Phase Transitions and Random Quantum Satisfiability

C. Laumann,1 R. Moessner,2 A. Scardicchio,1,3 and S. L. Sondhi1

1Department of Physics, Princeton University, Princeton, NJ 08544
2Max Planck Institut fur Physik Komplexer Systeme, 01187 Dresden, Germany
3Princeton Center for Theoretical Science, Princeton University, Princeton, NJ 08544

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Alongside the effort underway to build quantum computers, it is important to better understand which classes of problems they will find easy and which others even they will find intractable. We study random ensembles of the QMA1-complete quantum satisfiability (QSAT) problem introduced by Bravyi [1]. QSAT appropriately generalizes the NP-complete classical satisfiability (SAT) problem. We show that, as the density of clauses/projectors is varied, the ensembles exhibit quantum phase transitions between phases that are satisfiable and unsatisfiable. Remarkably, almost all instances of QSAT for any hypergraph exhibit the same dimension of the satisfying manifold. This establishes the QSAT decision problem as equivalent to a, potentially new, graph theoretic problem and that the hardest typical instances are likely to be localized in a bounded range of clause density.

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I. INTRODUCTION

The potential power of quantum computers motivates the intense effort in progress to understand and, eventually, build them. Much interest has, naturally, been focused on algorithms that outperform their classical counterparts. However the development of a complexity theory for quantum computers suggests that we already know problems, those shown to be QMA-complete, whose worst case solutions will take even quantum computers a time exponential in problem size. These appropriately generalize the class of NP-complete problems—those which are easy to check but (believed to be) hard to solve on classical computers—to quantum computers [2, 3].

The classic technique of complexity theory is assigning guilt by association, i.e. by exhibiting the polynomial time equivalence of a new problem to a particular problem believed not to be amenable to efficient solution. For NP-complete problems this leads to the celebrated Cook-Levin theorem which shows that the satisfiability problem with clauses in three Boolean variables (3-SAT) encapsulates the difficulty of the entire class [4]. While this is a powerful and rigorous approach it has two limitations. It does not directly tell us why a problem has hard instances and it does not tell us what general features they possess.

Over the past decade joint work by computer scientists and statistical physicists has produced an interesting set of insights into these two lacunae for 3-SAT (and related constraint satisfaction problems). These insights have come from the study of instances chosen at random from ensembles where the density of clauses \( \alpha \) acts as a control parameter. One can think of this as representing the study of typical, but not necessarily worst, cases with a specified density of clauses/constraints. Using techniques developed for the study of random systems in physics, it has been shown that the ensembles exhibit a set of phase transitions between a trivial satisfiable (SAT) phase at small \( \alpha \) and an unsatisfiable (UNSAT) phase at large \( \alpha \) (see e.g. [5, 6, 7, 8, 9]). These transitions mark sharp discontinuities in the organization of SAT assignments in configuration space as well a vanishing of SAT assignments altogether. This information has provided a heuristic understanding of why known algorithms fail on random SAT when the solution space is sufficiently complex and in doing so has localized the most difficult members of these ensembles to a bounded range of \( \alpha \).

In this work we begin an analogous program for quantum computation with the intention of gaining insight into the difficulty of solving QMA-complete problems. Specifically, we introduce and study random ensembles of instances of the quantum satisfiability (\( k \)-QSAT) prob-
lem formulated by Bravyi \[1\]. Like classical 2-SAT, 2-QSAT is efficiently solvable (i.e. it is in P), while for \(k \geq 4\), k-QSAT is QMA\(_1\)-complete \[19\].

After laying out some important definitions and background on our problem in Sec. \[I\] we attack the ("easy") \(k = 2\) problem in Sec. \[III\]. Here, we derive the complete phase diagram using transfer matrix techniques introduced by Bravyi \[1\]. In the process, we discover that the classical geometry of the 2-QSAT interaction graph determines the generic satisfiability of random instances without reference to the random quantum Hamiltonian imposed upon the graph. We further show that our random ensemble asymptotically satisfies a technically important constraint known as the promise gap with probability exponentially close to 1.

In Sec. \[IV\] we move on to establish rigorous bounds on the existence of SAT and UNSAT phases for the ("hard") \(k \geq 3\) problem. Although the proof is quite different from the analysis for \(k = 2\), we then directly recover the unexpected deduction that the satisfiability of a generic instance of k-QSAT reduces to a classical graph property. This allows us to more tightly bound the UNSAT phase transition. We comment briefly on the straightforward generalizations to some related ensembles and on the satisfaction of the promise gap. Finally, in Sec. \[V\] we lay out the goals for the continued study of random quantum problems and algorithms.

II. DEFINITION OF RANDOM QSAT

We consider a set of \(N\) qubits. We first randomly choose a collection of \(k\)-tuples, \(\{I_m, m = 1 \ldots M\}\) by independently including each of the \(\binom{N}{k}\) possible tuples with probability \(p = \alpha / \binom{N}{k}\). This defines an Erdős hypergraph with \(\alpha N\) expected edges; we exhibit simple examples of these for \(k = 2, 3\) in Fig. \[I\]. In classical \(k\)-SAT, the next step is to generate an instance of the problem by further randomly assigning a Boolean \(k\)-clause to each \(k\)-tuple of the hypergraph. In key contrast to the classical case, where the Boolean variables and clauses take on discrete values – true or false – their quantum generalizations are continuous: the states of a qubit live in Hilbert space, which allows for boolean combinations of \(|0\rangle\) ("false") and \(|1\rangle\) ("true"). Thinking of a Boolean clause as forbidding one out of \(2^k\) \(k\)-tuples, there is a continuous variable \(\alpha\) in each \(k\)-qubit system.

The sum of these projectors defines a positive semidefinite Hamiltonian \(H = \sum_{m=1}^{M} \Pi_{\phi_m}^{I_m}\) and the decision problem for a given instance is, essentially, to ask if there exists a state that simultaneously satisfies all of the projectors, i.e. to determine whether \(H\) has ground state energy, \(E_0\), exactly zero. The qualifier “essentially” is needed because \(E_0\) is a continuous variable and in order that the problem be checkable by a quantum verifier (and therefore in QMA\(_1\)), \(H\) must be accompanied by a promise that \(E_0\) is either exactly zero or exceeds an \(\epsilon \sim O(N^{-\alpha})\) where \(\alpha\) is a constant. The scaling function \(\epsilon\) is also known as the promise gap. For our random ensemble we wish to compute the statistics of this decision problem as a function of \(\alpha\). Specifically we would like to know if there are phase transitions, starting with a SAT-UNSAT transition, in the satisfying manifold as \(\alpha\) is varied. Additionally, we would like to check that the statistics in the large \(N\) limit are dominated by instances that automatically satisfy the promise gap.

We now note a few key properties of the hypergraph ensemble: At small clause density \(\alpha\), the size distribution of connected components (“clusters”) is exponential. Moreover, almost all of these clusters are treelike with a few, \(O(N^\alpha)\), containing a single closed loop (see Fig. \[I\]); the number of clusters with several closed loops vanishes as \(N \to \infty\). Above a critical value \(\alpha_{gc}(k) = 1 / (k(k-1))\), a giant component emerges on which a finite fraction of the qubits reside. Unlike the finite clusters, the giant component may contain a non-vanishing (hyper)core, defined as the subgraph remaining after recursively stripping away leaf nodes (i.e. nodes of degree 1). For \(k = 2\), an extensive hypercore emerges continuously at \(\alpha_{hc}(2) = \alpha_{gc}(2) = 1/2\); for \(k \geq 3\), the hypercore appears abruptly with a finite fraction of the nodes at \(\alpha_{hc}(k) = \alpha_{gc}(k)\) \[10\] \[11\].

III. 2-QSAT

Let us first consider our questions in the specific context of 2-QSAT (phase diagram in Fig. \[2\]). While this is a classically easy (P) problem, the random ensemble has much of the structure we also find for the harder \(k > 2\) case. An instance \(H\) of QSAT is satisfiable if its kernel \(\ker(H)\) has nonzero dimension; in the following, we will find the kernel of \(H\) explicitly by transforming the random projector problem almost surely into a simple Heisenberg ferromagnet, whose ground state space is well known.

A central tool in this analysis is Bravyi’s transfer matrix \(T_\phi\) which, given a vector \(|\psi_i\rangle\) of the state of qubit \(i\) yields \(|\psi_j\rangle = T_\phi^{ij}|\psi_i\rangle\) for qubit \(j\), such that the product state \(|\psi_{ij}\rangle = |\psi_i\rangle \otimes |\psi_j\rangle\) satisfies the projector onto \(\phi\): \(\langle \psi_{ij}|T_\phi^{ij}|\psi_{ij}\rangle = 0\). Concretely, if the projector penalizes a joint state of both qubits given by the complex vector \(\phi = (\phi_{00}, \phi_{01}, \phi_{10}, \phi_{11})\), we have \(T_\phi = \epsilon\phi^\dagger = (\epsilon \phi_{00}^* \phi_{11}^* - \epsilon \phi_{01}^* \phi_{10}^*)\). Here \(\epsilon\) is the standard antisymmetric matrix (Levi-Civita symbol) in two dimensions. We note that the transfer matrix \(T_\phi\) for any given link is almost surely invertible so this construction finds satisfying product states for any input \(|\psi_i\rangle\).
A. Trees are SAT

Consider the clusters that enter the 2-QSAT graph ensemble. Of these, a tree comprising $n$ qubits and $n-1$ edges has a satisfying product state $|\Psi\rangle = \bigotimes_{j=1}^{n} |\psi_j\rangle$, where $|\psi_j\rangle$ is obtained from an arbitrary reference qubit $i=1$ by repeated application of the $T$’s along the (unique) path joining $i$ with $j$. In fact, the satisfying subspace is, almost always, $n+1$ dimensional. To show this, we will map the random projector problem directly onto the ferromagnetic Hamiltonian on the same tree. In more mathematical terms, we construct a non-unitary action of the permutation group on the qubit Hilbert space $\mathcal{H}$ invariant. Thus, the ground state space is precisely the totally symmetric space $\text{Sym}^n\mathbb{C}^2$, which has dimension $n+1$.

We define a particular, non-orthogonal product basis for the Hilbert space $\mathcal{H} = \mathcal{H}_0 \otimes \mathcal{H}_1 \otimes \cdots \otimes \mathcal{H}^{n-1}$. First, choose any two linearly independent unit vectors $|\uparrow^0\rangle$ and $|\downarrow^0\rangle$ as a basis for $\mathcal{H}_0$. Second, use the transfer matrix $T_{ij}$ to transfer this basis of $\mathcal{H}_0$ to a normalized basis

$$
|\uparrow^j\rangle = \frac{T_{ij} |\uparrow^i\rangle}{\|T_{ij} |\uparrow^i\rangle\|}
$$

$$
|\downarrow^j\rangle = \frac{T_{ij} |\downarrow^i\rangle}{\|T_{ij} |\downarrow^i\rangle\|}
$$

(1)

for each of the neighbors $j$ of 0. Finally, recursively traverse the tree to produce a pair of vectors $|\uparrow^i\rangle, |\downarrow^i\rangle$ for each site $i$ in the tree. This produces a choice of basis for each of the individual qubit Hilbert spaces which we use to define a product basis we call the transfer basis. Although we use the notation of spin up and down, we emphasize that the states need not be orthogonal.

Finally, we show that the ground state space of the Hamiltonian $H$ is precisely the space of totally symmetric states defined with respect to the transfer basis. We consider the constraint that a generic vector $\Psi \in \mathcal{H}$ is annihilated by $\Pi^{01}$ by factoring $\Psi$: $|\Psi\rangle = |\uparrow^0 \downarrow^1\rangle |\uparrow^2 \ldots \uparrow^{n-1}\rangle + |\downarrow^0 \uparrow^1\rangle |\downarrow^2 \ldots \downarrow^{n-1}\rangle$.

To show this, we start with $\Pi^{01} |\Psi\rangle = |\Psi\rangle$. This follows trivially from the construction of the transfer basis for the first two terms while for the third term, keeping track of indices carefully (summation of the transfer basis for the first two terms while for the third term, keeping track of indices carefully (summation of the transfer basis for the first two terms while for the third term), we rewrite $|\Psi\rangle$ as $\sum_{\sigma} \phi_\sigma (|\uparrow^0 \downarrow^1\rangle |\downarrow^2 \ldots \downarrow^{n-1}\rangle + |\downarrow^0 \uparrow^1\rangle |\uparrow^2 \ldots \uparrow^{n-1}\rangle)$.

(2)

where we have exploited the antisymmetry of $\epsilon$. Thus, $|v_4\rangle$ must be zero and $\Psi$ clearly lies in the symmetric eigenspace for swaps $(01)$. A similar argument holds for each of the swaps $(ij)$ on links of the tree $B$. Since $B$ is connected, these swaps generate the full permutation group $S_n$ and $\Psi$ must be in the completely symmetric subspace of its action. In particular, this means that the ground state space is isomorphic to $\text{Sym}^n\mathbb{C}^2$ which is $n+1$ dimensional.
B. Loops

To understand hypergraphs with loops, we first look for satisfying product states. For a cluster $G$ comprising $nL$ independent closed loops \( i.e. \) $n$ qubits with $n + nL - 1$ edges, see Fig. 4, a product state is internally consistent only if the product of the $T$'s around each closed loop returns the state it started with. For a graph with a single closed loop, $O_1$, this imposes

\[
\left( \prod_{jk \in O_1} T^{jk} \right) |\psi_i\rangle = \lambda_1 |\psi_i\rangle,
\]

which in general allows two solutions, so that these graphs are SAT. By contrast, if a site $i$ is part of a second independent closed loop, $O_2$, the additional demand

\[
\left( \prod_{jk \in O_2} T^{jk} \right) |\psi_i\rangle = \lambda_2 |\psi_i\rangle
\]

already yields an overconstraint. Thus graphs with more than one closed loop lack satisfying product states with probability $1$.

In fact, we can generalize the above reasoning to show that it holds even for the entangled (non-product) states in the problem as follows. We choose a spanning tree on $G$ and starting qubit $0$ to define a transfer basis – the ground state space is necessarily a subspace of the associated symmetric space. In particular, if $G$ contains a single closed loop $O_1$, we take qubit $0$ on the loop. In this case, we choose the starting basis on qubit $0$ to satisfy the product state consistency condition for $O_1$

\[
\left( \prod_{m \in O_1} T^m \right) \psi^0 = \lambda \psi^0,
\]

\( i.e. \) we take the eigenbasis of the loop transfer matrix, which we denote $|↑0, ↓0\rangle$ with eigenvalues $\lambda_1, \lambda_↓$. These two states transfer to two linearly independent product ground states for the loop – all up and all down. A short calculation verifies that in fact there are no more: write a generic $|\Psi\rangle \in \ker(H) \subset \text{Sym}^n \mathbb{C}^2$ as

\[
|\Psi\rangle = |↑−1, ↓0\rangle |v_3\rangle + |↓−1, ↑0\rangle |v_3\rangle + \left(|↑−1, ↓0\rangle + |↓−1, ↑0\rangle\right) |v_3\rangle,
\]

where $-1$ is the qubit at the end of the loop $O_1$ and $v_i$ are vectors on the $n-2$ other qubits. The projection $\Pi^{(-1,0)}$ annihilates the first two terms by choice of the basis. However:

\[
\Pi^{(-1,0)} |\Psi\rangle = \left|\phi^{(-1,0)}\right\rangle \left(|↑−1, ↓0\rangle + |↓−1, ↑0\rangle\right) |v_3\rangle
\]

\[
= |\phi^{(-1,0)}\rangle \left(\lambda_↓^{|↑0\rangle ϵ |↓0\rangle} + \lambda_↑^{|↓0\rangle ϵ |↑0\rangle} \right) |v_3\rangle
\]

Since $\lambda_↑ \neq \lambda_↓$ w.p. $1$, $v_3$ must be zero and by symmetry, $\psi$ must be in the span of the all up and all down states.

An analogous calculation verifies that any additional closed loops introduce projectors that are violated on this two-dimensional subspace and thus graphs with more than one loop are unsatisfiable.

It is worthwhile to consider how the usual Heisenberg ferromagnet fits into the above calculations. In this case, the transfer matrices are all the identity and thus all of the loop constraint conditions are identically satisfied. From the point of view of Eq. [3], the ferromagnet has $\lambda_↑ = \lambda_↓ = 1$ and there is no reduction in the ground state degeneracy due to the introduction of loop-closing ferromagnetic bonds.

C. Phase diagram

In light of the above analysis, the existence of a SAT/UNSAT phase reduces to the presence of multiply connected components in the ensemble of 2-SAT graphs (rather than the combined ensemble of graphs and projectors). For $\alpha < \alpha_{gc} = 1/2$, the number of such clusters vanishes in the limit of large $N$, so any instance is SAT with probability that goes to $1$ as $N \to \infty$. At $\alpha_{gc} = 1/2$, closed loops proliferate as a giant component appears and thereafter all instances are UNSAT with probability $1$. A straightforward upper bound on the energy in the UNSAT phase, $E \leq O(\langle \alpha - \alpha_{gc}\rangle^2 N)$, follows from the fact that the fraction of sites in the core of the giant component grows as $(\alpha - \alpha_{gc})^2$.

Physical intuition suggests that the ground state energy should be extensive above the transition and thus likely saturate the given upper bound. In the next section, we adduce strong evidence for this by exhibiting a minimal lower bound on the energy $E > N^{-1-\epsilon}$ for any $\epsilon > 0$ which holds except for exponentially rare instances. This follows from the twin claims that (i) the energy of a figure eight comprising $n$ sites decays only polynomially with $n$, and (ii) that the number of disjoint figure eights in the random graph grows nearly linearly with $N$. Observe that this kind of lower bound is exactly what we need to establish that our ensemble keeps the promise that either $E = 0$ or $E > 1/N^a$ with probability ex-
We expect on physical grounds that the UNSAT phase of $k$-QSAT has extensive ground state energy with relatively vanishing fluctuations for any $k$. In this case, the promise that $E \geq O(N^{-a})$ fails to be satisfied only with exponentially small probability by Chebyshev’s inequality. More generally, so long as the average ground state energy is bounded below by a polynomially small scale $E \geq O(N^{-b})$ with relatively vanishing fluctuations, the promise will be violated with only exponentially small probability for $a > b$.

Placing rigorous lower bounds on the expected quantum mechanical ground state energy is generally difficult, but at least for $k = 2$, we argue as follows to find a nearly extensive bound scaling as $N^{1-\epsilon}$ for any $\epsilon > 0$. As the Hamiltonian for QSAT is a sum of nonnegative terms, we can bound the ground state energy from below by considering manageable subgraphs and ignoring the contribution from other terms. In particular, the average ground state energy of a figure eight graph, that is a loop of length $L$ with one additional crossing edge, is polynomially bounded below by $O(L^{-\delta})$. This already gives a polynomial lower bound on the expected energy simply by allowing $L$ to scale as $N$ and knowing that we will find at least one such subgraph in the giant component with probability exponentially close to 1. We will do better by finding a large number of disjoint such figure eight graphs.

The expected number of subgraphs $A$ in the random graph $G_{N,p}$ is given by the following formula:

$$E\#(A \subset G) = \frac{N!}{(N-|A|)! |Aut(A)|} e^{p\#(A)}$$  \hspace{1cm} (7)

where $|A|$ is the number of vertices in $A$, $e(A)$ is the number of edges in $A$ and $Aut(A)$ is the group of automorphisms of $A$ ($|Aut(A)|$ is the cardinality of this set). This formula simply counts the number of ways of finding permutations of $|A|$ nodes in $G$ and connecting them up into an $A$ subgraph. For the fixed clause density ensemble, we take $p = 2\alpha/N$.

A figure eight graph is uniquely specified by giving its size $L$ (we take $L$ even) and the distance $d = 2\ldots L/2$ between the two nodes that are connected by the crossing link. We let $A$ be the disjoint union of $K$ figure eight graphs with $L$ nodes each and cross bars at separation $L/2 - 1$. Such a graph has $|Aut(A)| = K!2^K$ from the $K!$ permutations of the disjoint subgraphs and the two-fold symmetry of each figure eight. Thus, the expected number of $K$-fold disjoint figure eights of size $L$ is

$$E\# = \frac{N!p^{K(L+1)}}{(N-KL)!K!2^K}$$ \hspace{1cm} (8)

We now allow $K,L$ to scale with $N$ such that $KL \ll N$ and use Stirling’s formula to find the asymptotic entropy

$$S \approx KL \left( \log 2\alpha - \frac{KL}{N} \right) + K (\log \alpha + 1 - \log K - \log N).$$ \hspace{1cm} (9)

This entropy is positive and growing with $N$ so long as $L \gg \log N$, $KL \ll N$ and $\alpha > \alpha_g = 1/2$. In this regime, assuming that the fluctuations in $S$ are relatively small, we find that there are $K \sim N^{1-\epsilon}$ disjoint figure eights of size $L \sim N^{\epsilon/2}$ with exponentially high probability. These lead to a nearly extensive lower bound on the expected ground state energy.

## IV. $k$-QSAT

We now turn to $k$-QSAT with $k > 2$. As noted above, for sufficiently small $\alpha$, the important hypergraphs have vanishing cores. In this regime, it is possible to construct the hypergraph by adding in edges one by one where each additional edge brings with it at least one new leaf node. Thus by a generalization of the transfer matrix arguments for $k = 2$, it is possible to construct a satisfying product state on the full hypergraph; the details are in Sec. [IV A]. Deducing the existence of the UNSAT phase is not as straightforward however for, unlike in the $k = 2$ problem, we do not know a set of unsatisfiable graphs which are present with finite probability as $N \to \infty$; indeed these are not known for classical SAT either. Thus, in Sec. [IV B] we produce an indirect proof that the dimension of the kernel of $H$ vanishes for sufficiently large $\alpha$.

Recall that for 2-QSAT the dimension of the satisfying manifold ended up depending, with probability 1, only on the topology of the graph and not on the choices of the projectors. This is in fact a very general result, which we prove directly in Sec. [IV C]. For any hypergraph for $k$-QSAT (regardless of its likelihood) the dimension of the satisfying manifold is independent of the choice of projectors with probability 1. Moreover, non-generic choices of projectors result in larger satisfying manifolds. Implicitly, this means that satisfiability for the random ensemble is a graph theoretic property and raises the extremely interesting possibility that random $k$-QSAT can be formulated without reference to quantum Hamiltonians at all.

Finally, we comment on the satisfaction of the promise gap in Sec. [IV D] and on the generalization of QSAT to higher rank projectors in Sec. [IV E].

### A. Existence of SAT phase

We prove the existence of a satisfiable phase at low clause density for rank 1 $k$-QSAT by constructing product states of zero energy for arbitrary hypergraphs containing a hypercore with a satisfying product state. Since

...
the core of a random graph vanishes for \( \alpha < \alpha_{hc}(k) \), this proves the existence of a satisfiable phase below the emergence of a hypercore.

**Lemma 1.** Suppose that \( H \) is an instance of QSAT on \( N \) qubits with a satisfying product state \( |\Psi\rangle \). Let \( H' = H + \Pi \) be an instance of QSAT on \( N' > N \) qubits where \( \Pi \) is a QSAT projector touching at least one qubit not among the original \( N \). Then \( H' \) has a satisfying product state \( |\Psi'\rangle \).

**Proof.** Assume first that the hyperedge \( \Pi \) adjoins only one dangling qubit (untouched by \( H \)), say qubit 0. Let the components of \( |\Psi\rangle \) be \( \Psi_{i_1 \cdots i_N} \). The key observation here is that the projector \( \Pi \) with \( k-1 \) of its qubits fixed in a product state of \( N \) qubits uniquely specifies the state of the dangling qubit 0. If \( \phi_{i_0 \cdots i_{k-1}} \) is the vector onto which \( \Pi \) projects, then the state of qubit 0 must be perpendicular to \( \phi_{i_0 \cdots i_{k-1}} \Psi_{i_1 \cdots i_{k-1}} \).

Thus, contraction with \( e^{i\alpha_0} \) defines a \( k-1 \) to 1 transfer matrix \( T_{i_1 \cdots i_{k-1}}^{i_0} \) generalizing the Bravyi transfer matrix of 2-QSAT. The new product state given by \( |\Psi'\rangle = |\psi^{0}\rangle \otimes |\Psi\rangle \) where \( \psi^{0} = T_{i_1 \cdots i_{k-1}}^{i_0} \Psi_{i_1 \cdots i_{k-1}} \) satisfies \( \Pi \). Moreover, since \( H \) does not act on qubit 0, \( H |\Psi'\rangle = |\psi^{0}\rangle \otimes H |\Psi\rangle = 0 \). Thus \( H' = H + \Pi \) is satisfied by the product state \( |\Psi'\rangle \) on \( N + 1 \) qubits.

Finally, if \( \Pi \) adjoins more than one dangling qubit, simply fix all but one of these to an arbitrary state and then apply the above transfer matrix procedure.

**Theorem 2.** Suppose that \( H \) is an instance of random QSAT and let \( H_C \) be the restriction to its hypercore. If there is a satisfying product state on \( H_C \), then there is almost surely a satisfying product state on \( H \).

**Proof.** Let \( \Pi_i, i = 1 \cdots M(G) - M(C) \) be the sequence of projectors removed in the process of stripping leaf hyperedges from the hypergraph of \( H \) to produce the hypercore. At each stage of this process, the removed projector \( \Pi_i \) has at least one dangling qubit. Thus, if we iteratively reconstruct \( H \) from \( H_C \) by adding back the \( \Pi_i \) in reverse order, we can apply the lemma at each stage to lift the product state on \( H_C \) to a product state on \( H \).

**B. The UNSAT phase exists**

For a given random hypergraph, let us construct \( H \) via the sequence \( H_m = \sum_{j=1}^{M} \Pi_{j}^{i_j} \), i.e. we add the projectors one at a time in some order. Clearly \( H = H_M \). Let \( D_m \) be the dimension of the satisfying subspace for \( H_m \); evidently \( D_0 = 2^{N} \). At each step, if the next added projector involves a set of qubits that were not previously acted upon, then \( D_{m+1} = D_m(1 - \frac{1}{2^k}) \) as we may simply implement the projection by reducing the size of a basis that can be factorized between the target qubits and all others. It is intuitively plausible that this is the best we can do—in cases where the target qubits are entangled with the rest of the system we should expect to lose even more states, i.e.

\[
D_{m+1} \leq D_m(1 - \frac{1}{2^k})
\]

in general. A proof that this bound holds with probability 1 for projectors randomly chosen according to the uniform Haar measure is contained in Appendix A. From this we conclude that

\[
D_{mN} \leq 2^N \left(1 - \frac{1}{2^k}\right)^{\alpha N}.
\]

Thus, for \( \alpha > -1/\log_2 (1 - \frac{1}{2^k}) \sim 2^k \) the problem is asymptotically almost always UNSAT.

**C. Geometrization theorem**

**Geometrization Theorem.** Given an instance \( H \) of random \( k \)-QSAT over a hypergraph \( G \), the degeneracy of zero energy states \( \dim(\ker(H)) \) takes a particular value \( D \) with probability 1 with respect to the choice of projectors on the edges of the hypergraph \( G \).

**Proof.** For a fixed hypergraph \( G \) with \( M \) edges, \( H = H_0 = \sum_{i=1}^{M} \Pi_i = \sum_{i=1}^{M} |\phi_i\rangle \langle \phi_i| \) is a matrix valued function of the \( 2^k M \) components of the set of \( M \) vectors \( |\phi_i| \). In particular, its entries are polynomials in those components. Choose \( |\phi\rangle \) such that \( H \) has maximal rank \( R \). Then there exists an \( R \times R \) submatrix of \( H \) such that \( \det(H|_{R\times R}) \) is nonzero. But this submatrix determinant is a polynomial in the components of \( |\phi\rangle \) and therefore is only zero on a submanifold of the \( |\phi\rangle \) of codimension at least 1. Hence, with probability 1, \( H \) has rank \( R \) and the degeneracy \( \dim(\ker(H)) = 2^N - R \).

The theorem holds for general rank \( r \) problems as well by a simple modification of the argument to allow extra \( \phi \)'s to be associated to each edge. A nice corollary of this result is a more stringent upper bound on the size of the SAT phase. Consider any assignment of classical clauses on a given hypergraph: we can also think of this as a special instance of \( k \)-QSAT where the projectors are all diagonal in the computational basis. As this is a non-generic choice of projectors, the dimension of its satisfying manifold is an upper bound on the dimension for generic choices. We conclude then that the classical UNSAT threshold is an upper bound on the quantum threshold. Indeed, if we can identify the most frustrated assignment of classical clauses on a given hypergraph, i.e. the assignment that minimizes the number of satisfying assignments, we could derive an even tighter bound.

**Corollary 3.** The zero state degeneracy of the random quantum problem is bounded above (w.p.1) by the number of satisfying assignments of the most constrained classical \( k \)-SAT problem on the same graph.
We have encountered many numerical examples in which the quantum problem has fewer ground states than the most frustrated classical problem on the same hypergraph. Thus, the bound of corollary 3 is not tight.

In the $k = 2$ case, corollary 2 gives us another way to show the critical point corresponds to the emergence of a two-core: once a two-core exists, it will contain a giant loop with two crossing bonds. This graph can be made unsatisfiable by setting each of the clauses around the loop to disallow the state 01 and the crossing bonds to disallow 00 and 11 respectively. Thus, the quantum problem is UNSAT w.p. 1. Ideally, we could extend this geometric characterization to the $k > 2$ problem, but we have thus far failed to identify a family of unsatisfiable hypergraphs that appear in the random hypergraph with non-vanishing probability.

D. Satisfying the promise

As in the case of 2-QSAT, physical experience suggests that the ground state energy for $k$-QSAT should be extensive with small fluctuations above the satisfiability transition. If this is true, the promise is satisfied with probability exponentially close to 1, as in the $k = 2$ case. Unfortunately, we do not know a set of unsatisfiable subgraphs analogous to the figure eight’s that might be used to provide a more rigorous bound.

E. QSAT at rank $r > 2$

We now briefly consider the extension of $k$-QSAT to $(k,r)$-QSAT, in which the projectors $\Pi^I$ have rank $r$, i.e. they penalize a uniformly chosen $r$-dimensional subspace of the $2^k$ dimensional $k$ qubit Hilbert space. The main results regarding $k$-QSAT generalize naturally: satisfiability is almost surely only dependent on the underlying graph and the weak bound on the ground state degeneracy becomes $D_{\alpha N} \leq 2^N \left(1 - \frac{r}{2^k}\right)^{\alpha N}$, implying a bound $\alpha^{(r)} \leq -1/\log_2 \left(1 - \frac{r}{2^k}\right)$. However, there need not be a SAT phase at all: if $r = 2^k$, the projectors $\Pi^I$ are each the identity and the ground state energy is $\alpha N > 0$ for any positive $\alpha$. More generally, there is some critical rank $r_c$ above which the SAT phase disappears.

We bound $r_c$ above by $2^k/2$ by exhibiting an unsatisfiable subgraph of the random hypergraph that arises asymptotically almost surely ($N \to \infty$) in the random graph ensemble even at small $\alpha$. Consider a chain with two clauses and one shared qubit. Classically, this corresponds to a two-clause problem on $2k - 1$ bits where we allow each clause to forbid $r = 2^k/2$ configurations. Now let the first clause forbid all configurations in which the shared bit is 0 and the second clause all configurations in which it is 1. This classical problem is clearly unsatisfiable and therefore so is any quantum problem on this subgraph w.p. 1. Indeed, since there are extensively many such small chain components in the hypergraph, each with an independently chosen $O(N^0)$ ground state energy, this provides an extensive lower bound on the total ground state energy.

The bound $r_c \leq 2^k/2$ is not tight however. For example, for $k = 3$, an open chain of length four with rank 3 $< 2^3/2 = 4$ is quantum mechanically unsatisfiable w.p. 1, as can be checked numerically, and therefore so is the ensemble for $(k,r) = (3,3)$. However, there is no classically unsatisfiable problem on this chain and it is harder to construct a rigorous bound that scales with $k$ using this starting point.

V. CONCLUSIONS AND OUTLOOK

As in the classical work that was our inspiration, we expect that the phase structure that we have identified also shows up in the actual performance of potential quantum algorithms. Specifically, the classical results and the general intuition from the study of quantum phase transitions [13, 14] both strongly suggest that the decision problem will be hardest near the SAT-UNSAT boundary and easy away from such a bounded range. This expectation is given further force by the graph theoretic formulation of typical instances whereby we may again expect that graphs with an intermediate clause density are the ones where it might prove hardest to decipher the relevant property. As a first step in testing this expectation, we have begun an investigation of the simplest adiabatic algorithm for this problem [15] and the relevant excitation spectra.

Looking ahead, the immediate challenge is to explicitly identify the graph theoretic property that characterizes satisfiable instances of random QSAT without reference to the quantum mechanical problem. If this can be accomplished it will shed light on the difficulty of deciding generic instances of random QSAT and provide a surprising connection to classical complexity classes. Separately, we intend to investigate whether there are analogs of the clustering transitions of classical $k$-SAT in the quantum problem and whether generalizations of the cavity method to quantum problems [16, 17, 18] can help locate them. Finally, we would like to translate this improved understanding of this ensemble of problems into new algorithms for solving them on the lines of the be-
lief/survey propagation algorithms for classical SAT.

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[1] S. Bravyi, arXiv (2006), quant-ph/0602108.
[2] A. Y. Kitaev, in AQIP’99 (DePaul University, 1999).
[3] A. Y. Kitaev, A. Shen, and M. N. Vyalyi, Classical and quantum computation, vol. 47 of Graduate studies in mathematics (American Mathematical Society, Providence, R.I., 2002).
[4] M. R. Garey and D. S. Johnson, Computers and Intractability: A Guide to the Theory of NP-Completeness, Series of Books in the Mathematical Sciences (W. H. Freeman & Co Ltd. 1979).
[5] S. Kirkpatrick, B. Selman, Science 264, 1297 (1994).
[6] R. Monasson, R. Zecchina, S. Kirkpatrick, B. Selman, and L. Troyansky, Nature 400, 133 (1999).
[7] M. Mezard, G. Parisi, and R. Zecchina, Science 297, 812 (2002).
[8] F. Krzakala, A. Montanari, F. Ricci-Tersenghi, G. Semerjian, and L. Zdeborová, Proceedings of the National Academy of Sciences 104, 10318 (2007).
[9] F. Altarelli, R. Monasson, G. Semerjian, and F. Zamponi, in Handbook of Satisfiability (IOS press, 2009), vol. 185 of Frontiers in Artificial Intelligence and Applications, arXiv:0802.1829.
[10] M. Mézard, F. Ricci-Tersenghi, and R. Zecchina, J. Stat. Phys. 111, 505 (2003).
[11] M. Molloy, Random Struct. Algorithms 27, 124 (2005).
[12] R. Monasson and R. Zecchina, Phys. Rev. E 56, 1357 (1997).
[13] S. L. Sondhi, S. M. Girvin, J. P. Carini, and D. Shahar, Rev. Mod. Phys. 69, 315 (1997).
[14] S. Sachdev, Quantum Phase Transitions (Cambridge University Press, Cambridge, UK, 2001).
[15] E. Farhi, J. Goldstone, S. Gutmann, J. Lapan, A. Lundgren, and D. Preda, Science 292, 472 (2001).
[16] C. Laumann, A. Scardicchio, and S. L. Sondhi, Phys. Rev. B 78, 134424 (2008).
[17] M. B. Hastings, Phys. Rev. B 76 (2007).
[18] M. Leifer and D. Poulin, Ann. Phys. 323, 1899 (2008).
[19] QMA1 is the one-sided version of QMA, in which “yes” instances of problems always have proofs that are verifiable with probability 1, while “no” instances need only have bounded false-positive error rates. QMA allows the verifier to occasionally produce false-negatives as well. This small difference is not believed to significantly influence the difficulty of the classes.

APPENDIX A: WEAK UNSAT BOUND

Weak UNSAT Bound. Given an instance of random $(k, r)$-QSAT with Hamiltonian $H$ over a hypergraph $G$ with $N$ qubits and $M$ clauses, the degeneracy of zero energy states (i.e. the number of satisfying assignments) $D = \dim(\ker(H))$ is bounded above by

$$D \leq 2^N \left(1 - \frac{r}{2^k}\right)^M$$

(A1)

with probability 1.

Proof. For a given random hypergraph $G$, we construct $H = H_M$ via the sequence $H_m = \sum_{j=1}^m \Pi_j$, where $m = 1 \ldots M$. That is, we add projectors one at a time in some fixed order. Let $D_m$ be the degeneracy of zero energy states after $m$ projectors have been added. Clearly,

$$D_0 = 2^N.$$

(A2)

The result will follow by induction after we show that each additional projector reduces the degeneracy by at least a factor $1 - \frac{r}{2^k}$ w.p. 1. That is,

$$D_{m+1} \leq D_m \left(1 - \frac{r}{2^k}\right).$$

(A3)

We consider adding a projector $\Pi = \sum_{n=1}^r |\phi^\alpha\rangle \langle \phi^\alpha|$, where $\langle \phi^\alpha| \phi^\beta\rangle = \delta_{\alpha\beta}$, to the Hamiltonian $H$ to construct $H'$. By appropriate reordering, we can always assume that $\Pi$ acts on the first $k$ qubits of the full $N$ qubit Hilbert space. We choose a basis for the zero energy subspace $\ker(H)$ prior to the addition of $\Pi$ as follows

$$|d\rangle = \sum_{i=1}^L |a^i_d\rangle |l^i\rangle$$

(A4)

where $d \in \{1, \ldots, D\}$ labels the basis states, $|a^i_d\rangle \in \mathbb{C}^k$ are vectors in the first $k$ qubit factor of the Hilbert space and $|l^i\rangle$ are $L$ linearly independent vectors on the remaining $(N-k)$ qubit factor.

The kernel of $H' = H + \Pi$ is the subspace of the kernel of $H$ which is annihilated by $\Pi$. We write a generic vector of $\ker(H)$ as:

$$|\psi\rangle = \sum_{d=1}^D \psi_d |d\rangle.$$

(A5)

Annihilation by $\Pi$ leads to the condition

$$0 = \Pi |\psi\rangle = \sum_{i=1}^L \sum_{\alpha=1}^r \langle \phi^\alpha| I^i \rangle \sum_d \langle \phi^\alpha| a^i_d\rangle \psi_d.$$

(A6)
which by linear independence of the $|\phi^\alpha\rangle |\tilde{t}\rangle$ requires, for any $\alpha, i$:

$$0 = \sum_{d=1}^{D} \langle \phi^\alpha | a^\dagger_i d \rangle \psi_d. \quad (A7)$$

That is, $\psi_d$ must lie in the kernel of the $rL \times D$ matrix $A_{(\alpha)d} = \langle \phi^\alpha | a^\dagger_d \rangle$. We now claim that $A_{(\alpha)d}$ has rank $R$ at least $rD/2^k$ with probability 1, which will prove our inductive step. This follows from two observations:

1. The rank $R$ of $A$ is a bounded random variable that takes on its maximal value over the choice of $|\phi^\alpha\rangle$ with probability 1. $A$ can be viewed as a matrix of monomials in the components of $|\phi^\alpha\rangle$. Choose an orthonormal frame $|\phi^\alpha\rangle$ maximizing the rank $R$ of $A$; with this choice, there exists some $R \times R$ submatrix of $A$ such that $\det(A_{R \times R})$ is nonzero. But this submatrix determinant is a homogenous polynomial in the components of $|\phi^\alpha\rangle$ and therefore is only zero on a submanifold of codimension at least 1 in the complex Grassmannian $\text{Gr}(r, 2^k)$ (i.e. the space of choices of rank $r$ projectors $\Pi$). Hence, almost every matrix $A$ will have maximal rank $R$.

2. We now need only exhibit one set of $|\phi^\alpha\rangle$’s such that $A$ has rank $R \geq rD/2^k$. This is easy: at least one $R$-tuple of the standard basis vectors will provide such an $A$. Consider the matrix of vectors $|a^\dagger_d\rangle$ and let $r_1, r_2, \cdots, r_{2^k}$ be the ranks of each of the $2^k$ component matrices (i.e. the matrices $A$ obtained by using the standard basis elements for $|\phi\rangle$). The $r_j$ are the number of linearly independent rows in each of these matrices. Now concatenate each of these component matrices vertically into a giant $2^k L \times D$ matrix. The row rank of this matrix can be no larger than $r_1 + r_2 + \cdots + r_{2^k}$ by construction, but there must be a full $D$ linearly independent columns: any linear relation on the columns of the giant matrix corresponds to a linear relation $\sum_d w_d |a^\dagger_d\rangle = 0$ which lifts trivially to a linear relation among the basis vectors $|d\rangle$. Hence,

$$r_1 + r_2 + \cdots + r_{2^k} \geq D.$$

From this relation, there must exist some collection $r_i$ of size $r$ such that $r_1 + \cdots + r_i \geq rD/2^k$. This collection of basis vectors provides the frame we desire.

\[\square\]