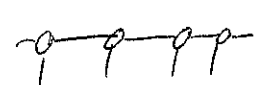


Notes of MERA Intro

Last week: see some tensor networks like MPS, PEPS.

These mimic the geometry of the system being modeled,

e.g.  for spin chain

and work well for gapped systems: $S_A \leq 2 \log X$ $\rightarrow X$ size of bonds being int. independent of size.

For gapless systems, a better alternative is the Multiscale ~~Ren~~Entanglement Renormalization Ansatz (MERA).

Basic idea is to add another dimension to the network, corresponding to RG flow. This should sound a bit like the radial direction in AdS/CFT, as we will make more precise later.

Basics of MERA

We want to implement RG in real-space, corresponding to a coarse-graining transformation. In more detail, take a 1D lattice system $\mathcal{H} = \bigotimes \mathbb{C}^k$ Hilbert space $\mathcal{H} = \bigotimes \mathbb{C}^k$ implement coarse-graining with an isometry,

$$w: \mathbb{C}^k \otimes \mathbb{C}^k \rightarrow \mathbb{C}^k \otimes \mathbb{C}^k, \quad w^\dagger w = \mathbb{1}_{\mathbb{C}^k}, \quad w w^\dagger = P$$



states mapped $w^\dagger |\psi\rangle = |\psi'\rangle \in \mathbb{C}^k$.
~~operator~~ operators, e.g. Hamiltonians ~~is~~ mapped as

$$H' = w^\dagger H w.$$

How faithful is this transformation? Look at the density matrix $\rho = \text{Tr}_{\mathbb{C}^k \otimes \mathbb{C}^k} |\psi\rangle\langle\psi| = \sum_{\alpha=1}^k p_\alpha |\varphi_\alpha\rangle\langle\varphi_\alpha|, \quad \sum p_\alpha = 1.$

Then we should choose w s.t.

$$P = w w^\dagger = \sum_{\alpha=1}^k |\varphi_\alpha\rangle\langle\varphi_\alpha|$$

In practice, we can neglect α with smallest p_α , accepting

an error s.t.

$$1 - \sum_{\alpha=1}^{\infty} p_{\alpha} \leq \epsilon \ll 1,$$


χ is a measure of entanglement between n_1, n_2 and the rest of the system:

$$S = - \sum_{\alpha=1}^{\infty} p_{\alpha} \log p_{\alpha} \leq \log \chi, \text{ saturated for } p_{\alpha} = 1/\chi.$$

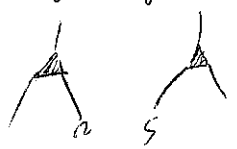
~~This is the same idea of DMRG~~

For critical systems, if we repeat this many times, we need larger and larger χ for accuracy. This reflects a failure of the simple coarse-graining; it is affected by entanglement at ~~regions~~ scales smaller than the graining we have, if the regions are at the boundary.

For example: Take ~~the~~ two sites n, s , in a state $\frac{|\uparrow_n \uparrow_s\rangle + |\downarrow_n \downarrow_s\rangle}{\sqrt{2}}$.

If coarse-graining does , don't contribute to entanglement? ~~between~~ between $n \oplus s$ and rest \rightarrow removed, i.e., s' is in a pure state.

If instead they belong to different blocks,

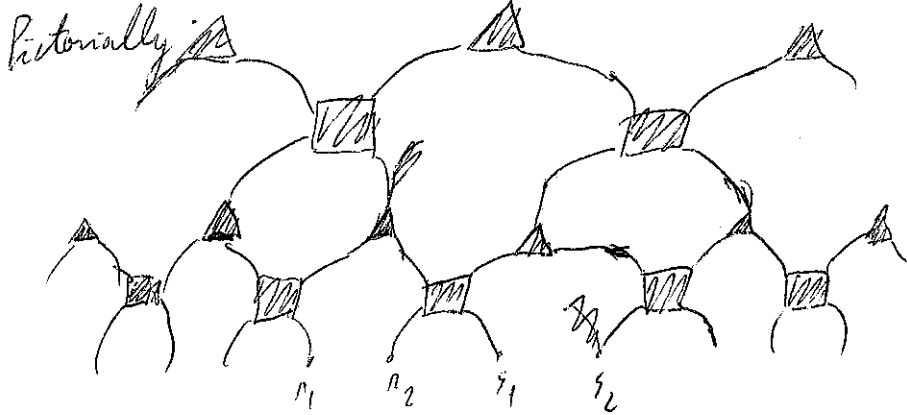


entanglement never removed!

So this retains short-range features: not a proper RG flow. (Still more, for 1D gapped systems, DMRG)

We can modify the procedure by adding disentangler, (3)
 $u: V^{\otimes 2} \rightarrow V^{\otimes 2}$, $u u^\dagger = u^\dagger u = \mathbb{1}$,

along the boundary of 2 blocks, before we coarse-grain them, such that u removes the entanglement between the boundary.



Again, map $H = W^\dagger u^\dagger H u W$

The density matrix now has a smaller rank, that in particular does not increase, even for 1D critical systems, as the procedure goes on.

Note we can map operators up ~~and down~~ the ~~chain~~ network,
 $\mathcal{O} \rightarrow \mathcal{O}' = W^\dagger u^\dagger \mathcal{O} u W$ as W goes up,
 and density

Our final tensor network MERA is the collection of all disentanglers and isometries. For a translation-inv. system, all u, w same for each layer, depending on $\mathcal{O}(X^4)$ parameters. Since there are $\mathcal{O}(\log N)$ layers, whole thing gives

$$\mathcal{O}(X^4 \log N).$$

Also true for 2D systems in the form $\mathcal{O}(X^9 \log N)$.

In practice, this is what is optimized to find $\langle \mathcal{O} \rangle$ of a given Hamiltonian.

On a critical system, since it is scale-invariant, the picture is improved: disentangles and isometries should have the same form for all layers, only two ~~states~~ ^{terms} u, v to optimize on. (4)

This method was used with $K=16$ to find scaling dimensions of operators of 1D Ising chain. Accurate up to 0.02%.

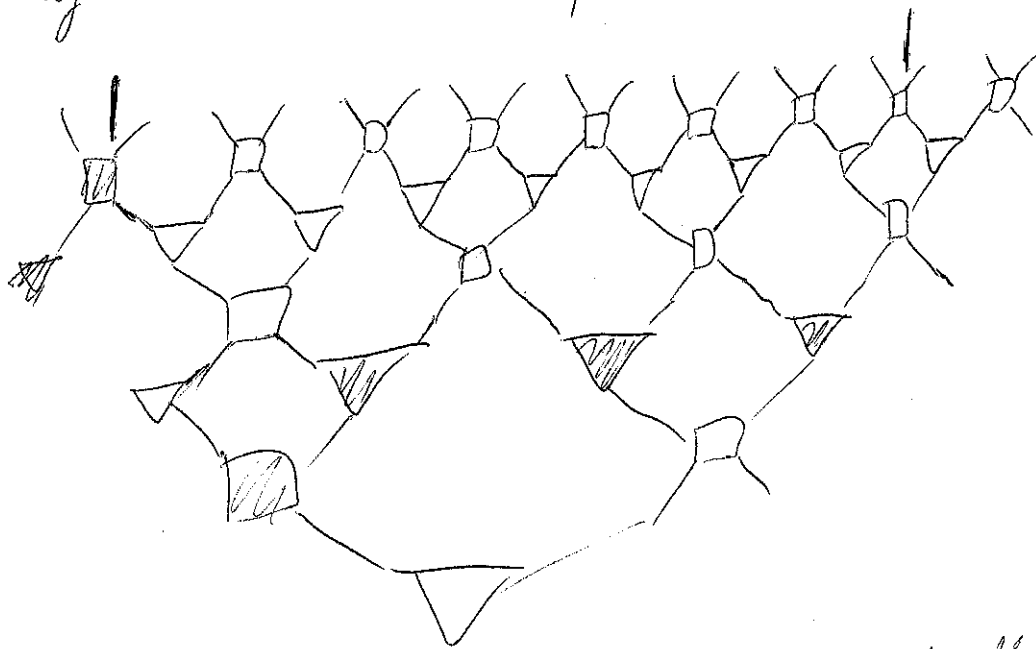
[One advantage of MERA is it makes meaning of scale invariance precise in a lattice, without taking continuous limit.]

• MERA can be seen as a quantum circuit. We will see more of this viewpoint next week.

• Support of an operator does not grow.

Behavior of ~~entanglement~~ entanglement entropy:

to compute S_A for a block of lattice sites, need the density matrix. How to compute it?



We get entropy from tracing out that occurs when we cut a bond. So we get an ^{upper} limit for the entanglement entropy, ~~given by~~ ^{interpretable as} $S_A \leq n \log X$ where n is a length/area (number of bonds cut) and X is the number of states per site.

Can we express this length as coming from a metric? (5)

~~For a cylindrical system, each layer is identical, and we cut~~
~~the cylinder~~

How many bonds do we break at each layer?

1D: 2.

Layers needed: $\log_2(L)$

$$S_A \leq 2 \log_2 L \log X$$

2D: $\left(\frac{L}{z^2}\right)^{d-1}$

Layers: $\log_d(L)$

$$S_A \leq \int_{z=0}^L L^{d-1} \left(\frac{L}{z^2}\right)^{d-1} dz \log X \approx \partial A \log X$$

Interpret this as area: for z labelling layer,

$$ds^2 = R^2 (dz^2 + e^{-2z} dx^i dx^i)$$
$$= R^2 \left(\frac{dy^2}{y^2} + \left(\frac{dx^i dx^i}{y^2} \right) \right)$$

$$y = e^z$$
$$dy = dz e^z = y dz$$

\Rightarrow spatial part of AdS metric! (state at fixed time)

for this part, count bonds you would cut if you sliced surface on that layer.

Note original system is at small z , small y ...
Boundary of AdS.

~~When to go from here~~

RT backwards: start from entanglement, build geometry from structure of entanglement in the tensor network.

Notes

- If correlation length is finite, entanglement is removed after $\textcircled{6}$ finite number of steps. Space ends.
- If we start ~~with~~ with a thermal state, at some point in the process we find a density matrix for a fully mixed state, proportional to $\mathbb{1}$. This looks like \propto temperature \Rightarrow horizon of a BH?
- Where do things like large N enter? For $S \propto \partial A$, need bonds to be maximally entangled. If not, this changes, and we need the geometric interpretation to change ...
- Use this to move away from large N , AdS?
- how do we use this to find time-component?
- Quantum circuit...

Refs

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