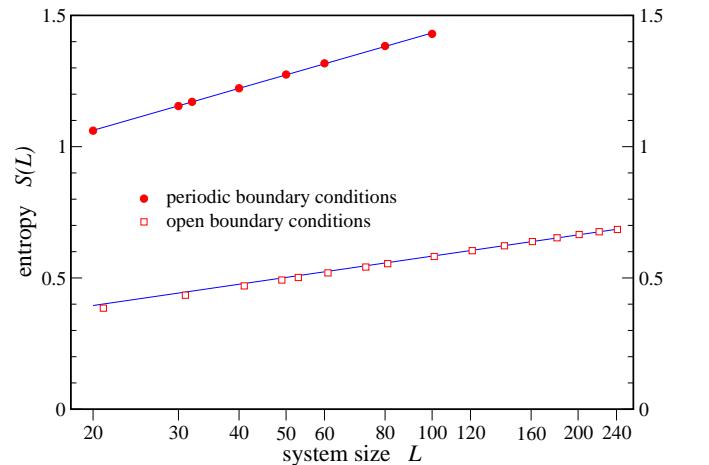


FIG. 2: (Color online) Entropy scaling for interacting Fibonacci anyons arranged along an open (open squares) or periodic chain (closed circles) versus the system size L . Logarithmic fits (solid lines) give central charge estimates of $c_{\text{PBC}} = 0.701 \pm 0.001$ and $c_{\text{OBC}} = 0.70 \pm 0.01$ respectively, where for the open boundary conditions only the values for the 5 largest systems have been taken into account due to large finite-size effects.

$|j_1, j_2, \dots\rangle$ where for odd-numbered sites $j_{2i+1} \in \{0, 1\}$, and for even numbered-sites $j_{2i} \in \{1/2, 3/2\}$. This re-labeling maps the matrix elements of \mathbf{X}_i into those of \mathbf{e}_i from Eq. (5).

We can now see that the Hamiltonian in Eq. (1) is that corresponding to a standard (integrable) lattice model description of the classical 2D tricritical Ising model, known as the RSOS model [14]. Specifically, the two-row transfer matrix $\mathbf{T} := \mathbf{T}_2 \mathbf{T}_1$ of this lattice model, shown in Fig. 3, is written in terms of Boltzmann weights $\mathbf{W}[i]$ assigned to a plaquette i of the square lattice

$$\mathbf{T}_1 := \prod_n \mathbf{W}[2n], \quad \text{and} \quad \mathbf{T}_2 := \prod_n \mathbf{W}[2n+1]$$

with

$$\mathbf{W}[i]_{\vec{j}}^{j'} = \frac{\sin[\frac{\pi}{k+2} - u]}{\sin \frac{\pi}{k+2}} \mathbf{1}_{\vec{j}}^{j'} + \frac{\sin u}{\sin \frac{\pi}{k+2}} \mathbf{e}[i]_{\vec{j}}^{j'}. \quad (6)$$

The parameter $u > 0$ is a measure of the lattice anisotropy, $\mathbf{1}$ is the identity operator, and

$$\mathbf{e}[i]_{\vec{j}}^{j'} := \left[\prod_{m \neq i} \delta_{j'_m, j_m} \right] \left(\mathbf{e}[i]_{j_{i-1}}^{j_{i+1}} \right)_{j_i}^{j'_i}. \quad (7)$$

The Hamiltonian of the so-defined lattice model is obtained from its transfer matrix by taking, as usual [15], the extremely anisotropic limit, $u \ll 1$,

$$\mathbf{T} = \exp\{-a(\mathbf{H} + c_1) + O(a^2)\}, \quad a = \frac{u\varphi}{\sin[\pi/(k+2)]} \ll 1$$

yielding $\mathbf{H} = -\sum_i \frac{1}{\varphi} \mathbf{e}_i$ (c_1 is an unimportant constant). Since the operators \mathbf{X}_i can be identified with \mathbf{e}_i , this demonstrates that the Hamiltonian of the Fibonacci chain is exactly that of the corresponding $k = 3$ RSOS model which is a lattice description of the tricritical Ising model at its critical point. The latter is a well-known (supersymmetric) CFT with central charge $c = 7/10$ [16, 17]. Analogously one obtains [18] for general k the $(k-1)^{\text{st}}$ unitary minimal CFT [19] of central charge $c = 1 - 6/(k+1)(k+2)$. A ferromagnetically coupled Fibonacci chain (energetically favoring the fusion along the τ -channel) is described by the critical 3-state Potts model with $c = 4/5$ and, for general k , by the critical Z_k -parafermion CFT [14, 18, 20] with central charge $c = 2(k-1)/(k+2)$.

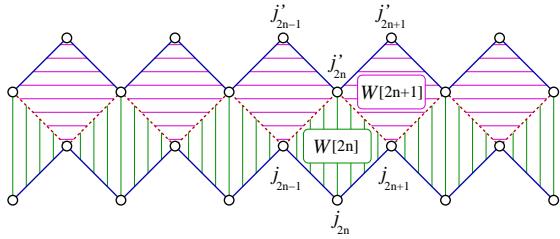


FIG. 3: Transfer matrix of the RSOS model.

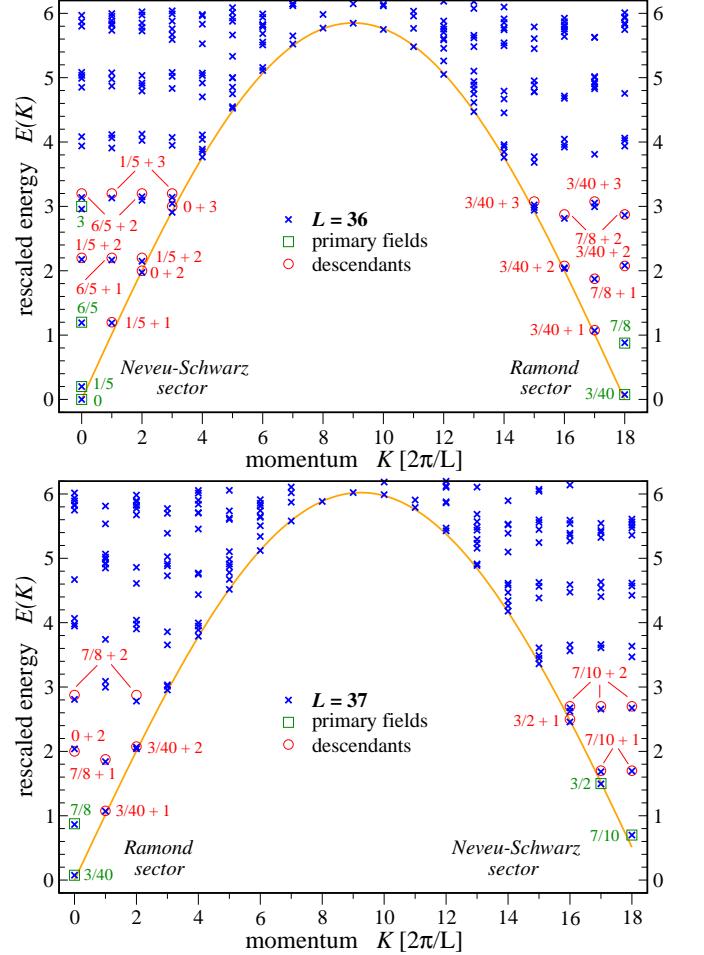


FIG. 4: Energy spectra for periodic Fibonacci chains of size $L = 36$ and $L = 37$. The spectra have been rescaled and shifted such that the two lowest eigenvalues match the conformal field theory assignments. The open boxes indicate the positions of the primary fields of the $c = 7/10$ conformal field theory. The open circles give the positions of multiple descendant fields as indicated. While we find excellent agreement in general, finite-size effects lead to small discrepancies for the higher energy states. The solid line is a cosine-fit of the dispersion which serves as a guide to the eye.

Excitation spectra We have calculated the excitation spectra of chains up to size $L = 37$ with open and periodic boundary conditions using exact diagonalization, as shown in Fig. 4. The numerical results not only confirm the CFT predictions but also reveal some important details about the correspondence between continuous fields and microscopic observables. In general, low-energy states on a ring are associated with local conformal fields [21], whose holomorphic and antiholomorphic parts belong to representations of the Virasoro algebra, described by conformal weights h_L and h_R . The energy levels are given by

$$E = E_1 L + \frac{2\pi v}{L} \left(-\frac{c}{12} + h_L + h_R \right), \quad (8)$$

eigenvalue	numerics $L = 31$	CFT assignment	numerics $L = 32$	CFT assignment
0	0.10	1/10	0	0
1	1.10	1/10 + 1	0.60	3/5
2	1.49	3/2	1.60	3/5 + 1
3	2.09	1/10 + 2	2.02	0 + 2
4	2.47	3/2 + 1	2.58	3/5 + 2
5	3.07	1/10 + 3	2.59	3/5 + 2
6	3.11	1/10 + 3	3.01	0 + 3
7	3.44	3/2 + 2	3.56	3/5 + 3
8	3.46	3/2 + 2	3.56	3/5 + 3

TABLE I: Lowest eigenvalues for open Fibonacci chains of size $L = 31$ and $L = 32$. The two lowest eigenvalues have been rescaled and shifted such that they match the conformal field theory assignments.

corresponding to states with a choice of momenta $K = h_L - h_R$ or $K = h_L - h_R + L/2$ in units of $2\pi/L$, where E_1 , v are non-universal constants. Here, $h_L = h_L^{(0)} + m_L$ and $h_R = h_R^{(0)} + m_R$, where $h_L^{(0)}, h_R^{(0)}$ correspond to weights of ‘primary’ fields and m_L and m_R are non-negative integers describing so-called ‘descendant’ fields. The numerical spectra for even values of L (see the first plot in Fig. 4) agree with Eq. (8), exhibiting primary fields with $h_L^{(0)} = h_R^{(0)} = 0, 1/10, 3/5, 3/2, 3/80, 7/16$, which are conventionally denoted by $I, \varepsilon, \varepsilon', \varepsilon'', \sigma, \sigma'$, respectively[24]. The momenta of the last two fields and their descendants are near $K = L/2$, as compared to the other four, indicating that the corresponding microscopic observables have alternating sign on the lattice. Such “staggered” fields must have nontrivial monodromy with respect to a space-time dislocation (i.e., the insertion or removal of a site at some particular time). Such a dislocation is characterized by a chiral ε'' field, say, $\varepsilon_L^{''}$ [22]. The log of the monodromy factor $\exp(2\pi i(h_L^{\psi \times \varepsilon''} - h_L^\psi - h_L^{\varepsilon''}))$ matches the momenta K in Fig. 4 [23]. Given this information, we may predict that the states of an odd size ring are associated with fields of the form $\xi_L \eta_R$, where $\eta = \xi \times \varepsilon''$. These include six primary fields, $\varepsilon_L^{''}, \varepsilon_R^{''}, \varepsilon_L' \varepsilon_R, \varepsilon_L \varepsilon_R', \sigma_L \sigma_R, \sigma_L' \sigma_R'$, as well as their descendants. Integrality of the momentum K dictates the choice of K (see below Eq. (8)), as in Fig. 4.

For open boundary conditions the spectra are known to be described by, say, the holomorphic sector only [26]. To explain the numerical data, we need to assume that the ends of the chain are characterized by a boundary field ε (or equally well ε'). Thus, for an even number of sites the spectrum is described by $\varepsilon \times \varepsilon = I + \varepsilon'$ (plus descendants). For an odd number of sites, this result is to be modified by fusion with ε'' , yielding $\varepsilon'' + \varepsilon$. The numerical low-energy spectra agree excellently with these predictions as shown in Table I.

Hidden symmetries The critical behavior of our model is not just a peculiarity of the exact solution but rather has topological origin. In general, an effective low-energy Lagrangian admits perturbations of the form $\int \psi(x, \tau) dx d\tau$, where ψ may be any local field that is consistent with all applicable

symmetries. Such terms are relevant if $h_L^\psi + h_R^\psi < 2$, in which case they may open a spectral gap or induce crossover to different critical behavior at large distances. In the tricritical Ising model, there are four relevant fields: $\varepsilon_L \varepsilon_R, \varepsilon'_L \varepsilon'_R, \sigma_L \sigma_R, \sigma'_L \sigma'_R$. Some explanation is in order as to why these fields do not appear in the effective Lagrangian of our model. The fields σ and σ' are staggered and thus prohibited by translational symmetry. Excluding ε and ε' requires a more subtle argument. The Fibonacci ring has a topological symmetry, which corresponds to adding an extra τ -line parallel to the spine of the fusion diagram (Fig. 1b) and merging it with the diagram using the F -matrix. We denote this operator by Y .

$$\langle x'_0, \dots, x'_{L-1} | Y | x_0, \dots, x_{L-1} \rangle = \prod_{i=0}^{L-1} \left(F_{\tau x_i \tau}^{x'_{i+1}} \right)_{x_{i+1}}^{x'_i},$$

where the identification $L \equiv 0$ is used. We may think of the fusion diagram as a description of a process that generates a set of τ -anyons on a circle from the local vacuum. Then Y describes another particle moving inside or outside that circle (or on the circle itself — before those anyons were created). The operator Y is sensitive to a possible topological charge $y = 1, \tau$ located at the center of the circle. Thus Y has two eigenvalues, $S_{y\tau}/S_{y1} = \varphi, -\varphi^{-1}$. We conjecture that the low-energy states associated with fields $I, \varepsilon'', \sigma'$ are in the trivial ($y = 1$) sector, and the fields $\varepsilon', \varepsilon, \sigma$ are in the $y = \tau$ sector. In fact, the topological fusion algebra (defined by the rule $\tau \times \tau = 1 + \tau$) is a quotient of the CFT fusion algebra.

We may imagine that the interaction between the anyons alters the topological liquid in which the anyons are excitations, producing an annulus of a different liquid. Some of the local fields correspond to the tunneling of a τ -anyon between the inner and outer edge of the annulus. Such a process is actually forbidden as it would change the topological charge y . Thus, only fields in the trivial topological sector $y = 1$ are allowed as perturbations. This excludes $\psi = \varepsilon_L \varepsilon_R$ and $\psi = \varepsilon'_L \varepsilon'_R$.

Outlook Extensions to chains of anyons in topological liquids described by $SU(2)_k$ Chern-Simons theory with $k > 3$ corresponding (up to phases) to the non-abelian statistics of higher members of the Read-Rezayi series have been mentioned below Eq. (7), but their topological stability is an open issue. In analogy to quantum spin chains, additional interactions such as dimerization or coupling of two Fibonacci chains in a ladder geometry should lead to gapped quantum liquids. Disordered anyonic chains are currently being investigated [25]. For 2D anyonic structures gapless phases of non-Fermi liquid type might potentially also emerge.

We thank E. Ardonne, N. Bonesteel, P. Fendley, C. Nayak, G. Refael, S. H. Simon, and J. Slingerland for stimulating discussions. Some of our numerical simulations were based on the ALPS libraries [27].

-
- [1] G. Moore and N. Read, Nucl. Phys. B **360**, 362 (1991).
[2] N. Read and E. Rezayi, Phys. Rev. B **59**, 8084 (1999).

- [3] A. Yu Kitaev, Ann. Phys. **303**, 3 (2003).
- [4] See e.g. N. Read and A.W.W. Ludwig, Phys. Rev. B **63**, 024404 (2001).
- [5] J. S. Xia *et al.*, Phys. Rev. Lett. **93**, 176809 (2004).
- [6] For a pedagogical introduction we refer to J. Preskill, Lecture notes on quantum computation; available online at <http://www.theory.caltech.edu/~preskill/ph229>
- [7] C. Holzhey et al., Nucl. Phys. B **424**, 44 (1994); P. Calabrese and J. Cardy, J. Stat. Mech. P06002 (2004).
- [8] S. R. White, Phys. Rev. Lett. **69**, 2863 (1992); U. Schollwöck, Rev. Mod. Phys. **77**, 259 (2005).
- [9] The DMRG calculations were done for system sizes up to $L = 240$, keeping up to $m = 2000$ states. The constrained Hilbert space was implemented by explicitly adding a nearest-neighbor repulsion $V = U \sum_{i=0}^{L-1} (1 - n_i)(1 - n_{i+1})$, with $U \gg 1$.
- [10] N. Temperley and E. Lieb, Proc. Roy. Soc. Lond. A **322**, 251 (1971).
- [11] D. Gepner and E. Witten, Nucl. Phys. B **278**, 493 (1986); A. B. Zamolodchikov and V. A. Fateev, Sov. J. Nucl. Phys. **43**, 657 (1986).
- [12] V. Jones, C. R. Acad. Sci. Paris Sér. I Math. **298**, 505 (1984); A. Kuniba, Y. Akutsu, and M. Wadati, J. Phys. Soc. Jpn. **55**, 3285 (1986); V. Pasquier, Nucl. Phys. B **285**, 162 (1987); H. Wenzl, Invent. Math. **92**, 349 (1988). For a recent discussion see: P. Fendley, J. Phys. A **39**, 15445 (2006).
- [13] The matrix $S_j^{j'}$ is the modular S-matrix of $SU(2)_k$ CFT.
- [14] G. E. Andrews, R. J. Baxter, and P. J. Forrester, J. Stat. Phys. **35**, 193 (1984).
- [15] See for example: R. J. Baxter, *Exactly solved models in statistical mechanics*, Academic Press, London (1982).
- [16] D. Friedan, Z. Qiu, and S. Shenker, Phys. Lett. B **151**, 37 (1985).
- [17] A. A. Belavin, A. M. Polyakov, and A. B. Zamolodchikov, Nucl. Phys. B **241**, 333 (1984).
- [18] D. A. Huse, Phys. Rev. B **30**, 3908 (1984).
- [19] D. Friedan, Z. Qiu and S. Shenker, Phys. Rev. Lett. **52**, 1575 (1985).
- [20] Here the sign of u in Eq. (6) is reversed [14] as compared to the antiferromagnetic case.
- [21] J. L. Cardy, J. Phys. A **17**, L385 (1984); Nucl. Phys. B **270**, 186 (1986).
- [22] ε'' is the superpartner of the stress tensor in this $N = 1$ superconformal theory [16].
- [23] We used the fact that the fusion with ε'' corresponds the mapping: $I \leftrightarrow \varepsilon'', \varepsilon \leftrightarrow \varepsilon', \sigma \leftrightarrow \sigma, \sigma' \leftrightarrow \sigma'$. The complete set of fusion rules can be found in [17], see e.g. also: P. Di Francesco, P. Mathieu, and D. Sénéchal, *Conformal Field Theory*, Springer, New York (1997).
- [24] The last two fields are the order operators, i.e. those coupling to applied magnetic fields in the tricritical Ising model.
- [25] N. Bonesteel, private communication
- [26] J. L. Cardy, Nucl. Phys. B **240**, 514 (1984).
- [27] F. Alet *et al.*, J. Phys. Soc. Jpn. Suppl. **74**, 30 (2005).