Recent development in bootstrap numerics

Ning SU (University of Pisa)

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Conformal bootstrap and Critical Phenomena

Problems in critical phenomena

↔ Questions on CFT data

 \leftrightarrow Conformal bootstrap constraints

↔ semidefinite program (SDP)

 \leftrightarrow high efficiency SDP numerics (bottleneck!)

In practice : the efficiency of the bootstrap numerics crucially limits our ability to access certain CFT data.

This talk : Latest examples of large scale bootstrap problem. New algorithm for SDP numerics (10+ times faster).

Sophisticated CFTs \leftrightarrow Large scale bootstrap constraints

Bootstrapping $\langle \phi \phi \phi \phi \rangle$ in 2D. Plot lowest singlet Δ_0 v.s. Δ_{ϕ} .



Why 2D Ising CFT sits at the kink of the bounds? How to make other $\mathcal{M}(m, m+1)$ sit on kinks?

Sophisticated CFTs ↔ Large scale bootstrap constraints

Why 2D Ising CFT sits on the kink of the bounds?

[Connor Behan 2017]

There is a series of solutions that saturate $\langle \phi \phi \phi \phi \rangle$: $\Delta_{\epsilon} = \frac{8}{3} \Delta_{\phi} + \frac{2}{3}$.

They smoothly connect $m = 3, 4, \dots \infty$ for $\mathcal{M}(m, m+1)$.

As $m \to 3$, an operator in spin 2 channel T' decouple from spectrum : $\lambda_{\phi\phi T'} \to 0$.

T' : level 2 null state in Virasoro primary

Sharp features in the allowed region (kinks, pikes) ↔ operators decoupling/disappear/recombine in the primary spectrum

[Belavin, Polykov, Zamolodchikov 1984] For c<1 to be unitary, there must be certain null states in the Virasoro primary

Sophisticated CFTs \leftrightarrow Large scale bootstrap constraints

In general, to access more and more sophisticated CFTs, we have to mix more and more operators.

Our setup must be able to access enough gap features in the spectrum that discriminate the target CFT from all other theories (solutions) with the same symmetry.

more correlators, more parameters to scan \rightarrow large scale SDP and challenging numerics

2D Potts and tricritical Potts model

Consider 2D square lattice of random spins with Hamiltonian for $s_i \in \{1, 2, ..., q\}$ [Potts 1952]

$$Z = \sum_{\{s_i\}} e^{-\beta \sum_{\langle ij \rangle} \delta_{s_i, s_j}}$$

Symmetry : S_q

Tuning β : at low temperature (small β), ordered phase with broken S_q . At high temperature, disordered phase. At $\beta = \beta_{crit}$: Second order phase transition, described by critical Potts $\subset \mathcal{M}(5, 6)$

Assuming some sites are vacant, turning both β and chemical potential of vacancies : Second order phase transition, described by tricritical Potts $\subset \mathcal{M}(6, 7)$

$2+\epsilon$ dimensional q=3 Potts and tricritical Potts model

Symmetry : S_3 has 3 irreps : singlet 1, sign rep 1 ', standard rep 2

Critical Potts : two relevant rep 2 operators σ , σ' , one relevant singlet ϵ . 2D values: $\Delta_{\sigma} = \frac{2}{15}$, $\Delta_{\sigma'} = \frac{4}{3}$, $\Delta_{\epsilon} = \frac{4}{5}$

Tricritical Potts : two relevant rep **2** operators σ , σ' , two relevant singlet ϵ , ϵ' . 2D values: $\Delta_{\sigma} = \frac{2}{21}$, $\Delta_{\sigma'} = \frac{20}{21}$, $\Delta_{\epsilon} = \frac{2}{7}$, $\Delta_{\epsilon'} = \frac{10}{7}$

In $2 + \epsilon$ dimension, q = 3 critical Potts and tricritical Potts get closer, merge and become complex CFT at d_{crit} (the merger/annihilation scenario [Gorbenko, Rychkov, Zan 2018]) Note : for q = 2 (Ising), $d_{\text{crit}} = 4$; For q = 4, $d_{\text{crit}} = 2$.

Two predictions around $d_{
m crit}$:

(1), Potts & tricritical Potts merges with $\Delta_{\epsilon} = \Delta_{\epsilon'} = 3$ at d_{crit}

(2), $\Delta_O - \Delta_{O, \text{crit}} \propto \sqrt{d_{\text{crit}} - d}$

Consider correlator $\langle\sigma\sigma\sigma\sigma\rangle$ in 2D, bound $\Delta_{\sigma'} \ v.s. \ \Delta_{\sigma}$:



[Junchen Rong, NS, 2017]

[Shai Chester, NS, to appear]

Consider correlator $\langle \sigma\sigma\sigma\sigma\rangle$ in 2D, assuming only σ, σ' are relevant operators



Consider all correlator of $\sigma, \sigma', \epsilon$:

 $\begin{array}{l} \langle \sigma\sigma\sigma\sigma\sigma\rangle, \langle \sigma'\sigma'\sigma'\sigma'\rangle, \langle \sigma\sigma'\sigma\sigma'\rangle, \langle \sigma\sigma\sigma'\sigma'\rangle, \langle \sigma\sigma\sigma\sigma'\rangle, \langle \sigma\sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\sigma\rangle, \langle \epsilon\epsilon\epsilon\epsilon\rangle, \langle \epsilon\sigma\epsilon\sigma\rangle, \langle \epsilon\sigma\epsilon\sigma'\rangle, \langle \epsilon\epsilon\sigma\epsilon\sigma'\rangle, \langle \epsilon\epsilon\sigma\sigma\rangle, \langle \epsilon\epsilon\sigma\sigma\rangle, \langle \epsilon\sigma\sigma\sigma\rangle, \langle \epsilon\sigma\sigma\sigma\rangle, \langle \epsilon\sigma\sigma\sigma\rangle, \langle \epsilon\sigma\sigma\sigma'\rangle, \langle \epsilon\sigma\sigma\sigma\rangle, \langle \sigma\epsilon\sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\rangle, \langle \sigma\sigma\sigma\rangle, \langle \sigma$

Totally 39 crossing equations.

Consider all correlators of σ , σ' , ϵ . Assume they are the only relevant scalars in 2 and 0. Bootstrap at 2D with truncation order : $\Lambda = 11$:



 $2 + \epsilon$ bootstrap :

For Potts : all correlators of σ , σ' , ϵ . Scan

 $\Delta_{\sigma}, \Delta_{\sigma'}, \Delta_{\epsilon}, \lambda_{\epsilon\epsilon\epsilon}, \lambda_{\sigma\sigma\epsilon}, \lambda_{\sigma\sigma'\epsilon}, \lambda_{\sigma\sigma\sigma}, \lambda_{\epsilon\sigma'\sigma'}, \lambda_{\sigma\sigma'\sigma'}, \lambda_{\sigma'\sigma'\sigma'}$ (10D)

For tri-Potts : all correlators of σ , σ' . Scan

 $\Delta_{\sigma}, \Delta_{\sigma'}, \Delta_{\epsilon}, \Delta_{\epsilon'}, \lambda_{\sigma\sigma\sigma}, \lambda_{\sigma\sigma'\sigma'}, \lambda_{\sigma'\sigma'\sigma'}, \lambda_{\sigma\sigma\epsilon}, \lambda_{\sigma\sigma'\epsilon}, \lambda_{\epsilon\sigma'\sigma'}, \lambda_{\sigma\sigma\epsilon'}, \lambda_{\sigma\sigma'\epsilon'}, \lambda_{\sigma'\sigma'\epsilon'}$ (13D)

Gigantic SDP and scan problem. We need the navigator method [Reehorst, Rychkov, Simmons-Duffin, Sirois, SN, van Rees 2021]. See Reehorst Marteen's talk.

In both setups, we minimize Δ_{σ} in the allowed region, i.e. go to the tip of the dagger.



Conclusion : $d_{\rm crit} \lesssim 2.5$. Approximately correct square root behavior $\Delta_{\epsilon'} - d_{\rm crit} \propto \sqrt{d_{\rm crit} - d}$, but not conclusive. Two possibilities in higher Λ :



We need much higher Λ bootstrap study (not possible with current techniques) (Higher Λ might even turn the dagger plot into an island)

O(N) vector model bootstrap

Lagrangian : $\mathcal{L} = \partial \phi_i \partial \phi_i + m^2 (\phi_i \phi_i) + \lambda (\phi_i \phi_i)^2$ Bootstrapping all 4pt involves $\{v = \phi_i, s = \phi^2, t = \phi_i \phi_j\}$:



[Chester, Landry, Liu, Poland, Simmons-Duffin, <mark>SN</mark>, Vichi 2019, 2020] Gigantic bootstrap problem. We used XSEDE super-cluster.

3D O(3) model

O(3) model : $\mathcal{L} = \partial \phi_i \partial \phi_i + \lambda_1 (\phi_i \phi_i)^2$ Cubic model : $\mathcal{L} = \partial \phi_i \partial \phi_i + \lambda_1 (\phi_i \phi_i)^2 + \lambda_2 \sum_i \phi_i^4$ (symmetry: $S_3 \times Z_2^3$)



A final verdict : $\Delta_4 \leq 2.99056$. $\mathit{O}(3)$ is unstable against cubic perturbation!

 $({\rm LaAlO}_3$ has structural phase transition in Cubic universality, not O(3) universality)

3D cubic model : conformal perturbation theory

(Junchen Rong, NS, to appear. See Junchen's talk for more details)

 $S_{ ext{cubic}} = S_{O(3)} + g \int d^3 x \ t_4 \ \text{where} \ \Delta_{t_4} \approx 2.988.$ Perturbation parameter : $\delta = 3 - 2.988 = 0.012$

$$\begin{split} \beta_g &= -\delta \ g - \frac{1}{2} \ S_{d-1} \ \lambda_{t_4 \ t_4} \ g^2 \\ \Delta_O &= \Delta_0 + \delta \ \frac{2 \lambda_{OO \ t_4}}{\lambda_{t_4 \ t_4}} \end{split}$$

 $\lambda_{OO \ t_4} / \lambda_{t_4 \ t_4}$ can be accessed by bootstrap correlators of $\{v, \ s, \ t, \ t_4\}$

 $\lambda_{OO t_4} / \lambda_{t_4 t_4 t_4} \approx 0.8$ from one allowed point at $\Lambda = 11$.

The O(3) rank 2 tensor (dim=5) ($\Delta_t = 1.2096$) split to cubic 2 \oplus 3 $\Delta_2 \approx 1.2200$, $\gamma_2 \approx 0.01043$ $\Delta_3 \approx 1.2091$, $\gamma_3 \approx -0.00044$

We couldn't know error bar of $\lambda_{OOX} / \lambda_{XXX} \approx 0.8$. Need larger Λ bootstrap study (not possible with current techniques).

large charge operators in O(3)



Non-abelian currents bootstrap

[Yin-Chen He, Junchen Rong, NS, Alessandro Vichi, ongoing work]

Consider $\langle J_{\mu} \; J_{\nu} \; J_{\rho} \; J_{\sigma}
angle$ where $(J_{\mu})^b_a$ is the current of global symmetry $\mathrm{SU}(N_f)$

 $S \overline{S} : A^{(a \ b)}_{(c \ d)}$, an example of decoupling operator that can detect color group (See Yinchen's discussion talk last week). $\Delta_{S \ \overline{S}, L=2} \approx 4$ for $U(N)_{N \ge 2}$ gauge theory, $\Delta_{S \ \overline{S}, L=2} \approx 6$ for QED.

Non-abelian currents bootstrap





Non-abelian currents bootstrap



To shrink the QED3 island and get precise CFT data, we might need to mix J^f with J^t and/or a monopole operator and do calculation at larger Λ . (not possible with current techniques)

A common challenge

To get sharp results for some problems in critical phenomena, we often have to bootstrap a large scale problem that beyond current limits.

Current bootstrap numerics

The old numerical method: a section of SDPs over a manifold. We solve SDPs one by one.



Typical Newtonian iterations per SDP : 100 to 300

(A well-known issue in SDP : for optimality runs, hotstart is not efficient. Best solution before : save a middle checkpoint.)

[Aike Liu, David Simmons-Duffin, NS, Balt van Rees, ongoing work]

The new numerical method: treat the section of SDPs over a manifold as a single optimization problem. Hopping to a new SDP without completely solving current SDP.



Simultaneously solves the optimization in the parameter manifold (maximize Δ_{σ}) and solves the optimization of SDP ($\mu \rightarrow 0$)

A novel algorithm:

- 1, Determine a good hopping point
- 2, Determine a good hopping direction in both SDP internal variables (x, y, X, Y) and external parameters ($\Delta_{\sigma}, ...$)

Key feature : only a few Newtonian iterations per SDP.

Example : O(3) bootstrapping correlators of v, s, t at $\Lambda = 19$. Maximize Δ_{t_4} in the allowed region of $\{\Delta_v, \Delta_s, \Delta_t, \Delta_{t_4}, \lambda_{vtv}, \lambda_{tts}, \lambda_{ttt}, \lambda_{sss}\}$



Example : O(3) bootstrapping correlators of v, s, t at $\Lambda = 19$. Maximize Δ_{t_4} in the allowed region of $\{\Delta_v, \Delta_s, \Delta_t, \Delta_{t_4}, \lambda_{vtv}, \lambda_{tts}, \lambda_{ttt}, \lambda_{sss}\}$ Average Newtonian iterations per SDP (excluding 1st SDP): New algorithm : ~4 Old algorithm : ~ 52

As the scale of the SDP get larger, the different is even bigger.

We invented a bag of tricks. Some of our tricks works for SDP itself. For example, in some cases, in one Newtonian iteration, we could make μ go from 10^{-15} to 10^{-33} .

(I will give a technical talk regarding the algorithm in the last week)

Future goals

Problems in critical phenomena

 $\begin{array}{c} \leftrightarrow \text{ Questions on CFT data} \\ \leftrightarrow \text{ Bootstrap constraints} \\ \leftrightarrow \text{ semidefinite program (SDP) (bottleneck)} \\ \leftrightarrow \text{ high efficiency SDP numerics} \end{array}$ Right now : Parameters $\xleftarrow{\text{iterate Casimir equ.}}$ numerical conformal blocks $\xleftarrow{\text{some algebra}}$ SDP
If we have a iterative process : Parameters $\xleftarrow{\text{iterate}}$ SDP and $d(\text{Parameters}) \leftrightarrow d(\text{SDP})$ We may make the two iterative processes play against each other.

Future goals

Given a set of operators { ϕ_1 , ϕ_2 , ϕ_3 , ...}, we assign different derivative truncation Λ to $\langle \phi_1 \phi_1 \phi_1 \phi_1 \rangle$, $\langle \phi_2 \phi_2 \phi_2 \phi_2 \rangle$, $\langle \phi_1 \phi_1 \phi_1 \phi_2 \phi_2 \rangle$. (Already implemented in simpleboot package)

Adding more derivatives to $\langle \phi_1 \phi_1 \phi_1 \phi_1 \rangle$ or adding a new operator ϕ_n ? Should be decided by their contribution to the navigator function.

Technical issues : non-uniform polynomial degrees. Create some degeneracies in SDP.

Thank you