

Entanglement Distribution on Complex Networks

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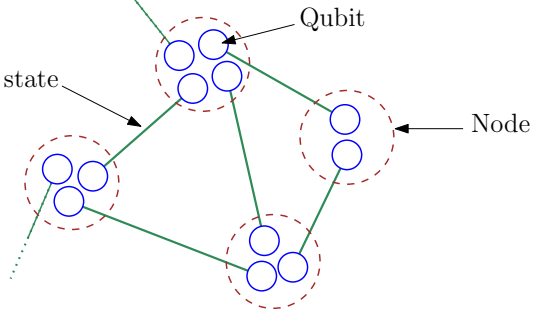
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Bi-partite entangled state



- *qubits* (two-level quantum systems) (spin $1/2$, e.g. photon polarization)
- Multiple qubits and classical resources at each *node* (vertex)
- *links* (edges): bi-partite entangled (pure/mixed) two-qubit states.
- Goal: entangle pairs of qubits between distant nodes
- Quantum operations local: within nodes. Classical can be global.
Local Operations and Classical Communication **LOCC**

Entanglement: Two entangled qubits



Two entangled qubits: four-dimensional Hilbert space.

Bi-partite pure state

All such states LOCC equivalent to unique state in Schmidt basis.

$$|\alpha\rangle = \sqrt{\alpha_0} |00\rangle + \sqrt{\alpha_1} |11\rangle$$

$$\alpha_0 > \alpha_1 \quad \alpha_0 + \alpha_1 = 1 \quad \alpha_1 \in [0, 1/2]$$

Pure, partially entangled, bipartite state

$\alpha_1 = 0$: no entanglement, $\alpha_1 = 1/2$: max. entanglement

Bell State: Singlet Conversion



Partially Entangled: $|\alpha\rangle = \sqrt{\alpha_0} |00\rangle + \sqrt{\alpha_1} |11\rangle$

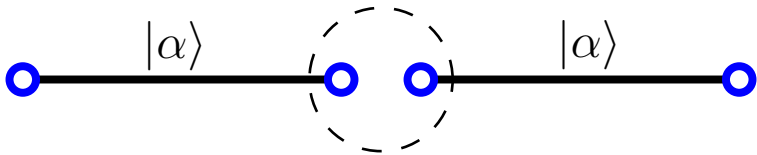
Local operations (and classical communication): qubits not allowed to interact

Maximally Entangled: $|\Psi\rangle = \frac{1}{\sqrt{2}} |00\rangle + \frac{1}{\sqrt{2}} |11\rangle$

Singlet, Bell State, Maximally Entangled State **Singlet Conversion Probability** $p = 2\alpha_1$, for $\alpha_0 > \alpha_1$

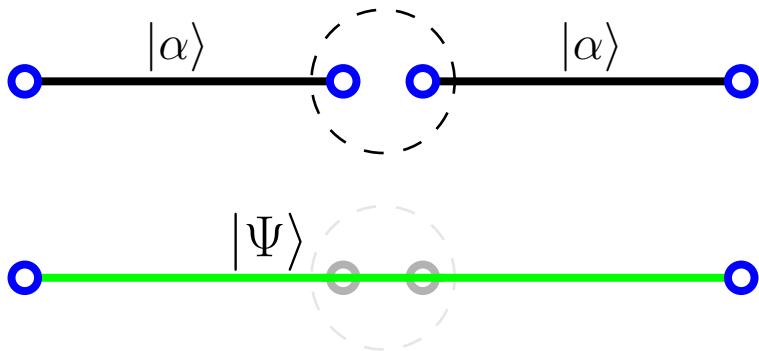
Otherwise: product state (failure)

Entanglement Swapping



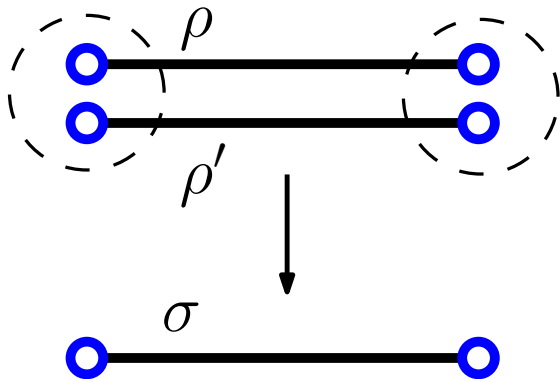
We can entangle the two outermost qubits, using only local operations and classical communication: *i.e.* without interacting outermost qubits. Using **entanglement swapping**.

Entanglement Swapping



Entanglement Swapping. Get **Bell state** with same probability as in singlet conversion $p = 2\alpha_1$! (product state otherwise) Note: if $\alpha_1 = 1/2$, then $p = 1$.

Mixed (or pure) states: Entanglement Purification

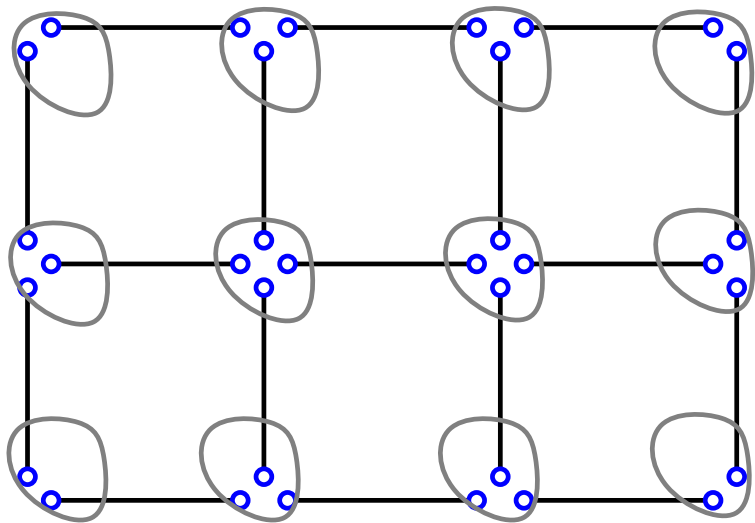


Obtain σ with entanglement greater than ρ, ρ' using LOCC
(local operations and classical communication)

Technical motivation: Generalize one-dimensional networks

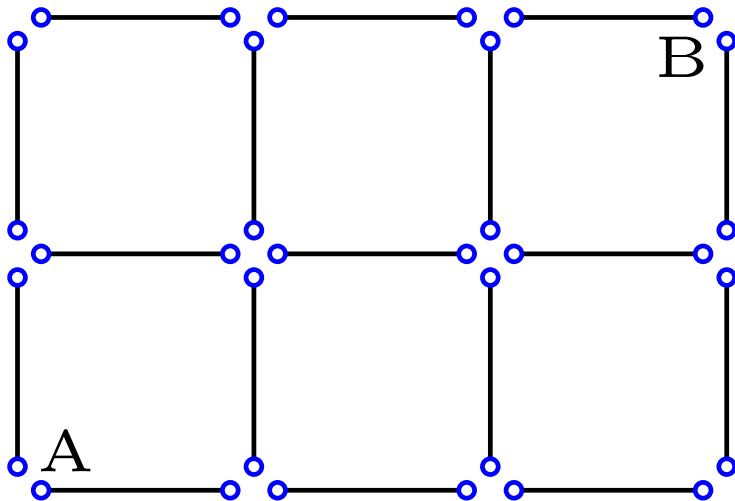
- Quantum Information: Entanglement is a resource for tasks: teleportation, key distribution, fault tolerant computation
 - Creating entanglement requires local interaction. Noise increases with distance. Depolarization. Absorption. **Can't distribute entanglement over long distance in a single stage!**
- Long range entanglement via Network of stations or nodes that store and purify a state.
 - Generalization of **quantum repeater** schemes.
Dür, Briegel, Cirac, Zoller, PRA 1999
 - Nodes share partially entangled states of qubits
 - Nodes(stations)/channels, Vertices/edges, Sites/bonds
 - Quantum operations **probabilistic**
 - Large number of random components \Rightarrow **Complex Networks, Percolation, Phase transition**

Quantum Network



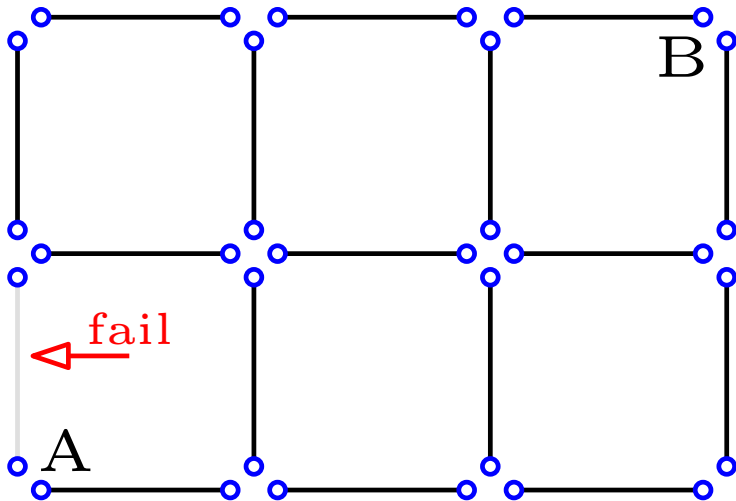
Concrete: Square lattice. Each bond is an entangled pair with amount of entanglement α_1 .

“Classical” Entanglement Percolation

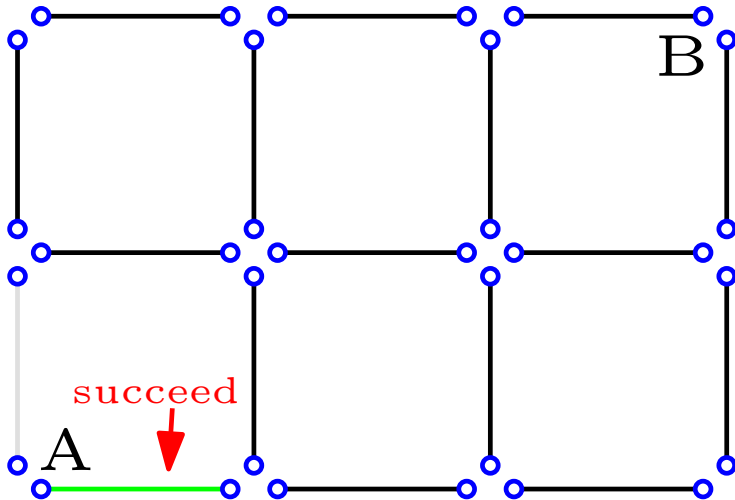


Entangle nodes A and B

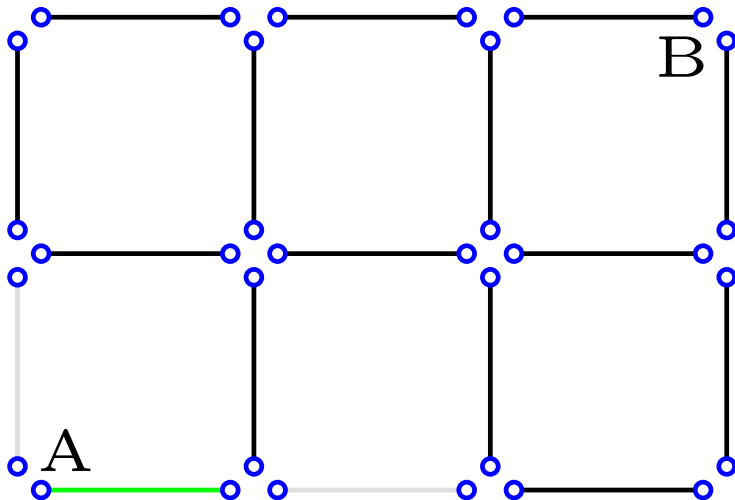
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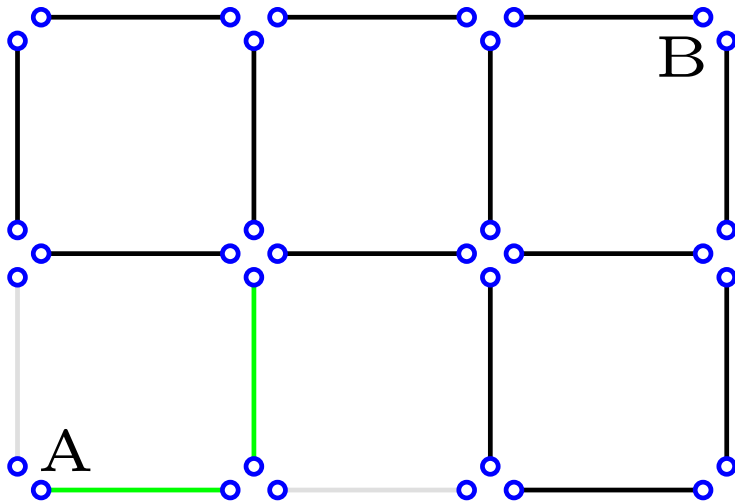
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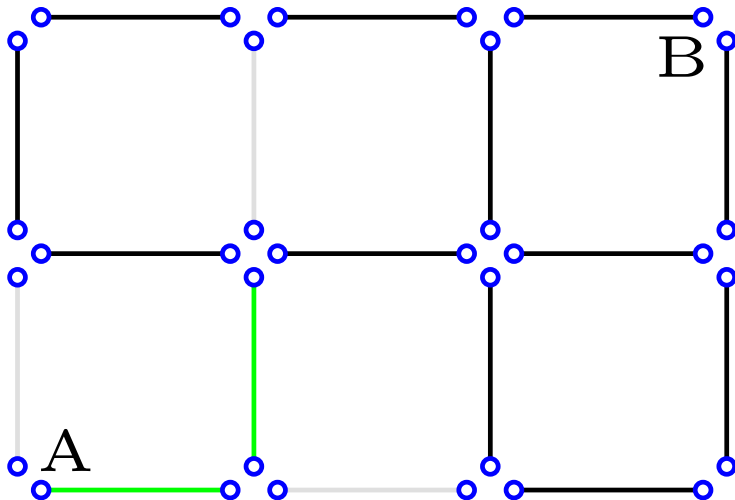
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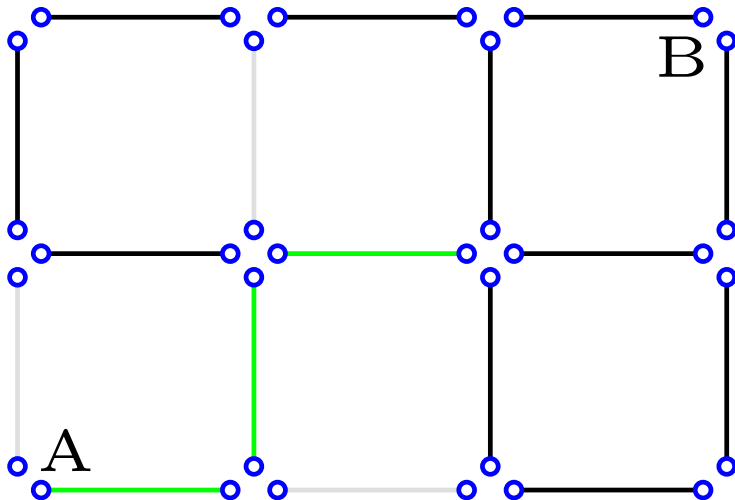
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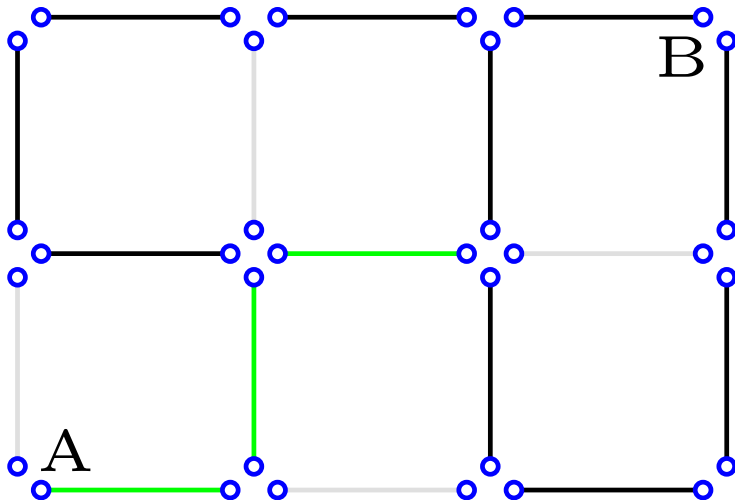
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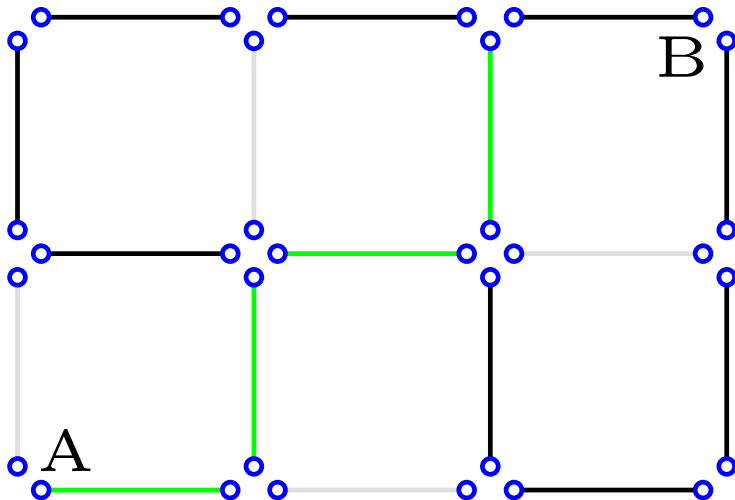
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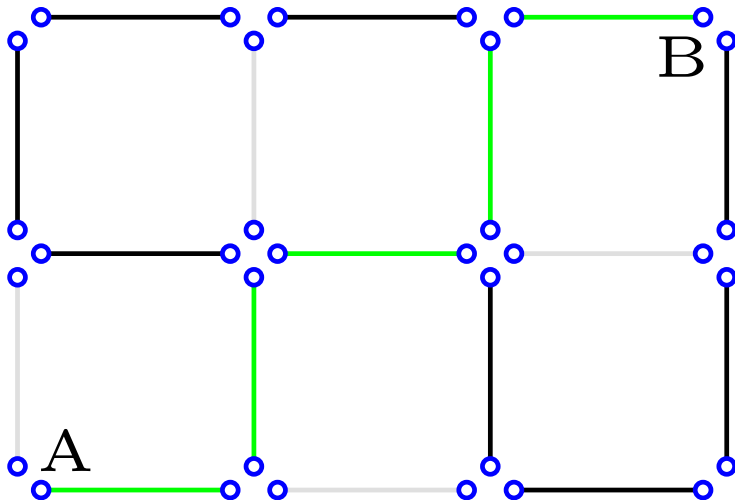
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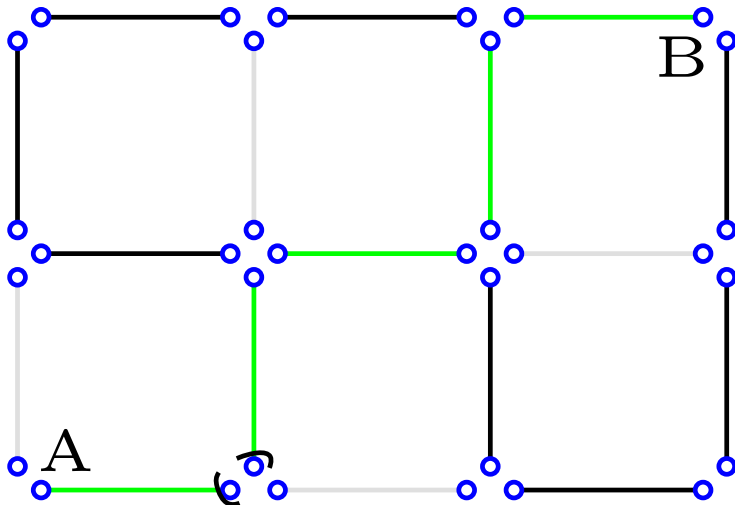
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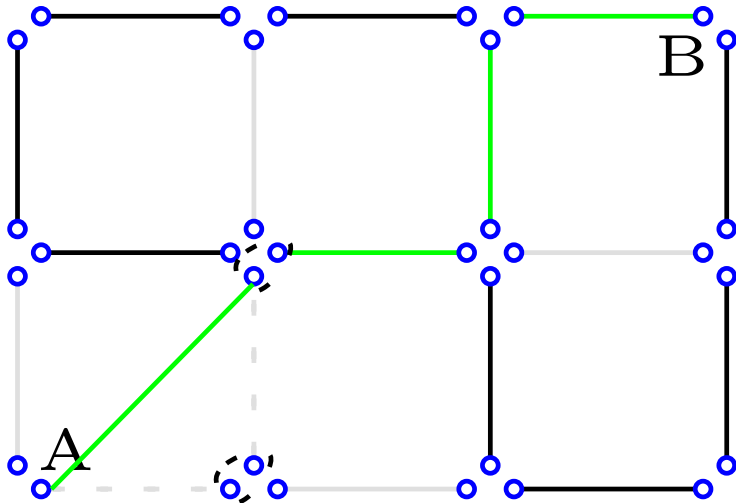
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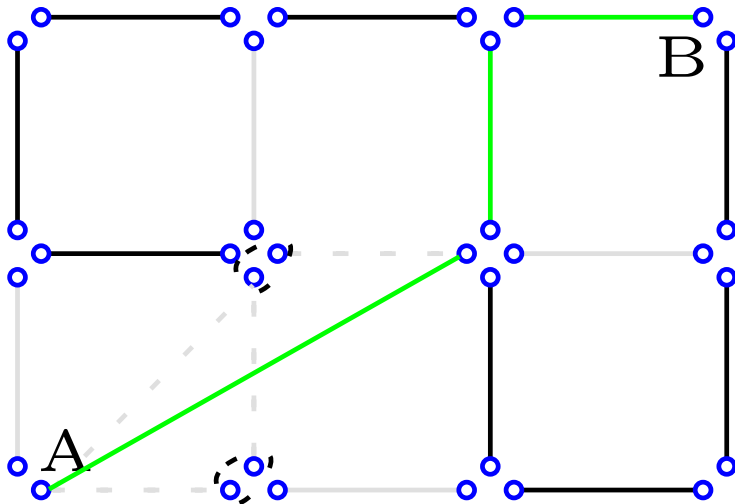
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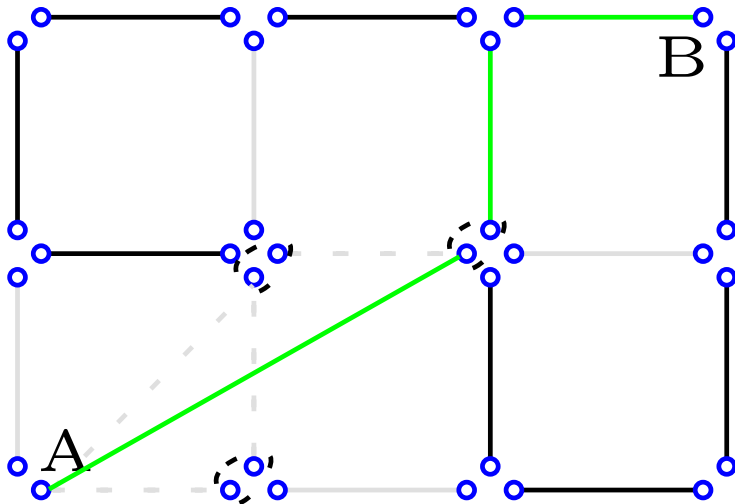
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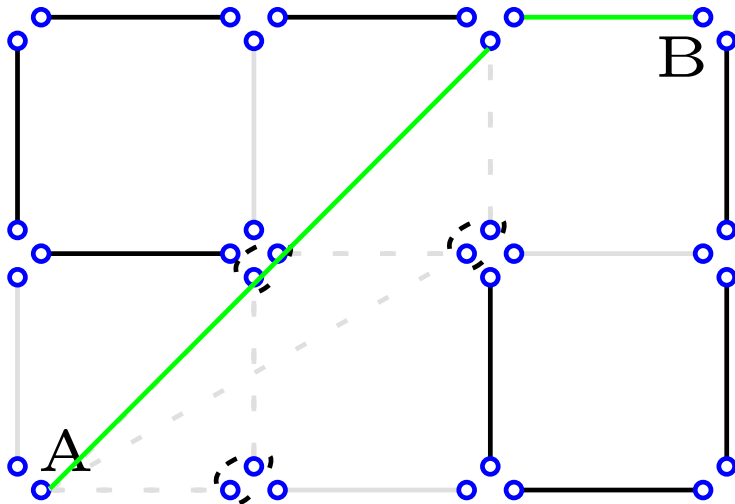
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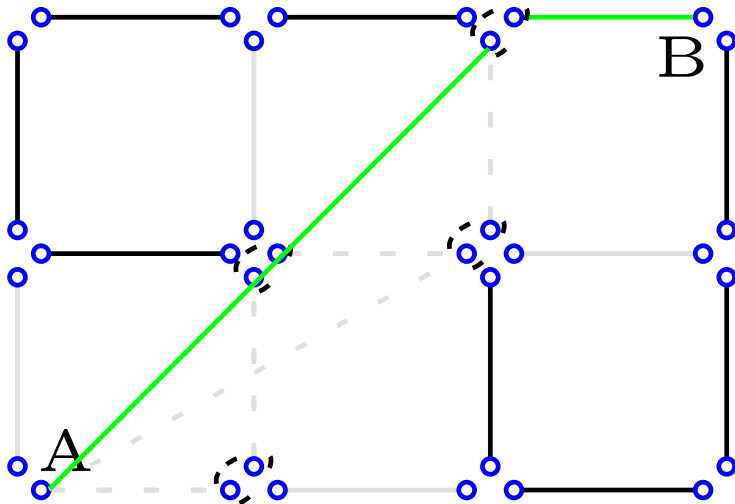
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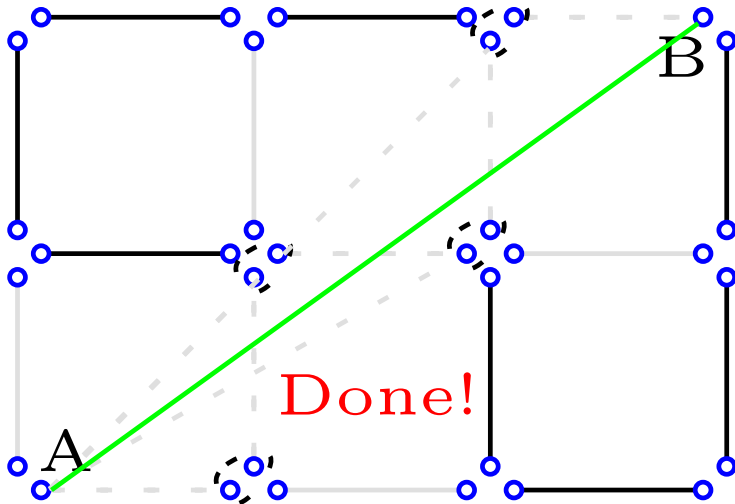
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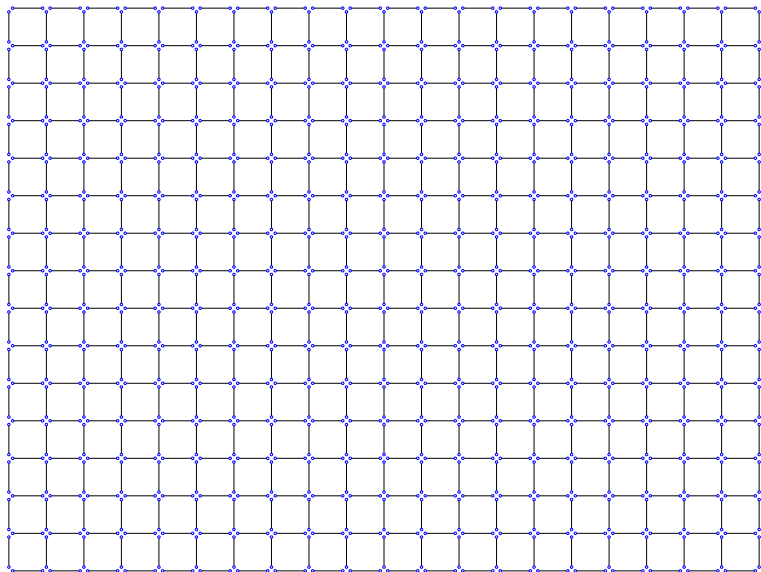
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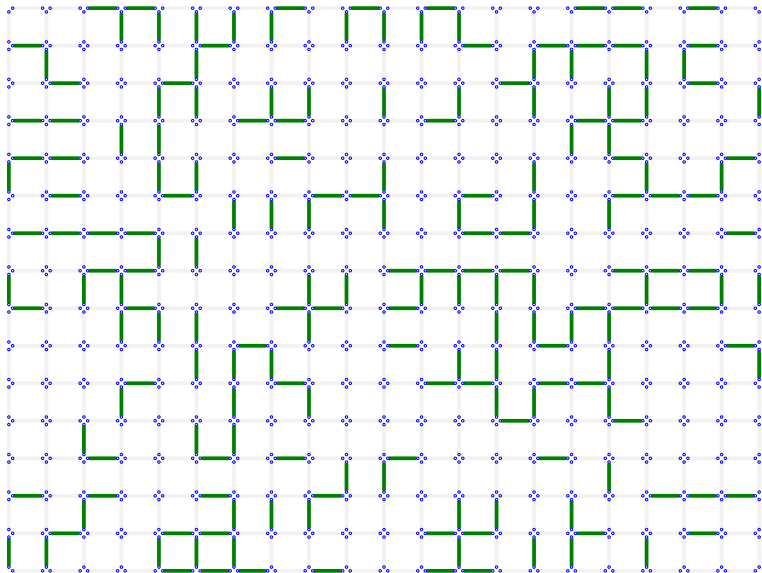
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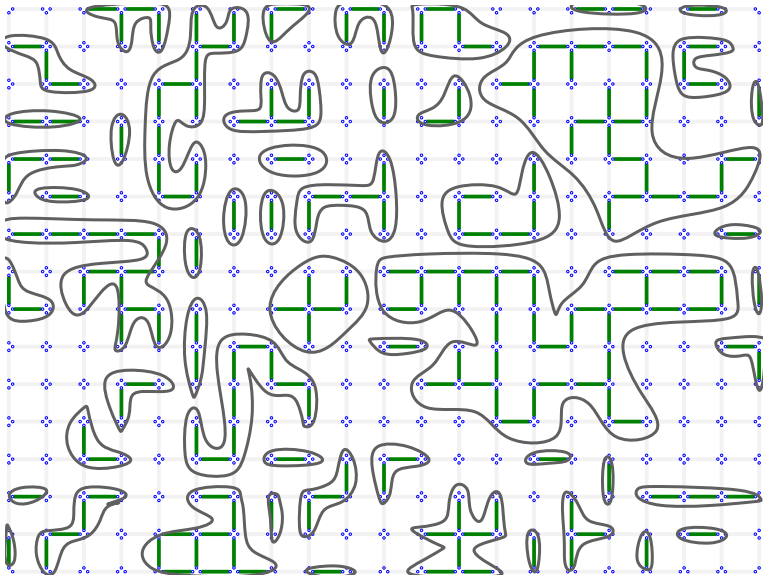
Big Network: $\alpha_1 = 0.175$ $p = 0.35$



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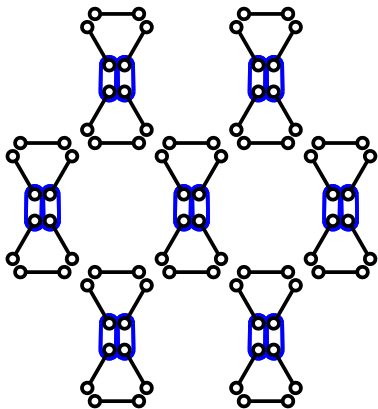


Big Network: $\alpha_1 = 0.175$ $p = 0.35$

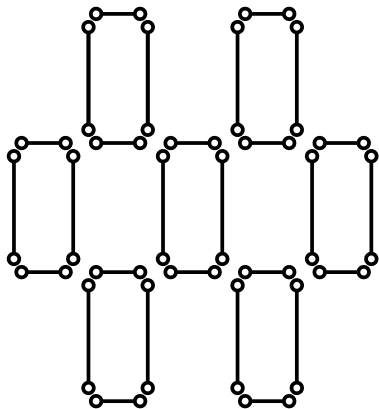


Percolation structure. Infinite cluster iff $p > p_c = 0.5$ on square lattice.

Kagome lattice to Square lattice



$p_c \approx 0.52$ Kagome



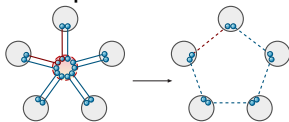
$p_c = 0.5$ Square lattice

Acín, Cirac, Lewenstein, Nature Phys (2007)

Perseguers, Cirac, Acín, Lewenstein, Wehr, PRA (2008)

JL, Wehr, Lewenstein, PRA (2009)

- Multipartite (GHZ) states \Rightarrow percolation non-planar graphs
Perseguers, Cavalcanti, JL, Lewenstein, and Acín, PRA (2010)
- Swapping: Project onto larger subspace. Conditionally complete swapping.
- Mixed states of rank ≤ 3 *Broadfoot, Dorner, Jaksch, PRA (2010), EPL (2009)*
- Q-star transformation on complex networks. increase



entanglement distance.

Cuquet,

Calsamiglia, PRL (2009), PRA (2011)

- Review: *Perseguers, JL, Cavalcanti, Lewenstein, Acín, Rep. Prog. Phys. 76 (2013)*

Let's leave these and **move to full-rank mixed states and complex networks.**

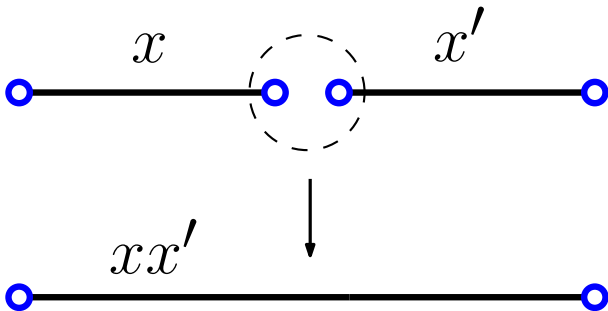
Two-qubit Mixed States

- Full-rank mixed state. All four eigenvalues positive.
- Realistic noise. But, Cannot purify finite number of states to Bell pair , *Jané QIC (2002)*
- Two-qubit **Werner** really (*isotropic*) state, useful for theory and experiment: One parameter x

$$\rho_W(x) = x |\Phi_{00}\rangle\langle\Phi_{00}| + \frac{1-x}{4} \mathbb{1}_4, \quad 0 \leq x \leq 1$$

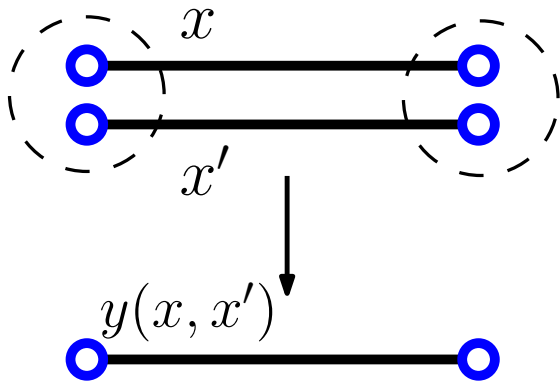
- Werner state is a full-rank state (for $x < 1$).
- Separable for $x \leq 1/3$.
- Make in lab: depolarization
- **Concurrence**: $C(x) = \max\{0, (3x - 1)/2\}$. **Linear in x** ,
 $C(\text{separable, i.e. not entangled}) = 0$, $C(\text{Bell pair}) = 1$

Swap Werner states. Get another Werner state.



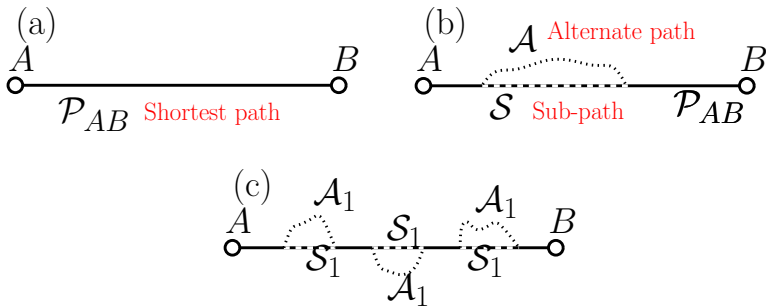
Entanglement increases with x ; Exponential decay of entanglement with length of chain. Swapping: lose entanglement, Purification: gain entanglement.

Purify Werner states. Get another Werner state.



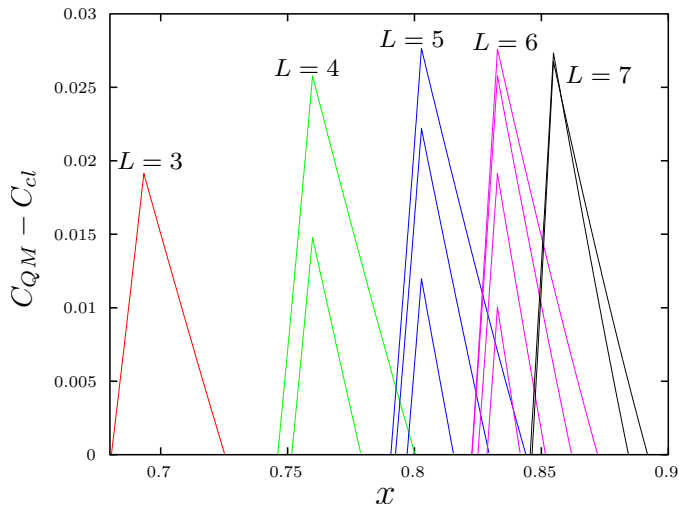
$$y(x, x') = \frac{x + x' + 4xx'}{3 + 3xx'}, \text{ with probability } \frac{1 + xx'}{2}$$

More generic network. Combine swapping and purification.



Swap first, or purify first? We call single-path purification **SPP**: 1) swap along sub-path; 2) swap along alternate path; 3) purify resulting states.

What is average concurrence? (over quantum outcomes)



Advantage of purify-swap depends on shortest path length L , sub-path length n , alternate path length m , Werner parameter x . Optimizing formula for gain in average concurrence is messy.

Poisson Random Graph (or Erdős–Rényi Graph)

Apply **SPP** between random pairs of nodes.

- Graph with N vertices. Zero or one edge between each pair. Each of the $N(N - 1)/2$ edges is present with probability p .
- Density of shortest paths of length L , σ_L

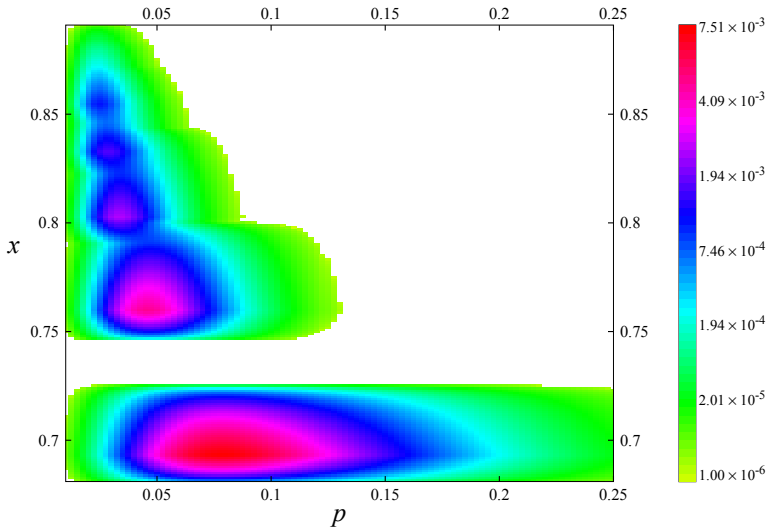
$$\sigma_1 = p,$$

$$\sigma_2 = (1 - p)[1 - (1 - p^2)^{N-2}] \approx (1 - p) \left(1 - e^{-p^2 N}\right),$$

$$\begin{aligned} \sigma_3 &\approx \left(1 - e^{-p^3(1-p)^5(N-2)(N-3)}\right) (1 - p^2)^{N-2}(1 - p), \quad \text{large } p \\ &\approx \left(1 - e^{-p^3(1-p)^5 N^2}\right) e^{-p^2 N} (1 - p) \end{aligned}$$

$$\sigma_L = p^L \frac{(N - 2)!}{(N - L - 1)!} + \mathcal{O}(p^{L+1}), \quad \text{small } p$$

$$\sigma_L \approx \frac{1}{N} \quad \text{for } pN = 1, \quad L < \text{radius}$$



Advantage of purify-swap depends on Werner parameter x and bond density of random graph p .

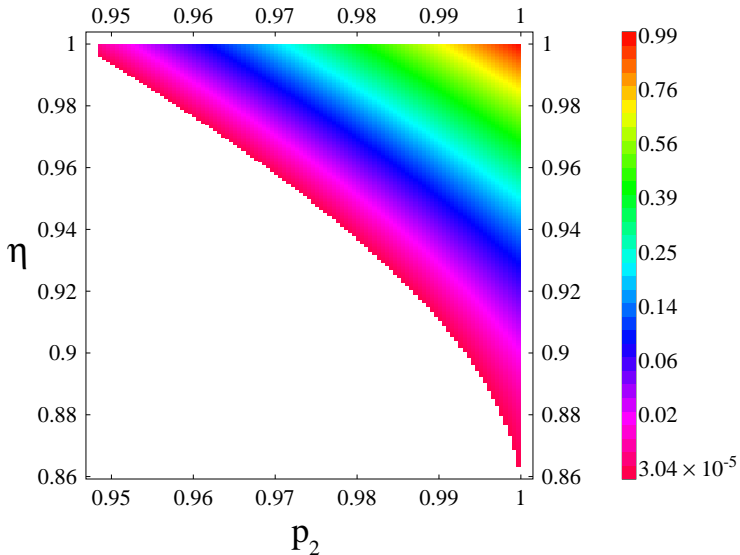
Monte Carlo $N = 200$, $L < 8$

Poisson Random Graph at critical point $pN = 1$

Choose random pair of vertices. Entangle pair via purify-swap, or direct swap. What is average gain in final entanglement ?

- Giant cluster of mass $N^{2/3}$
- Density of shortest paths independent of L . So, as Werner param. $x \rightarrow 1$ long paths dominate.
- “Good” ranges of x overlap more for large L : \Rightarrow integrate
- Each path-subpath occurs with probability $\approx 1/N^2$
- At fixed x , contributions are from $L \approx 1/(1-x)$. Four factors
 - $\approx L$ paths contribute near x
 - $\approx L$ sub-path lengths per path
 - $\approx L$ alternate paths per sub-path
 - $\approx L$ positions along path for sub,alt-path pair.
- Advantage of purify-swap over swap, averaged over network is

$$\Delta \bar{C} \sim \frac{K}{N^2(1-x)^4} \text{ for large } N \text{ small } 1-x, \quad (K \approx 6.5 \times 10^{-5})$$

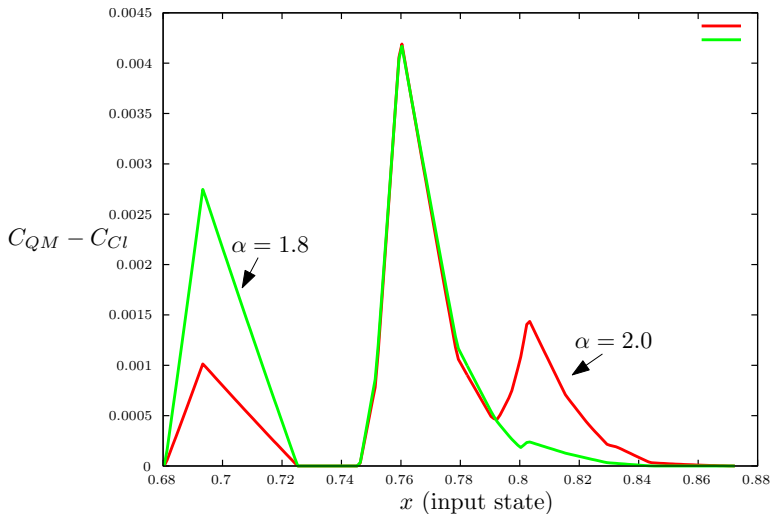


Noisy operations: $y = y_{\max}$, $a = a_{\max}(y_{\max})$, gives $\Delta C = 1/36$. η : reliability of measurement. p_2 : reliability of two-qubit operator.

Poisson Random Graph at critical point $pN = 1$

But wait, . . . there's more. Return to perfect operations.

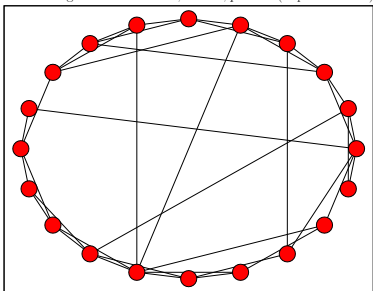
- For $Np = 1$, Radius grows like $N^{1/3}$ Nachmias, Peres, Ann. Prob. (2008)
- Our MC shows radius of largest cluster $\approx 3N^{1/3}$.
- Since $L \approx 1/(1-x) \Rightarrow \Delta \bar{C} < 81AN^{-2/3}$
- Purification protocols always give modest results. They must be used iteratively.
- *But*, Choose bond density to favor $L = 2, 3$: $p^2N = c$. Then $\sigma_2 \rightarrow (1 - e^{-c})$ and $\sigma_3 \rightarrow e^{-c}$. Now for Werner parameter around 0.7, we have many subgraphs for purify-swap.



Advantage of purify-swap for Werner states on scale-free network with $p(k) \propto k^{-\alpha}$. Monte Carlo with $N = 200$.

Watts-Strogatz model has parameter p that tunes between lattice-like model with regular local connections, and an Erdős–Rényi-like model. We compute distribution of shortest paths on Watts-Strogatz model for $p = 0$. One link to each neighbor at $\pm 1, \pm 2$, ($K = 4$).

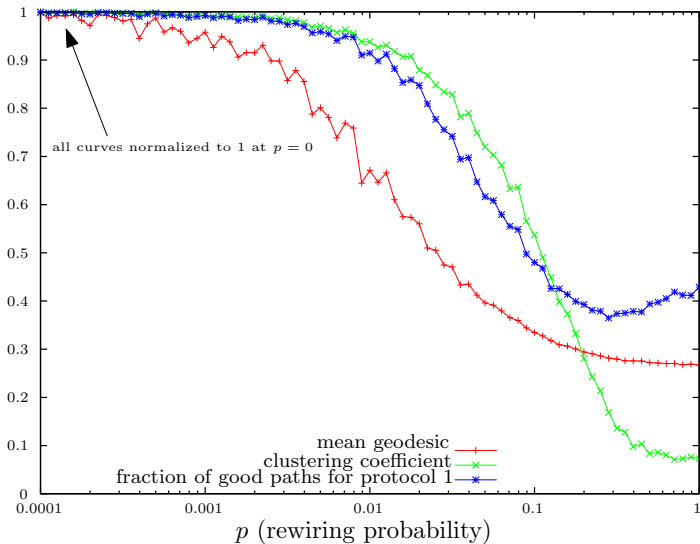
Watts-Strogatz model $N = 20$, $K = 4$, $p = 0.2$ (Arpad Horvath)



- $N(N - 1)/2$ shortest paths (SPs)
- Number of SPs of each length L from 1 through $N/4 - 1$ is $2N$ (and $3N/2$ for boundary case $L = N/4$.)
- Density of SPs of length L is then $\sigma_L = 4/(N - 1)$. Flat. (except for boundary case.)

Number of shortest paths admitting SPP (single purify-swap)

$$\frac{N^2 - N}{2} - 4N - \frac{1}{2}2N(5) = \frac{N(N - 19)}{2}.$$



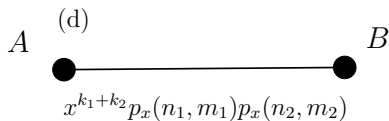
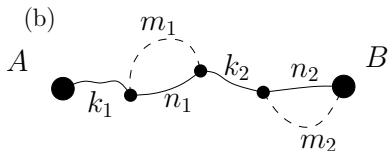
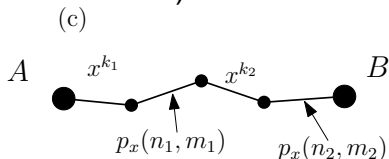
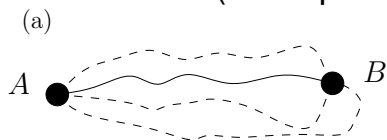
Watts-Strogatz small world (2×2 neighbors). Fraction of paths admitting purify-swap. ($N = 100$). The separation of the red and green lines is the main point of the model.

Summary

- Long-range entanglement by using short-range entangled links.
- Geometry, topology, etc. often determines optimal protocol
- Dynamics ?
- Coupled lattices ?
- Lower bound on entanglement in 2-d with local connections ?

Mixed states on complex network

Werner state on each link. What is average concurrence?
(over quantum outcomes)



$p_x(n, m)$ is average Werner parameter after **SPP** with subpath length n , alternate path length m .