

European Tensor Network School, San Sebastian
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An Introduction to MERA

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www.heptnseminars.org

(online seminar series on tensor networks)



MPS

Natural ansatz for
1d gapped systems

In **practice**, can be
applied to 1d critical
systems and also to
2d systems

Somewhat RG based

PEPS

Higher dimensional
extension of MPS

Applied to both
gapped and critical
systems in 2d

Not RG based

MERA

Natural ansatz for
1d critical systems

Possible to extend
to higher
dimensions

RG based

THIS TALK: MERA for 1d systems

Outline

PART 1: What is the MERA?

- (1) Efficient representations of quantum many-body states from **coarse-graining** the lattice
- (2) MERA as a numerical ansatz for 1d ground states

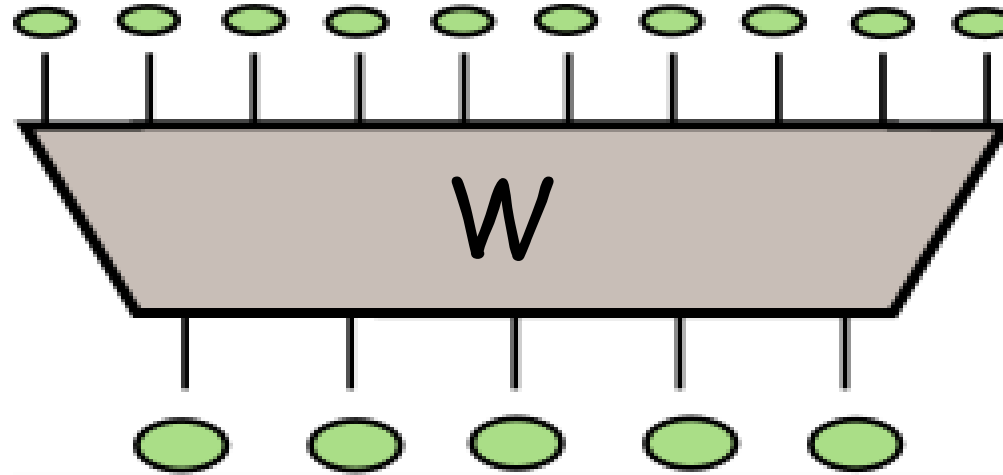
PART 2: MERA and critical systems

- (3) Entanglement in the MERA
- (4) infinite MERA as a lattice version of 2d CFTs
- (5) MERA and holography

Efficient representations of many-body states
from **coarse-graining** the lattice

Definition of a coarse-graining or RG transformation

A **linear operator (isometry) W** that maps states/operators on a lattice to states/operators on a coarse-grained lattice



RG flow of the Hamiltonian: $H \rightarrow H_1 \rightarrow H_2 \rightarrow H_3 \rightarrow \dots \rightarrow H_F \rightarrow H_F \rightarrow \dots$

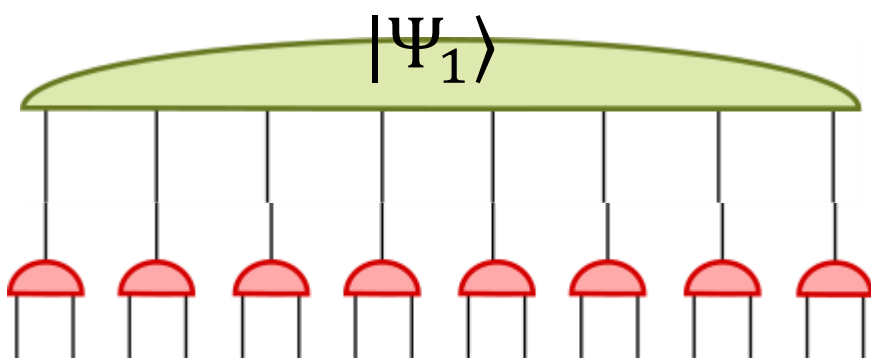
- 1) Must preserve the **low energy subspace**
- 2) Must reach a **fixed-point** (which captures universal low-energy properties)

The map W will be a tensor network

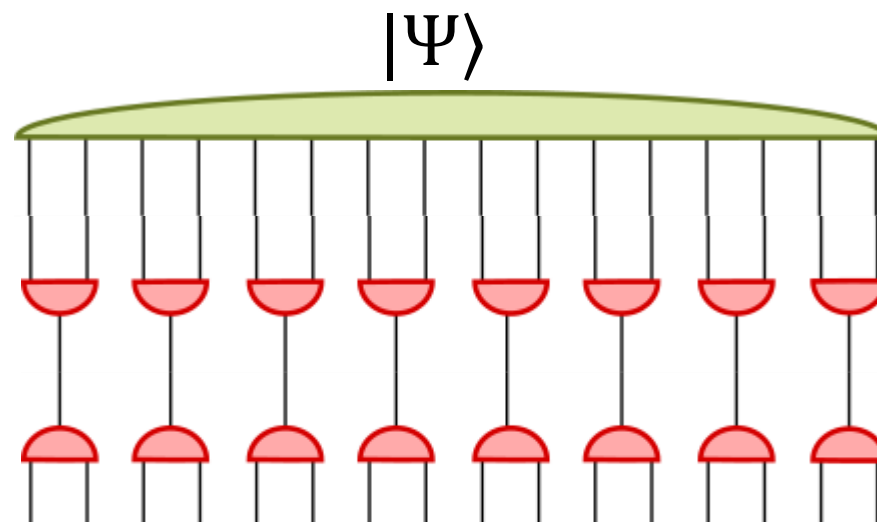
$$\rho \text{ (purple oval)} = D \begin{matrix} w^\dagger \\ \text{small purple oval} \\ w \end{matrix} \quad \text{(eigenvalue decomposition)}$$

Size of this index = rank of ρ

$$\begin{matrix} w \\ \text{red semi-circle} \\ w^\dagger \end{matrix} = \text{vertical line} \quad \text{(isometry)}$$



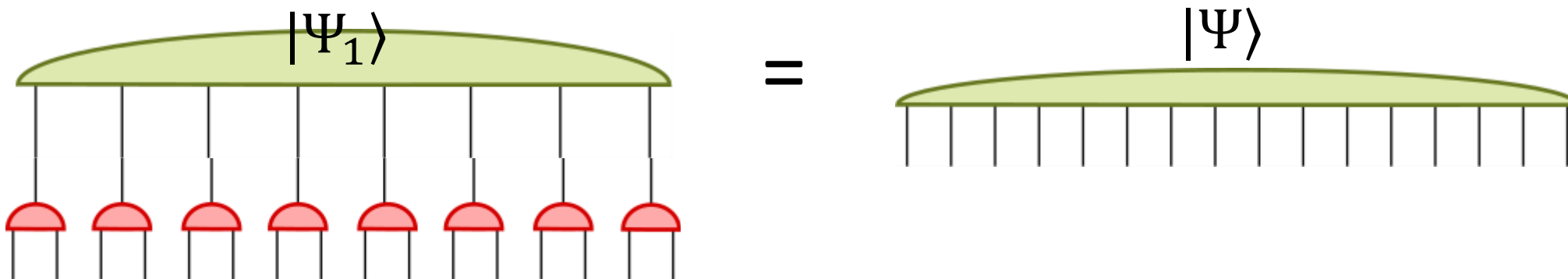
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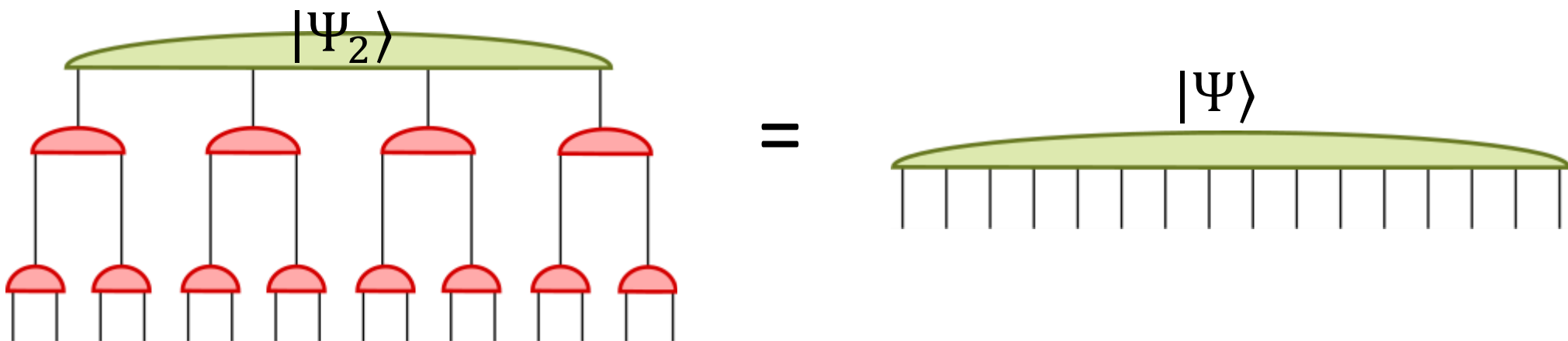


Alternative representation of $|\Psi\rangle$

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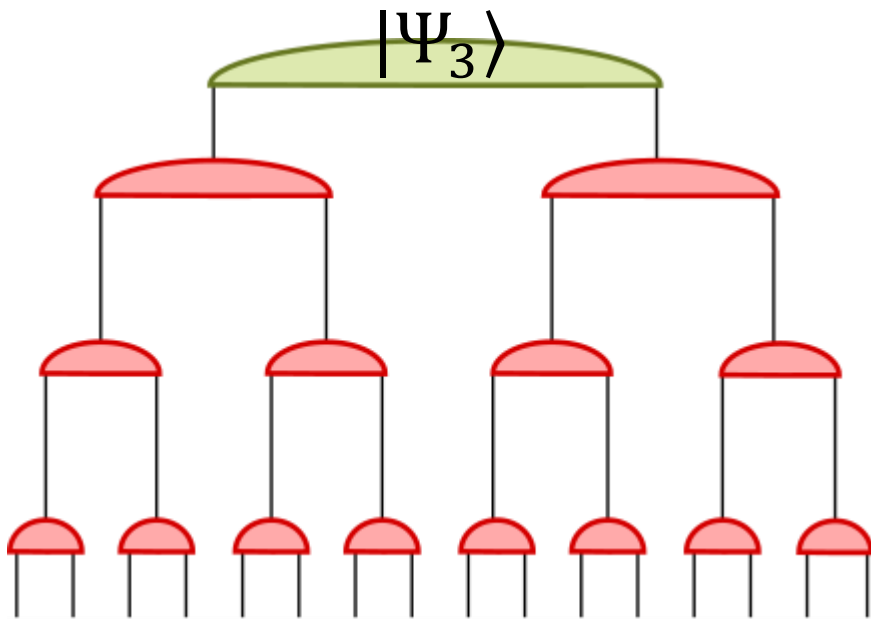
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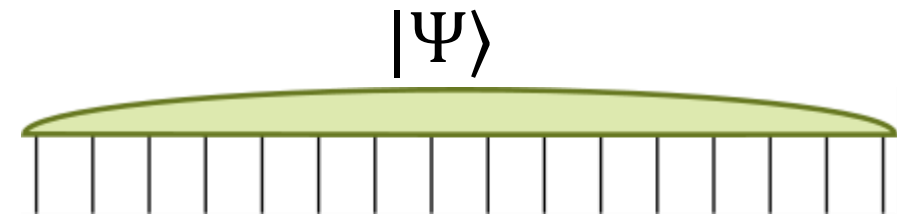
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Tree tensor network



=



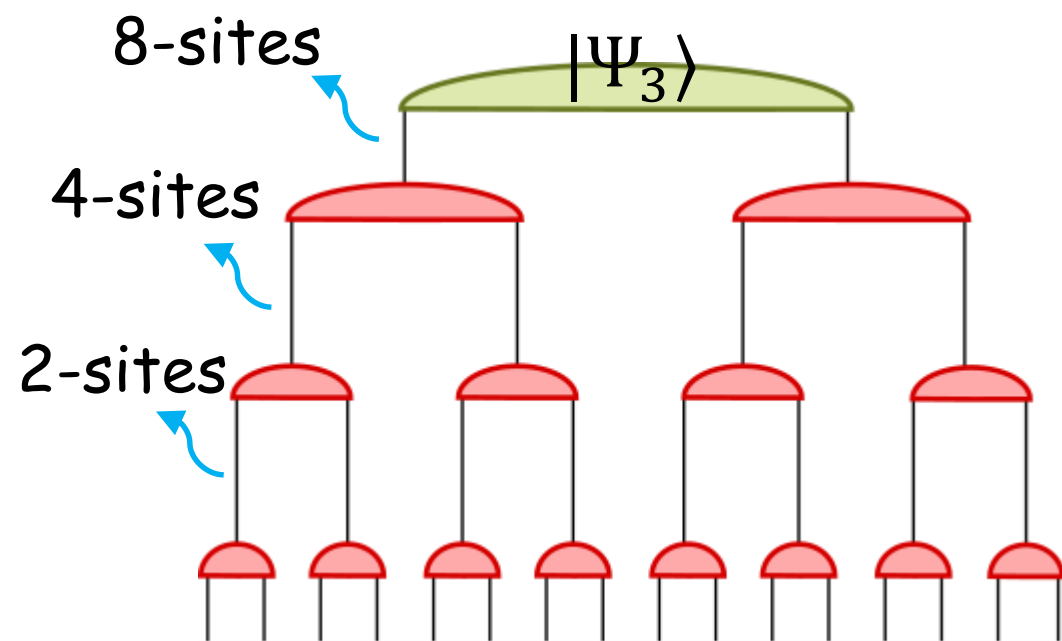
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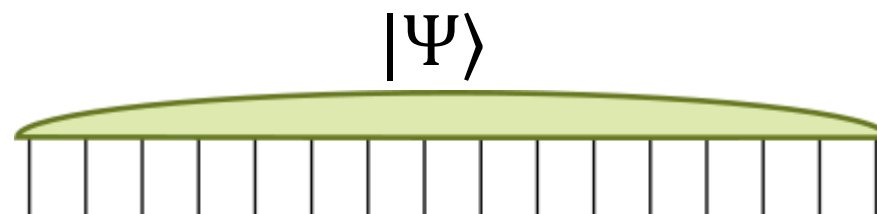
Size of this index = rank of ρ

$$\begin{matrix} w \\ \text{red semi-circle} \\ w^\dagger \end{matrix} = \text{vertical line} \quad \text{(isometry)}$$

Tree tensor network



=



Alternative representation of $|\Psi\rangle$

So far we have not assumed anything about the state

The point is only that we have an alternative representation of the state as a tree tensor network

Let's look at gapped states first ... We know that

$$S(\rho_\ell) \text{ saturates} \quad \Rightarrow \quad \text{Rank}(\rho_\ell) \text{ saturates}$$

A bond index in the tree tensor network corresponds to a block of sites on the original lattice.

This means that the bond dimension saturates (say to χ) after a finite number of coarse-graining steps

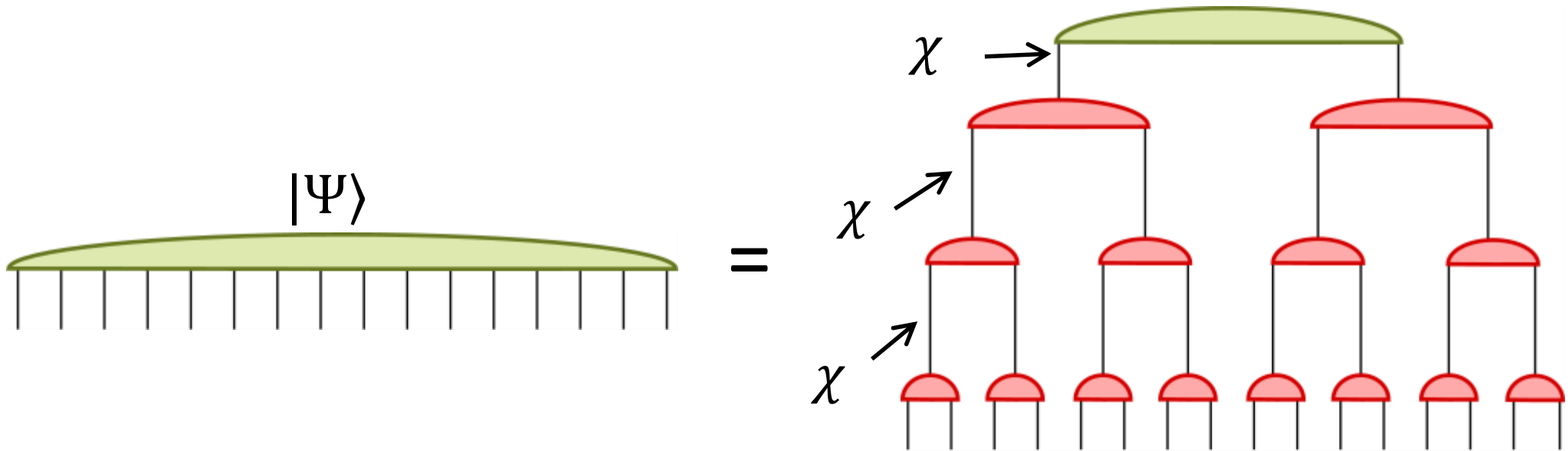
Quiz!

How many complex numbers in the tree representation of a 1d gapped ground state?

(i) $O(N^2)$

(ii) $O(\log N)$

(iii) $O(N)$



Quiz!

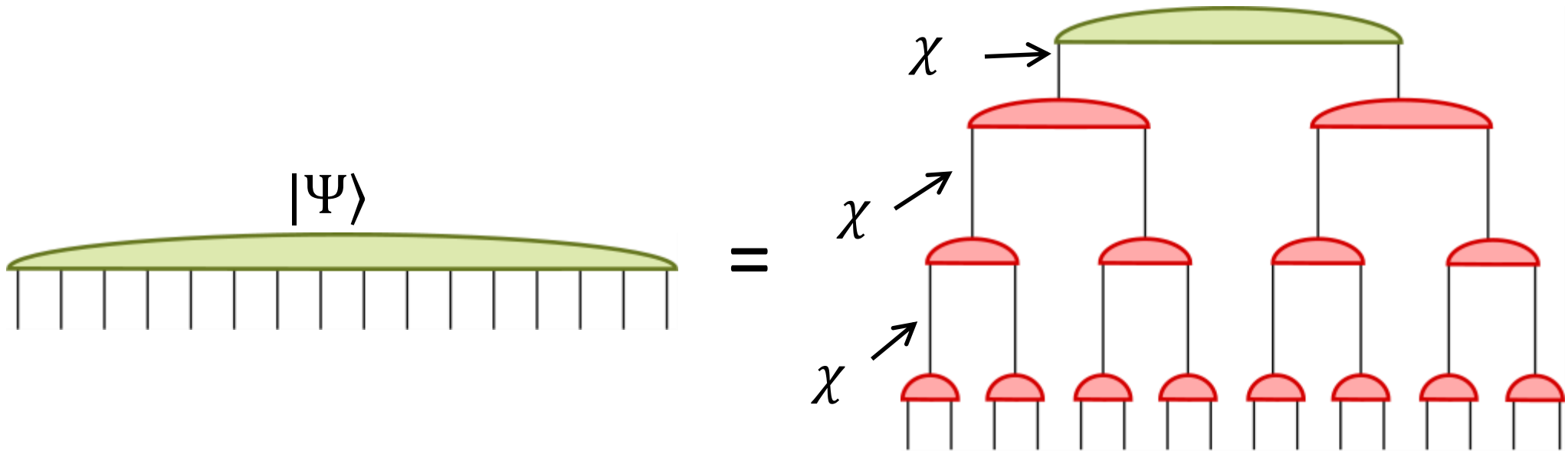
How many complex numbers in the tree representation of a 1d gapped ground state?

(i) $O(N^2)$

(ii) $O(\log N)$

(iii) $O(N)$

$O(N \chi^3)$



Tree tensor network can efficiently **parameterize** 1d gapped ground state

But being an efficient parameterization is not enough.

Can we also compute expectation values from the TTN efficiently (polynomial time)?

It turns out : yes.

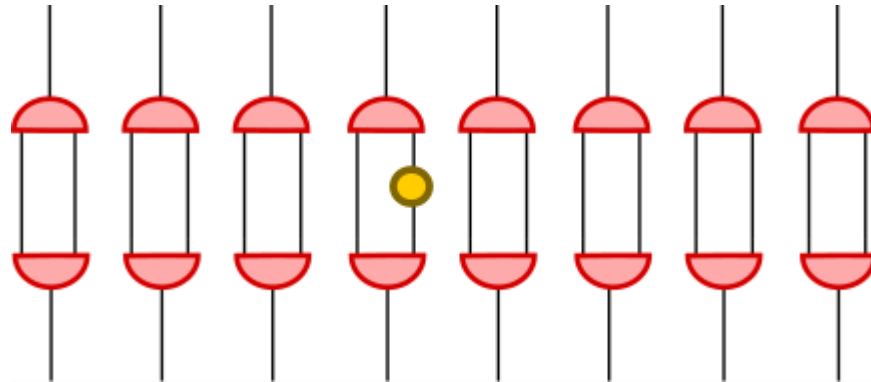
This is thanks to a particular feature of this coarse-graining transformation:

Local operators coarse-grain to local operators

1-site operators

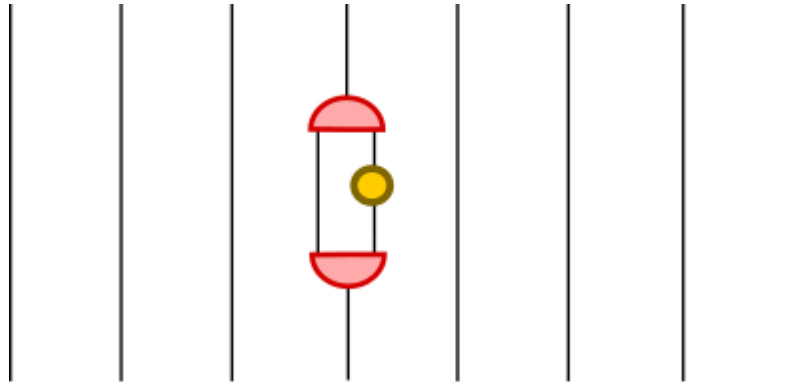
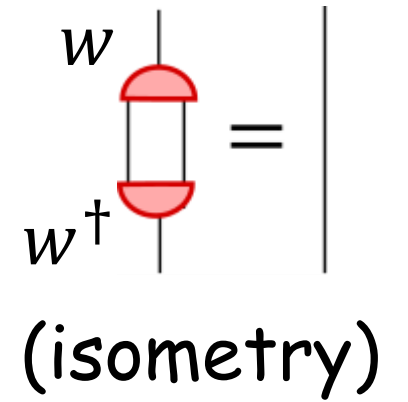
$$\begin{array}{c} w \\ \text{---} \text{---} \\ \text{---} \text{---} \\ w^\dagger \end{array} = \text{---}$$

(isometry)



1-site operators

Cost : $O(\chi^4)$

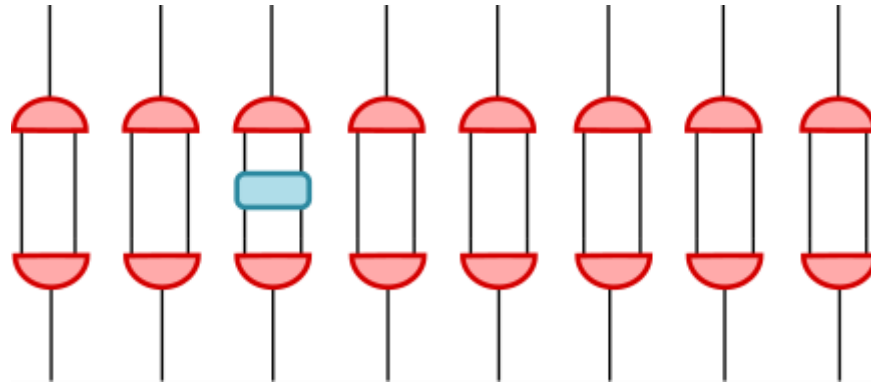


This also means that the cost of coarse-graining 1-site operators is independent of N

2-site operators

$$\begin{array}{c} w \\ \text{---} \text{---} \text{---} \\ \text{---} \text{---} \text{---} \\ w^\dagger \end{array} = \text{---}$$

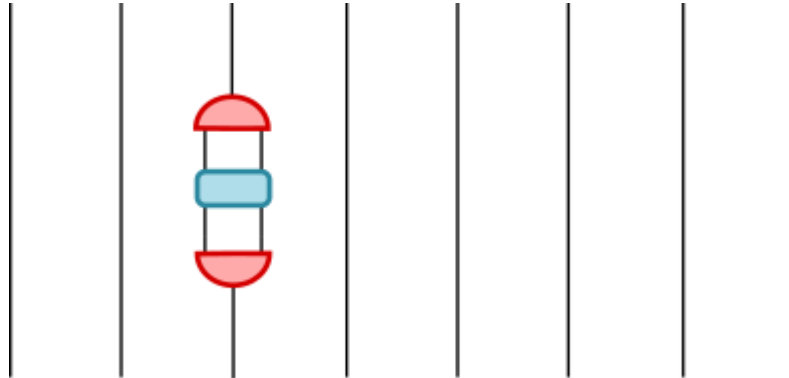
(isometry)



2-site operators

$$\begin{array}{c} w \\ \text{---} \text{red semi-circle} \\ \text{---} \text{white rectangle} \\ \text{---} \text{red semi-circle} \\ w^\dagger \end{array} = \text{---} \text{vertical line}$$

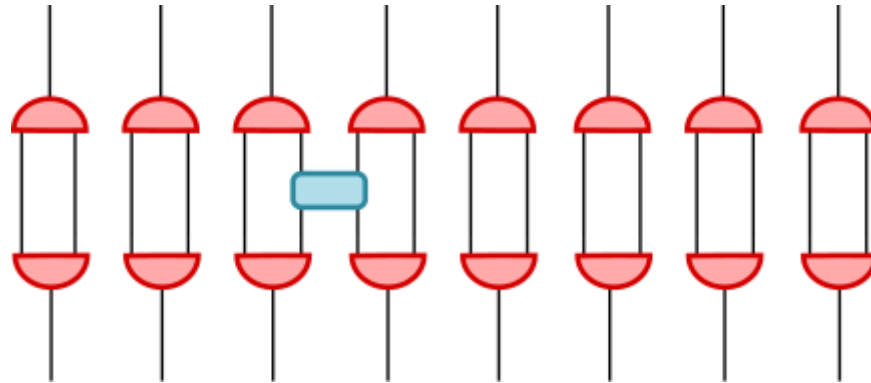
(isometry)



2-site operators

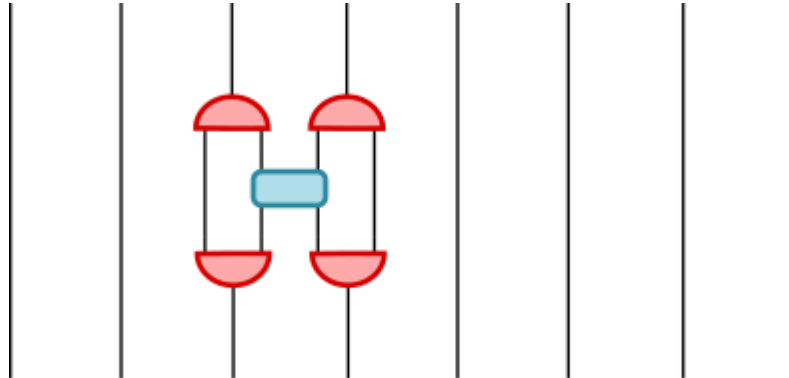
$$\begin{array}{c} w \\ \text{---} \text{red semi-circle} \\ \text{---} \text{red semi-circle} \\ w^\dagger \end{array} = \text{---} \text{vertical line}$$

(isometry)



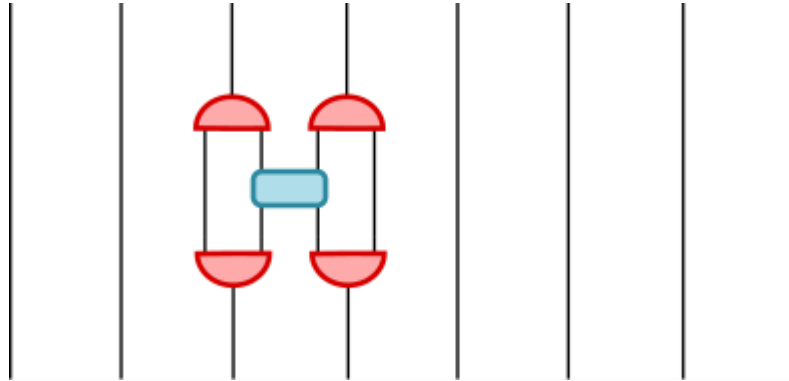
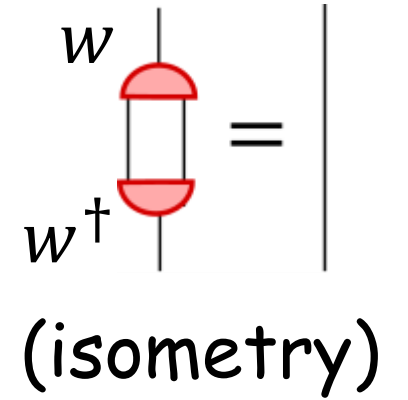
2-site operators

$$\begin{array}{c} w \\ \text{---} \text{red semi-circle} \\ w^\dagger \\ \text{---} \text{red semi-circle} \end{array} = \text{---} \text{vertical line} \quad \text{(isometry)}$$



2-site operators

Cost : $O(\chi^6)$

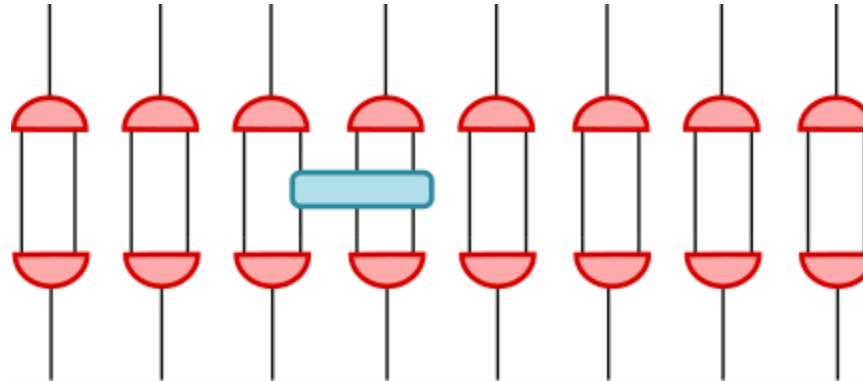


Cost of coarse-graining 2-site operators also independent of N

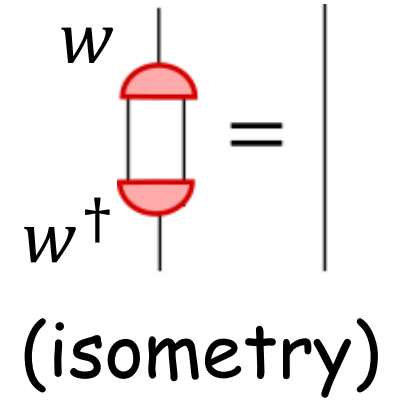
3-site operators

$$\begin{array}{c} w \\ \text{---} \text{---} \text{---} \\ \text{---} \text{---} \text{---} \\ w^\dagger \end{array} = \text{---}$$

(isometry)

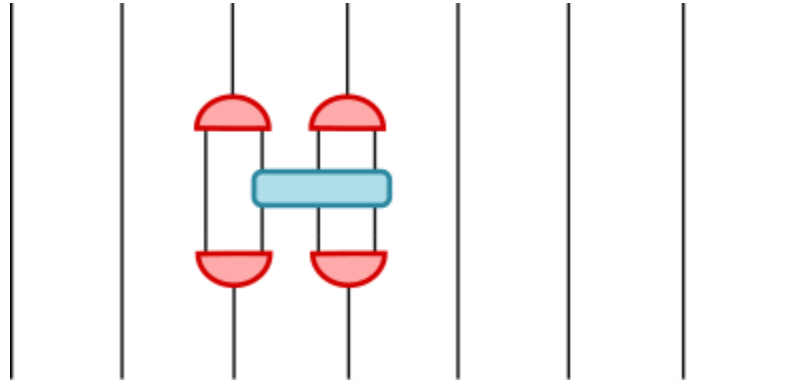


3-site operators



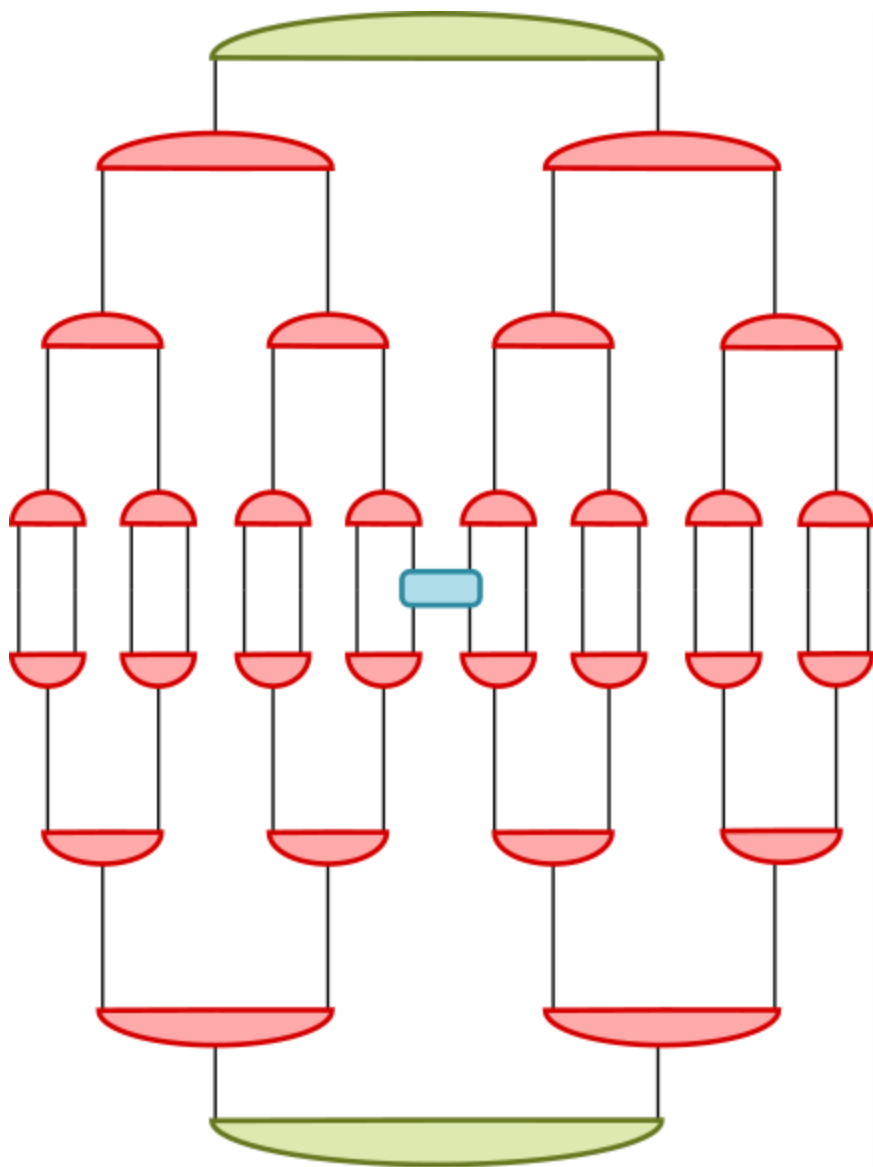
A diagram showing a single vertical line representing a site. Two red semi-circles are attached to the line, one above the other. The top semi-circle is labeled w and the bottom semi-circle is labeled w^\dagger . To the right of these semi-circles is an equals sign followed by another single vertical line. Below the entire diagram is the text "(isometry)".

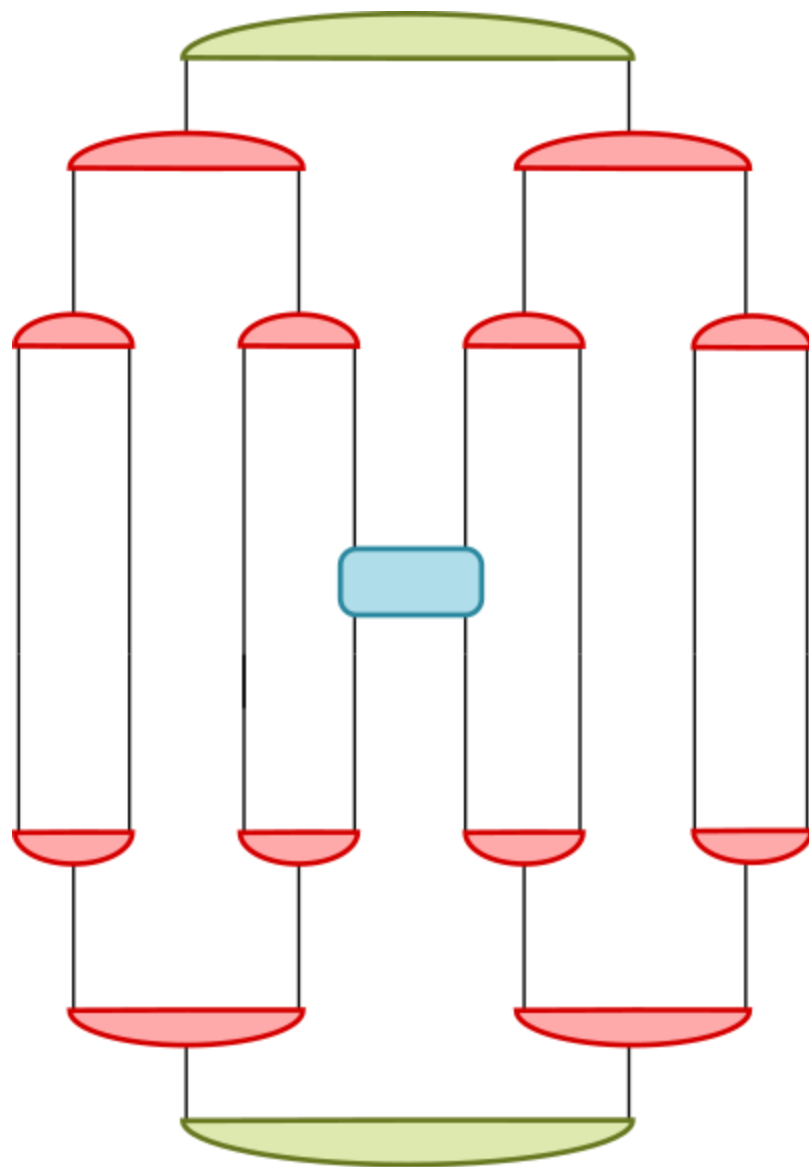
(isometry)

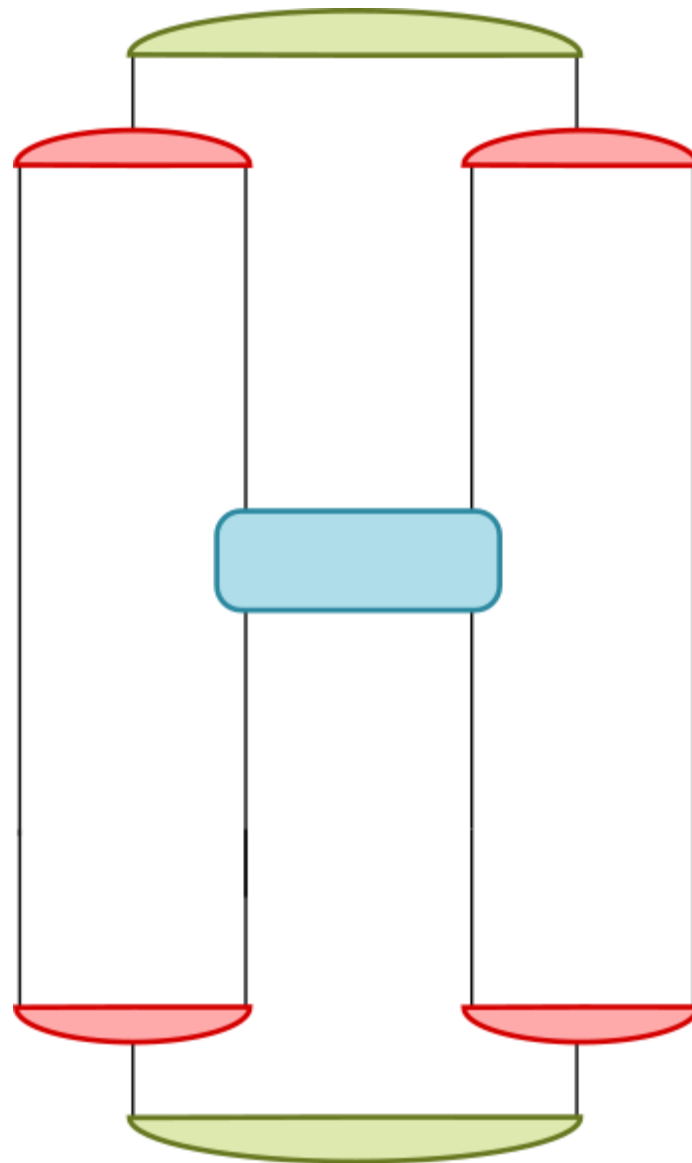


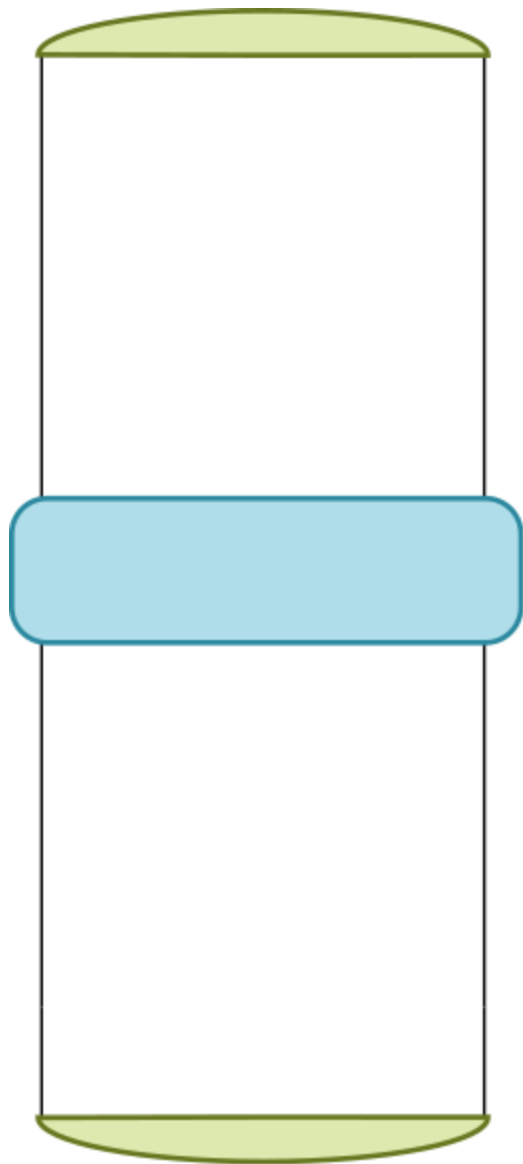
3 or more site operators **shrink** under coarse-graining (thus, also remain local)

Preservation of locality **implies** that expectation values of local operators can be computed efficiently

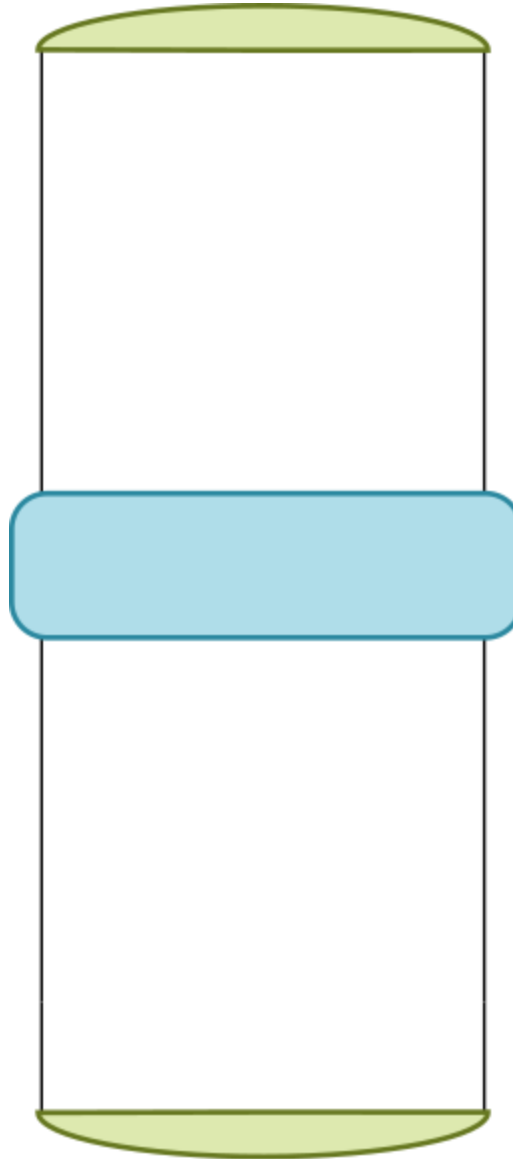






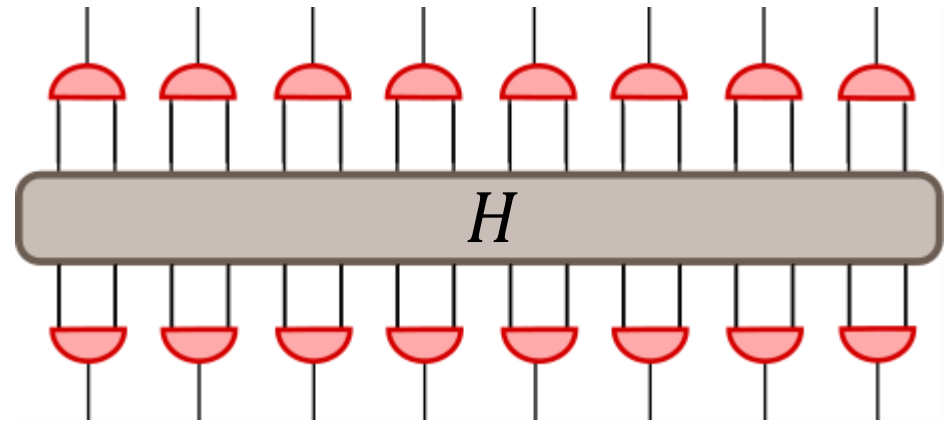
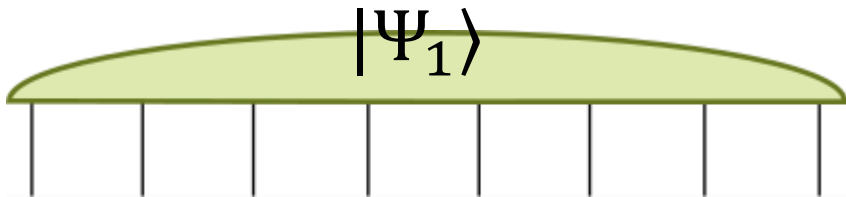
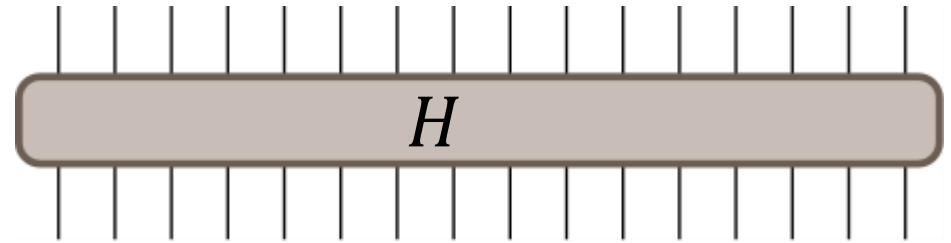
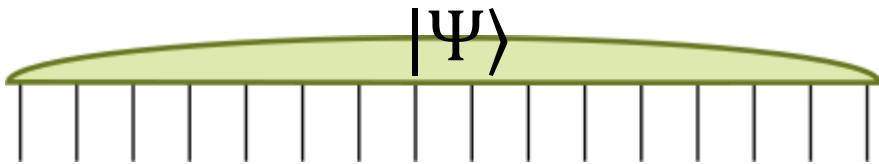


A TTN is an efficient ansatz
for 1D **gapped** ground states.

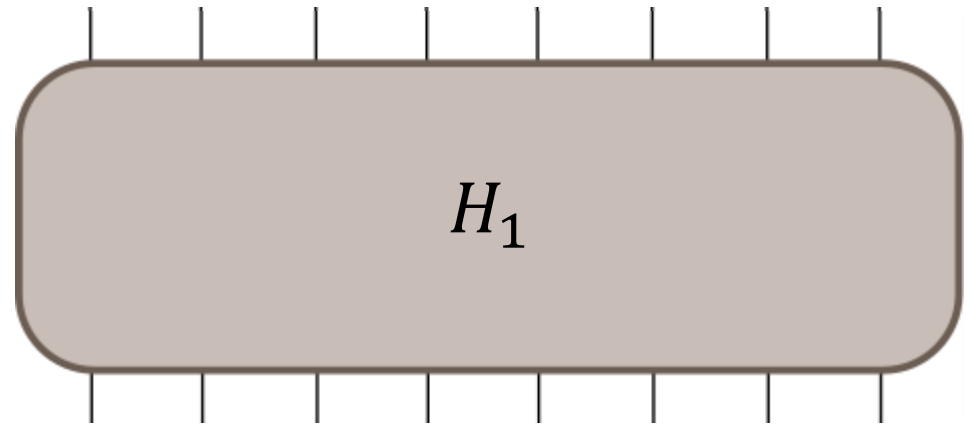
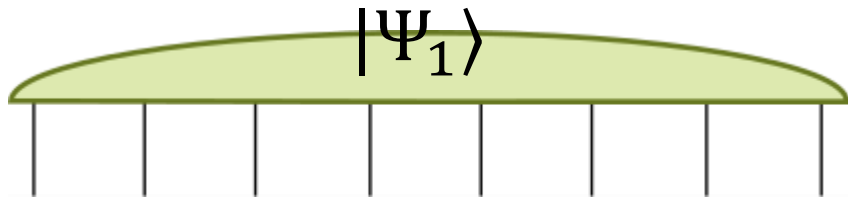
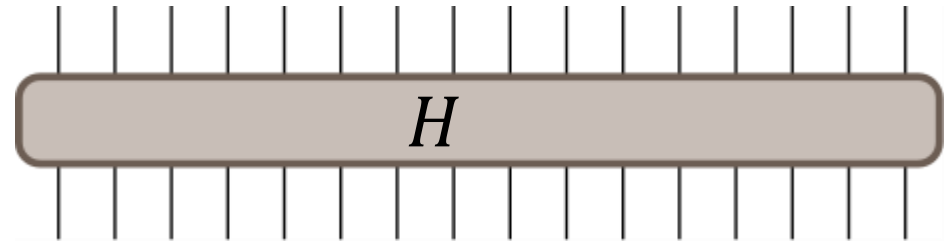
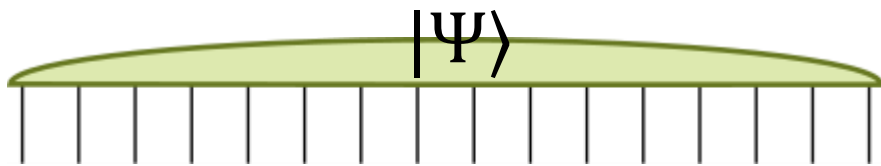


Cost per step: $O(\chi^6)$
Total number of steps: $O(\log N)$

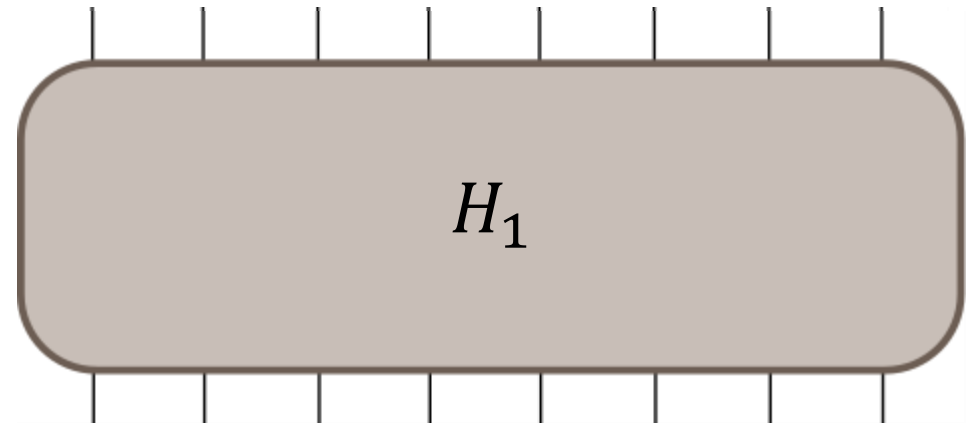
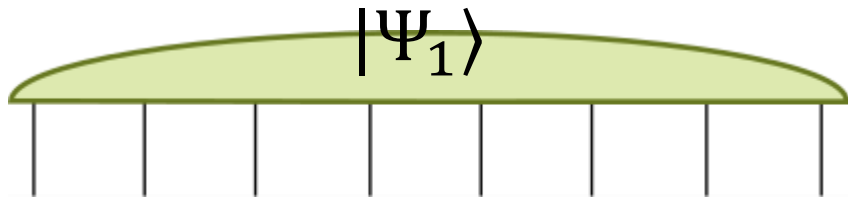
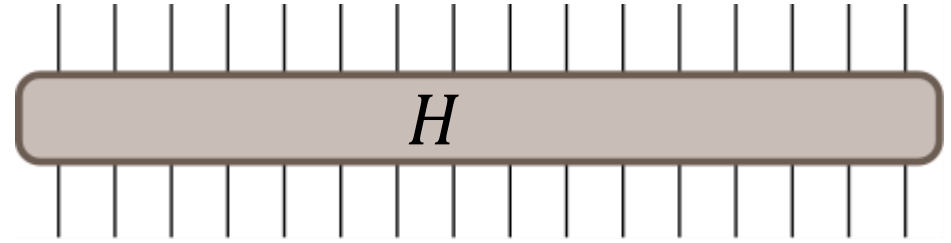
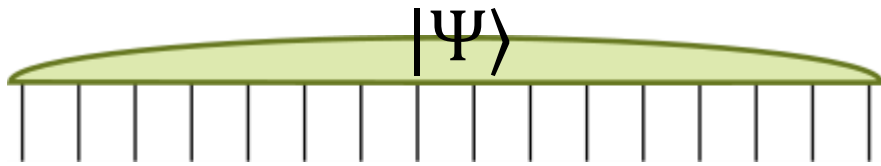
The tree representation of a gapped ground state also defines an **RG flow**



The tree representation of a gapped ground state also defines an **RG flow**

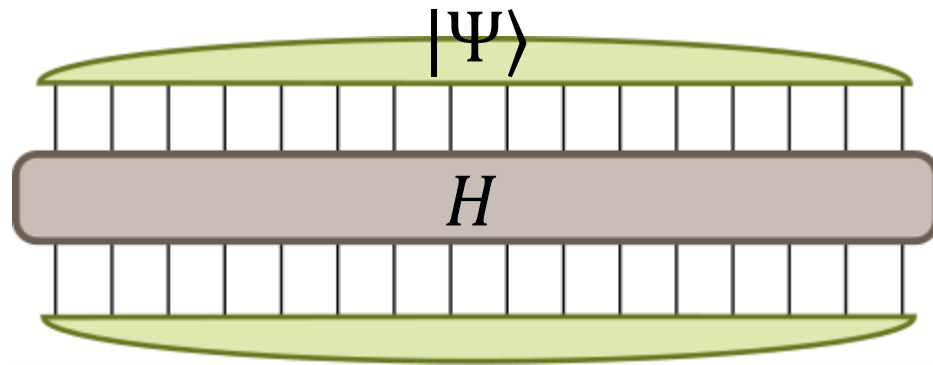


The tree representation of a gapped ground state also defines an **RG flow**

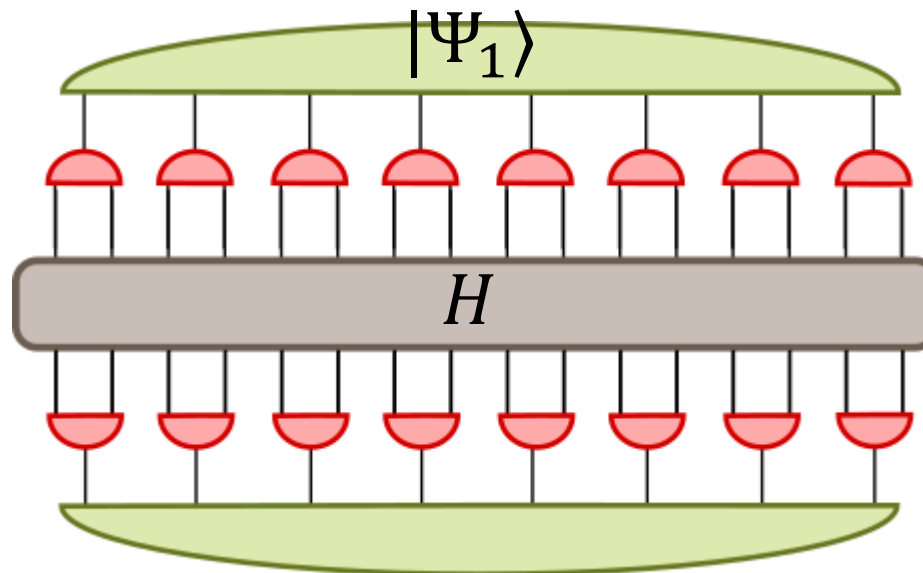


- (i) $|\Psi_1\rangle$ is the **ground state** of H_1
- (ii) $|\Psi_1\rangle$ has the **same** energy as $|\Psi\rangle$

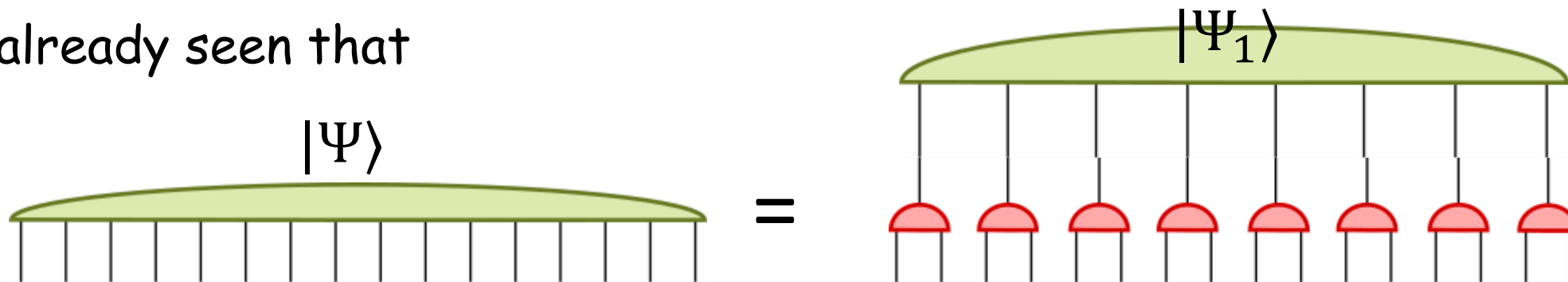
ground state energy of H =

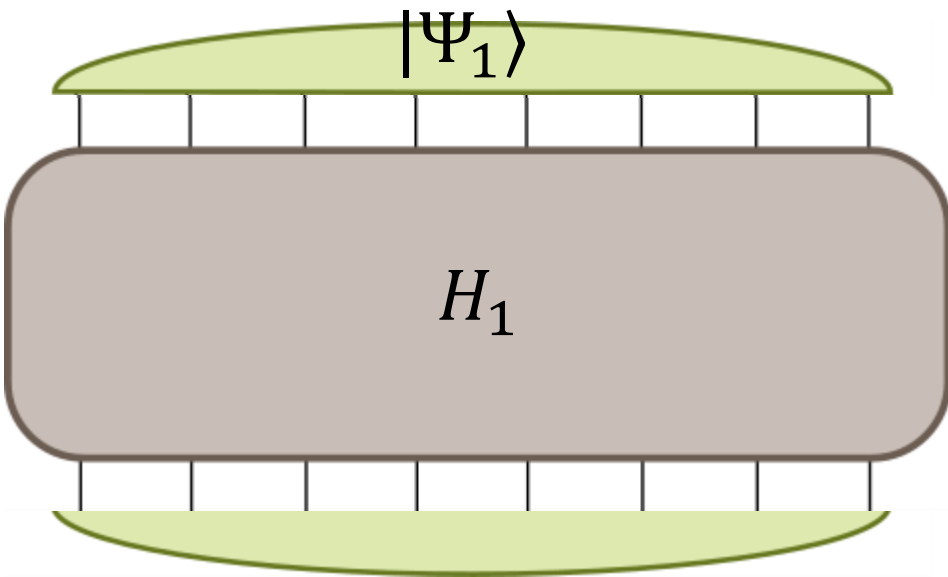


ground state energy of H =



We have already seen that

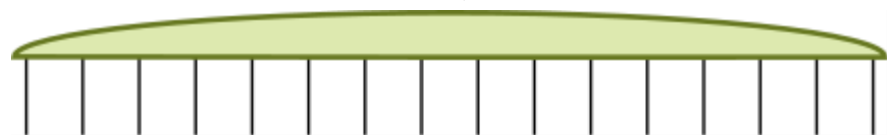
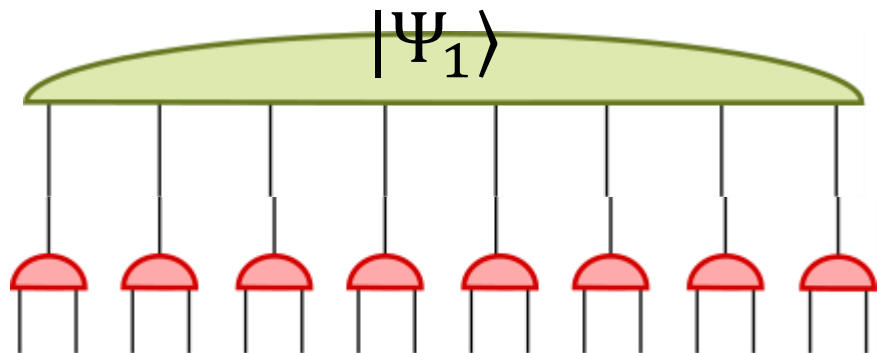


ground state energy of H =  = Energy of $|\Psi_1\rangle$

Also, clearly $|\Psi_1\rangle$ is the ground state of H_1 . Else this tensor network value can be minimized by replacing $|\Psi_1\rangle$ with the ground state.

But this cannot happen since $|\Psi\rangle$ is already the ground state of H .

We have already seen that

 = 

Not only energy match but also other expectation values are preserved

This means that the coarse-graining transformation preserves the ground state properties.

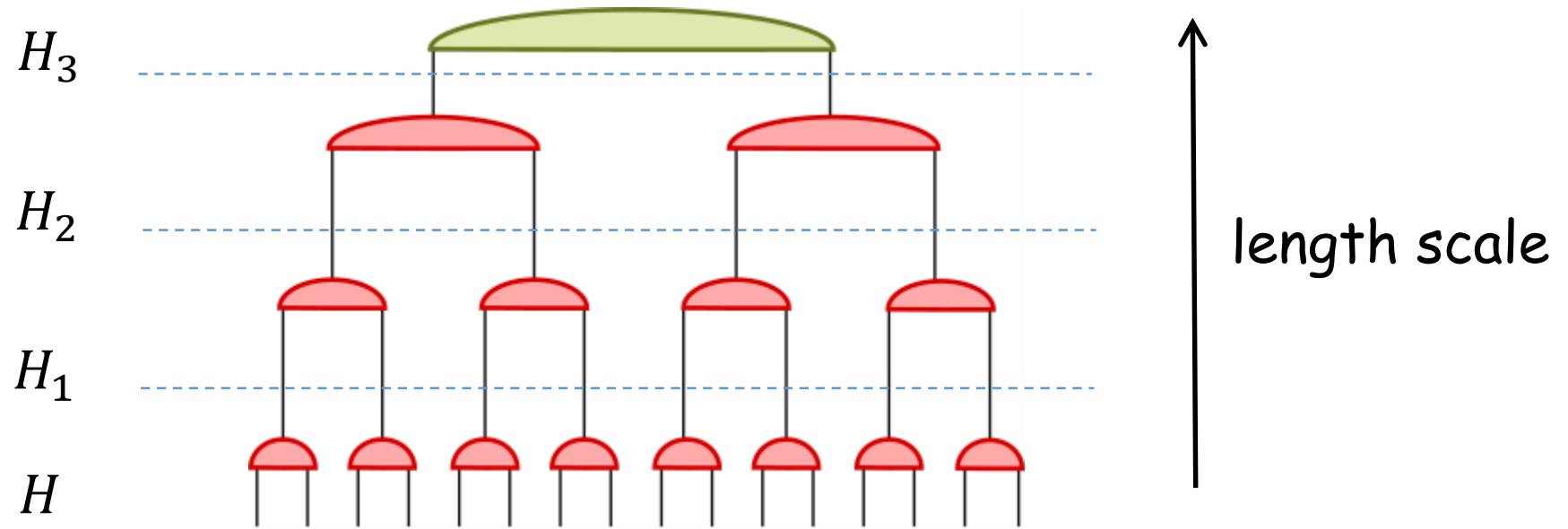
But H_1 is smaller than H .

So it must discard high energy states.

In fact, at the top we only have the ground state!

So the tree tensor network defines a RG flow in the space of Hamiltonians

The tree tensor network of a gapped ground state defines an **RG flow**



Summary: tree tensor network not only gives an efficient description of the gapped ground state but also defines an **RG flow** in the space of Hamiltonians.

What about **critical** ground states?

What about critical ground states?

$S(\rho_\ell)$ grows as $\log \ell \quad \Rightarrow \quad \text{Rank}(\rho_\ell) \text{ grows } O(\text{poly}(\ell))$

Bond dimension grows polynomially with number of coarse-graining steps

Unsustainable RG flow. **Worse:** does not reach a fixed point

Not reproducing a fixed point at criticality not just a failure of aesthetics

Obtaining the correct fixed point will allow contact with 2d CFTs that describe the critical system in the continuum

What about critical ground states?

$$S(\rho_\ell) \text{ grows as } \log \ell \quad \Rightarrow \quad \text{Rank } (\rho_\ell) \text{ grows } O(\text{poly}(\ell))$$

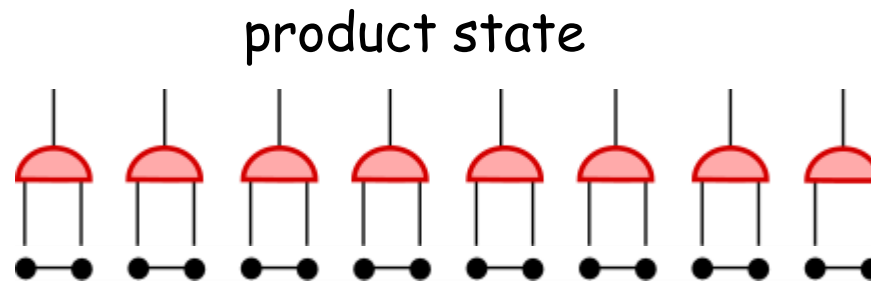
Bond dimension grows polynomially with number of coarse-graining steps

Unsustainable RG flow. **Worse:** does not reach a fixed point

Crux of the problem:

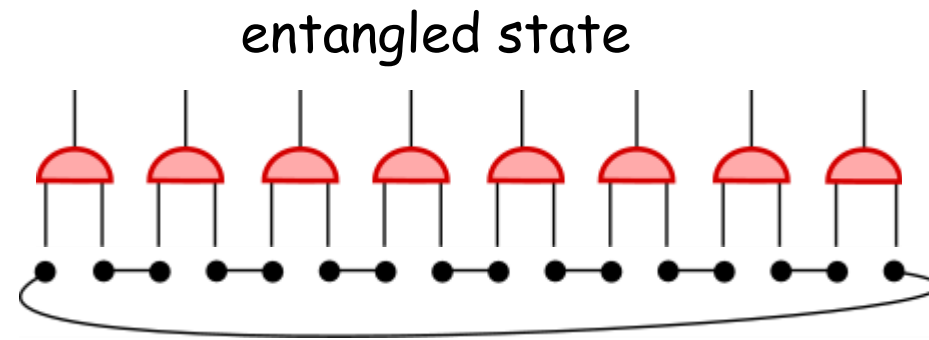
- (i) In a critical state entanglement appears at all length scales (as you block more and more sites, new **short-range** entanglement keeps appearing)
- (ii) The coarse-graining transformation does not remove **short-range** correlations properly

RG flow in the tree can accumulate of short-range correlations



maximally entangled
state

RG flow in the tree can accumulate of short-range correlations

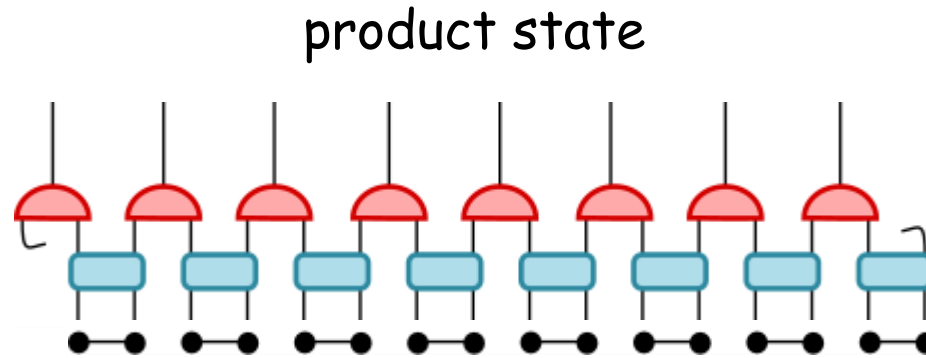


maximally entangled
state

In a critical state, where correlations appear at all length scales, some local correlations will not get removed properly

One fix: Introduce unitary gates that remove correlations locally

Disentanglers



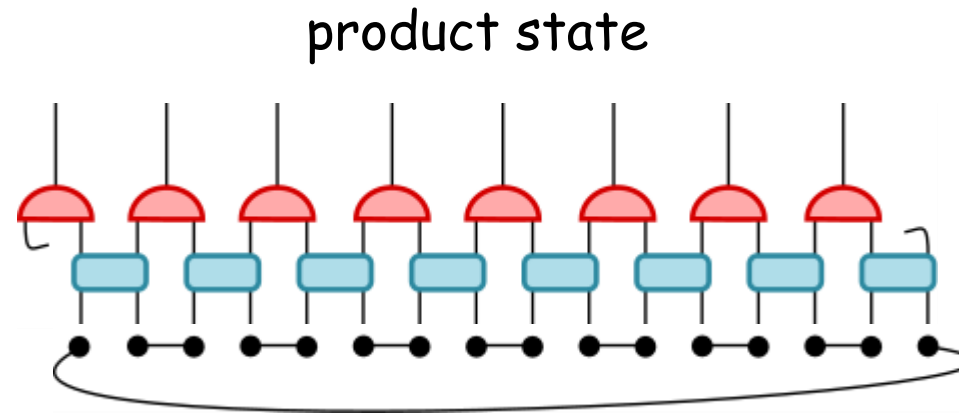
Disentanglers remove the entanglement

Isometries are trivial

One fix: Introduce unitary gates that remove correlations locally

Disentanglers

Entanglement
renormalization

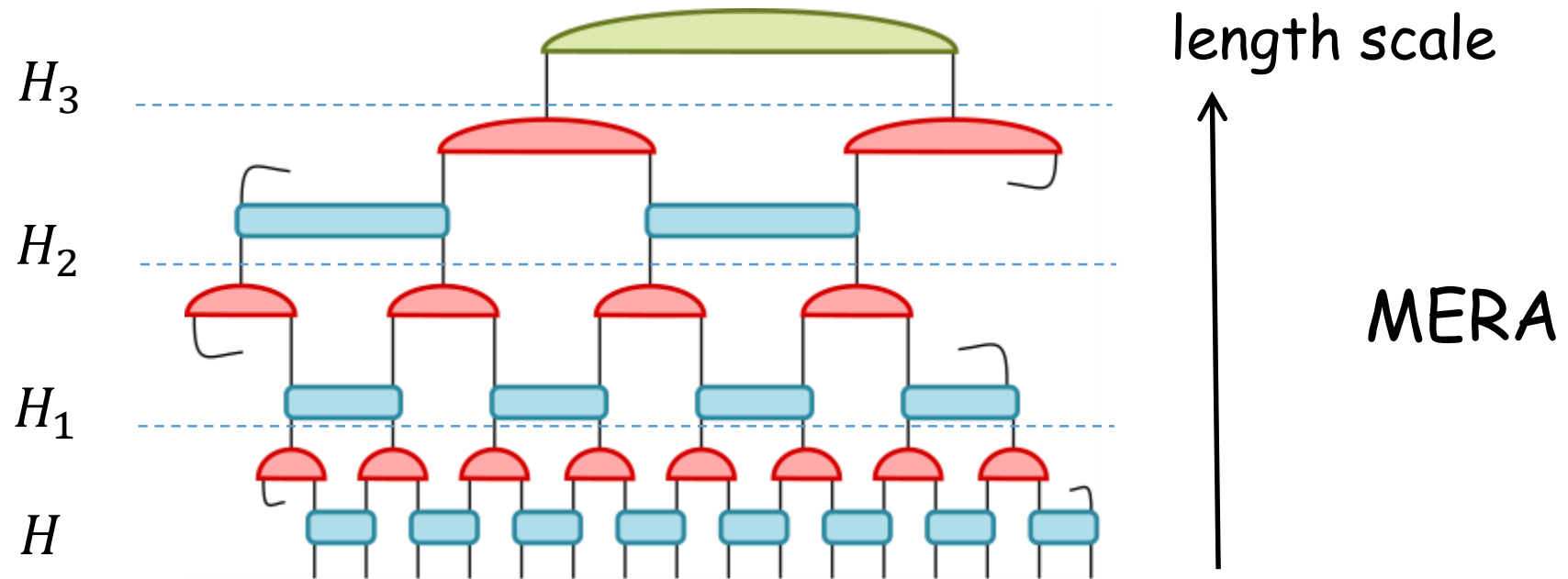


Disentanglers are trivial

But isometries remove the entanglement

Isometries + disentanglers are more **effective** at removing short range correlations

Tensor network generated by entanglement renormalization



Retains all the useful features of the tree: $O(N)$ parameters, preserves locality.

Also defines an RG flow, which can reach a fixed point at criticality.

MERA as a numerical **ansatz** for 1d ground states

So far we have discussed how to **translate** an exponentially large vector to a tree tensor network or MERA.

The isometries and disentanglers are determined from reduced density matrices in the state

In practice, we do not know the ground state to begin with.

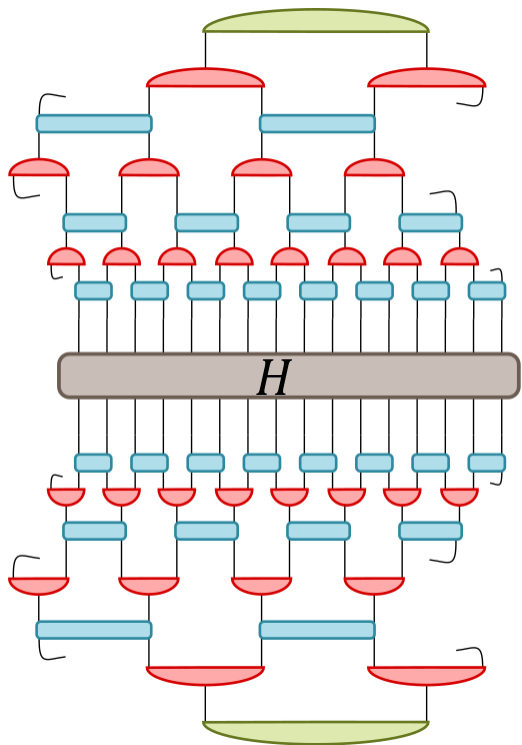
In this case, we can use the MERA as an **ansatz** for 1d ground states.

What we have argued so far is that it is **capable** of efficiently representing both gapped and critical ground states.

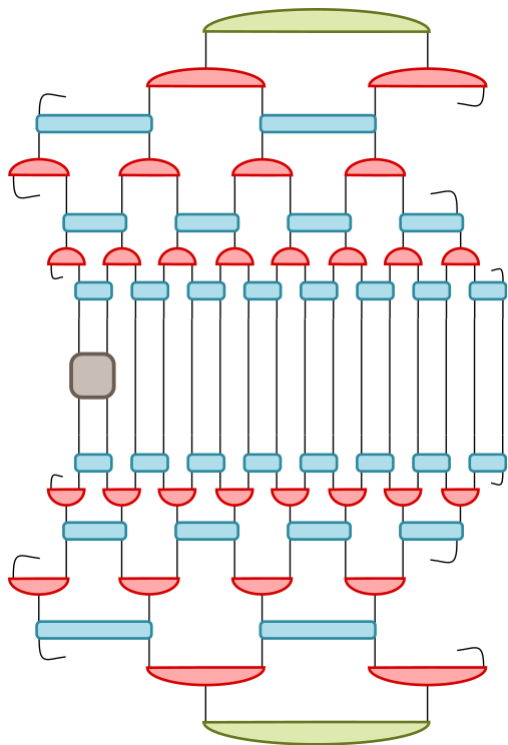
So how do we determine the ground state tensors in practice?

One method: variational energy minimization

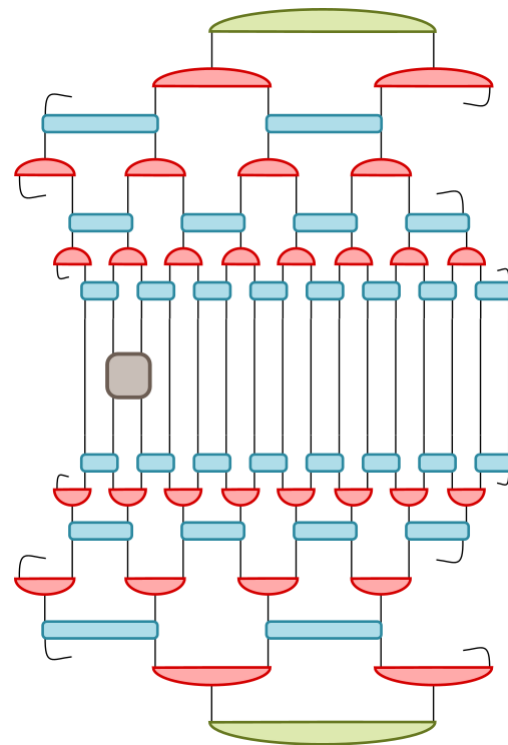
Hamiltonian is local



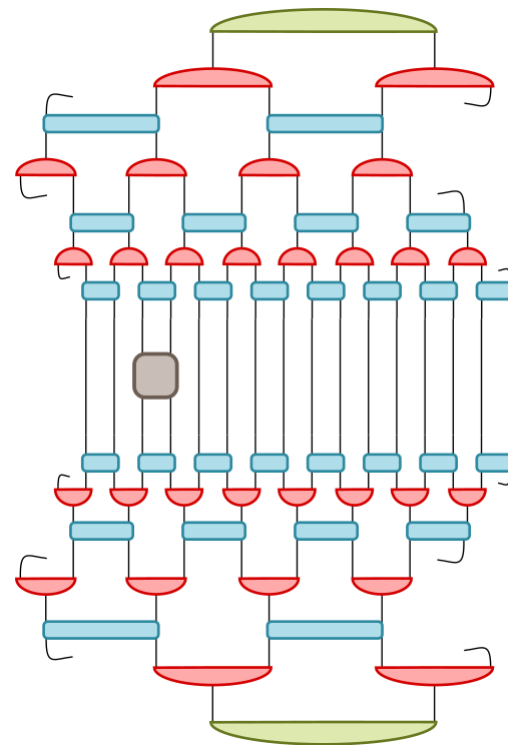
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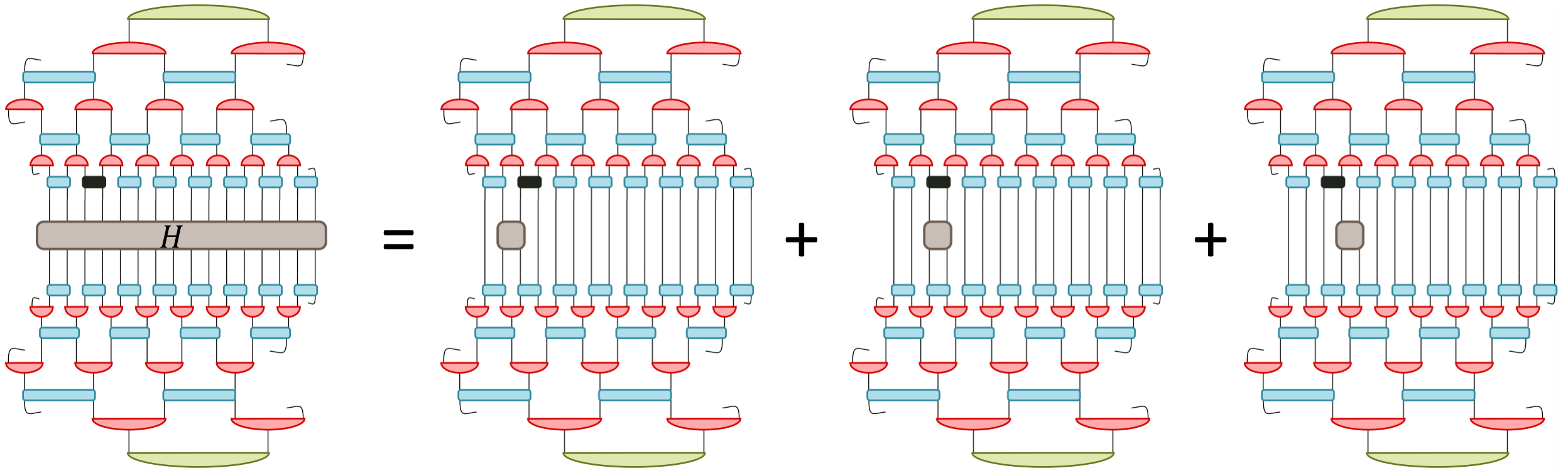


+ ... (N terms)

Minimize:
(over all tensors)

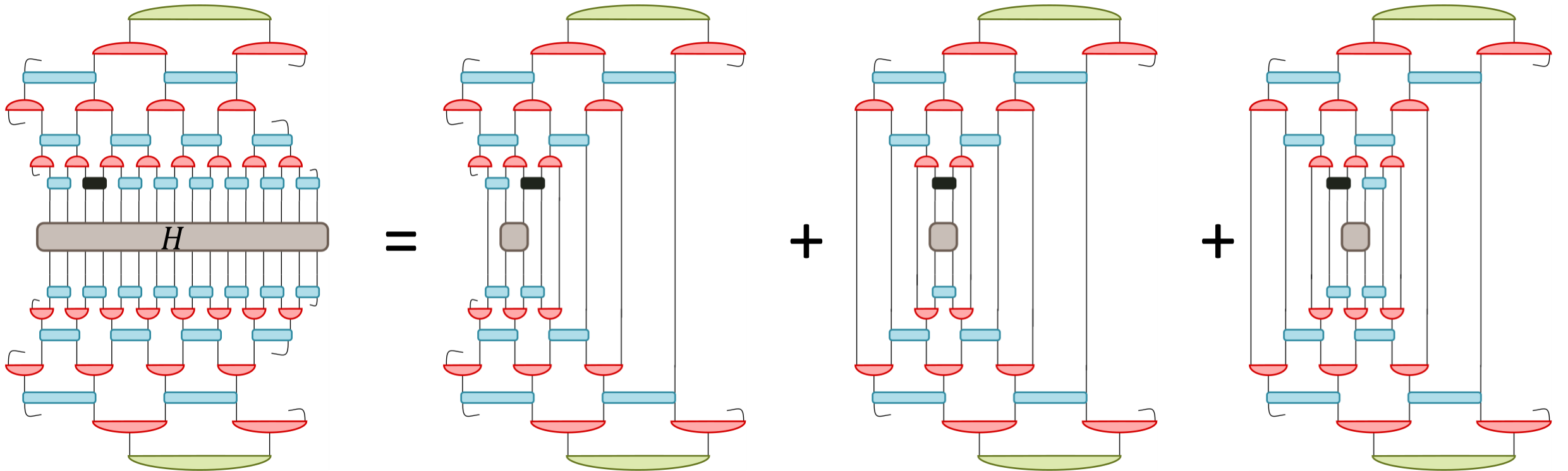
Optimize one tensor at time

Suppose we want to optimize the black disentangler



Only 3 terms contribute. (The black disentanglers cancels out in all other terms)

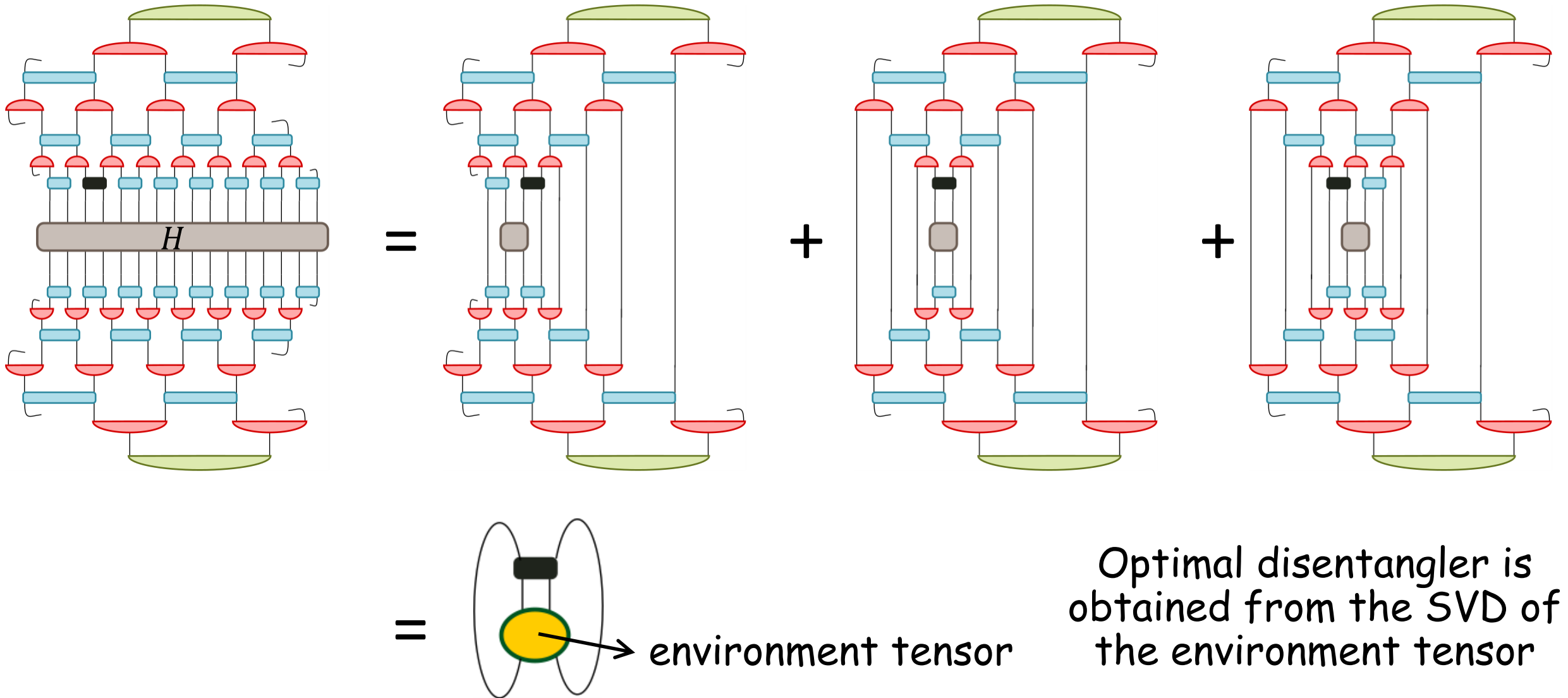
Suppose we want to optimize the black disentangler



Only 3 terms contribute. (The black disentanglers cancels out in all other terms)

Furthermore, these terms also simplify significantly

Suppose we want to optimize the black disentangler



In this way we can optimize one tensor at a time, say sweeping the tensor network from bottom to top

Repeat sweeps until e.g. energy converges

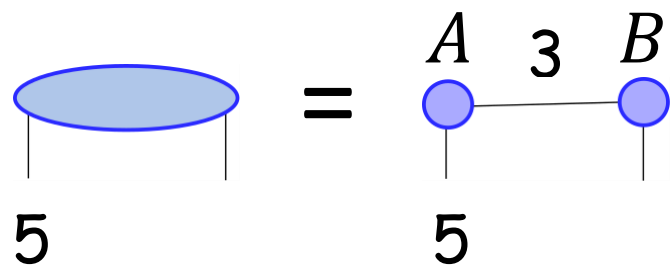
The numerically optimized MERA should describe an approximate RG flow as described earlier

PART 2

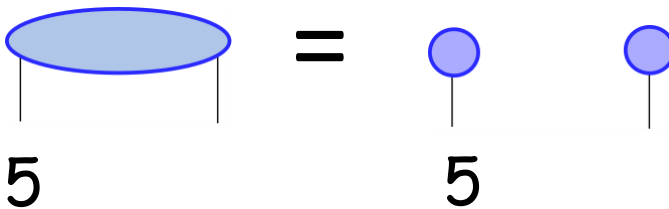
MERA and critical systems

Entanglement in the MERA

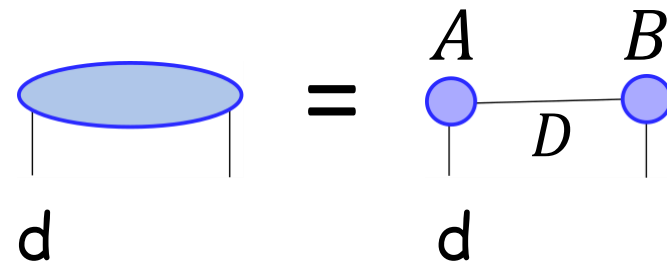
A simple observation ...



$$\text{Rank}(\rho_A) \leq 3$$

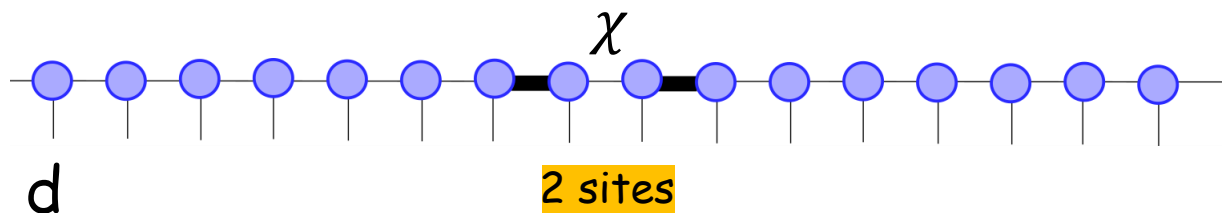


$$\text{Rank}(\rho_A) = 1$$



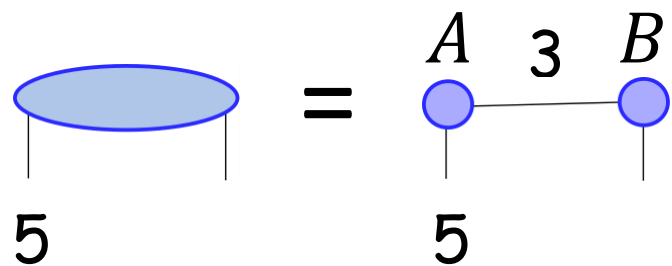
$$\text{Rank}(\rho_A) \leq \min(d, D)$$

iMPS

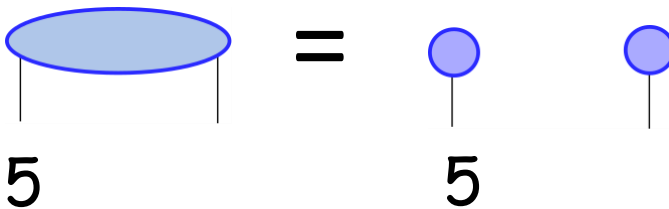


$$D = \chi^2$$

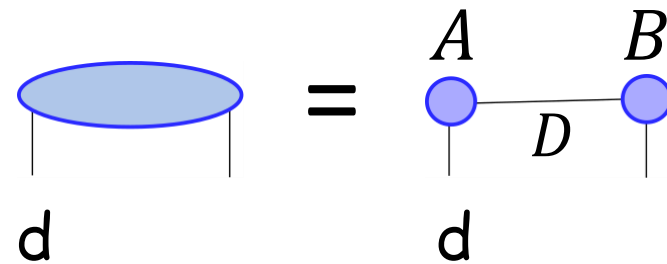
A simple observation ...



$$\text{Rank}(\rho_A) \leq 3$$

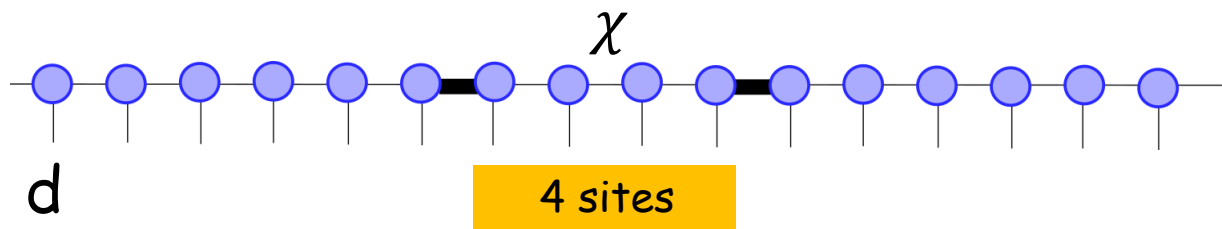


$$\text{Rank}(\rho_A) = 1$$



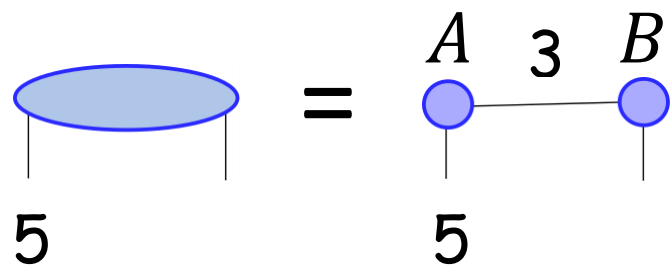
$$\text{Rank}(\rho_A) \leq \min(d, D)$$

iMPS

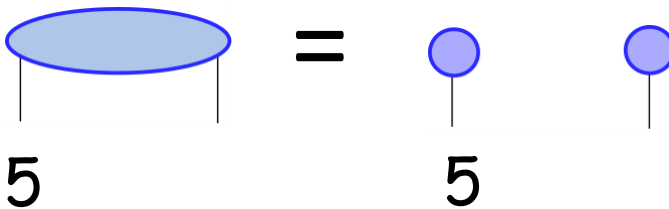


$$D = \chi^2$$

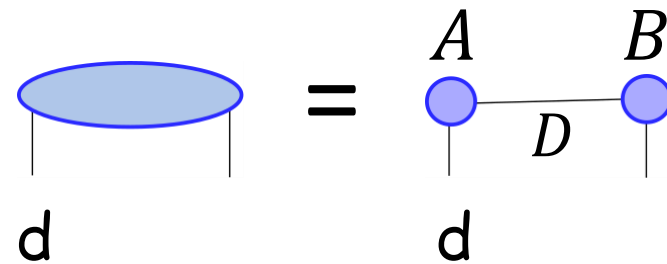
A simple observation ...



$$\text{Rank}(\rho_A) \leq 3$$

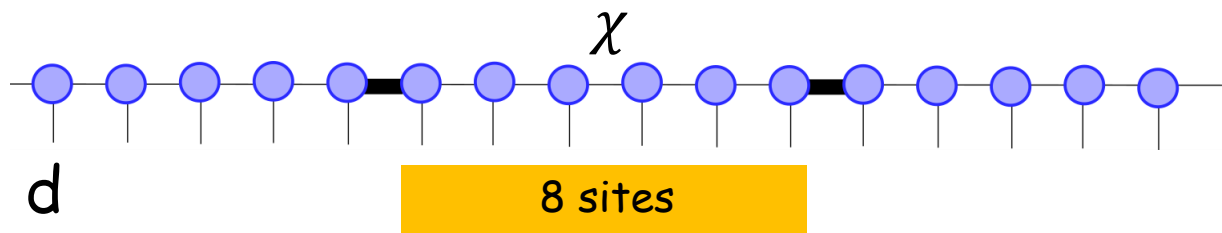


$$\text{Rank}(\rho_A) = 1$$



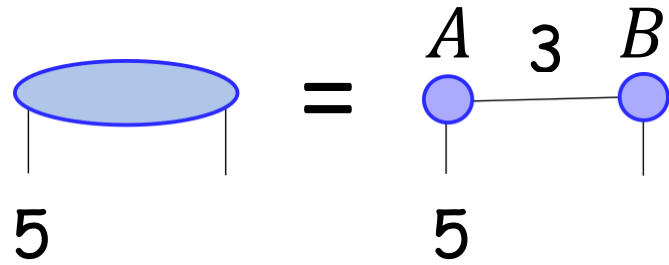
$$\text{Rank}(\rho_A) \leq \min(d, D)$$

iMPS

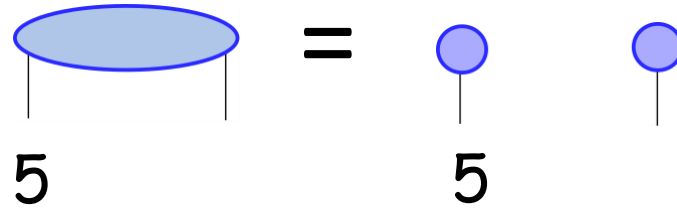


$$D = \chi^2$$

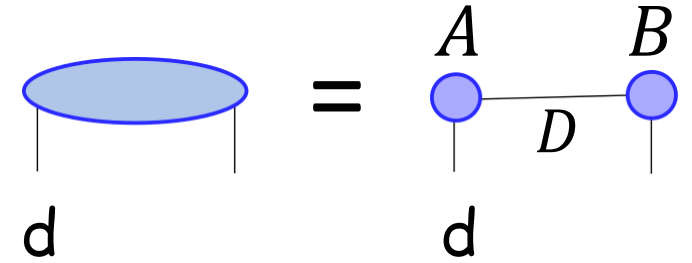
A simple observation ...



$$\text{Rank}(\rho_A) \leq 3$$

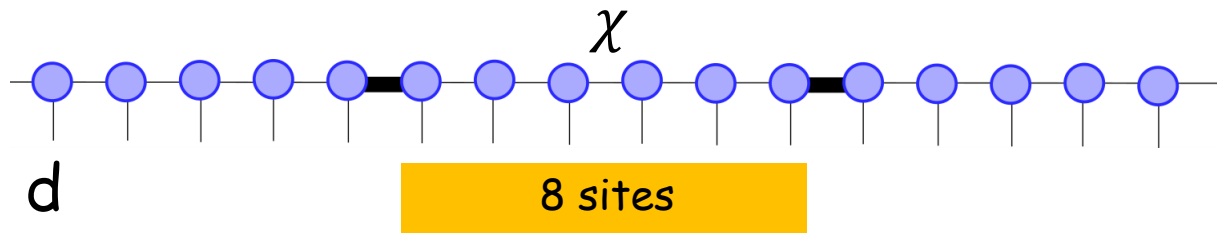


$$\text{Rank}(\rho_A) = 1$$



$$\text{Rank}(\rho_A) \leq \min(d, D)$$

iMPS

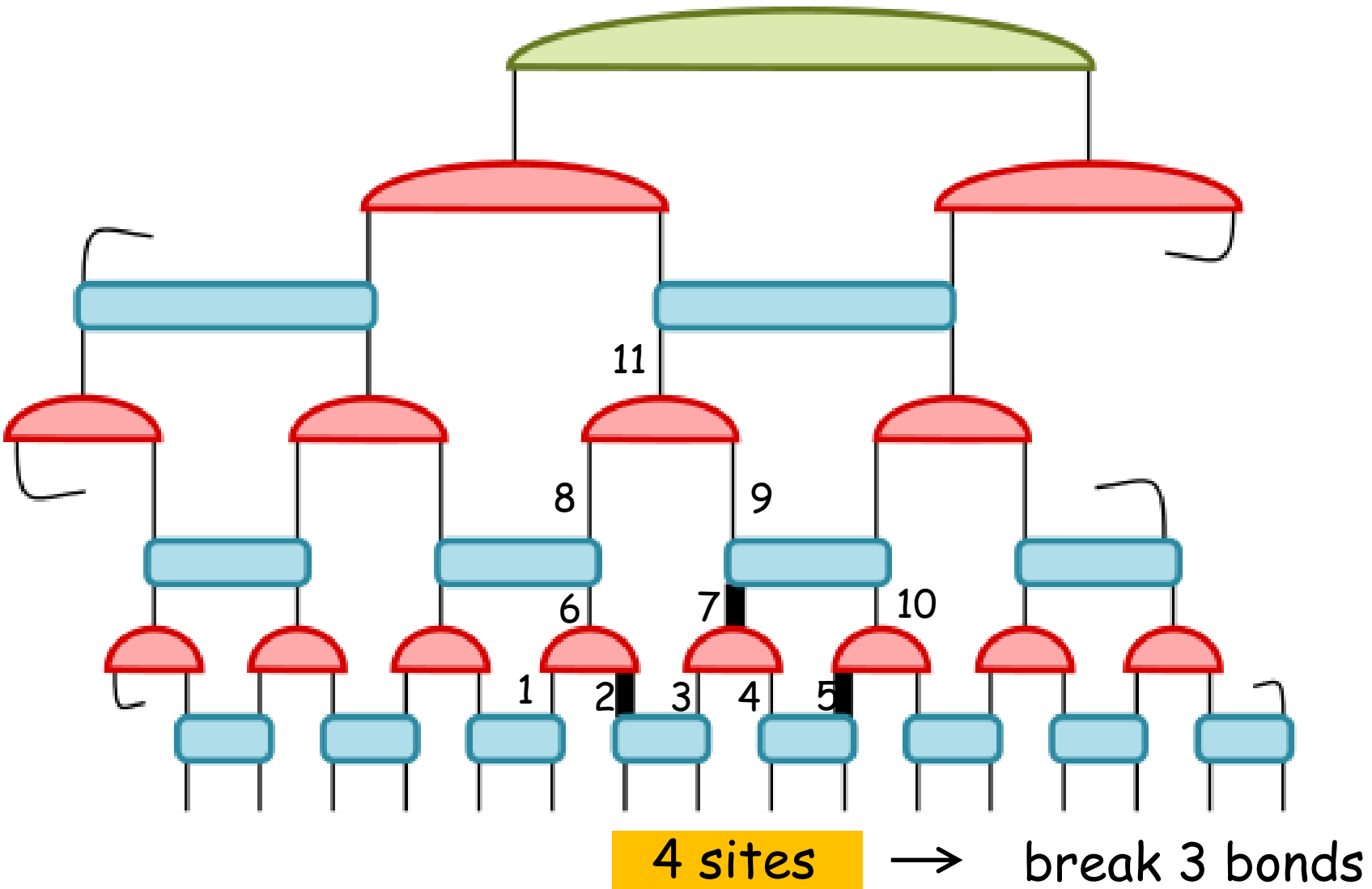


At some ℓ , $d^\ell > \chi^2$, thus $\text{Rank}(\rho_\ell) \leq \chi^2$

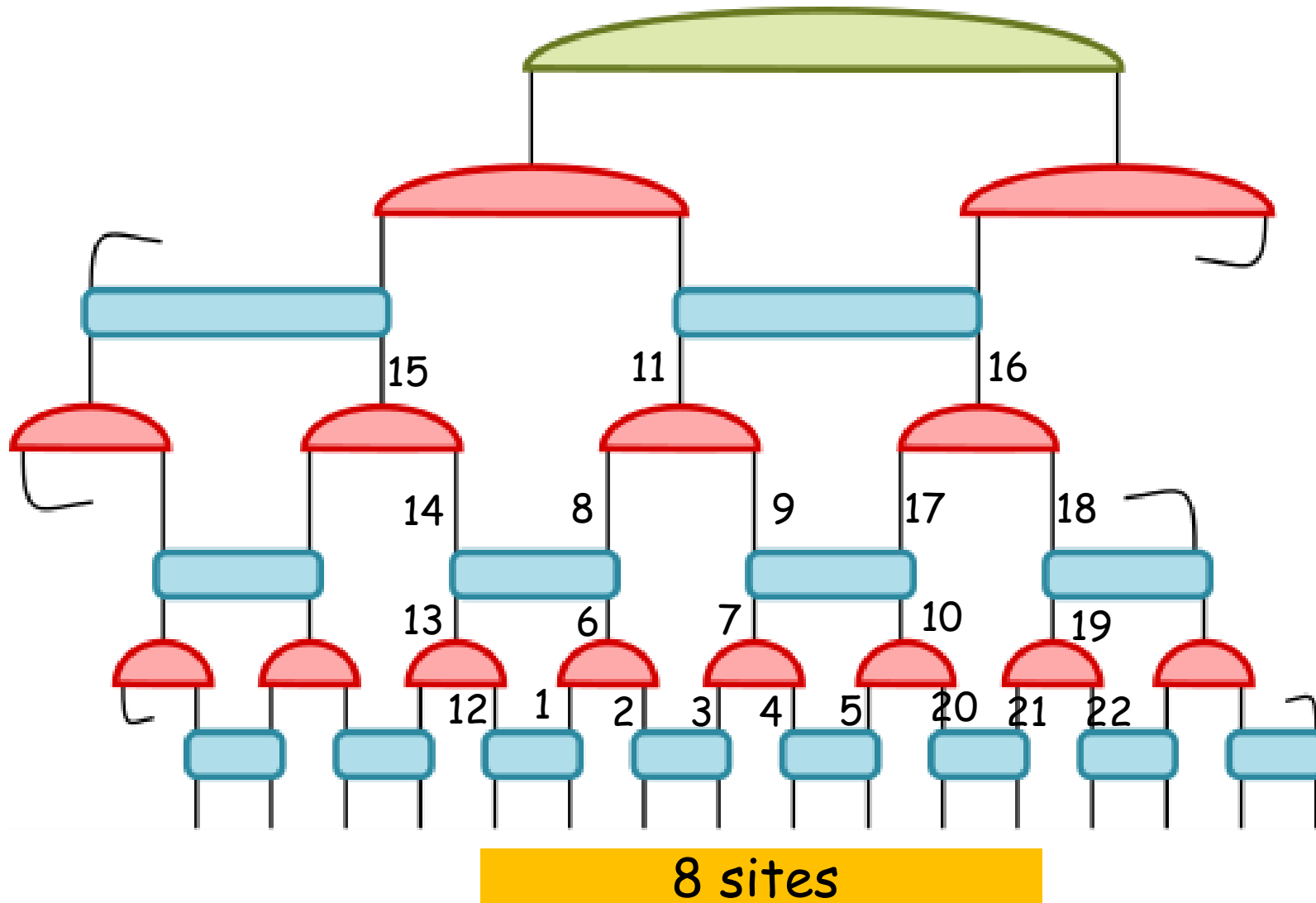
$$D = \chi^2$$

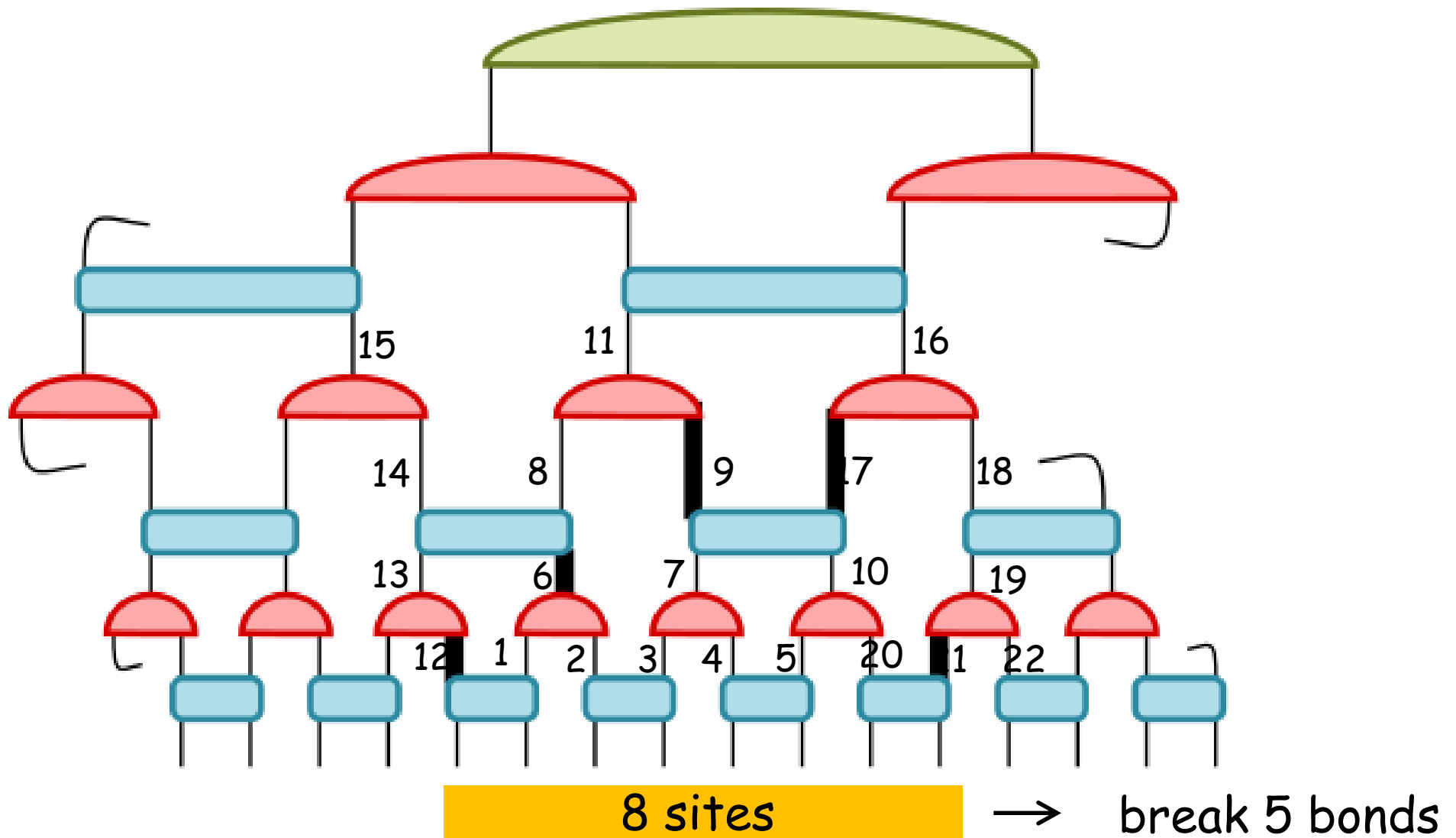
Matches the entanglement profile of gapped ground states

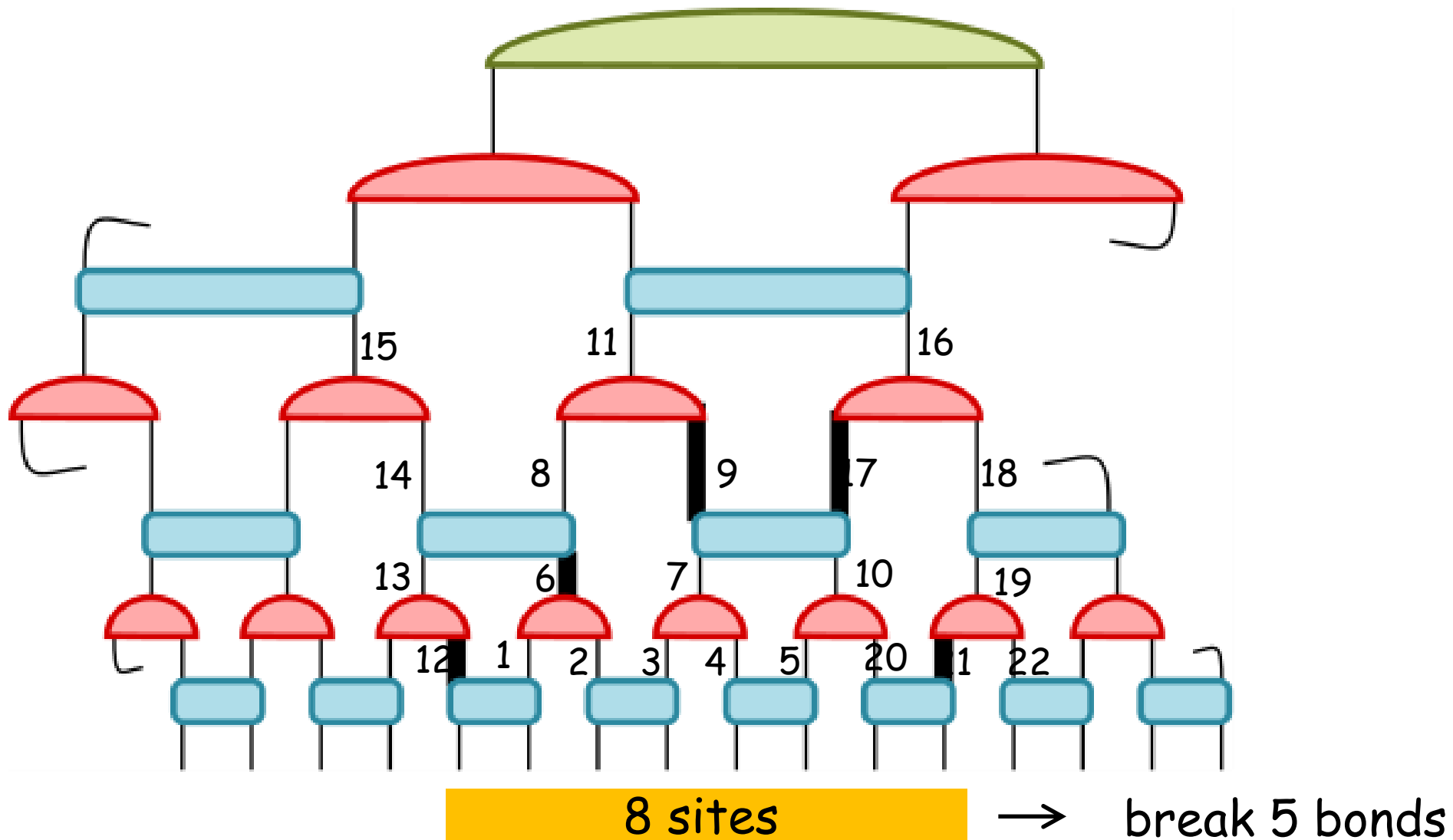
Quiz!



Quiz!







ℓ sites \rightarrow break $O(\log \ell)$ bonds \Rightarrow Entropy $\sim O(\log \ell)$

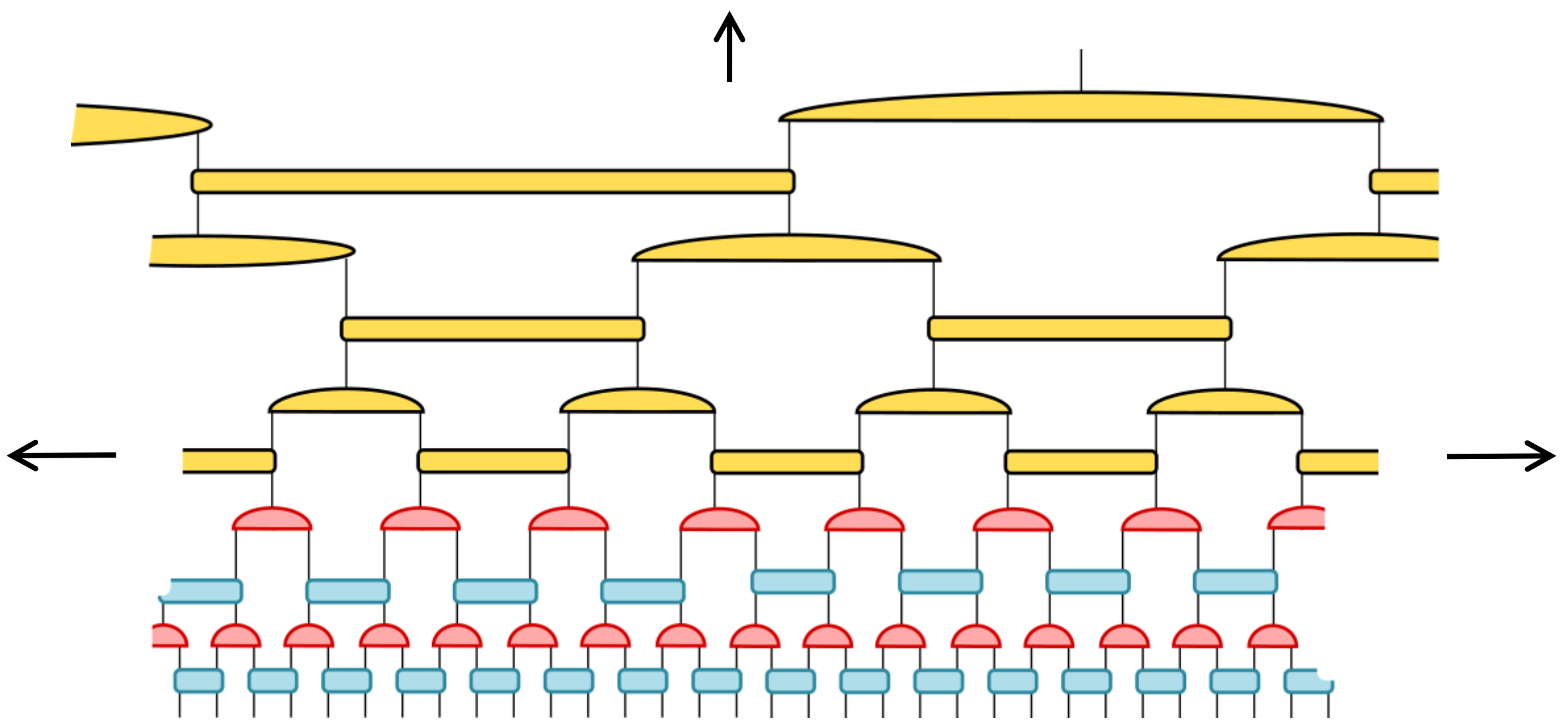
A random MERA state has logarithm scaling of entanglement entropy (similar to critical states)

One reason to hope that the MERA might be a **natural** ansatz for critical states

As another reason -- Next let us see how the MERA accurately captures the RG fixed point at criticality.

The *infinite* MERA:

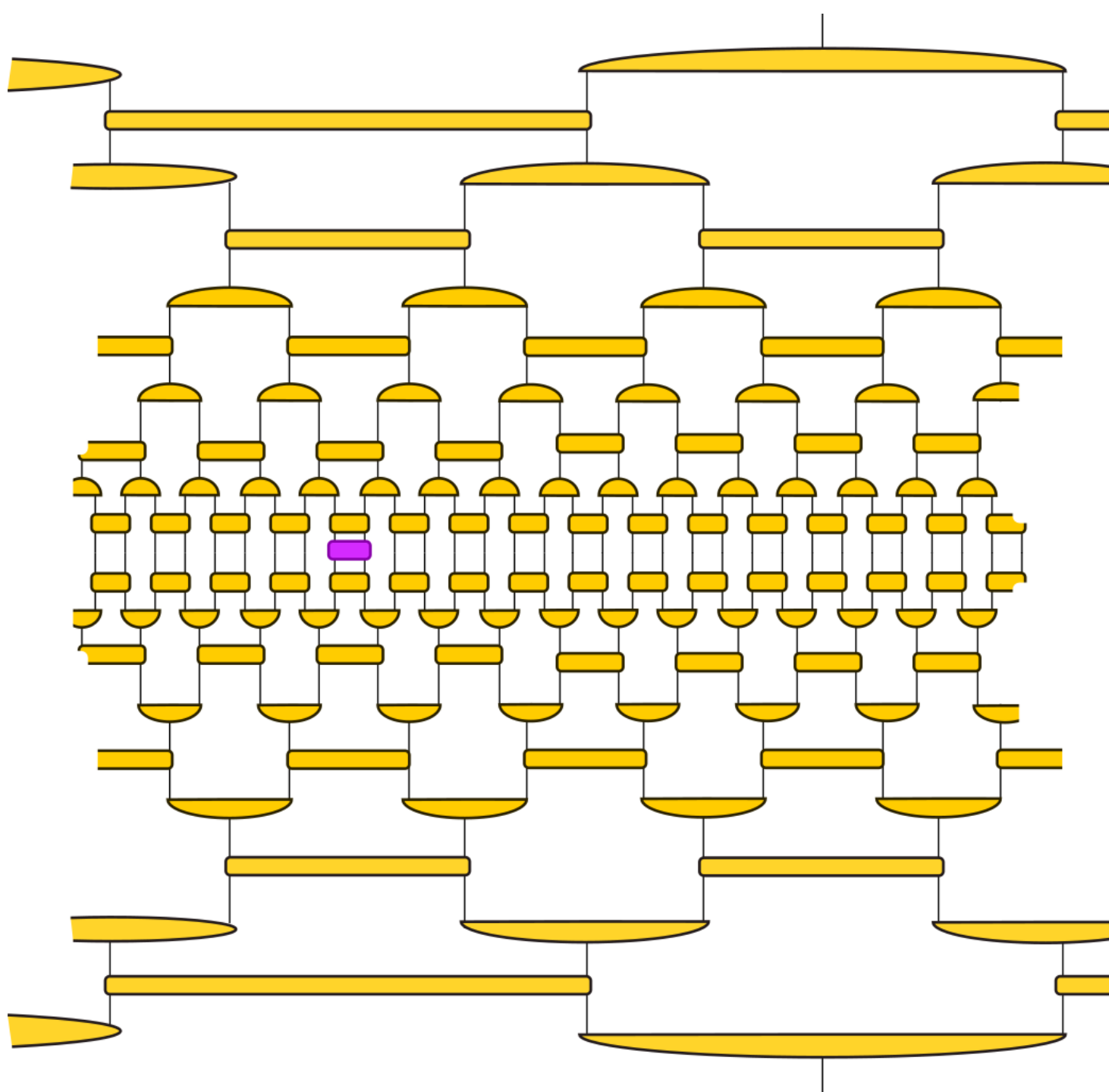
a lattice approximation of a 2d CFT

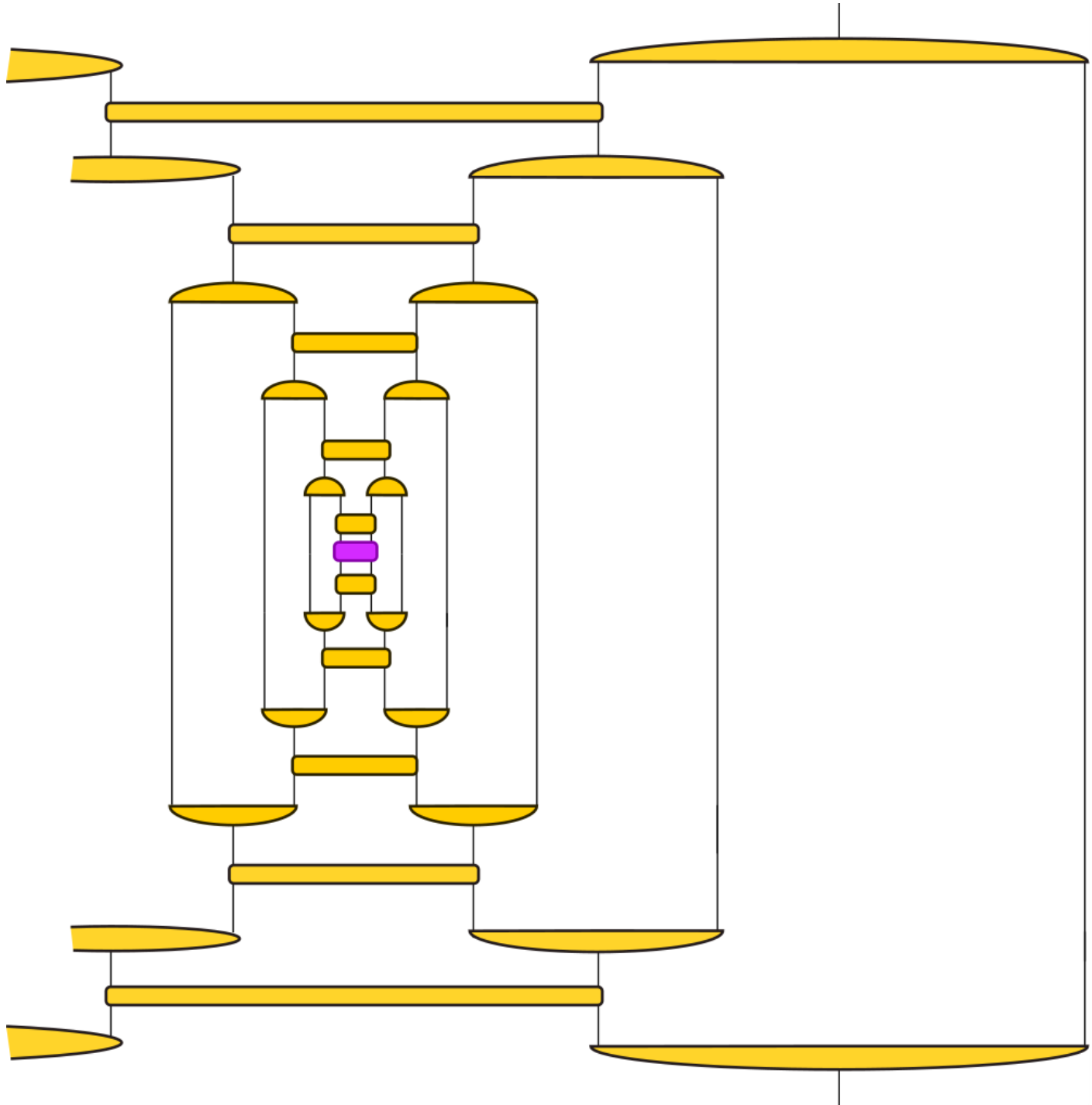


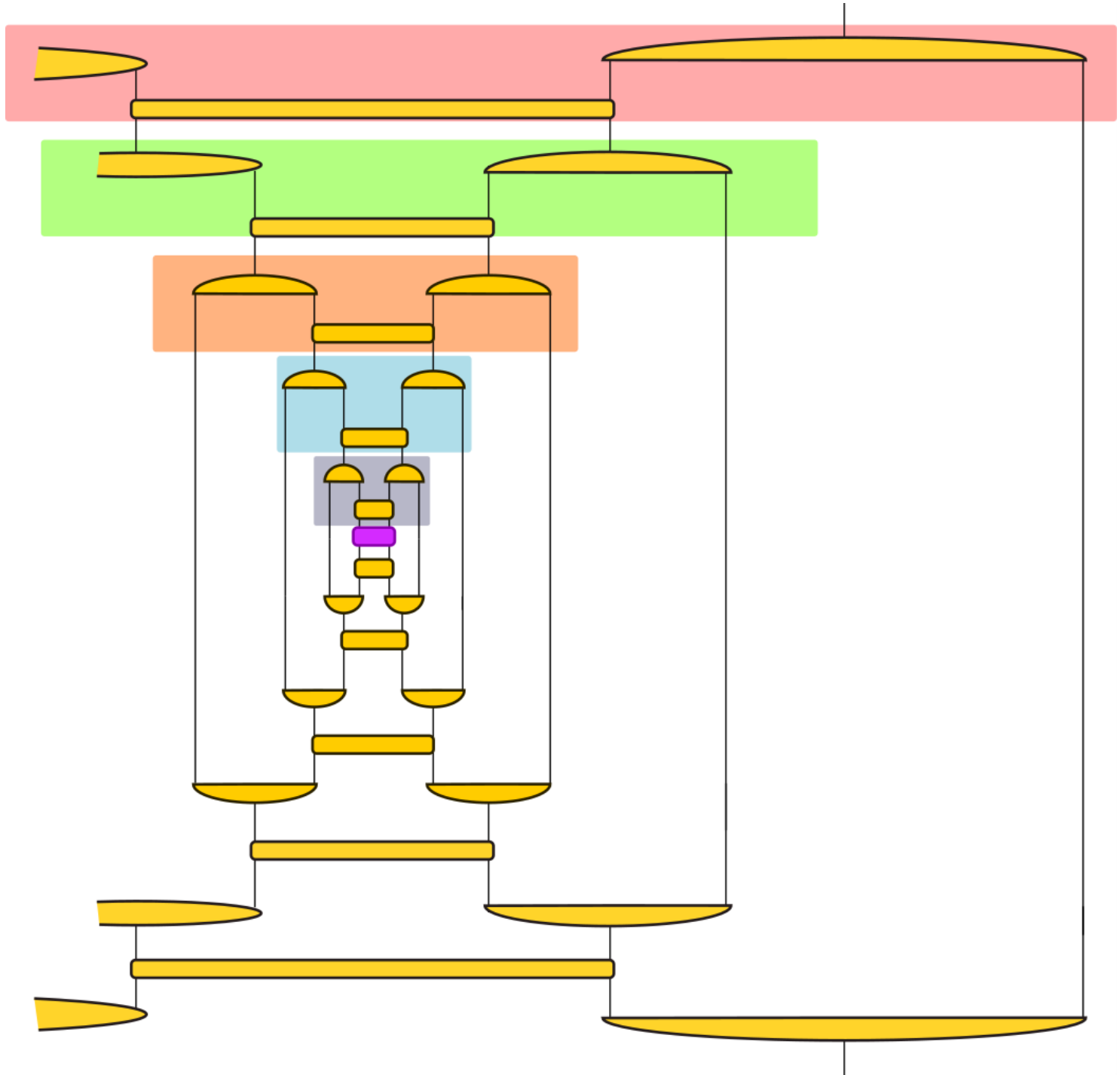
Choose identical tensors in each layer: **approximate** translation-invariance.
Choose identical tensors after some scale: **exact** scale-invariance.

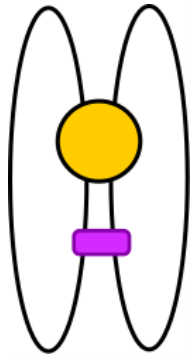
So the infinite MERA can be efficiently stored.

But how can we efficiently compute expectation values from such an infinite tensor network?

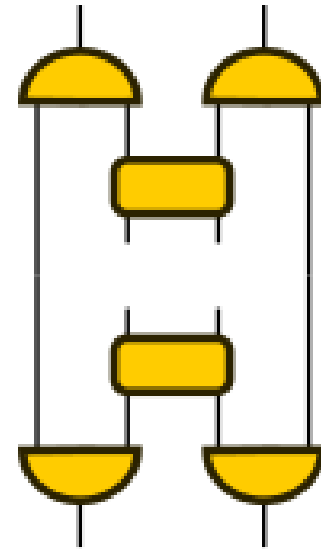








dominant eigenvector



"descending superoperator"

So the infinite MERA is also an efficient ansatz.

How do we use it in practice?

Number of **transitional layers** is an input parameter in numerics

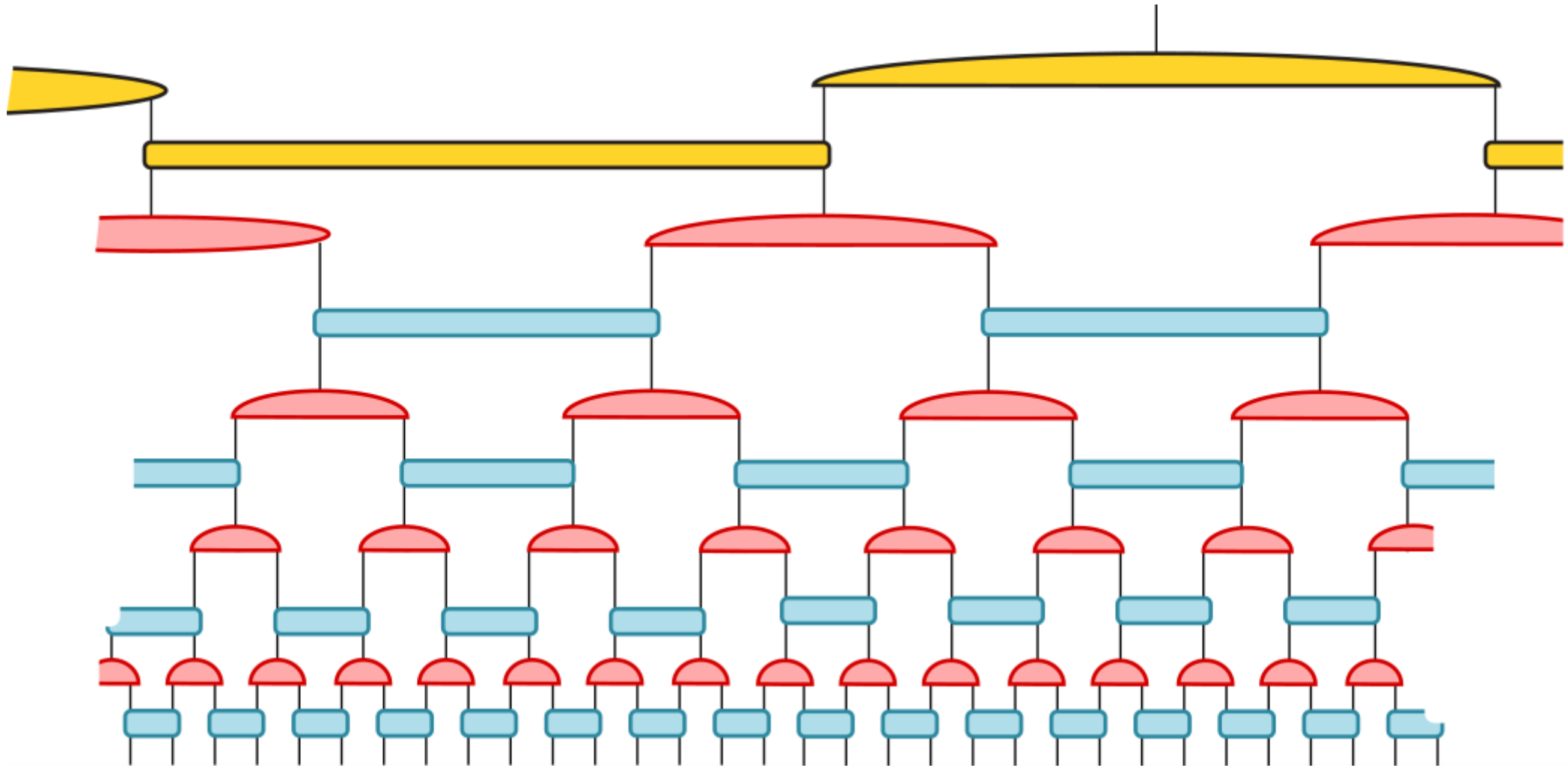
Converge energy for a small number of transitional layers

Increase the number of layers and check if energy lowers

Keep adding layers till the energy has converged (say L layers)

This numerically optimized MERA encodes a RG flow where the Hamiltonian reaches an approximate fixed point after L steps

$$H \rightarrow H_1 \rightarrow H_2 \rightarrow \cdots \rightarrow H_L \rightarrow H_L \rightarrow H_L \rightarrow \cdots$$



The **yellow tensors** encode a lot of important information about the 2d CFT that describes the critical system in the continuum limit

Conformal Field Theory in two dimensions

In 2d, conformal symmetry described by the infinite dimensional Virasoro algebra

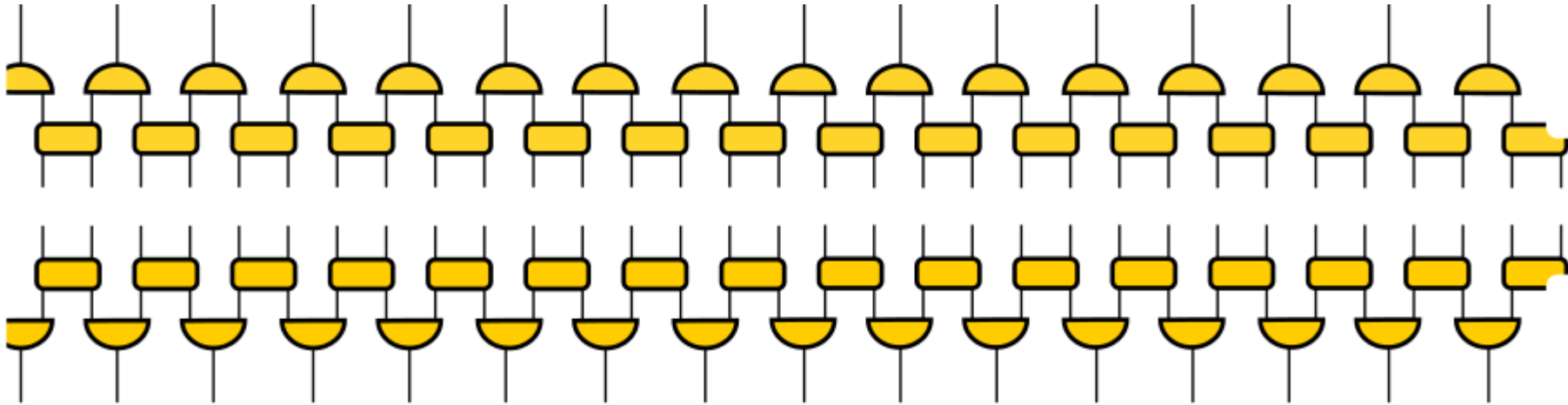
Scaling field operators $\phi_1, \phi_2, \phi_3, \dots$ with scaling dimensions $\Delta_1, \Delta_2, \Delta_3, \dots$

Scaling fields are eigenoperators of the dilation operator, and the scaling dimensions are the corresponding eigenvalues

2-point correlators $\langle \phi_\alpha(x) \phi_\beta(y) \rangle \sim \frac{C_{\alpha\beta}}{(x-y)^{\Delta_\alpha + \Delta_\beta}}$ (polynomial decay)

3-point correlators $\langle \phi_\alpha(x) \phi_\beta(y) \phi_\gamma(z) \rangle \sim \frac{C_{\alpha\beta\gamma}}{(x-y)^\dots (y-z)^\dots (x-z)^\dots}$ (OPE coefficients)

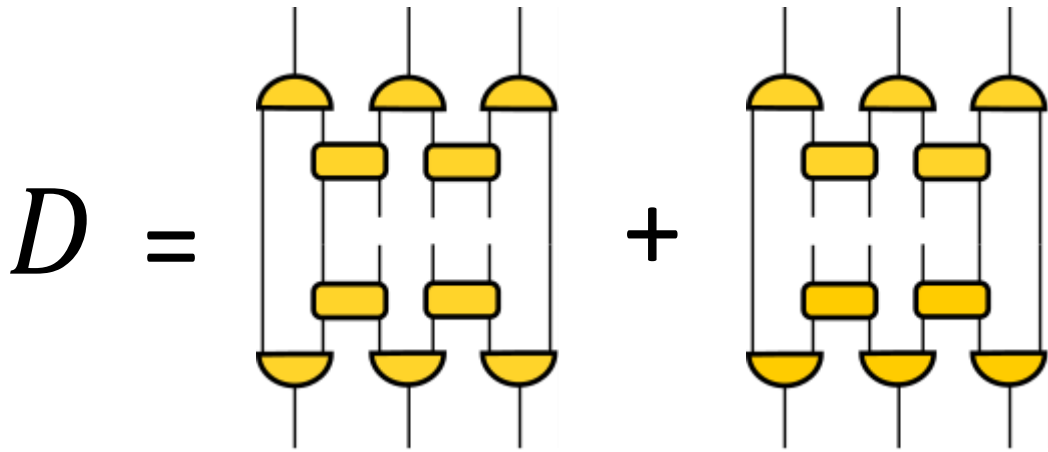
Dilation operator in the MERA (the scaling superoperator)



But we have seen that operators generally **shrink** under coarse-graining
(But eigenoperators cannot shrink!)

3-site operators remain 3-site, so we can find 3-site eigenoperators

3-site scaling superoperator



Note: This data can also be extracted from the MPS, but its easier and more intuitive with the MERA

Numerical observations

- Eigenvalues of D match some of the scaling dimensions
- Eigenvectors are some lattice representation of the scaling fields
- Can compute the 3-point correlator of these scaling operators to determine the OPE coefficients
- Patches of the fixed point tensors also give lattice representations of conformal transformations
- Fixed point tensors encode essential CFT data and also know about the conformal symmetry

iMERA \approx 2d CFT on the lattice

iMERA [?] = 2d CFT on the lattice

Open question

Does there exist an iMERA that encodes the CFT data **exactly**?

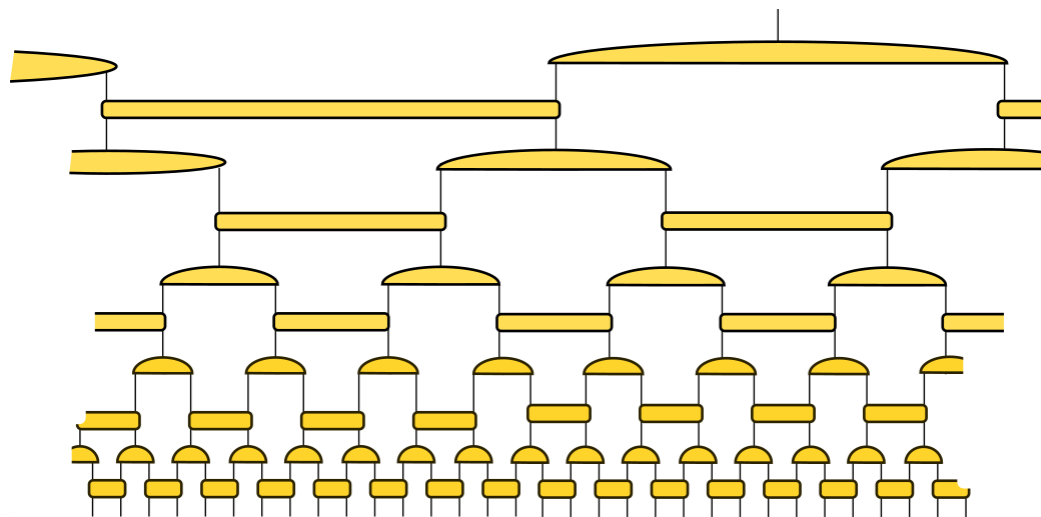
Of course, only a finite part of the infinite CFT data.

One requirement: incorporate the vacuum conformal symmetry in the MERA.

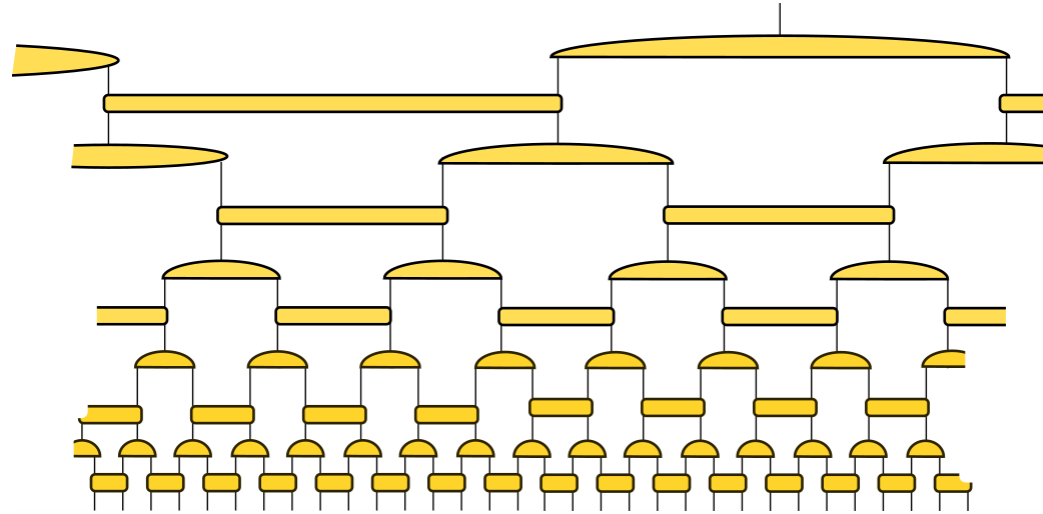
Vacuum conformal symmetry: scale invariance, translation invariance, special conformal transformations (those generated by the positive generators)

Fantasy/Folklore: If the MERA is scale + translation invariance and it is the ground state of a local Hamiltonian then it might possibly be an **exact** description of a 2d CFT on the lattice

Scale invariance

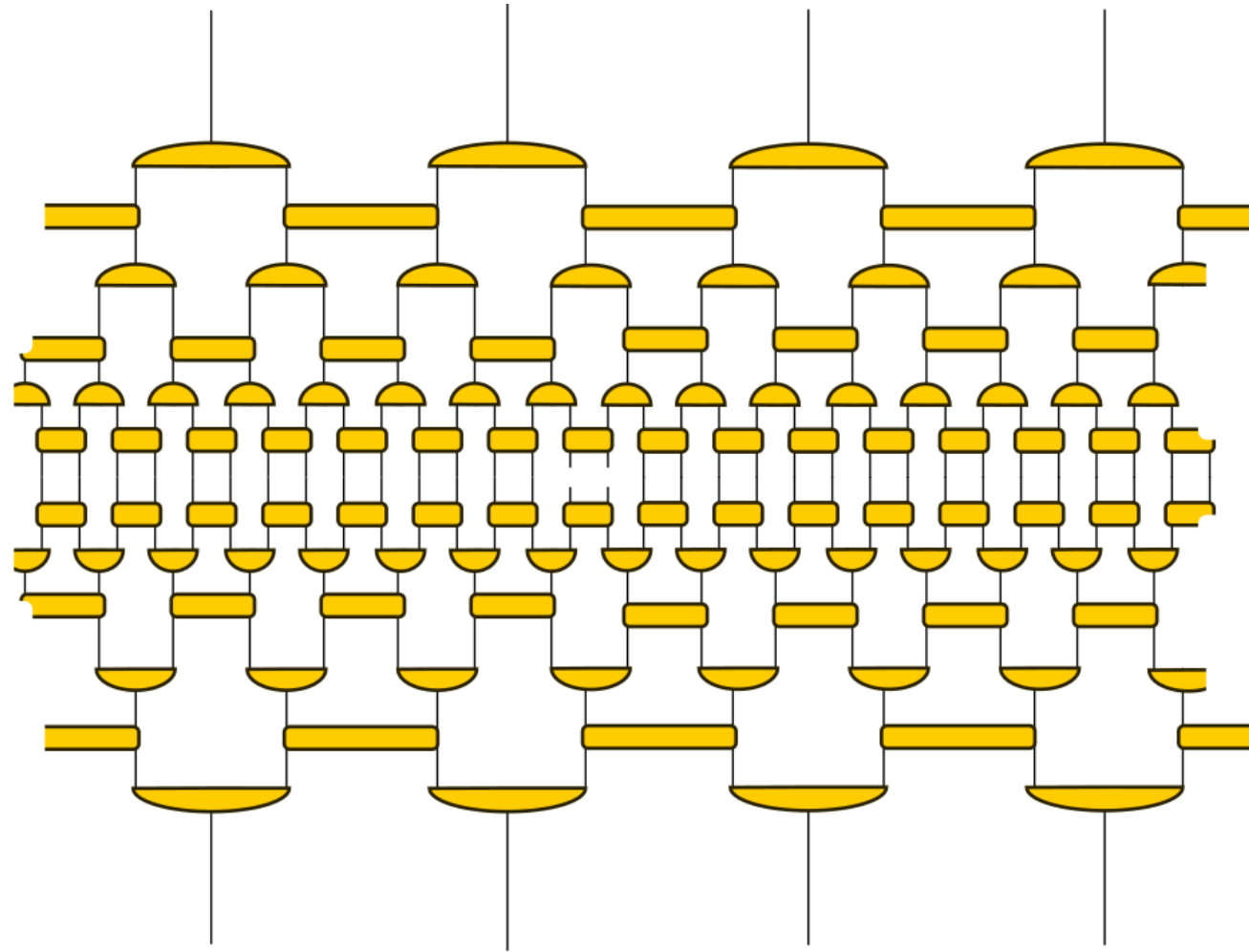


Scale invariance

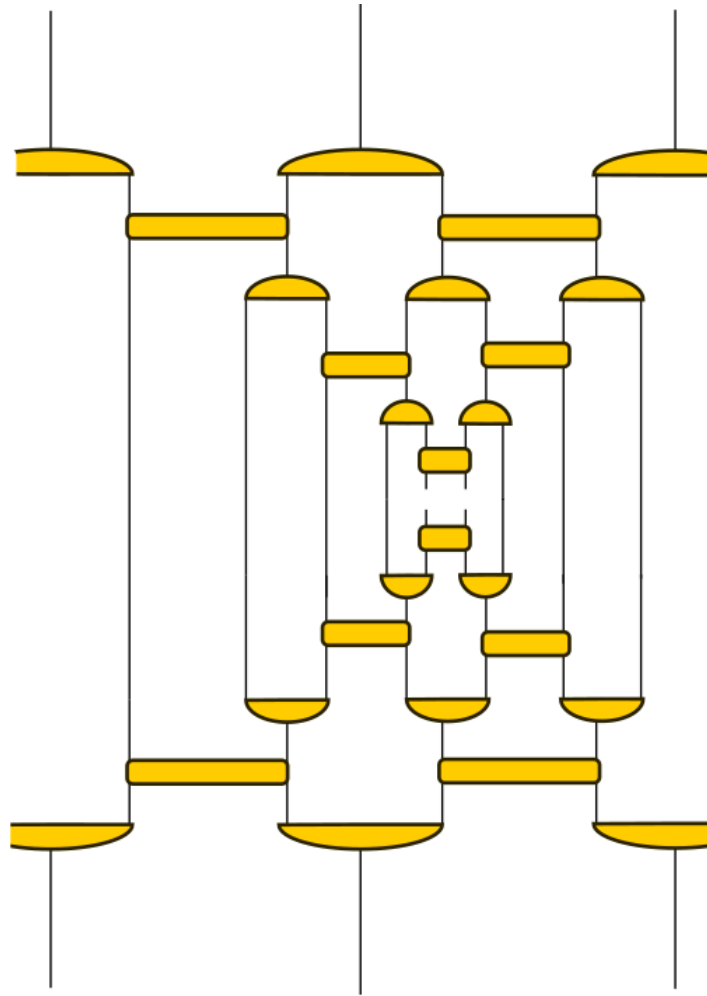


But most of these states are **NOT**
translation invariant

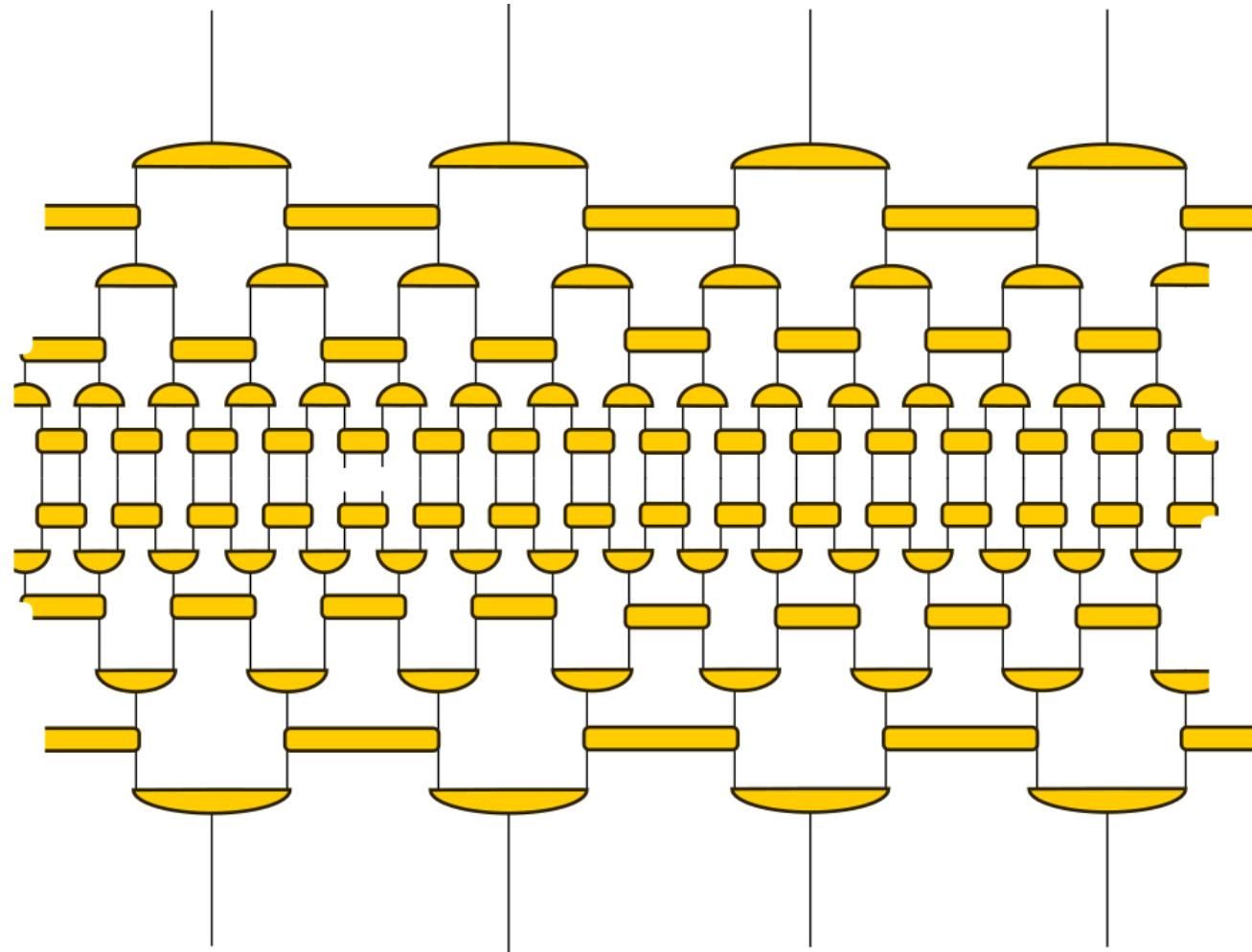
Translation invariance



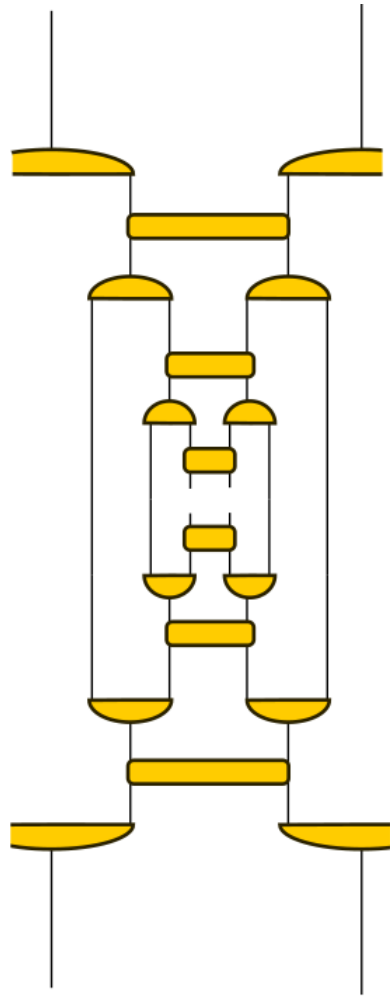
Translation invariance



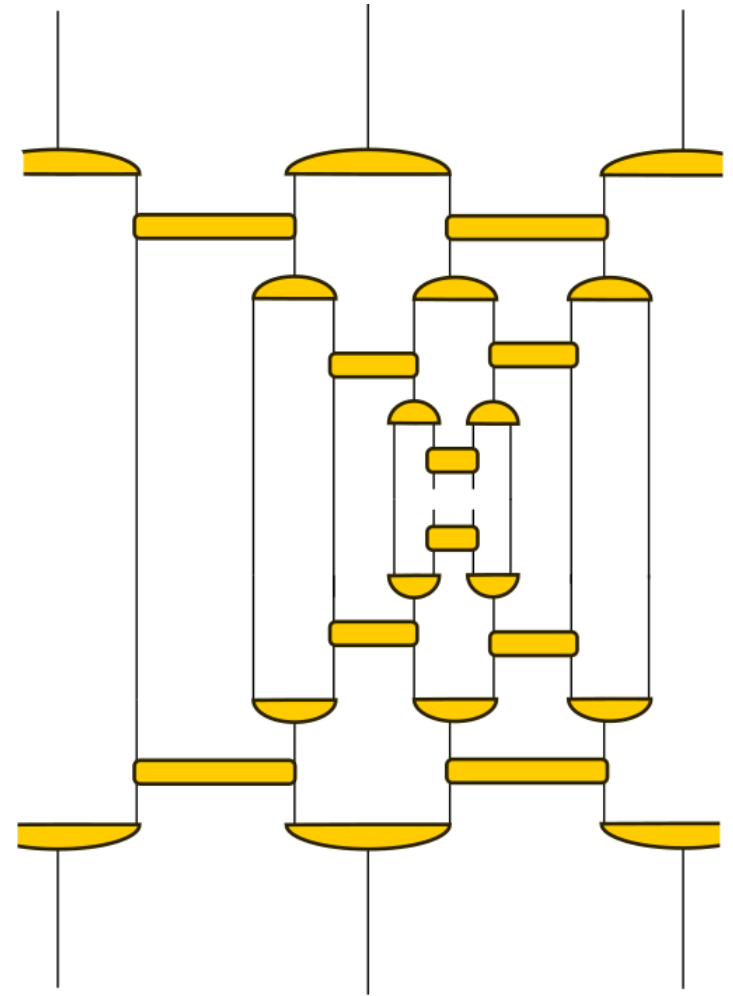
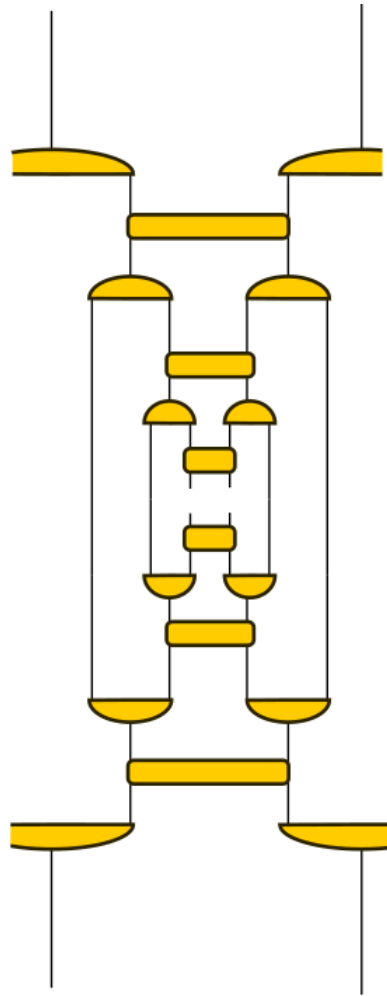
Translation invariance



Translation invariance



Translation invariance



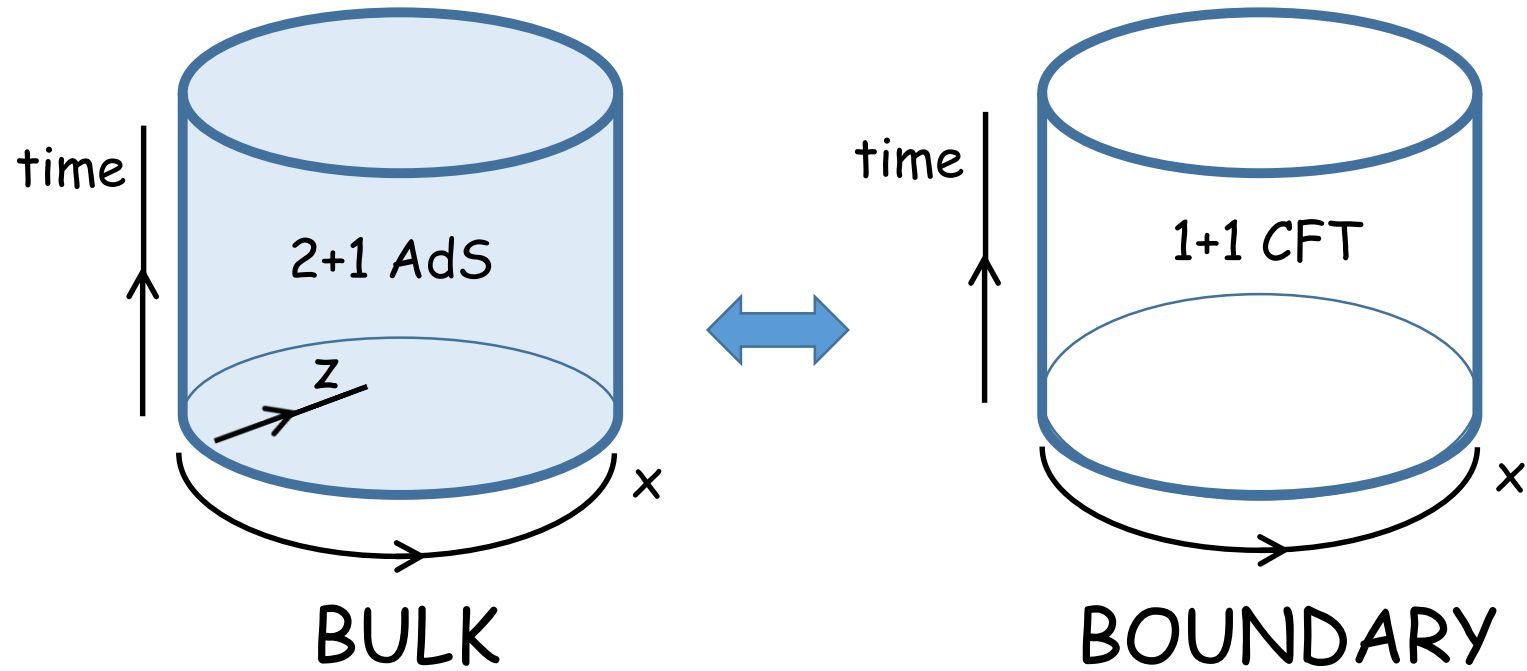
We can get approximate translation invariance in numerical simulations (by e.g. averaging environment tensors across the lattice)

Increasing bond dimension usually improves translation invariance and also the estimation of the CFT data

Can we make translation symmetry exact. (Open Problem)

MERA and holography: basic intro

The AdS/CFT correspondence



Gravity in 2+1 anti-deSitter spacetime



1+1 CFT

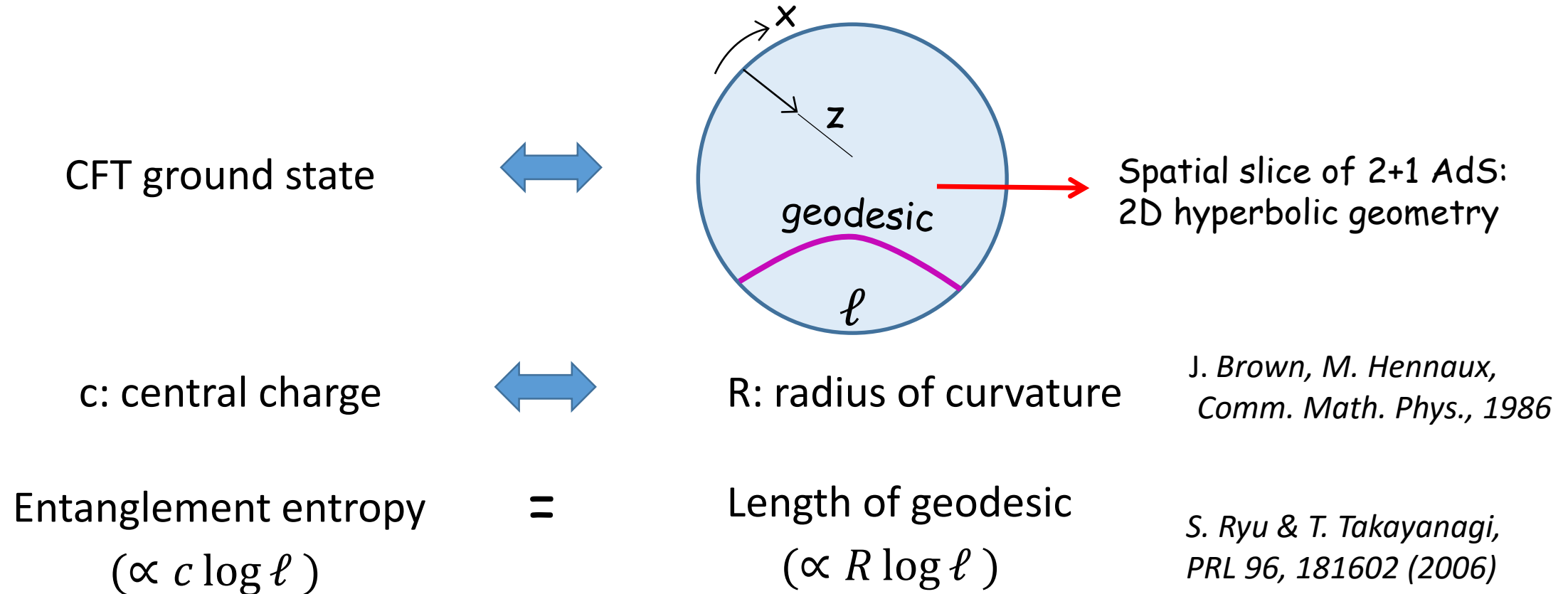
Extra dimension on the gravity side

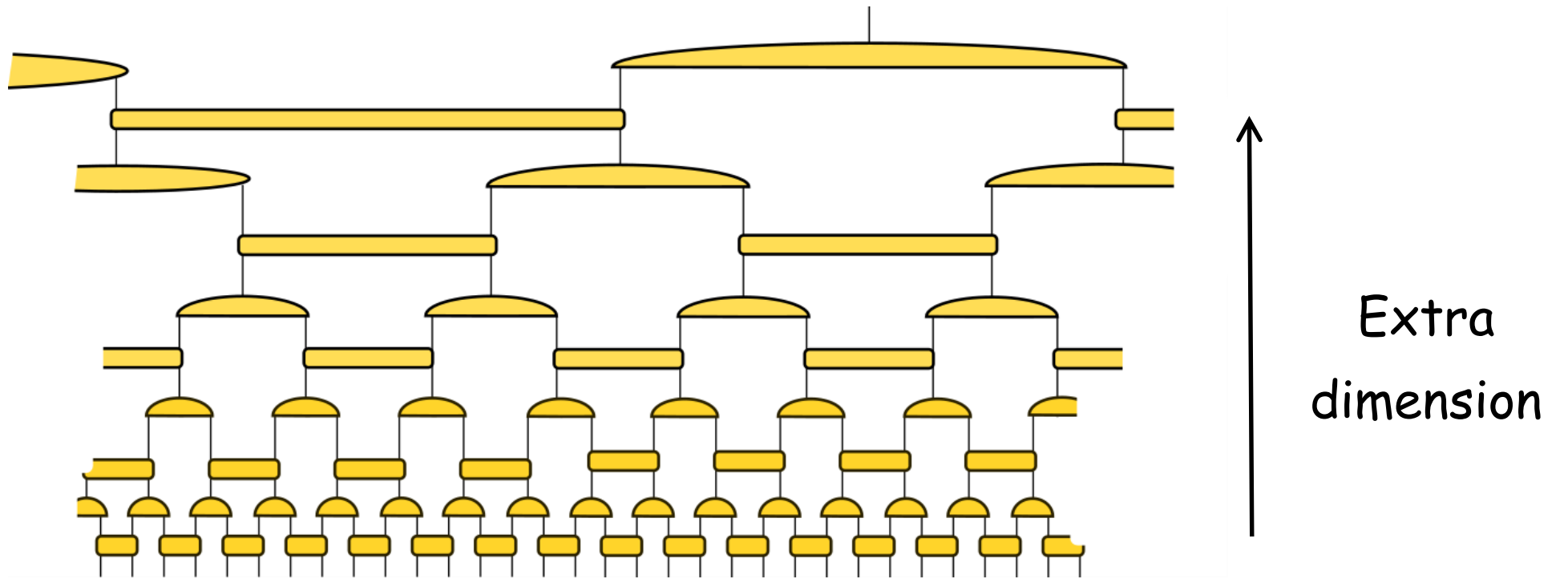


RG direction

The AdS/CFT correspondence

Certain 'Large' N QFTs correspond to (semi)-classical bulk gravity.



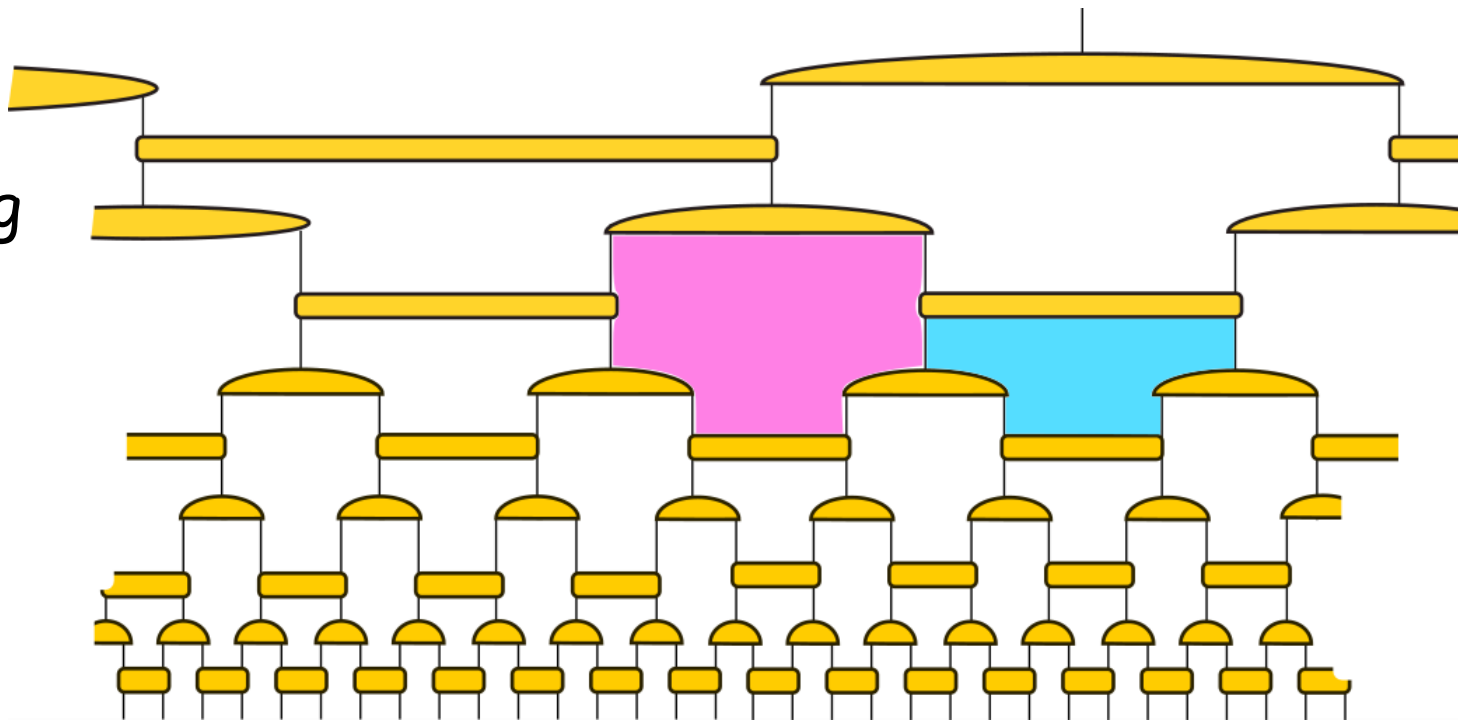


MERA is a good approximation of a 2d CFT

Has an extra dimension that corresponds to length scale

Is based on a negatively curved geometry

MERA as a tiling
of the plane
(tiles are flat
and regular)

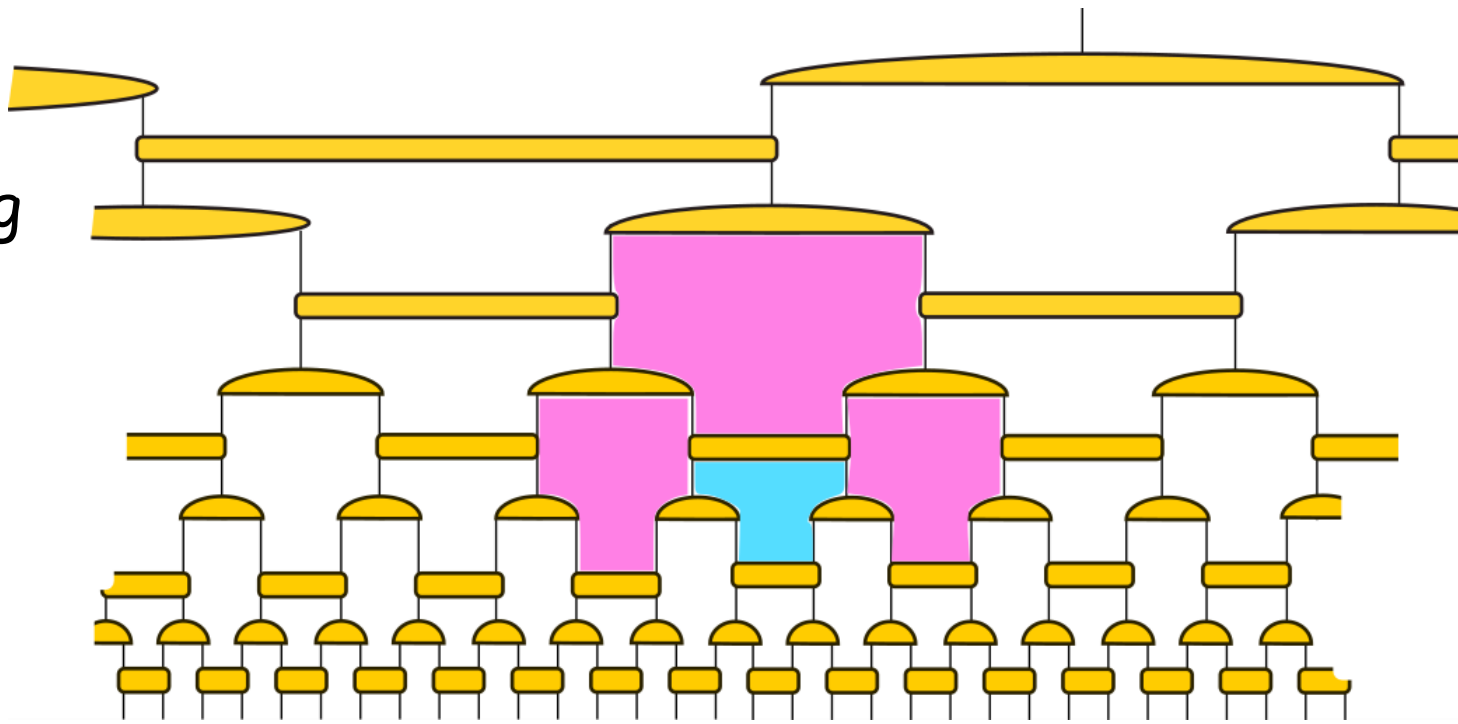


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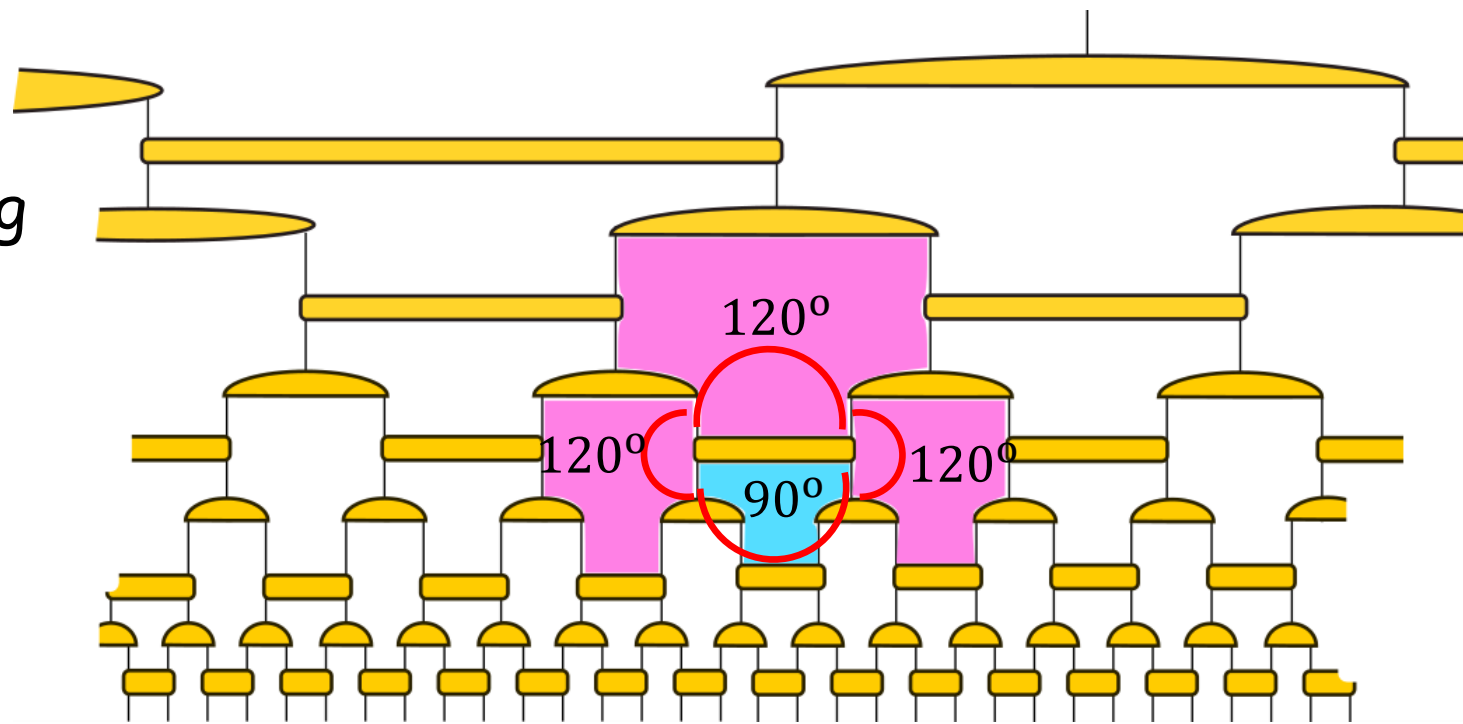
Extra
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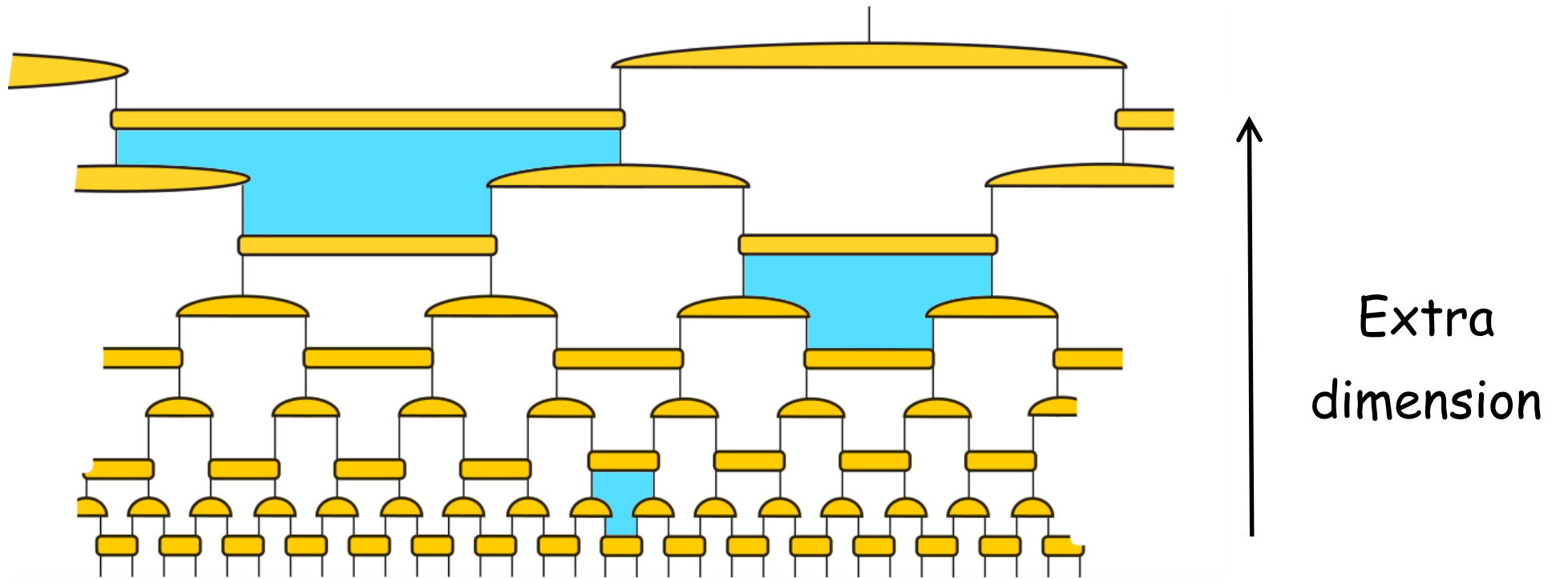


Extra
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MERA is a good approximation of a 2d CFT

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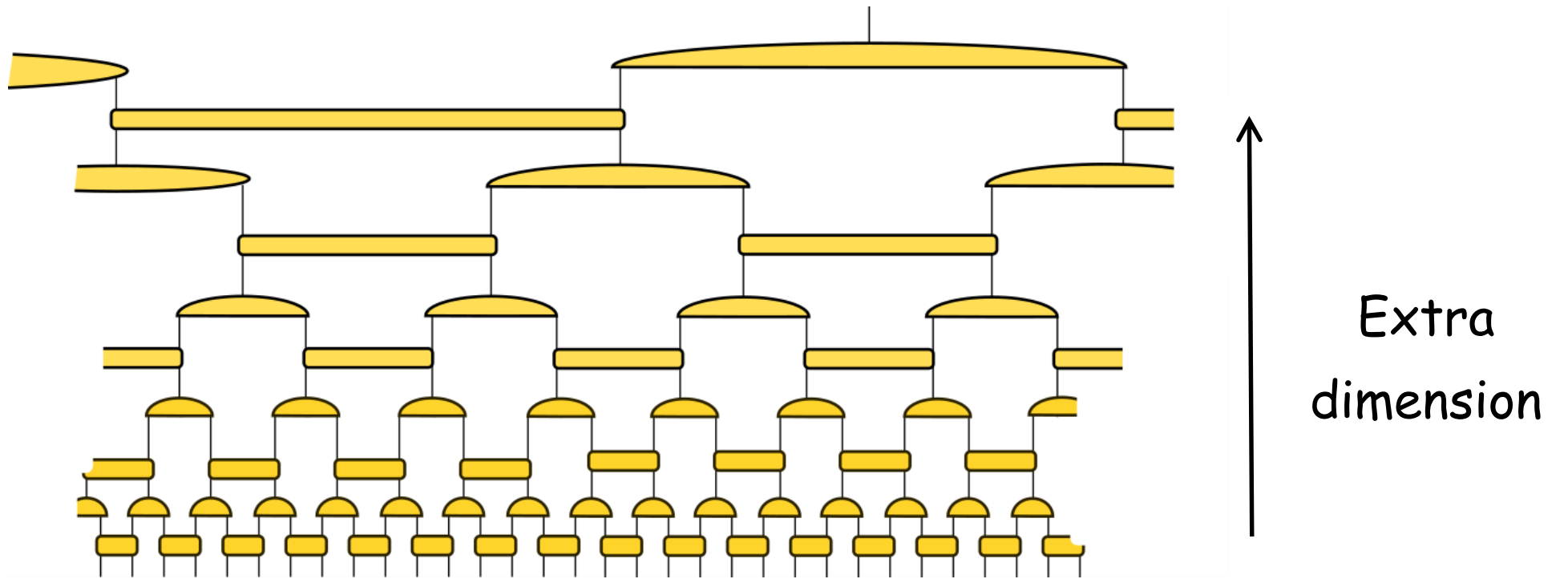
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MERA is a good approximation of a 2d CFT

Has an extra dimension that corresponds to length scale

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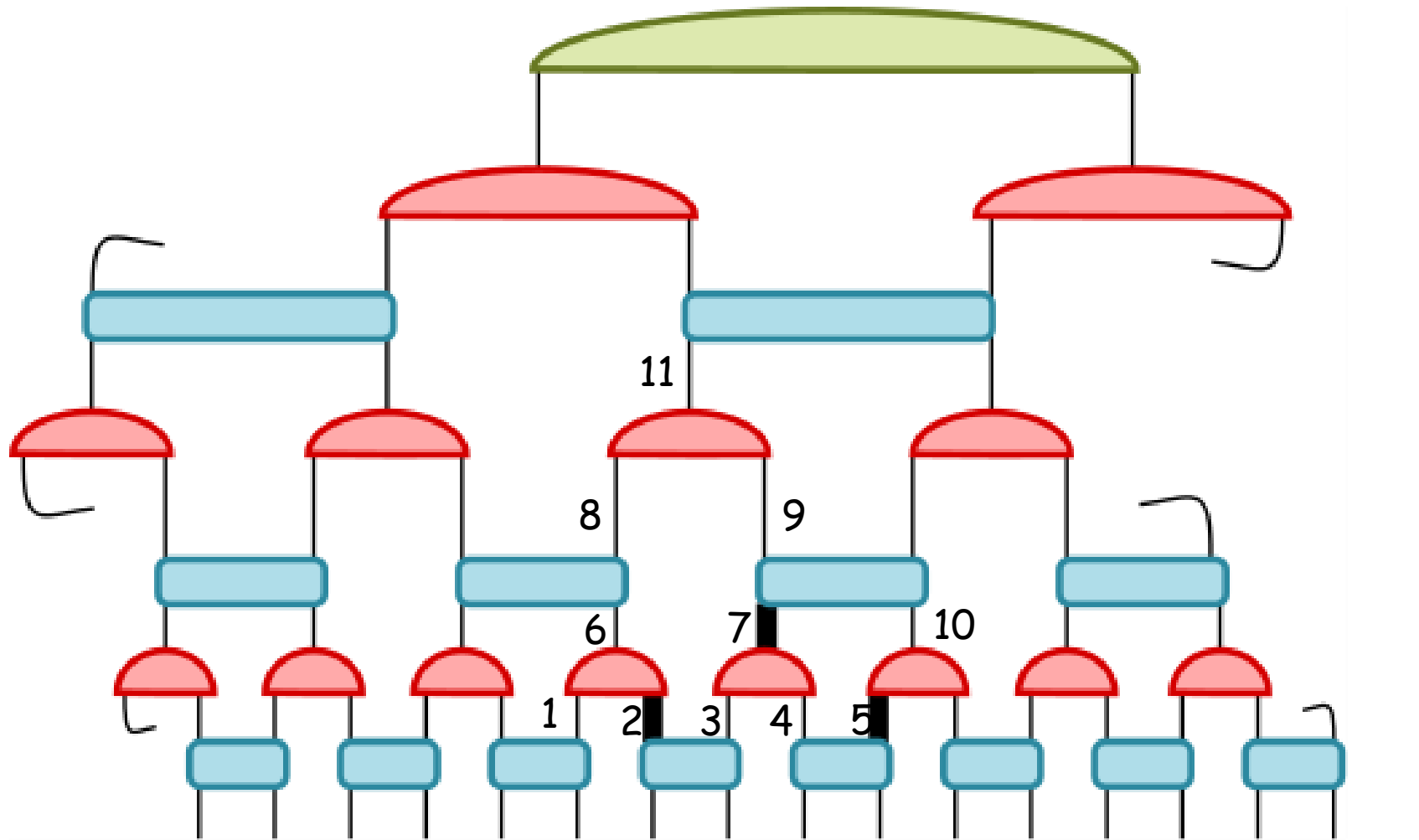


MERA is a good approximation of a 2d CFT

Has an extra dimension that corresponds to length scale

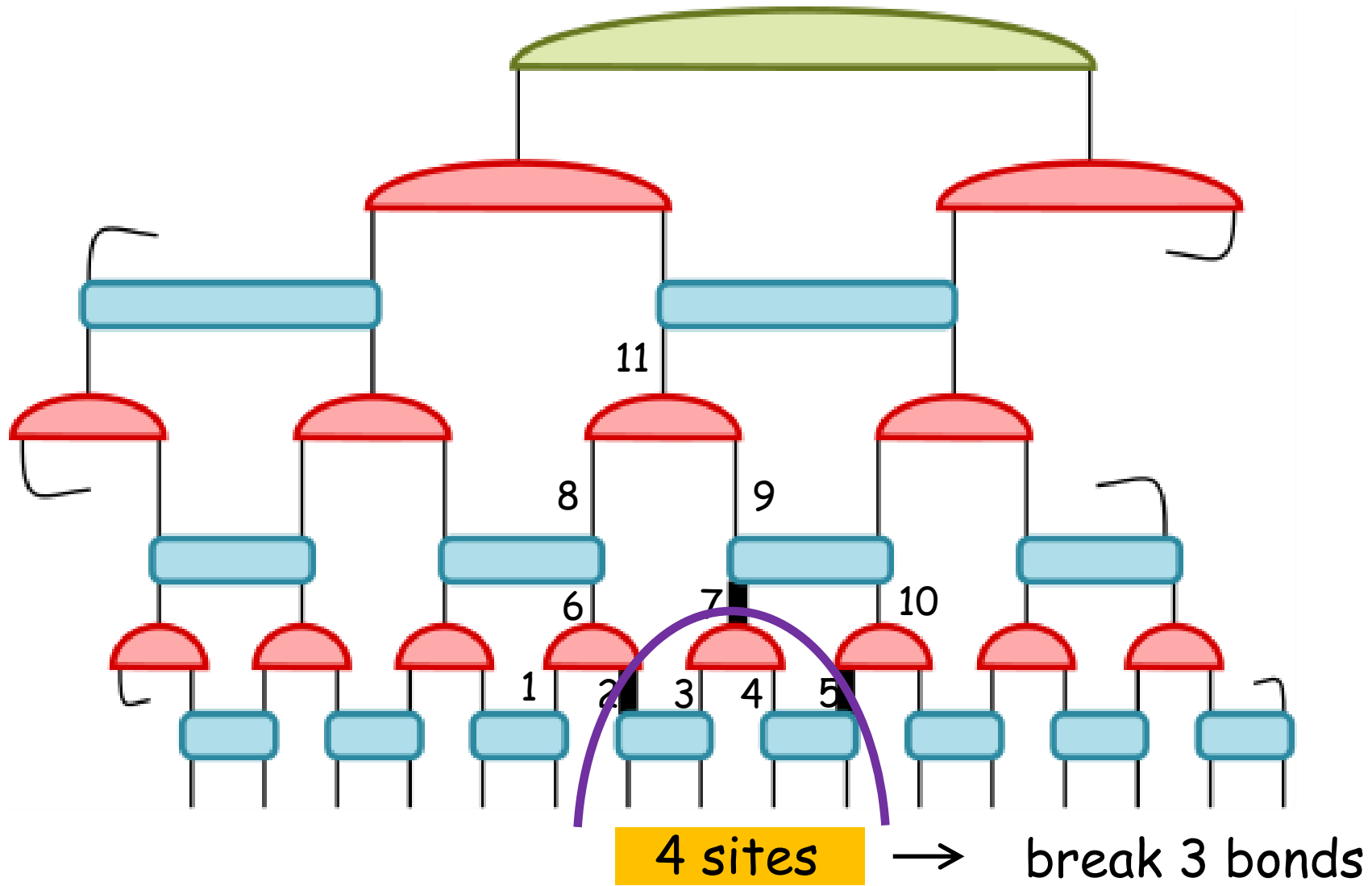
Is based on a negatively curved geometry

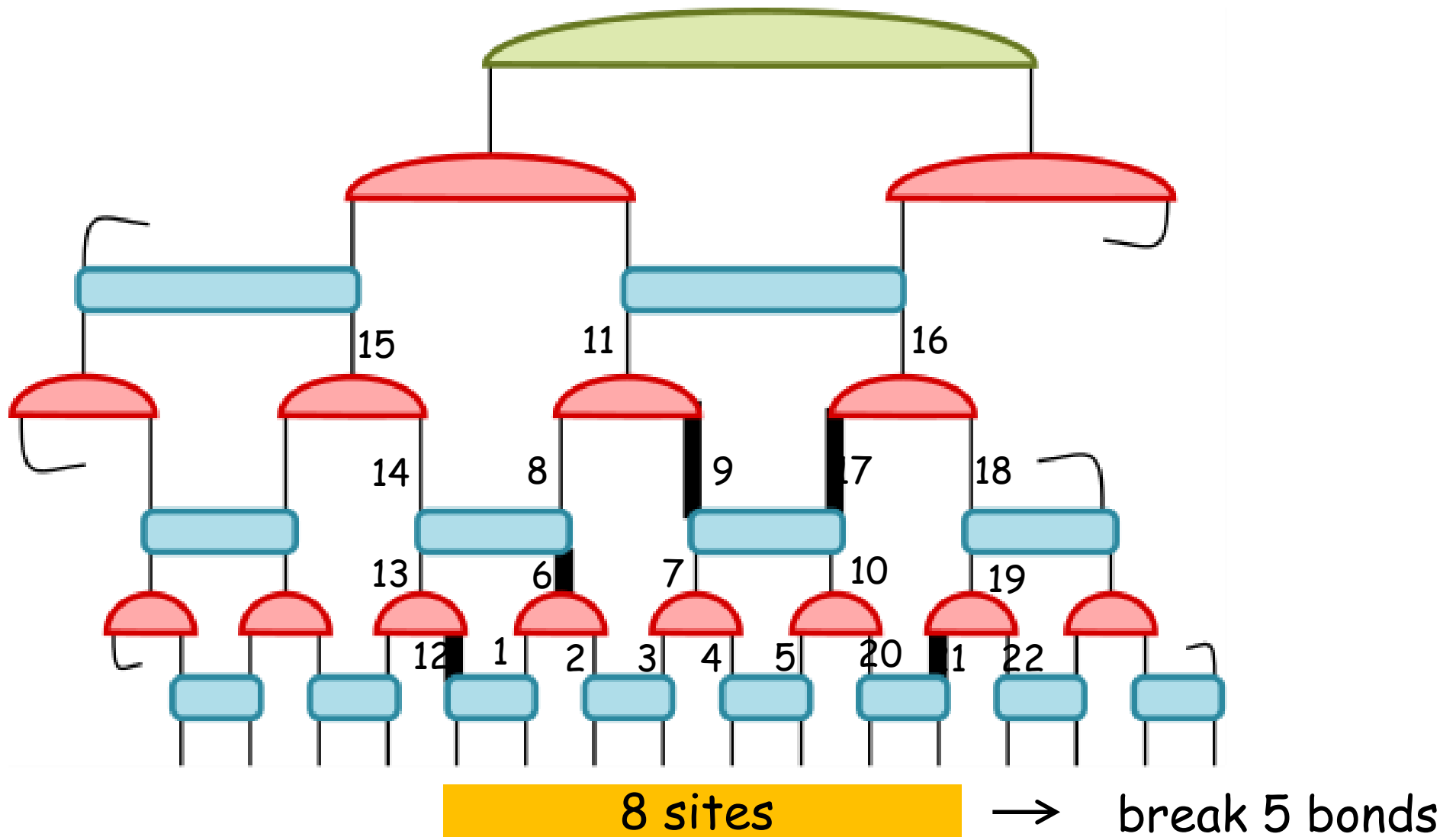
Bulk geodesics give bound on entanglement of intervals

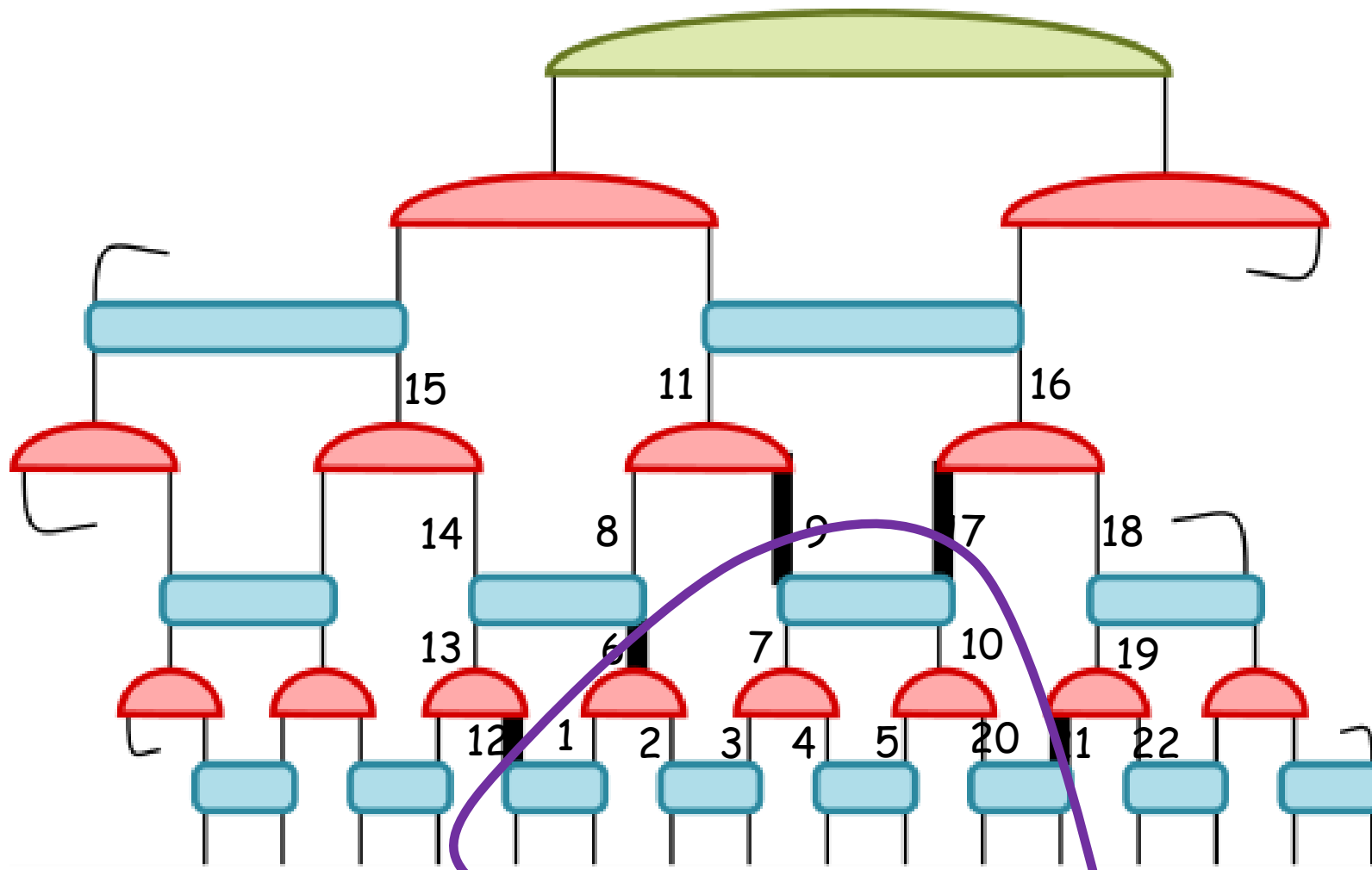


4 sites

→ break 3 bonds







8 sites

→ break 5 bonds

But these are mostly qualitative comparisons

In particular, the emergent geometry should come from the tensors (state-dependent). (Recent proposal by Vidal in this direction)

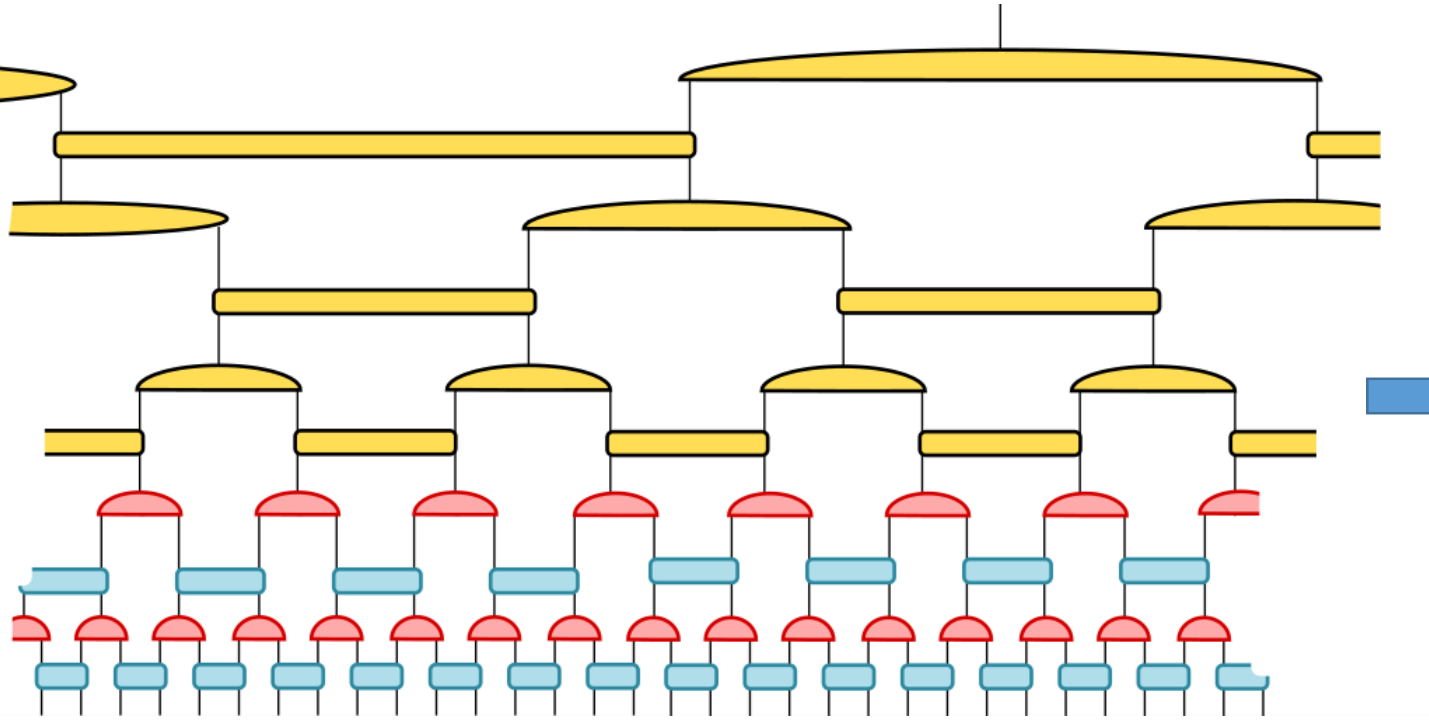
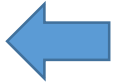
Personal philosophy: Make the CFT side of the MERA exact by e.g. adding conformal symmetry. If there is any holography, it should pop out.

Adding translation symmetry may be a good first step (in the tiling view, lack of translation invariance comes from the fact that MERA is an **aperiodic** tiling of the plane).

Summary

"Natural" ansatz for 1d critical ground states

RG flow



Lattice version
of 2d CFTs



May have something
to do with AdS/CFT

THANKS!

References

- 1) G. Vidal, PRL (2007), arXiv:cond-mat/0512165
- 2) G. Vidal, PRL (2008), arXiv:quant-ph/0610099
- 3) G. Evenbly and G. Vidal, PRB (2009), arXiv:0707.1454
- 4) R. N. C. Pfeifer, G. Evenbly, and G. Vidal, PRA (R) (2009), arXiv:0810.0580
- 5) Swingle, PRD (2012), arXiv:0905.1317