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#### Homework 2

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### 1 2D CFT and the 1D Transverse-Field Ising Chain

(a) We want to study the 2D CFT that underlies the quantum critical behavior at the 1D transverse field Ising model (TFIM) phase transition. Putting this 2D CFT on the plane, we have a natural parametrisation of the fields with the complex variable z=x+iy and its conjugate  $\bar{z}$ . Each primary field is caracterized by the way it transforms under a dilation  $z\to \lambda z$  with factor  $\lambda\in\mathbb{R}^+$  and rotation  $z\to e^{i\theta}z$  at angle  $\theta\in[0,2\pi)$ . For a primary field  $\varphi_a$ , these transformations respectively have the effects  $\varphi_a(z)\to\lambda^{\Delta_a}\varphi_a(\lambda z)$  and  $\varphi_a(z)=e^{-i\theta S_a}\varphi_a(e^{i\theta}z)$  where  $\Delta_a$  is the scaling dimension and  $S_a$  is the spin of the primary. Using the properties of the Virasoro algebra, we can construct conformal towers from each primary by regrouping them with their descendants.

It is possible to use quantization to extract the CFT data  $\{\Delta_\alpha, S_\alpha, c\}$  where c is the central charge. We start by applying a Weyl transformation to the complex plane to bring it to a cylinder: each concentric circle with z=0 at its center is mapped to a circle of length L at differents heights on the cylinder. We then quantize the CFT by associating an Hilbert space to each circle. The Hamiltonian  $H^{\text{CFT}}$  connecting these Hilbert spaces at different heights (times) is connecting circles of different radius in the original space and is related to the Dilation operator through an affine transformation. In this quantization, there is a correspondence between each CFT operator and states on a circle Hilbert space. The states  $|\varphi_\alpha\rangle$  corresponding to operators with well defined spins and scaling dimensions (more general than primaries and descendants, the stress energy tensor is included but is qasi-primary in 2D) are accessible through the spectrum of the dilation operator and inderictly through the spectrum  $\{E_\alpha = \frac{2\pi}{L}(\Delta_\alpha - \frac{c}{L})\}$  of  $H^{\text{CFT}}$ .

We can construct a sequence of finite size Hamiltonians (finite dimensionnal Hilbert space) that approach  $H^{\text{CFT}}$  for the 2D Ising CFT as size increases. The most natural family of Hamiltonians acheiveing this is the TFIM Hamiltonian given by

$$H = -\sum_{i=1}^{L} Z_i Z_{i+1} - \sum_{i=1}^{L} X_i$$

where  $Z_i, X_i$  are Pauli operators acting at site i of a system with  $L \in \mathbb{N}$  sites associated to a two-dimensionnal Hilbert space. We use the identification  $L \sim 0$  to have the model defined on a circle and eventually have its Hilbert space match the CFT Hilbert space in the  $L \to \infty$  limit. We note that the values of transverse field and ferromagnetic coupling are set equal to 1 so that the Hamiltonian is associated with the quantum critical point of the TFIM. From H, we can build operators that produce (in the  $L \to \infty$  limit) states corresponding to descendants when acted on states corresponding to primaries. These operators are Fourier modes of H and are expressed at

$$H_{n} = -\frac{N}{2\pi} \sum_{j=1}^{L} \exp\left(i(j+1/2)n\frac{2\pi}{N}\right) Z_{j} Z_{j+1} - \frac{N}{2\pi} \sum_{j=1}^{L} \exp\left(ijn\frac{2\pi}{N}\right) X_{j}$$

where *N* is the number of modes. how many modes are there This operator is affected by finite size effects and the connection between finite size approximations of descendants and primaries states is made stronger with the operator

$$O_n = \frac{H_n + H_{-n}}{2} \rightarrow_{L \rightarrow \infty} \frac{L_n^{\text{CFT}} + \overline{L}_{-n}^{\text{CFT}} + L_n^{\text{CFT}} + \overline{L}_n^{\text{CFT}}}{2} \quad \text{ for } n > 0$$

where  $\{L_n^{\text{CFT}}, \bar{L}_n^{\text{CFT}}\}$  of Virasoro algebra operators. For a state  $|\varphi\rangle$  corresponding to a primary, the Virasoro operators  $L_n^{\text{CFT}}, \bar{L}_n^{\text{CFT}}$  for n>0 act as  $\bar{L}_n^{\text{CFT}}|\varphi\rangle = L_n^{\text{CFT}}|\varphi\rangle = 0$ . This is the defining property of primary states. Going in the opposite direction, every descendant state can be reached from primary states by acting  $L_{-n}^{\text{CFT}}, \bar{L}_{-n}^{\text{CFT}}$ . We note that  $L_{-1}^{\text{CFT}}, L_{-2}^{\text{CFT}}$  (resp.  $\bar{L}_{-1}^{\text{CFT}}, \bar{L}_{-2}^{\text{CFT}}$ ) generate the subalgebra  $L_{-n}^{\text{CFT}}$  (resp.  $\bar{L}_{-n}^{\text{CFT}}$ ) implying that n=1,2 operators are sufficient to generate all the conformal towers. The scaling dimension  $\Delta_{\varphi'}$  of the descendant state  $|\varphi'\rangle$  of the primary state  $|\varphi\rangle$  with scaling dimension  $\Delta_{\varphi}$  is such that  $\Delta_{\varphi'} - \Delta_{\varphi} \in \mathbb{N}$ .

In what follows we describe a criterion to identify *approximate* primaries and associate them to descendants in the spectrum of *H*. We start by defining the projector

$$\Gamma_{\varphi} \equiv \sum_{\varphi_{\alpha}: E_{\alpha} < E_{\varphi}} \left| \varphi_{\alpha} \right\rangle \left\langle \varphi_{\alpha} \right|$$

which projects on the subspace of spaned by energy eigenstates below  $E_{\varphi}$  (the energy of the primary corresponding to the projector). Then we use " $\bar{L}_n^{\text{CFT}} | \varphi \rangle = L_n^{\text{CFT}} | \varphi \rangle = 0 \iff | \varphi \rangle$  is primary" to write

$$\begin{split} |\varphi\rangle \text{ is primary} &\Longrightarrow \Gamma_{\varphi}O_{n}\,|\varphi\rangle = \sum_{\varphi_{\alpha}:E_{\alpha}< E_{\varphi}} |\varphi_{\alpha}\rangle\,\langle\varphi_{\alpha}|\, \frac{L_{n}^{\text{CFT}} + \overline{L}_{-n}^{\text{CFT}} + L_{-n}^{\text{CFT}} + \overline{L}_{n}^{\text{CFT}}}{2}\,|\varphi\rangle \\ &= \sum_{\varphi_{\alpha}:E_{\alpha}< E_{\varphi}} |\varphi_{\alpha}\rangle\,\langle\varphi_{\alpha}|\, \frac{\overline{L}_{-n}^{\text{CFT}} + L_{-n}^{\text{CFT}}}{2}\,|\varphi\rangle\,, \quad |\varphi\rangle \text{ is primary} \\ &= \sum_{\varphi_{\alpha}:E_{\alpha}< E_{\varphi}} |\varphi_{\alpha}\rangle\,\langle\varphi_{\alpha}|\, \frac{|\varphi'\rangle + |\varphi''\rangle}{2}, \quad |\varphi'\rangle = L_{-n}^{\text{CFT}}\,|\varphi\rangle\,, \\ &= 0 \end{split}$$
 (descendants with  $E_{\varphi'} = E_{\varphi''} > E_{\varphi}$ )

where the energies of the descendant states  $|\varphi'\rangle$  and  $|\varphi''\rangle$  are both equal to  $E_{\varphi'} = E_{\varphi''} = \frac{2\pi}{L}(\Delta_{\alpha} + n - \frac{c}{12}) > E_{\varphi}$  (but their spins are different making them distinguishable add reference to paper).

The reverse direction of this implication is established with

$$0 = \Gamma_{\varphi} O_n \implies 0 = \sum_{\varphi_\alpha : E_\alpha < E_\omega} |\varphi_\alpha\rangle \, \langle \varphi_\alpha| \, \frac{L_n^{\rm CFT} + \overline{L}_n^{\rm CFT}}{2} \, |\varphi\rangle \implies |\varphi\rangle \, \text{ is primary}$$

where we removed the  $L_{-n}^{\rm CFT}$  and  $\overline{L}_{-n}^{\rm CFT}$  terms because they increase the energy beyond the projector scope (this is true if  $|\varphi\rangle$  is a primary or a descendant). The last implication is supported by the fact acting  $L_n^{\rm CFT}$ ,  $\overline{L}_n^{\rm CFT}$  on a descendant will lower its energy (possibly to a primary state) to the scope of the projector leading to a non-zero contribution. Only primaries escape the scope of the projector because they are mapped to zero directly by  $L_n^{\rm CFT}$ ,  $\overline{L}_n^{\rm CFT}$ .

In the finite size spectrum of H, the approximate primaries can be found by numerically computing the norm  $\varepsilon_{\varphi}^{(n)} \equiv |\Gamma_{\varphi}O_n|\varphi\rangle|$ . This norm is zero in the limit  $L \to \infty$  iff we have a primary state. For finite L, we relax the zero norm condition to a small norm condition. More precisely, we introduce a threshold  $\varepsilon_{\max}$  and declare an eigenstate of H "primary candidate" iff  $\varepsilon_{\varphi}^{(1)} + \varepsilon_{\varphi}^{(2)} \leq \varepsilon_{\max}$ .

- (b) Using the criterion described in (a) we can extract primary candidates for the TFIM CFT. We use a sparse exact diagonalization python code to find the 22 lowest energy states in the spectrum of H at size L = 16 and look for the three lowest primary states.
- (c) The TFIM Hamiltonian is related in the  $L \to \infty$  limit to the CFT Hamiltonian by the affine transformation  $H^{\text{CFT}} = aH + b$  for real constants a,b. At finite size, we can determine the values of a,b leading to an optimal agreement between the eigenvalues of the low energy states and expected scaling dimensions. We first note that the ground state  $|0\rangle$  (with  $H|0\rangle = E_0|0\rangle$ ) has scaling dimension  $\Delta_I = 0$  since it corresponds to the identity I. We choose  $b = -aE_0$  to shift the eigenvalue of  $|0\rangle$  to 0. At this point we have  $H^{\text{CFT}} \approx a(H E_0)$ . To fix a we use the fact  $L_{-2}^{\text{CFT}} |0\rangle$  (resp.  $\overline{L}_{-2}^{\text{CFT}} |0\rangle$ ) is a state  $|I,2\rangle$  (resp.  $|\overline{I},\overline{2}\rangle$ ) with scaling dimension  $\Delta_2^I = 0 + 2$  (shifted by 2 from  $\Delta_I$ ). To produce an approximation of the effect of  $L_{-2}^{\text{CFT}}$  at finite size, we use  $O_2$  and write

$$O_2\left|0\right> \approx \frac{L_2^{\mathrm{CFT}} + \overline{L}_{-2}^{\mathrm{CFT}} + L_{-2}^{\mathrm{CFT}} + \overline{L}_2^{\mathrm{CFT}}}{2}\left|0\right> = \frac{\overline{L}_{-2}^{\mathrm{CFT}} + L_{-2}^{\mathrm{CFT}}}{2}\left|0\right> = \frac{\left|I,2\right> + \left|\overline{I,2}\right>}{2}.$$

Since  $O_2|0\rangle$  is a superposition of two eigenstates with the same scaling dimension (and same energy), it is also an eigenstate of H. Acting H on it yields the constraint

$$(H - E_0)O_2 \mid 0 \rangle \approx \frac{1}{a} \times H^{\text{CFT}}O_2 \mid 0 \rangle = \frac{1}{a} \times \frac{2\Delta_2^I}{2}(O_2 \mid 0 \rangle) \implies \frac{1}{a} = \frac{\left| (H - E_0)O_2 \mid 0 \rangle \mid \Delta_2^I \mid O_2 \mid 0 \rangle}{\Delta_2^I \mid O_2 \mid 0 \rangle}$$

where whe chose a > 0 to ensure positive scaling dimensions (the spectrum is bounded from below implying the shifted spectrum is positive). This way to calculate a can be implemented numerically.

- (d)
- (e)
- (f)
- (g)

### 2 Acknowledgement

## References

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