

Exact three and four-point correlation functions in the $O(n)$ loop model

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Tale of two loop models

Q-state Potts model

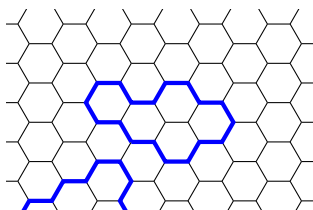
- Q-state spins; interactions have S_Q permutation symmetry.
- Equivalent loop model on medial lattice [Baxter-Kelland-Wu 1976].
- Respects fixed orientation of lattice edges: $U(n)$ symmetry.
- Related to integrable 6-vertex model and Temperley-Lieb algebra.
- S_Q commutes with partition algebra $\mathcal{P}_L(Q)$, descending to Potts–Temperley–Lieb algebra $P\mathcal{T}\mathcal{L}_{2L}(\sqrt{Q})$ in $d = 2$.

$O(n)$ model

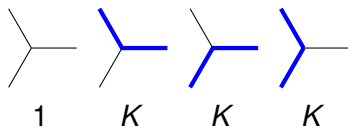
- Vector spins $\in \mathbb{R}^n$; interactions have $O(n)$ symmetry.
- Equivalent loop model in $d = 2$ after modification [Nienhuis 1982].
- Related to integrable 19-vertex model and Motzkin algebra.
- $O(n)$ commutes with Brauer algebra $\mathcal{B}_L(n)$, descending to unoriented Jones–Temperley–Lieb algebra $u\mathcal{J}\mathcal{T}\mathcal{L}_L(n)$ in $d = 2$.

$O(n)$ model [Nienhuis 1982]

loop weight n



Vertex weights 1 and K



All configurations can be built by a transfer matrix:

$$\check{R}_k = \begin{array}{c} \diagup \diagdown \\ \diagdown \diagup \end{array} + K \begin{array}{c} \diagup \text{---} \\ \diagdown \diagup \end{array} + K \begin{array}{c} \diagdown \diagup \\ \text{---} \diagdown \end{array} + K^2 \begin{array}{c} \diagup \text{---} \\ \text{---} \diagup \end{array} + K^2 \begin{array}{c} \text{---} \diagup \\ \diagdown \text{---} \end{array} + K^2 \begin{array}{c} \text{---} \diagdown \\ \diagup \text{---} \end{array} + K^2 \begin{array}{c} \diagup \text{---} \\ \text{---} \diagdown \end{array} + K^2 \begin{array}{c} \text{---} \diagup \\ \diagdown \text{---} \end{array}$$

Define the partition function

$$Z(K, n) = \sum_{\text{loops}} K^{\#\text{monomers}} n^{\#\text{loops}} .$$

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Monomer fugacity at the critical point:

$$K_c = \left(2 \pm \sqrt{2-n}\right)^{-1/2} ,$$

where $-2 \leq n \leq 2$. Plus (minus) sign for the dilute (dense) phase.

Special cases:

- $n = 1$ dense: Site percolation
- $n = 1$ dilute: Ising model
- $n = 0$ dilute: Self-avoiding walks
- $n = 2$ either: Gaussian free field, XY model

Most of there are really *logarithmic* CFTs.

Our first objective is to understand the case of 'generic' n .

As a warm-up, let us start by a simpler question



Conformal Field Theory of the $O(n)$ model

Central charge

$$c = 13 - 6\beta^2 - 6\beta^{-2} \quad \text{with} \quad \begin{cases} \Re\beta^2 > 0, \\ \beta^2 \notin \mathbb{Q}. \end{cases}$$

Conformal weight Δ and momentum P :

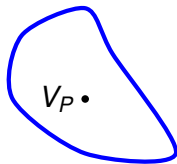
$$\Delta = P^2 - P_{(1,1)}^2, \quad \Delta_{(r,s)} = P_{(r,s)}^2 - P_{(1,1)}^2, \quad P_{(r,s)} = \frac{1}{2} \left(-\beta r + \beta^{-1} s \right).$$

Field content, with left- and right-moving conformal weights $(\Delta, \bar{\Delta})$:

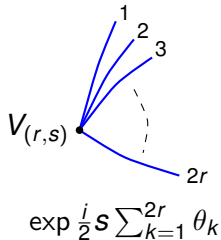
Name	Notation	Parameters	$(\Delta, \bar{\Delta})$
Degenerate	$V_{\langle r,s \rangle}^d$	$r = 1; s \in 2\mathbb{N} + 1$	$(\Delta_{(r,s)}, \Delta_{(r,s)})$
Diagonal	V_P	$P \in \mathbb{C}$	$(P^2 - P_{(1,1)}^2, P^2 - P_{(1,1)}^2)$
Non-diagonal	$V_{(r,s)}$	$r \in \frac{1}{2}\mathbb{N}^*; s \in \frac{1}{r}\mathbb{Z}$	$(\Delta_{(r,s)}, \Delta_{(-r,s)})$

Interpretation of fields within the loop model:

Diagonal and non-diagonal fields



$w(P)$

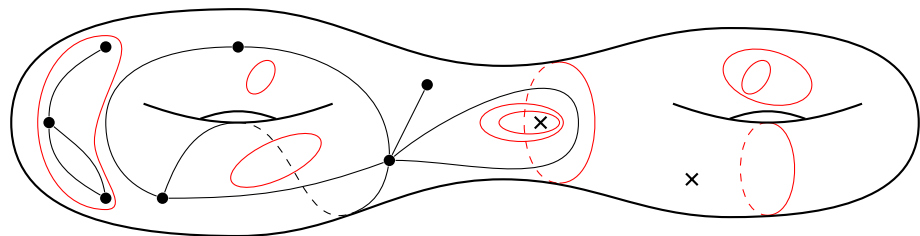


$V_{\langle 1,3 \rangle}^d$ is the energy operator.

The dense $O(n)$ model has a CFT limit iff $V_{\langle 1,3 \rangle}^d$ is irrelevant:

$$\Re \Delta_{(1,3)} > 1 \iff \Re \beta^{-2} > 1 .$$

Dream about correlation functions



Here \bullet are $V_{(r,s)}$ insertions, and \times are V_P insertions.

Open curves define a *combinatorial map* on a Riemann surface.

Segal's axioms: Three basic building blocks

1) Annulus with one insertion, 2) Disk with two insertions, 3) Pants.

The blocks are glued by integrating over eigenstates.

Progress this far

- Fields related to irreps of affine Temperley-Lieb algebra, $ATL_L(n)$.
- Bijection between correlation functions and combinatorial maps.
- Conformal symmetry enhanced to *interchiral symmetry* via $V_{\langle 1,3 \rangle}^d$.
- Global $O(n)$ symmetry in interplay with conformal symmetry.

First goal is to understand $N \leq 4$ points on the sphere.

- $N = 2$ understood from critical exponents.
- $N = 3$ conjecturally understood in all cases.
- $N = 4$ from conformal bootstrap. Partial analytical control.

The talk summarises this progress.

From this we can construct the TL generator:

$$e_i = s_{i+\frac{1}{2}} s_i s_{i+1} s_{i+\frac{1}{2}} = \begin{array}{c} \dots \quad \dots \quad i \quad i+1 \quad \dots \quad \dots \\ | \quad | \quad \cup \quad \cap \quad | \quad | \\ \dots \quad \dots \quad \cap \quad \cup \quad \dots \quad \dots \end{array}$$

To get the periodic algebra $\mathcal{ATL}_L(n)$ we add:

$$e_L = \begin{array}{c} \cup \quad \cup \quad \dots \quad \cup \quad \cup \\ | \quad | \quad \dots \quad | \quad | \\ \cap \quad \cap \quad \dots \quad \cap \quad \cap \end{array}, \quad u = \begin{array}{c} / \quad / \quad / \quad \dots \quad / \quad / \\ \backslash \quad \backslash \quad \backslash \quad \dots \quad \backslash \quad \backslash \end{array}$$

Define also the pseudo-translation t of the $2r \in \mathbb{N}^*$ through-lines:

$$\begin{array}{c} \cup \quad \cup \quad | \quad \cup \quad \cup \\ \cap \quad \cap \quad \cup \quad \cap \quad \cap \end{array} \xrightarrow{t} \begin{array}{c} \cup \quad \cup \quad \cup \quad \cup \quad \cup \quad \cup \\ \cap \quad \cap \quad \cap \quad \cap \quad \cap \quad \cap \end{array}$$

$\mathcal{AFL}_L(n)$ is ∞ -dimensional. A finite-dimensional quotient, the unoriented Jones–Temperley–Lieb algebra $u\mathcal{JTL}_L(n)$, is obtained by replacing non-contractible loops by n and imposing

$$t^{2r} \underset{u\mathcal{JTL}_L(n)}{=} 1 .$$

The standard modules $W_{(r,s)}^{(L)}$ are irreps of $u\mathcal{JTL}_L(n)$, spanned by link patterns with $2r$ defects. E.g. for $W_{(1,s)}^{(10)}$:



We have

$$(t - e^{\pi i s}) W_{(r,s)}^{(L)} = 0 .$$

The labels (r, s) carry over to the CFT.

Conformal partition function on the torus

Obtained by Di Francesco-Saleur-Zuber in 1987.

Let $q = e^{2\pi i\tau}$ with τ the modulus, and $\eta(q)$ is the Dedekind function.

$$Z^{O(n)}(q) = \sum_{s \in 2\mathbb{N}+1} \chi_{\langle 1,s \rangle}(q) + \sum_{r \in \frac{1}{2}\mathbb{N}^*} \sum_{s \in \frac{1}{r}\mathbb{Z}} L_{(r,s)}(n) \chi_{(r,s)}^N(q)$$

with the diagonal degenerate characters

$$\chi_{\langle r,s \rangle}(q) = \left| \frac{q^{P_{(r,s)}^2} - q^{P_{(r,-s)}^2}}{\eta(q)} \right|^2,$$

and the non-diagonal characters

$$\chi_{(r,s)}^N(q) = \frac{q^{P_{(r,s)}^2} \bar{q}^{P_{(r,-s)}^2}}{|\eta(q)|^2}.$$

$$Z^{O(n)}(q) = \sum_{s \in 2\mathbb{N}+1} \chi_{\langle 1, s \rangle}(q) + \sum_{r \in \frac{1}{2}\mathbb{N}^*} \sum_{s \in \frac{1}{r}\mathbb{Z}} L_{(r,s)}(n) \chi_{(r,s)}^N(q)$$

We have the Virasoro representations:

$\mathcal{R}_{\langle 1, s \rangle} =$ diagonal level- s degenerate rep. with character $\chi_{\langle 1, s \rangle}(q)$,

$\mathcal{W}_{(r,s)} =_{r,s \in \mathbb{N}^*}$ indecomposable rep. with character $\chi_{(r,s)}^N(q) + \chi_{(r,-s)}^N(q)$,

$\mathcal{W}_{(r,s)} =_{r \notin \mathbb{Z}^* \text{ or } s \notin \mathbb{Z}^*}$ Verma module with character $\chi_{(r,s)}^N(q)$.

The multiplicities $L_{(r,s)}(n)$ were obtained by Read-Saleur in 2001:

$$L_{(r,s)}(n) = \delta_{r,1} \delta_{s \in 2\mathbb{Z}+1} + \frac{1}{2r} \sum_{r'=0}^{2r-1} e^{\pi i r' s} x_{(2r) \wedge r'}(n),$$

with polynomials $x_d(n)$ defined by

$$x_0(n) = 2 \quad , \quad x_1(n) = n \quad , \quad n x_d(n) = x_{d-1}(n) + x_{d+1}(n) .$$

Global $O(n)$ symmetry

This looks like Schur-Weyl duality.

Indeed we have both CFT and $O(n)$ symmetry.

$O(n)$ can be defined for $n \in \mathbb{C}$ (Deligne category).

Under the global $O(n)$ symmetry, primary operators transform in irreps:

$$[] : \bullet, \quad [2] : \square\square, \quad [11] : \begin{array}{|c|} \hline \square \\ \hline \square \\ \hline \end{array}, \quad [5421] : \begin{array}{|c|c|c|c|c|} \hline \square & \square & \square & \square & \square \\ \hline \square & \square & \square & \square & \\ \hline \square & \square & & & \\ \hline \square & & & & \\ \hline \end{array}.$$

Known dimensions and tensor products (Newell-Littlewood numbers).

Loop-model interpretation:

Each loop carries $[1]$, the fundamental (defining) representation.

Empty space corresponds to $[\]$, the trivial representation.

$$Z^{O(n)}(q) = \sum_{s \in 2\mathbb{N}+1} \chi_{\langle 1, s \rangle}(q) + \sum_{r \in \frac{1}{2}\mathbb{N}^*} \sum_{s \in \frac{1}{r}\mathbb{Z}} L_{(r,s)}(n) \chi_{(r,s)}^N(q)$$

The proper way to understand it is that the $O(n)$ CFT has a space of states (spectrum)

$$\mathcal{S}^{O(n)} = \bigoplus_{s \in 2\mathbb{N}+1} [] \otimes \mathcal{R}_{\langle 1, s \rangle} \oplus \bigoplus_{r \in \frac{1}{2}\mathbb{N}^*} \bigoplus_{s \in \frac{1}{r}\mathbb{Z}} \Lambda_{(r,s)} \otimes \mathcal{W}_{(r,s)}$$

acted upon by $O(n) \times \mathfrak{C}$, where \mathfrak{C} is conformal symmetry.

So $\dim_{O(n)} \Lambda_{(r,s)} = L_{(r,s)}(n)$. And of course $\dim_{O(n)} [] = 1$.

Introduce the formal alternating hook representations

$$\Lambda_t = \delta_{t \equiv 0 \pmod{2}}[\] + \sum_{k=0}^{t-1} (-1)^k [t-k, 1^k].$$

We find then

$$\Lambda_{(r,s)} = \delta_{r,1} \delta_{s \in 2\mathbb{Z}+1}[\] + \frac{1}{2r} \sum_{r'=0}^{2r-1} e^{\pi i r' s} x_{(2r) \wedge r'} \left(\Lambda_{\frac{2r}{(2r) \wedge r'}} \right).$$

There exists an equivalent formula which makes clear that the expansion coefficients of Young tableaux $\in \mathbb{N}$.

Let us have a closer look:

$$\Lambda_{(\frac{1}{2},0)} = [1] ,$$

$$\Lambda_{(1,0)} = [2] ,$$

$$\Lambda_{(1,1)} = [11] ,$$

$$\Lambda_{(\frac{3}{2},0)} = [3] + [111] ,$$

$$\Lambda_{(\frac{3}{2},\frac{2}{3})} = [21] ,$$

$$\Lambda_{(2,0)} = [4] + [22] + [211] + [2] + [] ,$$

$$\Lambda_{(2,\frac{1}{2})} = [31] + [211] + [11] ,$$

$$\Lambda_{(2,1)} = [31] + [22] + [1111] + [2] .$$

We also have e.g. $[1] \otimes [1] = [2] + [11] + []$.

This tells us how to decompose two loop lines on $O(n)$ irreps.

Consequences for correlation functions

Two-point functions are given by the conformal dimensions, up to normalisation of the field.

Three-point functions are also fixed by global conformal invariance, up to structure constants.

Four-point functions could be determined by differential equations, if both $V_{\langle 1,s \rangle}^d$ and $V_{\langle r,1 \rangle}^d$ were present, but we only have the former!

Therefore we need the *conformal bootstrap*.

But we can do better than usual for two reasons:

- $V_{\langle 1,3 \rangle}^d$ generates an *interchiral symmetry*.
- We can exploit the global $O(n)$ symmetry.

Consider a four-point function of non-diagonal primary fields, and its s-channel decomposition into conformal blocks:

$$\left\langle \prod_{i=1}^4 V_{(r_i, s_i)} \right\rangle = \sum_{s \in 2\mathbb{N}+1} D_s \mathcal{G}_{\langle 1, s \rangle}^D + \sum_{r \in \frac{1}{2}\mathbb{N}^*} \sum_{s \in \frac{1}{r}\mathbb{Z}} D_{(r, s)} \mathcal{G}_{(r, s)} .$$

The blocks are known from Zamolodchikov's recursion relation.

Degenerate shift equations using $V_{\langle 1, 3 \rangle}^d$ determine $\frac{D_{(r, s+1)}}{D_{(r, s-1)}}$ and $\frac{D_{s+1}}{D_{s-1}}$.

So rewrite

$$\left\langle \prod_{i=1}^4 V_{(r_i, s_i)} \right\rangle = D_{s_0} \mathcal{H}_{s_0} + \sum_{r \in \frac{1}{2}\mathbb{N}^*} \sum_{s \in \frac{1}{r}\mathbb{Z} \cap (-1, 1]} D_{(r, s)} \mathcal{H}_{(r, s)} ,$$

in terms of interchiral blocks

$$\mathcal{H}_{s_0} = \sum_{s \in s_0 + 2\mathbb{N}} \frac{D_s}{D_{s_0}} \mathcal{G}_{\langle 1, s \rangle}^D , \quad \mathcal{H}_{(r, s)} = \sum_{j \in 2\mathbb{N}} \frac{D_{(r, s+j)}}{D_{(r, s)}} \mathcal{G}_{(r, s+j)} .$$

Solve then the crossing equations

$$\sum_{V \in \mathcal{S}^{(s)}} D_V^{(s)} \text{ (s-channel diagram)} = \sum_{V \in \mathcal{S}^{(t)}} D_V^{(t)} \text{ (t-channel diagram)} = \sum_{V \in \mathcal{S}^{(u)}} D_V^{(u)} \text{ (u-channel diagram)}$$

s-channel
t-channel
u-channel

We know the spectrum. And we can constrain the solution space by fixing the $O(n)$ symmetry of the exchanged fields V .

In favourable cases this gives a unique (numerical) solution.

Conjecture: Each solutions to the crossing equations gives a valid correlation function in the $O(n)$ CFT.

We have computed the 30 correlation functions with $\sum_{i=1}^4 r_i = 2, 3, 4$.

We have two ways to prove

$$\Lambda_{(r,s)} = \delta_{r,1} \delta_{s \in 2\mathbb{Z}+1} [[]] + \frac{1}{2r} \sum_{r'=0}^{2r-1} e^{\pi i r' s} \chi_{(2r) \wedge r'} \left(\Lambda_{\frac{2r}{(2r) \wedge r'}} \right).$$

1st proof: Compute the torus partition function twisted by a non-trivial group element of $O(n)$. This produces the character $\Lambda_{(r,s)}$, not just its dimension $L_{(r,s)}(n)$ as in Read-Saleur (2001).

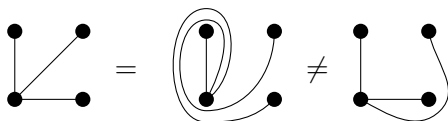
2nd proof: The commutant of $O(n)$ on $\mathcal{S}_L^{O(n)}$ is the Brauer algebra, $\mathcal{B}_L(n)$, generated by e_i and p_i . But in $d = 2$, it reduces to $u\mathcal{JTL}_L(n)$.

Hence we must compute the branching rules $\mathcal{B}_L(n) \downarrow u\mathcal{JTL}_L(n)$.

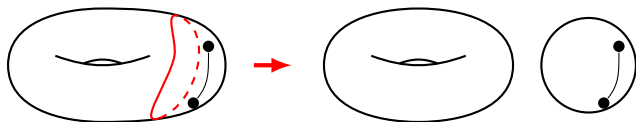
This is a solvable combinatorial problem.

Combinatorial maps

A (connected) *combinatorial map* is a (connected) graph, together with a cyclic permutation of the half-edges around each vertex. Monogons are forbidden.



A map is *weakly connected* if it cannot be split into two non-trivial maps (a sphere with 0 or 1 vertex).

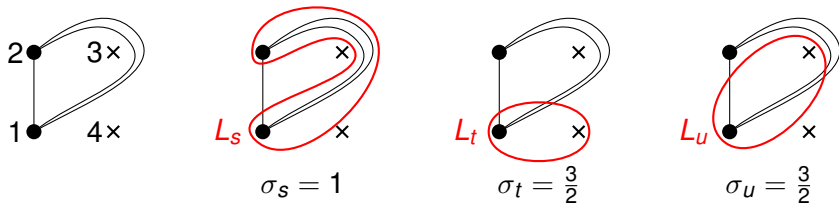


This map is not weakly connected (it should have ‘used’ the handle).

Number of maps $|\mathcal{M}_{g,N}(r_i)|$ and of weakly connected maps $|\mathcal{M}_{g,N}^c(r_i)|$.
 Genus g , number of points N , vertex valencies $2r_j$.

$$|\mathcal{M}_{0,4}^c| = \left[\sum_{i=1}^4 r_i^2 - \frac{1}{2} \right].$$

Signature of a planar map with four vertices:



A map M is weakly connected iff $\forall x \in \{s, t, u\}, \sigma_x(M) > 0$.

For any N -point function of diagonal and non-diagonal fields, the dimension of the space of solutions of conformal bootstrap equation with spectra made only of non-diagonal fields is $\left| \mathcal{M}_{g,N}^C(r_i) \right|$.

The critical limit of a loop model correlation function is a solution of the conformal bootstrap equations.

The set of correlation functions is a basis of solutions of the corresponding conformal bootstrap equations.

Digression on the Barnes double gamma function

Recall $c = 13 - 6\beta^2 - 6\beta^{-2}$ and set $Q = \beta + \beta^{-1}$.

For $\Re x > 0$ define $\Gamma_\beta(x)$ through

$$\log \Gamma_\beta(x) = \int_0^\infty \frac{dt}{t} \left[\frac{e^{-xt} - e^{-Qt/2}}{(1 - e^{-\beta t})(1 - e^{-t/\beta})} - \frac{(Q/2 - x)^2}{2e^t} - \frac{Q/2 - x}{t} \right]$$

and the shift equations

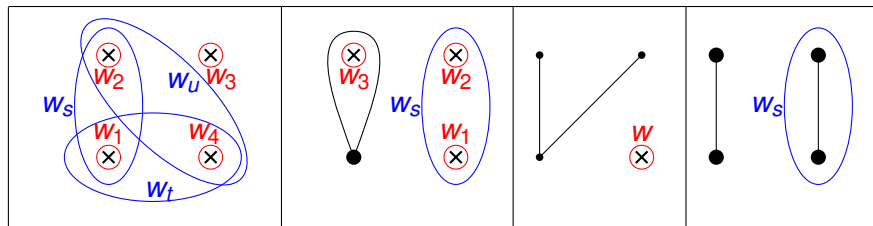
$$\frac{\Gamma_\beta(x + \beta)}{\Gamma_\beta(x)} = \sqrt{2\pi} \frac{\beta^{\beta x - \frac{1}{2}}}{\Gamma(\beta x)} \quad , \quad \frac{\Gamma_\beta(x + \beta^{-1})}{\Gamma_\beta(x)} = \sqrt{2\pi} \frac{\beta^{\frac{1}{2} - \beta^{-1}x}}{\Gamma(\beta^{-1}x)} .$$

Sometimes one defines also the upsilon function

$$\Upsilon_\beta(x) = \frac{1}{\Gamma_\beta(x)\Gamma_\beta(Q - x)} .$$

4-point functions of diagonal and non-diagonal fields

In addition to monomer weight K and bulk loop weight n , define *vertex weights* w_i (with $i = 1, 2, 3, 4$) and *channel weights* w_x (with $x = s, t, u$):



At most one of the loop types w_x can exist in a given configuration.

In the lattice model, define $C^{\text{loop}}(L, \ell | K, n, w_i, w_x)$, with L the size and ℓ the separation between z_1, z_2 and z_3, z_4 (in the s -channel).

We find the s-channel decomposition

$$C^{\text{loop}}(L, \ell | K, n, w_i, w_x) = \sum_{\omega \in S(L)} A_{\omega}(L | K, n, w_i, w_x) \left(\frac{\Lambda_{\omega}(L | K, n, w_s)}{\Lambda_{\max}(L | K, n, w_s)} \right)^{\ell},$$

with $(\Lambda_{\omega})_{\omega \in S(L)}$ the $ATL_L(n)$ spectrum of transfer matrix eigenvalues.

Remarkable that only $ATL_L(n)$ eigenvalues participate here!

Define ratios wrt different values of the weights: $f(x : x') = \frac{f(x)}{f(x')}$.

Even more remarkably, we find that

$$A_{(r,s),\rho}(L | K, n, w_i, w_x : w'_x) = D_{(r,s)}^{(s)}(n, w_i, w_x : w'_x).$$

Here $\omega = (r, s), \rho$, where ρ labels states in the same module (r, s) . There is no dependence on ρ, L and K . Hence the amplitudes have **nothing to do with CFT** and should be computable from $ATL_L(n)$.

Looks like a Wigner-Eckart theorem, but lifted from QM to $ATL_L(n)$.

Reference 2- and 3-point structure constants

Omitting the known coordinate dependence, define:

$$\langle V_1 V_2 \rangle = \delta_{12} B_1 \quad , \quad \langle V_1 V_2 V_3 \rangle = C_{123} .$$

For non-diagonal fields, set:

$$B_{(r,s)}^{\text{ref}} = \frac{(-)^{rs}}{2 \sin(\pi(\text{frac}(r) + s)) \sin(\pi(r + \beta^{-2}s))} \prod_{\pm, \pm} \Gamma_{\beta}^{-1} \left(\beta \pm \beta r \pm \beta^{-1} s \right) ,$$

$$C_{(r_1, s_1)(r_2, s_2)(r_3, s_3)}^{\text{ref}} = \prod_{\epsilon_1, \epsilon_2, \epsilon_3 = \pm} \Gamma_{\beta}^{-1} \left(\frac{\beta + \beta^{-1}}{2} + \frac{\beta}{2} |\sum_i \epsilon_i r_i| + \frac{\beta^{-1}}{2} \sum_i \epsilon_i s_i \right) .$$

For diag fields, set $V_P = V_{(0, 2\beta P)}$, so $C_{(0, 2\beta P_1)(0, 2\beta P_2)(0, 2\beta P_3)}^{\text{ref}} = C_{P_1, P_2, P_3}$.

When $w_i \equiv 0$, C_{P_1, P_2, P_3} gives the probability that three points belong to the same FK cluster [Delfino-Viti, 2013; Ikhlef-J-Saleur, 2016].

Normalised 4-point structure constants

$$D_{(r,s)}^{(x)} = \frac{C_{(r_1,s_1)(r_2,s_2)(r,s)}^{\text{ref}} C_{(r,s)(r_3,s_3)(r_4,s_4)}^{\text{ref}}}{B_{(r,s)}^{\text{ref}}} d_{(r,s)}^{(x)}$$

Combining analytical arguments with numerical bootstrap and transfer matrices, we find that $d_{(r,s)}^{(x)}$ is a polynomial in $n = -2 \cos(\pi\beta^2)$, with β -independent coefficients and $\deg_n d_{(r,s)}^{(x)} \leq r(r-1)$.

If the x -channel decomposition involves a diagonal field V_{P_x} , then $d_{(r,s)}^{(x)}$ is also polynomial in $w(P)$.

If some $V_i = V_{P_i}$ is diagonal, then $d_{(r,s)}^{(x)}$ is polynomial in $w_i = w(P_i)$.

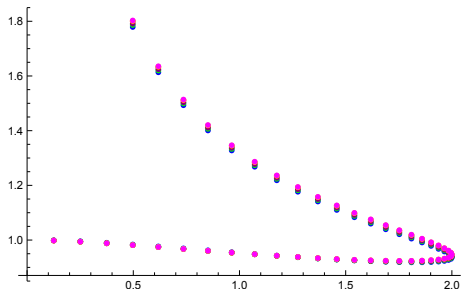
The dependence on $w_x = w(P_x)$ becomes polynomial after we subtract a rational term that is needed for the 4-point function to be holomorphic in P_x .

Results on 3-point structure constants

Not clear how to factorise 4-point structure constants on 3-point ones, since several fields can have the same dimension.

But we find that 3-point structure constants of combinatorial maps are simply given by $C_{(r_1, s_1)(r_2, s_2)(r_3, s_3)}^{\text{ref}}$. TM check for a dozen of cases.

E.g. $C_{(1,0)(1,0)(1,0)}^{\text{ref}}$ gives the probability that 3 points \in same loop.



Xin Sun et al. have an unpublished proof of this one case (using CLE_{κ}).