

Irregular Conformal States and Spectral Curve: Irregular Matrix Model Approach

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Abstract. We present recent developments of irregular conformal states. Irregular vertex operators and their adjoint in a new formalism are used to define the irregular conformal states and their inner product instead of using the colliding limit procedure. Free field formalism can be augmented by screening operators which provide more degrees of freedom. The inner product is conveniently given as the partition function of an irregular matrix model. (Deformed) spectral curve is the loop equation of the matrix model at Nekrasov–Shatashvili limit. We present the details of analytic structure of the spectral curve for Virasoro symmetry and its extensions, W -symmetry and super-symmetry.

Key words: irregular state; irregular conformal block; random matrix model; spectral curve

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1 Introduction

Virasoro symmetry is typically represented by Virasoro module which consists of primary state and its descendants. An irregular conformal state is a different representation of the Virasoro symmetry and is a simultaneous eigenstate of a certain positive modes of conformal generators. Therefore, the irregular state is not a primary or descendant state and is rather a coherent state, combination of primary and its descendants. For example, the simplest irregular state is called Whittaker state [17], the eigenstate of Virasoro L_{+1} mode, and later generalized by Gaiotto [18]. We use the definition of the Virasoro irregular state of rank m as the simultaneous eigenstate of Virasoro generators L_k with $m \leq k \leq 2m$. This definition can be similarly extended to W -symmetry; the simultaneous eigenstate of W -generators of spin s ; $W_k^{(s+1)}$ with $sm \leq k \leq (s+1)m$. One can also construct the irregular states with the super-symmetry.

The irregular state is constructed as the superposition of a primary state and its descendants in [6, 22, 26]. However, there appear some ambiguities for the rank greater than 1. The superposition of states is not completely fixed by the defining relations, simultaneous eigenstate of a certain positive Virasoro generators. One of the main reason for this insufficiency is that other positive generators of the rank m , $W_k^{(s+1)}$ with $0 \leq k < sm$ are non-commuting each other and require consistency condition when applied to the irregular state. It is also noted in [11, 12] that the consistency allows more degrees of freedom other than the simultaneous eigenvalues. Therefore, a non-trivial systematic tool is needed to fix the consistency conditions.

The irregular state is introduced in physics community in relation with AGT conjecture [1]. AGT connects Nekrasov partition function of $N = 2$ super Yang–Mills theory in 4 dimensions [28] with the Liouville regular conformal block in 2 dimensions. At IR limit, one can have a different class of conformal states, so called Argyres–Douglas theory of $N = 2$ super Yang–Mills theory [2],

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which is not the deformation of UV conformal theory. It was pointed out in [18] that the Argyres–Douglas theory can be described in terms of the irregular state. In fact, a simple way to obtain the irregular state is suggested in [16, 19], which uses ‘colliding limit’ of the Liouville conformal block or equivalently from Penner-type matrix model [15, 21].

During the last few years, even though many of the technical developments are achieved for the evaluation of irregular conformal block (ICB) in a significant way, the clear understanding of the irregular conformal block is still missing. The role of the colliding limit is not well understood. In this paper, instead of directly using the ‘colliding limit’, we try a new approach to find ICB. We hope this new alternative approach provides a more intuitive way to understand the structure of ICB and Argyres–Douglas theory.

To do this, we adopt the irregular state of bosonic free fields, which is directly obtained from the definition of the irregular vertex operator (IVO) [20, 27, 32]. The next step is to find the adjoint of the state, a necessary ingredient to define its inner product. However, any explicit form of the adjoint of the irregular state has never been defined in the literature. We present here the definition of the adjoint, which uses the conformal transformation $z \rightarrow 1/z$ as used in the usual definition of the adjoint of the regular primary state. There needs, however, a proper rescaling of the state for the irregular one, which is the main difference from the usual definition of the regular one. After this setup, the inner product is easily evaluated, and turns out to be fixed by coherent coordinates only. This shows that the free field formalism is too restrictive to describe the Argyres–Douglas theory, since more degrees of freedom are needed. In fact, the necessary degrees of freedom are provided by the screening operators. As a result, the inner product is effectively given in terms of the partition function of a beta-deformed Penner-type matrix model with the additional potential of the polynomial and inverse polynomial type, which is called the irregular matrix model in short. According to this set-up, the necessary building blocks of ICB are the irregular states, their adjoints and screening operators. It is demonstrated that all the previous results appeared for ICB are reproduced in this approach without using the colliding limit.

This paper is organized as follows. Section 2 is the new set up of the formalism of ICB. Using the one bosonic free field formalism, the main idea is presented in a formal way. The irregular vertex operator of the free bosonic field is employed and its adjoint are newly defined. The inner product between the irregular state and its adjoint is explicitly evaluated. This idea can be easily extended to many field formalism including fermionic fields. After the set up, we include the screening operators to provide the inner product with more degrees of freedom. It is demonstrated that the effect of including screening operator in the inner product is conveniently represented as the partition function of an irregular matrix model. The result is equivalent to the colliding limit of the Penner-type matrix model which describes the regular conformal block.

Some explicit examples are given in Section 3 and 4. The Virasoro symmetry is considered in Section 3. The inner product of the irregular conformal states coincides with the colliding limit of the Liouville regular conformal block. Therefore, one can use the technique developed in the random matrix formalism, namely, the loop equation which encodes the conformal symmetry of the irregular matrix model. For simplicity, we are considering the Nekrasov–Shatashvili (NS) limit [29], where the loop equation is truncated to the one-point resolvent only. The resulting loop equation is viewed as the (deformed) spectral curve. The analytic structure of the spectral curve provides all the information of the system, including the partition function known as tau function. We discuss the analytic properties of the spectral curve, nature of the cut structure/pole structure, and the flow equation which plays the important role of finding the partition function. One may extend the same method to W - and super-symmetry, which is shortly described in Section 4. Section 5 is the conclusion.

We skip many of technical details since the purpose of this paper is to describe the structure of the irregular conformal block in a more intuitive way. Section 2 and 3 are enough to read

for understanding the mainstream of ICB. Dedicated readers may find more technical details in [10, 12] for W -symmetry, and [33, 34] and references therein for supersymmetry.

2 Irregular vertex operator, irregular state and its inner product

2.1 free field formalism

General formalism of free field representation of the irregular vertex operator (IVO) $I^{(m)}(z)$ is considered in [20, 27, 32] using bosonic field ϕ . In this section, we consider IVO of rank m using one field:

$$I^{(m)}(w) = e^{2\Phi^{(m)}/\hbar(w)}, \quad \Phi^{(m)}(w) = \sum_{k=0}^m \frac{c_k}{k!} \frac{\partial^k \phi(w)}{\partial w^k}, \quad (2.1)$$

where \hbar is a convenient parameter and the bosonic field has the holomorphic normalization

$$\langle \phi(z)\phi(w) \rangle = -\frac{1}{2} \log(z-w), \quad \langle e^{\alpha_1 \phi(z)} e^{\alpha_2 \phi(w)} \rangle = (z-w)^{-2\alpha_1 \alpha_2}.$$

The energy momentum tensor and its moments are given as

$$T(z) = -(\partial\phi)^2, \quad L_k = \oint \frac{dz}{2\pi} z^{1+k} T(z) \quad (2.2)$$

and IVO satisfies the operator product

$$\oint_{w=0} \frac{dz}{2\pi} z^{1+k} T(z) I^{(m)}(w) = \Lambda_k I^{(m)}, \quad (2.3)$$

where $\hbar^2 \Lambda_k = -\sum_{t+r=k} c_t c_r$. This shows that $\Lambda_k = 0$ when $k > 2m$.

Irregular state $|I^{(m)}\rangle$ is obtained if IVO is applied on the vacuum $|0\rangle$:

$$|I^{(m)}\rangle = \lim_{z \rightarrow 0} I^{(m)}(z)|0\rangle = I^{(m)}|0\rangle. \quad (2.4)$$

The condition (2.3) shows that the irregular state is the simultaneous eigenstate of positive generators

$$[L_k, I^{(m)}] = \Lambda_k I^{(m)}.$$

Therefore, $|I^{(m)}\rangle$ in (2.4) can be the simultaneous eigenstate of positive mode Virasoro generators L_a with $m \leq a \leq 2m$ since $L_{a \geq 0}$ satisfies the commutation relation $[L_a, L_b] = (a-b)L_{a+b}$.

The field derivative terms in IVO represent the descendant contributions. And the coefficients c_k correspond to the coherent coordinates of the Heisenberg positive modes a_k with $1 \leq k \leq m$. This can be seen if one expands the bosonic field in terms of the Heisenberg modes $\partial\phi(z) = -\sum_k a_k/z^{k+1}$ and

$$\oint_{w=0} \frac{dz}{2\pi} z^{1+k} \partial\phi(z) I^{(m)}(w) = c_k I^{(m)}$$

or $[a_k, I^{(m)}] = c_k I^{(m)}$. Therefore, it is obvious that

$$a_k |I^{(m)}\rangle = c_k |I^{(m)}\rangle, \quad 1 \leq k \leq m.$$

To define an inner product we need an adjoint state. The corresponding adjoint operator is defined at infinity

$$\langle I^{(n)} | = \lim_{\zeta \rightarrow 0} \langle 0 | \hat{I}^{(n)}(\zeta),$$

where conformal-transformation from the infinity to 0, $z \rightarrow 1/\zeta$ is used. The adjoint operator is equivalent to the same one (2.1) but is defined at $1/\zeta$

$$\hat{I}^{(n)}(\zeta) = R_0^{(n)} e^{2\hat{\Phi}^{(n)}/\hbar}(1/\zeta), \quad \hat{\Phi}^{(n)} = \sum_{\ell=0}^n \frac{\hat{c}_\ell}{\ell!} \frac{\partial^\ell \phi(1/\zeta)}{\partial \zeta^\ell}.$$

$R_0^{(n)}$ is the scaling factor

$$R_0^{(n)} = e^{2\frac{\hat{c}_0}{\hbar^2} \sum_{\ell=0}^n \frac{\hat{c}_\ell}{\ell!} \frac{\partial^\ell \log(\zeta)}{\partial \zeta^\ell}} = (\zeta)^{2\frac{\hat{c}_0^2}{\hbar^2}} \prod_{\ell=1}^n e^{2\frac{\hat{c}_0 \hat{c}_\ell}{\hbar^2 \ell!} \frac{\partial^\ell \log(\zeta)}{\partial \zeta^\ell}}. \quad (2.5)$$

Using the state and its adjoint we can define the inner product $\langle I^{(n)} | I^{(m)} \rangle$ which is given as two-point correlation

$$\langle I^{(n)} | I^{(m)} \rangle_0 = \lim_{w, \zeta \rightarrow 0} \langle 0 | \hat{I}^{(n)}(\zeta) I^{(m)}(w) | 0 \rangle_0.$$

We put the subscript 0 to emphasize that the inner product is defined in terms of the free field representation. It is noted that the primary operator product has the non-vanishing expectation value if $\hat{c}_0 + c_0 = 0$. This will be called the neutrality condition for the inner product. The inner product need be symmetric

$$\langle I^{(n)} | I^{(m)} \rangle_0 = \langle I^{(m)} | I^{(n)} \rangle_0.$$

Explicit evaluation shows the symmetric property

$$\langle I^{(n)} | I^{(m)} \rangle_0 = \lim_{w, \zeta \rightarrow 0} \exp \left(-2 \sum_{k, \ell \geq 1} \frac{\hat{c}_\ell c_k}{\ell! k!} \partial_\zeta^\ell \partial_w^k \log(1 - w\zeta) \right) = e^{\zeta_{n,m}/\hbar^2},$$

$$\zeta_{n,m} = \sum_{\ell \geq 1} 2\hat{c}_\ell c_\ell / \ell, \quad (2.6)$$

where the summation in $\zeta_{n,m}$ holds until $\ell \leq \min(m, n)$. The scaling factor $R_0^{(n)}$ compensates the infinite contribution at $\zeta = 0$ so that the inner product is finite. In addition, $\langle I^{(0)} | I^{(m)} \rangle_0 = \langle I^{(n)} | I^{(0)} \rangle_0 = 1$. This is consistent with the fact that $|I^{(m)}\rangle$ consists of primary state $|I^{(0)}\rangle$ and its descendants where $\langle I^{(0)} | I^{(0)} \rangle_0$ is normalized to 1.

The inner product is the function of coherent coordinates c_k and \hat{c}_ℓ . We may use c_k and \hat{c}_ℓ as the coordinates to represent the positive Virasoro mode acting on the state. For the case with $0 \leq a \leq m-1$, we have

$$L_a |I^{(m)}\rangle = v_a |I^{(m)}\rangle, \quad v_a = \sum_k k c_{a+k} \frac{\partial}{\partial c_k} \quad (2.7)$$

and v_a satisfies the commutation relation

$$[v_a, v_b] = (b-a)v_{a+b}, \quad (2.8)$$

which corresponds to right action of the Virasoro commutation relation. One can check the commutation relation holds for the inner-product

$$[v_a, v_b] \langle I^{(n)} | I^{(m)} \rangle = (b - a) v_{a+b} \langle I^{(n)} | I^{(m)} \rangle \quad \text{or} \quad [v_a, v_b] (\zeta_{n,m}) = (b - a) v_{a+b} (\zeta_{n,m}).$$

Likewise, the adjoint state introduces the Virasoro representation in \hat{c}_ℓ space

$$\langle I^{(n)} | L_a = \hat{v}_a \langle I^{(n)} |, \quad \hat{v}_a = \sum_{\ell} \ell \hat{c}_{a+\ell} \frac{\partial}{\partial \hat{c}_\ell}, \quad (2.9)$$

and \hat{v}_a satisfies the commutation relation

$$[\hat{v}_a, \hat{v}_b] = (b - a) \hat{v}_{a+b}.$$

It is obvious that the commutation relation also holds for the inner-product

$$[\hat{v}_a, \hat{v}_b] (\zeta_{n,m}) = (b - a) \hat{v}_{a+b} (\zeta_{n,m}).$$

Finally, it is noted that the two representations \hat{v}_a and v_a are commuting each other

$$[v_a, \hat{v}_b] = 0.$$

It is noted that the first-order differential representation of the positive Virasoro generators allows the calculation much easier. On the other hand, if one wants to consider the full representation of the positive and negative Virasoro generators, then one inevitably needs to consider the second-order differential representation.

2.2 Screening operator and irregular matrix model

Whittaker state, the eigenstate of L_1 and annihilated by $L_{k>1}$ was constructed explicitly in [17] as the superposition of a primary state and its descendants. In [18, 26], the irregular state of rank 1 is also constructed. And later, the irregular state of arbitrary rank $m > 1$ is suggested in [22] with unknown coefficients

$$|G_{2m}\rangle = \sum_{\ell, Y, \ell_p} \Lambda^{\ell/m} \left\{ \prod_{i=1}^{m-1} a_i^{\ell_{2m-i}} b_i^{\ell_i} \right\} \times \mu^{\ell_m} Q_{\Delta}^{-1} (1^{\ell_1} 2^{\ell_2} \dots (2m-1)^{\ell_{2m-1}} (2m)^{\ell_{2m}}; Y) L_{-Y} |\Delta\rangle, \quad (2.10)$$

where $|\Delta\rangle$ is the primary state with conformal dimension Δ . $L_{-Y} = L_Y^+$ represents the product of lowering operators and $L_Y = L_1^{\ell_1} L_2^{\ell_2} \dots L_s^{\ell_s}$. The summation ℓ and Y run from 0 to ∞ , maintaining $\ell = |Y| = \sum_p p \ell_p$. $Q_{\Delta}(Y; Y')$ is the Shapovalov form, $Q_{\Delta}(Y; Y') = \langle \Delta | L_{Y'} L_{-Y} | \Delta \rangle$. The coefficients a_i and μ are closely related with eigenvalues; $\Lambda_{2m} = \Lambda^m$, $\Lambda_k = \Lambda^{k/m} a_{2m-k}$ for $m < k < 2m$ and $\Lambda_m = \Lambda \mu$. However, b_i 's are not fixed by the eigenvalues but still gives the contribution to the expectation values of the lower positive generators. The coefficients are to be fixed with the consistency conditions given in (2.8). This suggests that free field IVO cannot be a unique solution. In addition, considering AGT which relates CFT with $N = 2$ super gauge theory, we need more degrees of freedom to connect the irregular states with Argyres–Douglas theory.

More degrees of freedom can be added when homogeneous solution O satisfying $[L_k, O] = 0$ for $k \geq 0$ is considered. The solution O is called screening operator and is given as a contour integral of primary operator $\psi(z)$ with holomorphic dimension 1

$$O = \oint dz \psi(z).$$

Then, the screening operator allows a general solution which will be the free field solution (2.1) with any number of screening operators multiplied.

Liouville field theory has the scaling dimension 1 operator in the presence of background charge $Q = b + 1/b$. The vertex operator $V_\alpha(z) = e^{2\alpha\phi}(z)$ has the scaling dimension $\Delta_\alpha = \alpha(Q - \alpha)$ since the background charge modifies the energy-momentum tensor from the free field case (2.2)

$$T(z) = -(\partial\phi)^2 + Q\partial^2\phi.$$

The background charge makes the conformal dimension of $V_b(z)$ be 1 and changes the eigenvalue Λ_k of IVO appeared in (2.1)

$$\hbar^2\Lambda_k = \epsilon(k+1)c_k - \sum_{r+s=k} c_r c_s, \quad (2.11)$$

where we use the notation $\epsilon = \hbar Q$.

Due to the insertion of the screening operators, the inner product becomes more complicated

$$\langle I^{(0)}|I^{(m)}\rangle_N = \langle I^{(0)}|O^N|I^{(m)}\rangle_0 = Z_N^{(0,m)}[c_0; \mathbf{c}],$$

where N denotes that there are N number of screening operators are inserted. $Z_N^{(0,m)}$ is the partition function of irregular matrix model [16, 30] and depends on the coherent coordinates $\mathbf{c} = \{c_1, \dots, c_n\}$:

$$\begin{aligned} Z_N^{(0,m)}[c_0; \mathbf{c}] &= \left[\int \prod_{I=1}^N d\lambda_I \right] \prod_{I < J} (\lambda_I - \lambda_J)^{2\beta} e^{\frac{\sqrt{\beta}}{g} V^{(0,m)}(\{\lambda_I\})}, \\ V^{(0,m)}(\{\lambda_I\}) &= \sum_{I=1}^N \left\{ c_0 \log(\lambda_I) - \sum_{k=1}^m \frac{c_k}{k\lambda_I^k} \right\}. \end{aligned} \quad (2.12)$$

Here we redefine $\beta = -b^2$ or $b = i\sqrt{\beta}$ and $g = i\hbar/2$ for later convenience.

When no screening operator is included, the inner product $\langle I^{(0)}|I^{(m)}\rangle_N$ reduces to $\langle I^{(0)}|I^{(m)}\rangle_0 = 1$ as in free field theory formalism in (2.6). However, it should be emphasized that $\langle I^{(0)}|I^{(m)}\rangle_N$ is not normalized as 1 due to the screening operator effect and the parameter dependence of the inner product is of the primary concern.

The inner product $\langle I^{(0)}|I^{(m)}\rangle_N$ should be symmetric

$$\langle I^{(0)}|I^{(m)}\rangle_N = \langle I^{(m)}|I^{(0)}\rangle_N,$$

where the proper scaling factor in the definition of the adjoint definition (2.5) is modified due to the background charge

$$R_\epsilon^{(m)} = e^{2\frac{(\hat{c}_0 - \epsilon)}{\hbar^2} \sum_{\ell=0}^m \frac{\hat{c}_\ell}{\ell!} \frac{\partial^\ell \log(\zeta)}{\partial \zeta^\ell}} = (\zeta)^{2\frac{(\hat{c}_0 - \epsilon)\hat{c}_0}{\hbar^2}} \prod_{\ell=1}^m e^{2\frac{(\hat{c}_0 - \epsilon)\hat{c}_\ell}{\hbar^2 \ell!} \frac{\partial^\ell \log(\zeta)}{\partial \zeta^\ell}}. \quad (2.13)$$

Note that the symmetric property of the inner-product is related with the conformal invariance. After the conformal transformation $\lambda_I \rightarrow 1/\lambda_I$, the partition function changes to $Z_N^{(m,0)}[\hat{c}_0; \mathbf{c}]$ whose potential is now given as

$$V^{(m,0)}(\{\lambda_I\}) = \sum_{I=1}^N \left\{ \hat{c}_0 \log(\lambda_I) - \sum_{k=1}^m \frac{\hat{c}_k \lambda_I^k}{k} \right\},$$

where $\hat{c}_k = c_k$ and \hat{c}_0 is a quantity hidden at infinity and is defined through the neutrality condition

$$\hat{c}_0 + \hbar N b + c_0 = \epsilon. \quad (2.14)$$

Conformal invariance of the partition function shows that

$$Z_N^{(0,m)}[c_0; \mathbf{c}] = Z_N^{(m,0)}[\hat{c}_0; \hat{\mathbf{c}}]$$

and therefore, the symmetric property of the inner product is assured.

The inner product of two irregular states is similarly given as

$$\langle I^{(n)} | I^{(m)} \rangle_N = \mathcal{Z}_N^{(n,m)}[c_0; \hat{\mathbf{c}}; \mathbf{c}] = \lim_{w, z \rightarrow 0} \langle 0 | I^{(n)}(1/w) O^N I^{(m)}(z) | 0 \rangle = e^{\zeta_{n,m}/\hbar^2} Z_N^{(n,m)}[c_0; \hat{\mathbf{c}}; \mathbf{c}],$$

where $\zeta_{n,m}$ is the free field contribution and $Z_N^{(n,m)}$ is the screening field effect [9]

$$Z_N^{(n,m)}[c_0; \hat{\mathbf{c}}; \mathbf{c}] = \left[\int \prod_{I=1}^N d\lambda_I \right] \prod_{I < J} (\lambda_I - \lambda_J)^{2\beta} e^{\frac{\sqrt{\beta}}{g} V^{(n,m)}(\{\lambda_I\})}, \quad (2.15)$$

$$V^{(n,m)}(\{\lambda_I\}) = \sum_{I=1}^N \left\{ c_0 \log(\lambda_I) - \sum_{k=1}^m \frac{c_k}{\lambda_I^k} - \sum_{\ell=1}^n \frac{\hat{c}_\ell}{\ell} \lambda_I^\ell \right\}. \quad (2.16)$$

The conformal invariance

$$Z_N^{(n,m)}[c_0; \hat{\mathbf{c}}; \mathbf{c}] = Z_N^{(m,n)}[\hat{c}_0; \mathbf{c}; \hat{\mathbf{c}}]$$

is obvious with the neutrality condition (2.14) and so is the symmetric property of the inner product.

One may also define irregular conformal block $\mathcal{F}_{N_1, N_2}^{(n,m)}[c_0; \hat{\mathbf{c}}; \mathbf{c}]$ in terms of inner product [12]

$$\langle I^{(n)} | I^{(m)} \rangle_N = \sum_{\Delta} \mathcal{F}_{N_1, N_2}^{(n,m)}[c_0; \hat{\mathbf{c}}; \mathbf{c}] \langle I^{(n)} | \Delta \rangle_{N_1} \langle \Delta | I^{(m)} \rangle_{N_2},$$

where $N = N_1 + N_2$ and $|\Delta\rangle$ is an intermediate primary state of conformal dimension $\hbar^2 \Delta = (\hat{c}_0 + N_1 \epsilon)(c_0 + N_2 \epsilon)$. In this definition, the intermediate states are to be inserted in the definition of the inner product $\langle I^{(n)} | I^{(m)} \rangle_N$. The irregular conformal block is given in terms of the partition function as

$$Z_N^{(n,m)}[c_0; \hat{\mathbf{c}}; \mathbf{c}] = e^{-\zeta_{n,m}/\hbar^2} \sum_{N_1 + N_2 = N} \mathcal{F}_{N_1, N_2}^{(n,m)}[c_0; \hat{\mathbf{c}}; \mathbf{c}] Z_{N_1}^{(n,0)}[\hat{c}_0; \hat{\mathbf{c}}] Z_{N_2}^{(0,m)}[c_0; \mathbf{c}].$$

One practical way of evaluating ICB is to use the perturbative approach [11], where the potential of partition function $Z_N^{(n,m)}[c_0; \hat{\mathbf{c}}; \mathbf{c}]$ is splitted so that the reference one is given as $Z_{N_1}^{(n,0)}[\hat{c}_0; \hat{\mathbf{c}}] \times Z_{N_2}^{(0,m)}[c_0; \mathbf{c}]$ and the rest as the perturbative one. The explicit calculation demonstrates that the parameter b_k appeared in (2.10) is identified as $\Lambda^{k/m} b_k = \Lambda_k + v_k (-\hbar^2 \log Z_N^{(0,m)})$ with $\Lambda^m = \Lambda_{2m}$.

The irregular matrix model (2.15) can also be obtained using the colliding limit [9] of he regular Liouville conformal block, holomorphic correlation of vertex operators $V_\alpha = e^{2\alpha\phi}$ together with the screening operators. The regular conformal block with $(n + m + 2)$ primary operators

$$\mathcal{G}_N^{(n+m)} = \left\langle \prod_{A=1}^{n+m+2} V_{\alpha_A}(z_A) \left(\oint dz e^{b\phi(z)} \right)^N \right\rangle$$

satisfies the neutrality condition $\sum_{A=1}^{n+m+2} \alpha_A + bN = Q$. Colliding limit is to put $m + 1$ operators at the origin with c_k ($1 \leq k \leq m$) finite

$$c_k = \sum_{A=1}^m \hbar \alpha_A z_A^k,$$

where z_A is the coordinates of the operators approaching to 0. We rescaled α_A with \hbar so that c_k is identified with the coefficients in the potential in (2.16). At infinity, $n + 1$ primary operators is put to infinity with \hat{c}_ℓ ($1 \leq \ell \leq n$) finite

$$\hat{c}_\ell = \sum_{A=1}^n \hbar \alpha_A \zeta_A^\ell$$

with $z_A = 1/\zeta_A \rightarrow \infty$. After the proper scaling $R_\epsilon^{(n)}$ introduced in (2.13), $\mathcal{G}_N^{(n+m)}$ reduces to the inner product

$$\langle I^{(n)} | I^{(n)} \rangle_N = e^{\zeta_{n,m}/\hbar^2} \mathcal{Z}_N^{(n,m)}[c_0; \hat{\mathbf{c}}, \mathbf{c}].$$

It is to be noted that the free field contribution $e^{\zeta_{n,m}}$ is called $U(1)$ effect which appears from the factor

$$\lim_{\zeta_A, z_B \rightarrow 0} \prod_{A,B} (1 - \zeta_A z_B)^{-2\alpha_A \alpha_B} = e^{\zeta_{n,m}/\hbar^2}$$

at the colliding limit (z_B, ζ_A) $\rightarrow 0$.

3 Properties of inner product

3.1 (Deformed) spectral curve

The inner product (2.12) or (2.15) requires a systematic evaluation of the effect of the screening operators. We are using the salient feature of the partition function, the conformal invariance. Under the holomorphic transform $\lambda_I \rightarrow f(\lambda_I) = \lambda_I + \delta/(z - \lambda_I)$, the conformal invariance of the partition function is encoded in the loop equation

$$4W(z)^2 + 4V'(z)W(z) + 2\epsilon W'(z) - \hbar^2 W(z, z) = f(z). \quad (3.1)$$

Here $W(z)$ is the one-point resolvent

$$W(z) = g\sqrt{\beta} \left\langle \sum_I \frac{1}{z - \lambda_I} \right\rangle, \quad (3.2)$$

where the bracket $\langle \dots \rangle$ denotes the expectation value with respect to the irregular matrix model. Primed quantity denotes the derivative with respect to its argument. Two point resolvent $W(z, z)$ is given as the connected part

$$W(z_1, z_2) = \beta \left\langle \sum_I \frac{1}{z_1 - \lambda_I} \sum_J \frac{1}{z_1 - \lambda_J} \right\rangle_{\text{conn}}.$$

$f(z)$ has the potential dependence

$$f(z) = 4g\sqrt{\beta} \left\langle \sum_I \frac{V'(z) - V'(\lambda_I)}{z - \lambda_I} \right\rangle.$$

For simplicity, we will consider the classical/NS limit of the loop equation (3.1). Nekrasov–Shatashvili (NS) limit is obtained if $\hbar \rightarrow 0$ and $b \rightarrow \infty$ so that $\epsilon = \hbar Q = 2g\sqrt{\beta}$ is finite [29]. On the other hand, classical limit is obtained in the Liouville field theory when $b \rightarrow 0$ and $\epsilon = \hbar Q$ is finite. However, Liouville field theory has the duality $b \rightarrow 1/b$ so that two limits, $b \rightarrow \infty$ and $b \rightarrow 0$ are equivalent. At the classical/NS limit, one can check that multi-point resolvent vanishes and the loop equation is given in terms of one-point resolvent only

$$4W(z)^2 + 4V'(z)W(z) + 2\epsilon W'(z) = f(z). \quad (3.3)$$

This loop equation becomes a more informative form if $x = 2W + V'$ is used

$$x^2 + \epsilon x' + \xi_2(z) = 0. \quad (3.4)$$

This Riccati equation is regarded as a deformed (due to $\epsilon x'$ term) spectral curve.

Analytic structure of the spectral curve is specified by $\xi_2(z) = -V'^2 + \epsilon V'' - f$, contains details of the information of Virasoro symmetry. This fact is seen as follows: For the type of potential $V = V^{(n,m)}$, f has the analytic structure

$$f = \sum_{a=-n}^{m-1} \frac{d_a}{z^{a+2}},$$

where the prime in the summation denotes that $a = -1$ is missing since $d_{-1} = 0$. Therefore, $\xi_2(z)$ has irregular poles and zeros on the Riemann surface

$$\xi_2(z) = \sum_{k=-2n}^{2m} \frac{\hbar^2 \Lambda_k}{z^{k+2}} - \sum_{a=-n}^{m-1} \frac{d_a}{z^{a+2}},$$

where Λ_k is the same one given in (2.11) with the definition $c_{-k} = \hat{c}_k$. Due to the irregular singularity, the Riccati equation (3.4) shows the different analytic structure from the regular ones.

Suppose we ignore $f(z)$ (putting $d_a = 0$). Then, $\xi_2(z)$ is given in terms of Λ_k . The positive modes L_k for $m \leq k \leq 2m$ apply on the state $|I^{(m)}\rangle$, and the negative modes L_k for $-2n \leq k \leq -n$ on the state $\langle I^{(n)}|$, both resulting in the desired eigenvalues.

To understand the role of d_a , let us consider the case d_a with $a \geq 0$. One can check that from the definition of f one has

$$d_a = v_a(F^{(n;m)}) \quad \text{for } 0 \leq a \leq m-1, \quad (3.5)$$

where $F_N^{(n;m)} = -\hbar^2 \log Z_N^{(n;m)}$ and v_a is given in (2.7), Virasoro representation in $\{c_k\}$ space. For the case d_{-a} with $a \geq 2$ one has $d_{-a} = -2\epsilon \hat{c}_a + \hat{d}_a$. We have \hat{d}_a with $0 \leq a \leq n-1$ in terms of Virasoro representation

$$\hat{d}_a = \hat{v}_a(F_N^{(n;m)}), \quad (3.6)$$

where \hat{v}_a is given in (2.9), Virasoro representation in $\{\hat{c}_\ell\}$ space. Therefore, it is clear that $\xi_2(z)$ is identified with the expectation value of the energy momentum tensor (Virasoro current)

$$\xi_2 = \langle \hbar^2 T(z) \rangle = \sum_{k \in \mathbb{Z}} \frac{\langle \hbar^2 L_k \rangle}{z^{k+2}}.$$

Here, expectation is given in terms of irregular matrix model and is also considered as the normalized expectation value

$$\langle A \rangle = \frac{\langle I^{(n)} | A | I^{(m)} \rangle_N}{\langle I^{(n)} | I^{(m)} \rangle_N}.$$

It is also important to note that $\hbar^2 \Lambda_k$ and the free energy $F^{(n;m)}$ is regarded as finite at the classical/NS limit [37]. Moreover, since $Z_N^{(n;m)} = \exp(-F^{(n;m)}/\hbar^2)$, the flow equations (3.5) and (3.6) satisfies consistency conditions

$$\begin{aligned} v_a(d_b) - v_b(d_a) &= (a-b)d_{a+b}, & \hat{v}_a(\hat{d}_b) - \hat{v}_b(\hat{d}_a) &= (a-b)\hat{d}_{a+b}, \\ v_a(\hat{d}_b) - \hat{v}_b(d_a) &= 0. \end{aligned} \quad (3.7)$$

The consistency conditions are useful tools to construct the partition function. The major step is to find the values d_a and \hat{d}_a directly from the analytic property of the spectral curve (3.4). This procedure is presented in the following subsection.

3.2 Irregular spectral curve and polynomial equation

When $\epsilon = 0$, the spectral curve (3.4) becomes the large N limit of random matrix models which can be solve as in the usual approach [7, 8]. Therefore, one can solve the spectral curve keeping x is $O(1)$ so that $\epsilon x'$ is regarded as sub-dominant. In this case, the dominant term is simply given as $x = \pm\sqrt{-\xi_2}$ and solves the spectral curve (3.4) as the perturbation series in powers of ϵ [9, 30]. As a result, the solution has $(m+n)$ square-root branch cuts and provides the double covering of the Riemann surface [14, 15].

However, at the classical/NS limit, the derivative term x' survives in the spectral curve (3.4) and can change the analytic structure of the solution space. This is related with the well-known fact that one can transform the Riccati equation into a second-order linear differential equations. Putting $x = \epsilon(\log \Psi)'$ one has

$$\left(\epsilon^2 \frac{\partial^2}{\partial z^2} + \xi_2(z) \right) \Psi(z) = 0.$$

To understand the meaning of the transformation of the spectral curve, let us consider an expectation value [4, 36]

$$P(z) \equiv \left\langle \prod_I (z - \lambda_I) \right\rangle = \sum_{A=0}^N P_A z^A = \prod_{\alpha} (z - z_{\alpha}), \quad (3.8)$$

which is a monic polynomial of degree N ($P_N = 1$). z_{α} are zeros of the polynomial, which are assumed distinct. One can check that $P(z)$ is related with the resolvent $W(z)$ in (3.2) at the classical/NS limit. Using $\log \left(\prod_I (z - \lambda_I) \right) \propto \sum_I \int^z \frac{dz'}{z' - \lambda_I}$ one has

$$\log \left(\frac{P(z)}{P(z_0)} \right) = \frac{2}{\epsilon} \int_{z_0}^z dz' W(z'), \quad (3.9)$$

since the multi-point resolvent vanishes at the classical/NS limit [4, 25]. Taking the derivative of (3.9), we have

$$2W(z) = \epsilon P'(z)/P(z) = \epsilon \sum_{\alpha=1}^N \frac{1}{z - z_{\alpha}}. \quad (3.10)$$

Then, the monic polynomial $P(z)$ satisfies the second-order differential equation

$$\epsilon^2 P''(z) + 2\epsilon V'(z)P'(z) = f(z)P(z). \quad (3.11)$$

The solution of the differential equation shows that there are N zeros and therefore, the resolvent $W(z)$ has simple poles. As a result, one can conclude that there will appear N -simple poles in

the spectral curve rather than the branch cut present when $\epsilon = 0$. The branch cut disappears and only simple poles are present.

In addition, one can find d_a from the differential equation (3.11) in terms of coherent coordinates, $\{c_k\}$ and $\{\hat{c}_k\}$. Then, the consistency condition (3.7) is solved by noting that $v_a = \sum_k U_{ak}^{-1} \partial_k$ or $\partial_k = \sum_a U_{ka} v_a$, where $\partial_k = \partial/\partial c_k$. U is a $m \times m$ matrix for the rank m case and is simply given by the coherent coordinates. Then, a closed one form $d(F^{(n,m)}) = \sum_k dc_k \partial_k(F^{(n,m)})$ satisfies the flow equations $d(F^{(n,m)}) = \sum_{k,a} dc_k U_{ka} d_a$. According to the spectral curve (3.3), d_a is the residue of $z^{1+a}(4W^2(z) + 4V'(z)W(z) + 2\epsilon W'(z))$ so that

$$d_a = -4 \sum_{\ell \geq 1} \frac{c_{a+\ell}}{\ell!} \partial^\ell W(0) = 2\epsilon \sum_{\ell \geq 1} c_{a+\ell} \left(\sum_{\alpha=1}^N \frac{1}{z_\alpha^\ell} \right),$$

where the solution $W(0)$ is the function of coherent coordinates. Using the zeros of $P(z)$, one has the flow equation

$$d(F_N^{(n,m)}) = \sum_{k,a} dc_k U_{ka} d_a = 2\epsilon \sum_{k,a} dc_k U_{ka} \sum_{\ell \geq 1} c_{a+\ell} \left(\sum_{\alpha=1}^N \frac{1}{z_\alpha^\ell} \right). \quad (3.12)$$

The same form of the flow equation holds for the hat coordinates

$$\hat{d}(F_N^{(n,m)}) = \sum_{k,a} d\hat{c}_k \hat{U}_{ka} \hat{d}_a = 2\epsilon \sum_{k,a} d\hat{c}_k \hat{U}_{ka} \sum_{\ell \geq 1} \hat{c}_{a+\ell} \left(\sum_{\alpha=1}^N \frac{1}{\hat{z}_\alpha^\ell} \right).$$

where \hat{U} and \hat{z}_α are the adjoint expression which is obtained from the U and z_α by putting $c_a \leftrightarrow \hat{c}_a$. (The relation $\hat{d}_{a \geq 2}$ can be checked from the residue $z^{1-a}(4W^2(z) + 4V'(z)W(z) + 2\epsilon W'(z))$).

It is to be noted that the polynomial $P(z)$ is closely related with the degenerate primary operator expectation value. Let us consider $\langle I^{(n)} | V_+(z) | I^{(m)} \rangle$, where $V_+(z) \equiv V_{-1/(2b)}$ has the conformal dimension $\Delta_+ = -\frac{1}{2} - \frac{3}{4b^2}$ and has the null vector at level 2. Let us define $\Psi_+^{(n,m)}(z)$, normalized expectation value

$$\Psi_+^{(n,m)}(z) = \frac{\langle I^{(n)} | V_+(z) | I^{(m)} \rangle_{N_+}}{\langle I^{(n)} | I^{(m)} \rangle_N},$$

where $N_+ - N = 1/(2b^2)$ is assumed for the neutrality condition to hold. This requires that the number of screening operators used to evaluate the partition function is different from that used to evaluate the expectation value. At NS limit, however, this unpleasant feature disappears since $N_+ = N$. Then it is easy to find

$$\Psi_+^{(n,m)}(z) = P(z) e^{V^{(n,m)}(z)/\epsilon} = \exp\left(\frac{1}{\epsilon} \int^z x(z') dz'\right),$$

where $P(z)$ is the solution of (3.11). This show that the deformed spectral equation (3.4) or (3.11) leads to the second-order differential equation of $\Psi_+^{(n,m)}(z)$

$$\left(\epsilon^2 \frac{\partial^2}{\partial z^2} + \xi_2(z) \right) \Psi_+^{(n,m)}(z) = 0,$$

which is exactly the second-order differential equation obtained from the Riccati equation.

One may also use the null-vector appearing at level 2

$$\chi_+(z) = \left[\hat{L}_{-2}(z) - \frac{3}{2(2\Delta_+ + 1)} \hat{L}_{-1}^2(z) \right] V_+(z).$$

The null vector vanishes when inner product is evaluated with any state. Therefore, one has the null constraint $\langle I^{(n)} | \chi_+(z) | I^{(m)} \rangle = 0$ which is represented in terms of the second-order differential equation with respect to z . It is shown in [23, 31] that the equation is the Mathieu equation for the Whittaker case. If $m = n$, we have Generalized Mathieu equation on a circle with $z = e^{2ix}$ when the wave function $\Psi_+^{(n,m)}(z)$ is properly rescaled [36].

As a side remark, Liouville field theory has an equivalent screening operator

$$\tilde{O} = \oint dz e^{\phi/b}(z)$$

due to the dual symmetry under $b \rightarrow 1/b$. Therefore, one can equivalently use either O or \tilde{O} . We can find the effect of inserting a new screening operator \tilde{O} by evaluating the expectation value with respect to the partition function

$$\tilde{\Psi}^{(n,m)}(z) = \frac{\langle I^{(n)} | e^{\phi/b}(z) | I^{(m)} \rangle_{\tilde{N}}}{\langle I^{(n)} | I^{(m)} \rangle_N},$$

where $\tilde{N} = N + 1/b^2$ and at NS limit, one has $\tilde{N} = N$. Noting that $\tilde{\Psi}^{(n,m)}(z) = 1/(\Psi_+^{(n,m)}(z))^2$, one has

$$\langle \tilde{O} \rangle = \oint \frac{dz}{P^2(z)} e^{-\frac{2}{\epsilon} V^{(n,m)}(z)} = \sum_k \frac{2\pi i}{(dP/dz)^2} \frac{de^{-\frac{2}{\epsilon} V^{(n,m)}}}{dz} \Big|_{z=z_k},$$

where z_k are the zeros of $P(z)$. One may also find that the expectation value of higher level degenerate operators is given as the higher powers of $\Psi_+^{(n,m)}(z)$'s. However, simultaneous use of the two screening operators (or infinite number of \tilde{O} 's) will break the conformal symmetry as seen in the sine-Gordon case, which needs further investigation.

3.3 Filling fraction and branch cuts

The deformed spectral curve is written as the second-order differential equation (3.11) of a monic polynomial $P(z)$. For the type of potential $V^{(0,m)}$ one may multiply (3.11) by z^{m+1} to find $N + m - 1$ order of polynomial equation which provides $N + m$ independent relations. Noting the number of unknowns are $N + m$ (N number of pole positions and m number of d_a 's), the polynomial equations determine the unknowns completely. For the case $V^{(n,0)}$, the same conclusion arises if one uses the dual potential $V^{(0,n)}$ after the conformal transformation. For $V^{(n,m)}$, we have $N + m + n - 1$ order of polynomial equation and end up with $N + m + n$ independent relations. This is consistent with the fact that there are $N + m + n$ unknowns (N number of pole positions and m number of $d_{a \geq 0}$ and n number of $d_{a < 0}$). Therefore, the differential equation (3.11) completely fixes $d_{a \geq 0}$ and $\hat{d}_{a \geq 0}$.

Once the solution is known, the partition function can be found through the flow equation (3.12). On the other hand, the branch cut structure is absent in the spectral curve. This raises a question. In the regular case, the integration contour of the partition function (2.15) is defined so that the integration contour includes the branch cut. The partition function is fixed according to the distribution of integration contours around the branch cuts. On the other hand, how can one define the contour integral of the partition function if the branch cut disappears?

The hint lies on the degrees of freedom in the solution space. To see this, let us consider the case $n = 0$, $m = 2$ and $N = 1$. We have 3 unknowns, P_0 , d_0 and d_1 . d_0 is trivially given: $d_0 = 2\epsilon c_0$. However, $d_1 = 2\epsilon c_1 - d_0 P_0$ is fixed by a quadratic equation

$$d_1^2 - 2\epsilon c_1 d_1 + 2\epsilon c_2 d_0 = 0.$$

Therefore, d_1 has two solutions: $d_1^\pm = \epsilon c_1 (1 \pm \sqrt{1 - \eta})$ where $\eta = 4c_2 c_0 / c_1^2$. Each solution corresponds to $P_0^- \sim c_1 / c_0$ and $P_0^+ \sim c_2 / c_1$ which shows that the root lies near one of two stationary points of the potential. In conclusion, zeros of the polynomial (or the poles of the resolvent) may distribute differently around the different stationary points of the potential. Accordingly, the different solution results in the different partition function.

As N increases, the solution space of (3.11) rapidly becomes very complicated. Suppose we consider the potential $V^{(0,m)}$. One can easily convince that the coefficient P_{N-1} in (3.8) has $N + 1$ solutions. (For the case m with N zeros, one has $\frac{(N+m-1)!}{N!(m-1)!}$ solutions. One may view this solutions as the zero distribution with $N = \sum_{i=1}^m N_i$. N_i is the number of zeros around each stationary point of the potential.) Therefore, we need more efficient way to find d_a 's and \hat{d}_a 's.

Alternative approach is to use the filling fraction from the beginning. One can find the number N_a of inserted screening operators using the one point resolvent $W(z)$ in (3.2) or (3.10)

$$\epsilon N_a = \oint_{\mathcal{A}_a} \frac{dz}{2\pi i} 2W(z)$$

if the integration contour \mathcal{A}_a locates around the saddle point of the potential. Therefore, the integration contours in the partition function (2.15) are chosen among the (A-cycle) contour loops \mathcal{A}_a . This suggests that one can find d_a in perturbative power series of ϵ . Regarding $V(z) = O(1)$, we have $W(z) = O(\epsilon)$ and $f(z) = O(\epsilon)$ and the dominant contribution is found in the equation

$$4W(z)V'(z) \sim f(z).$$

This is consistent with the expectation that poles of $W(z)$ (zeros of the polynomial $P(z)$) (3.9) are accumulated around the stationary points of the potential $V(z)$. Therefore, we can put $W(z) = \sum_{k \geq 1} \epsilon^k W^{(k)}(z)$ and $f(z) = \sum_{k \geq 1} \epsilon^k f^{(k)}(z)$ and apply the ϵ expansion to the loop equation (3.3) directly.

According to the perturbation, the leading order contribution $f^{(1)}(z)$ is related with the filling fraction N_k

$$N_a = \oint_{\mathcal{A}_a} \frac{f^{(1)}}{2V'} dz, \quad (3.13)$$

where \mathcal{A}_a is the contour encircling around the stationary point ξ_a of the potential. Note that the maximum number of stationary point is the same as that of d_a . Therefore, N_a 's in (3.13) fix d_a 's to the order of ϵ . The sub-dominant terms of the filling fraction should vanish, which leads to the null identity. For example, at order of ϵ^2 one has

$$0 = \oint_{\mathcal{A}_a} dz \left\{ \frac{f^{(2)}}{2V'} - \frac{(f^{(1)})'}{4(V')^2} + \frac{(2V'' - f^{(1)})f^{(1)}}{8(V')^3} \right\}.$$

The flow equation (3.12) shows that to the lowest order in ϵ

$$d(F_N^{(n,m)}) = 2\epsilon \sum_{k,a} dc_k U_{ka} \sum_{\ell \geq 1} c_{a+\ell} \left(\sum_{a=1}^{n+m} \frac{N_a}{\xi_a^\ell} + O(\epsilon) \right).$$

Note that $\xi_a(\{c_k\}, \{\hat{c}_k\}) = 1/\hat{\xi}_a(\{\hat{c}_k\}, \{c_k\})$. Therefore,

$$\hat{d}(F_N^{(n,m)}) = 2\epsilon \sum_{k,a} d\hat{c}_k \hat{U}_{ka} \sum_{\ell \geq 1} \hat{c}_{a+\ell} \left(\sum_{a=1}^{n+m} N_a \xi_a^\ell + O(\epsilon) \right).$$

If one identifies the spectral curve with the Seiberg–Witten curve, then, $x(z)dz$ is the Seiberg–Witten one-form λ and the filling fraction of the deformed spectral curve corresponds to the Coulomb branch parameter a_k ,

$$a_a = \epsilon N_a + \epsilon/2.$$

According to the identification, the partition function has the relation

$$\frac{\partial d(F^{(n,m)})}{\partial a_a} = 2 \sum_{k,a} dc_k U_{ka} \sum_{\ell \geq 1} c_{a+\ell} \left(\frac{1}{\xi^\ell} + O(\epsilon) \right),$$

where $O(\epsilon)$ corresponds to $O(a_a)$.

4 Extension of symmetry

4.1 W -symmetry

$W^{(s+1)}$ -symmetry can be easily incorporated if one employs s -number of fields. The irregular state is generated by IVO which contains s bosonic fields and its finite number of derivatives. Using the bosonic free field with holomorphic normalization

$$\langle \phi^{(a)}(z) \phi^{(b)}(w) \rangle = -\delta^{ab} \log(z-w)$$

one has IVO [32]

$$I^{(s|m)}(w) = \exp \left\{ \sum_{a=1}^s \sum_{k=0}^m \frac{c_k^{(a)}}{\hbar k!} \partial_w^k \phi^{(a)}(w) \right\}, \quad |I^{(s|m)}\rangle = \lim_{w \rightarrow 0} I^{(s|m)}(w)|0\rangle.$$

Likewise, its adjoint is defined at $1/\zeta$

$$\hat{I}^{(s|n)}(\zeta) = R_0^{(s|n)} \exp \left\{ \sum_{a=1}^s \sum_{\ell=0}^n \frac{\hat{c}_\ell^{(a)}}{\hbar \ell!} \partial_\zeta^\ell \phi^{(a)}(1/\zeta) \right\},$$

where ∂_ζ is the derivative with respect to ζ and $R^{(s|n)}$ is the scale factor

$$R_0^{(s|n)} = \exp \left\{ \sum_{a=1}^s \sum_{k=0}^n \frac{\hat{c}_0^{(a)} \hat{c}_k^{(a)}}{\ell!} \partial_\zeta^\ell \log(\zeta) \right\}.$$

The adjoint state is defined as

$$\langle I^{(s|n)} | = \lim_{\zeta \rightarrow 0} \langle 0 | \hat{I}^{(s|n)}(\zeta).$$

Their inner product is defined as in Section 2

$$\langle I_0^{(s|n)} | I_0^{(s|m)} \rangle = \lim_{w, \zeta \rightarrow 0} \langle 0 | \hat{I}_0^{(s|n)}(\zeta) I_0^{(s|m)}(w) | 0 \rangle = e^{\zeta_{n,m}^{(s)}/\hbar^2},$$

where $\zeta_{n,m}^{(s)} = \sum_{\ell \geq 1}^{\min(m_a, n_a)} \frac{(\hat{\mathbf{c}}_\ell, \mathbf{c}_\ell)}{\ell}$ and $\hat{\mathbf{c}}_0 + \mathbf{c}_0 = 0$ is assumed. We use bold letters for vectors,

$\mathbf{c}_\ell = (c_\ell^{(1)}, \dots, c_\ell^{(s)})$ and $(\hat{\mathbf{c}}_\ell, \mathbf{c}_\ell)$ represents the inner product between two vectors.

Beyond the free theory, we introduce s -kind of screening operators $V_{b\mathbf{e}_k}$ ($k = 1, \dots, s$)

$$V_{b\mathbf{e}_k} = e^{b(\mathbf{e}_k, \Phi)}.$$

Φ is the scalar fields $\Phi = (\phi^{(1)}, \dots, \phi^{(s)})$ with s components and s -dimensional vector \mathbf{e}_k are the simple roots of Lie algebra. The bosonic fields with screening operators are represented as Toda field theory. (\mathbf{e}_k, Φ) denotes the scalar product. The scalar product $(\mathbf{e}_i, \mathbf{e}_j) = K_{ij}$ is the Cartan matrix (for A_s Lie algebra, $K_{ii} = 2$, $K_{ii+1} = -1$ and other components are zero). Vertex operator $V_\alpha = e^{(\alpha_i, \Phi)}$ has the holomorphic dimension $\Delta_\alpha = (\alpha, (\hat{Q} - \frac{1}{2}\alpha))$, where \hat{Q} is the background charge vector $\hat{Q} = Q\rho$ and ρ is the Weyl vector (half of the sum of all positive roots). Therefore, $V_{\mathbf{e}_k}$ has the holomorphic dimension $\Delta_{\mathbf{e}_k} = 1$.

Using the screening operators one constructs the inner product of the irregular states in terms of irregular matrix model with s -set of variables

$$\langle I^{(s|n)} | I^{(s|m)} \rangle_{\mathbf{N}} = e^{\zeta_{n,m}} Z_N^{(s|n,m)},$$

where $e^{\zeta_{n,m}}$ is the inner-product due to free field contribution and $Z_N^{(s|n,m)}$ is the screening operator contribution

$$Z_N^{(s|n,m)} = \left\{ \prod_{a=1}^s \prod_{I=1}^{N^{(a)}} \int d\lambda_I^{(a)} \right\} \prod_{a \geq b} (\Delta_{ab})^{\beta K_{ab}} e^{\frac{\sqrt{\beta}}{g} V^{(s|n,m)}},$$

where Δ_{ab} is the Vandermonde determinant

$$\Delta_{ab} = \delta_{a,b} \prod_{I < J} (\lambda_I^{(a)} - \lambda_J^{(a)}) + (1 - \delta_{a,b}) \prod_{I,J} (\lambda_I^{(a)} - \lambda_J^{(b)}),$$

and $V^{(s|n,m)}$ is the potential $V^{(s|n,m)} = \sum_a V^{(s|n,m)}(\lambda^{(a)})$, where

$$V^{(s|n,m)}(\lambda^{(a)}) = \sum_{I=1}^{N^{(a)}} \left\{ c_0^{(a)} \log(\lambda_I^{(a)}) - \sum_{k=1}^{m^{(a)}} \frac{c_k^{(a)}}{(\lambda_I^{(a)})^k} - \sum_{\ell=1}^{n^{(a)}} \frac{\hat{c}_\ell^{(a)}}{\ell} (\lambda_I^{(a)})^\ell \right\}.$$

Neutrality condition $\hat{\mathbf{c}}_0 + \mathbf{c}_0 + \epsilon \mathbf{N} = \epsilon \rho$ insures that conformal invariance of the system and symmetry of the inner product. The scaling factor in the adjoint definition is also modified due to the background charge

$$R_\epsilon^{(s|n)} = \exp \left\{ \sum_{a=1}^s \sum_{k=0}^n \frac{(\hat{c}_0^{(a)} - \epsilon \rho) \hat{c}_k^{(a)}}{\hbar^2 k!} \partial_\zeta^k \log \zeta \right\}.$$

In general, one has $(s+1)$ -th order spectral curves for the A_s type potential. For example, for A_2 case, one has cubic spectral curves [10, 12, 38] if one uses the conformal transformation $\lambda_I^{(1)} \rightarrow \lambda_I^{(1)} + \sum_J \frac{\delta}{(\lambda_I^{(1)} - z)(\lambda_I^{(1)} - \lambda_J^{(2)})}$ and $\lambda_J^{(2)} \rightarrow \lambda_J^{(2)} + \sum_I \frac{\delta}{(\lambda_J^{(2)} - z)(\lambda_I^{(1)} - \lambda_J^{(2)})}$. Its classical/NS limit is conveniently written in two symmetric forms

$$\begin{aligned} X_1^3 + \xi_2 X_1 + 3\epsilon X_1 X_1' + \epsilon^2 X_1'' &= +\frac{2}{3\sqrt{3}} \xi_3 - \frac{\epsilon}{2} \xi_2', \\ X_2^3 + \xi_2 X_2 + 3\epsilon X_2 X_2' + \epsilon^2 X_2'' &= -\frac{2}{3\sqrt{3}} \xi_3 - \frac{\epsilon}{2} \xi_2', \end{aligned} \quad (4.1)$$

where X_1 and X_2 are one-point resolvents (R_1 and R_2) shifted by potential

$$X_1 = 2 \left(R_1 + \frac{1}{3} (2V_1' + V_2') \right), \quad X_2 = 2 \left(R_2 + \frac{1}{3} (V_1' + 2V_2') \right),$$

where we use the abbreviation $V_a = V^{(\mathbf{n}, \mathbf{m})}(\lambda^{(a)})$.

There exists also the quadratic form of the spectral curve

$$X_1^2 + X_2^2 - X_1 X_2 + \epsilon(X_1' + X_2') + \xi_2 = 0, \quad (4.2)$$

which presents the Virasoro symmetry. In fact, (4.1) and (4.2) are not independent each other. Only two of the three are independent. $\xi_2(z)$ is the energy momentum tensor (Virasoro current) expectation value

$$\xi_2(z) = \sum_{k=-2m}^{2n} \frac{A_k}{z^{k+2}} - \sum_{k=-m}^{n-1} \frac{d_k}{z^{k+2}} = \frac{\langle I_m | \hbar^2 T(z) | I_n \rangle}{\langle I_m | I_n \rangle}.$$

Here A_k is a constant obtained from the potential with $\mathbf{c}^{(1)} = \mathbf{a}$ and $\mathbf{c}^{(2)} = \mathbf{b}$

$$A_k = 2\epsilon(k+1)(a_k + b_k) - \frac{4}{3} \sum_{r+s=k} (a_r a_s + b_r b_s + a_r b_s).$$

The mode d_k ($0 \leq k \leq m-1$) is related with the partition function

$$d_k = v_k(F_N^{(m;n)}), \quad v_k = \sum_{s>0} s \left(a_{s+k} \frac{\partial}{\partial a_s} + b_{s+k} \frac{\partial}{\partial b_s} \right)$$

and its dual form \hat{d}_k ($0 \leq k \leq n-1$) if one replaces (\mathbf{a}, \mathbf{b}) with $(\hat{\mathbf{a}}, \hat{\mathbf{b}})$.

The expectation of the \mathcal{W}_3 current $W(z)$ is given as $\xi_3(z)$

$$\xi_3(z) = \frac{\langle I^{(n)} | \hbar^2 W(z) | I^{(m)} \rangle}{\langle I_m | I_n \rangle} = \sum_{k=-3m}^{3n} \frac{B_k}{z^{k+3}} - \sum_{k=-2m}^{2n-1} \frac{e_k}{z^{k+3}},$$

where B_k comes directly from the potential coefficients

$$\begin{aligned} B_k &= \frac{4}{3\sqrt{3}} \sum_{r+s+t=k} (2(a_r a_s a_t - b_r b_s b_t) + 3(a_r a_s b_t - b_r b_s a_t)) \\ &\quad - \frac{\sqrt{3}}{2} \epsilon \sum_{r+s=k} (2(k+2)(a_r a_s - b_r b_s) + (r-s)(a_r b_s - b_r a_s)) \\ &\quad + \frac{\sqrt{3}}{2} \epsilon^2 (k+1)(k+2)(a_k - b_k). \end{aligned}$$

The moment e_k induces the flow equation

$$e_k = \mu_k(F_N^{(n,m)}),$$

where μ_k is the W -current. To find the partition function we need the mode with $n \leq k \leq 2n-1$ which is

$$\mu_k = \sum_{\substack{k=r+s-t; \\ t>0}} \sqrt{3}t \left(-(a_r a_s + 2a_r b_s) \frac{\partial}{\partial a_t} + (b_r b_s + 2a_r b_s) \frac{\partial}{\partial b_t} \right).$$

The flow equations show the consistent conditions

$$v_p(d_q) - v_q(d_p) = (p-q)d_{p+q}, \quad v_p(e_q) - \mu_q(d_p) = (2p-q)\mu_{p+q}$$

and its duals, replacing (\mathbf{a}, \mathbf{b}) with $(\hat{\mathbf{a}}, \hat{\mathbf{b}})$, realize the \mathcal{W}_3 symmetry.

It is noted in [12] that the spectral curve (4.1) can be put into a third-order differential equation of $\Psi_i(z)$, where $\Psi_i(z) = \exp\left(\frac{1}{\epsilon} \int^z X_i(z') dz'\right)$ with $i = 1, 2$:

$$\left(\epsilon^3 \frac{\partial^3}{\partial z^3} + \xi_2 \epsilon \frac{\partial}{\partial z} + U_i(z) \right) \Psi_i(z) = 0,$$

where $U_1(z) = +\frac{2}{3\sqrt{3}}\xi_3 - \frac{\epsilon}{2}\xi_2'$ and $U_2(z) = -\frac{2}{3\sqrt{3}}\xi_3 - \frac{\epsilon}{2}\xi_2'$. Note that $\Psi_i(z)$ corresponds to the normalized expectation value $\langle I^{(n)} | V_{-\mathbf{w}_i/b}(z) | I^{(m)} \rangle / \langle I^{(n)} | I^{(m)} \rangle$, where $\mathbf{w}_1 = (2\mathbf{e}_1 + \mathbf{e}_2)/3$ and $\mathbf{w}_2 = (2\mathbf{e}_2 + \mathbf{e}_1)/3$, $((\mathbf{e}_j, \mathbf{w}_k) = \delta_{jk})$. Therefore, the expectation value of $V_{++}(z) = V_{-(\mathbf{w}_1+\mathbf{w}_2)/b}(z)$ is given as

$$\frac{\langle I^{(n)} | V_{++}(z) | I^{(m)} \rangle}{\langle I^{(n)} | I^{(m)} \rangle} = \Psi_1(z) \Psi_2(z).$$

One may also put the spectral curves (4.1) into two coupled third-order differential equations of two monic polynomials $P(z)$ and $Q(z)$ of the degree N : $P(z) = \langle \prod_{i=1}^N (z - x_i) \rangle$ and $Q(z) = \langle \prod_{j=1}^M (z - y_j) \rangle$. (4.2) becomes a second-order differential equation of two polynomials. Putting $2R_1(z) = \epsilon P'(z)/P(z)$ and $2R_2(z) = \epsilon Q'(z)/Q(z)$, one has for (4.2)

$$\epsilon^2 (P''Q - P'Q' + PQ'') + 2\epsilon (V_1'P'Q + V_2'PQ') = FPQ$$

and for (4.1)

$$\begin{aligned} \epsilon^3 P''' + 2\epsilon^2 (2V_1' + V_2')P'' + \epsilon (4V_1'(V_1' + V_2') + 2\epsilon V_1'' - F)P' &= G_1P, \\ \epsilon^3 Q''' + 2\epsilon^2 (V_1' + 2V_2')Q'' + \epsilon (4V_2'(V_1' + V_2') + 2\epsilon V_2'' - F)Q' &= G_2Q. \end{aligned}$$

G_1 and G_2 are given in terms of e_k and d_k with irregular poles and zeros

$$\begin{aligned} G_1 &= \sum_{k=-2n}^{2m-1} \frac{1}{z^{k+3}} \left\{ -\frac{2}{3\sqrt{3}}e_k + \frac{2}{3} \sum_{r+s=k} d_r(2b_s + a_s) \right\} - \frac{\epsilon}{2} \sum_{k=-n}^{m-1} \frac{(k+2)d_k}{z^{k+3}}, \\ G_2 &= \sum_{k=-2n}^{2m-1} \frac{1}{z^{k+3}} \left\{ \frac{2}{3\sqrt{3}}e_k + \frac{2}{3} \sum_{r+s=k} d_r(2a_s + b_s) \right\} - \frac{\epsilon}{2} \sum_{k=-n}^{m-1} \frac{(k+2)d_k}{z^{k+3}}. \end{aligned}$$

4.2 Supersymmetry

One may have different type of irregular states if one adopts supersymmetric theory. Considering that $N = 1$ super Liouville conformal field theory is related with the instanton partition function of $N = 2$ quiver gauge theories on the ALE space $\mathcal{C}^2/\mathcal{Z}_2$ [3, 5], we expect that the supersymmetric irregular matrix model will provide the useful information on the Argyres–Douglas theory.

The supersymmetric irregular vertex operator is constructed in [33, 34]. We present the operator in the superfield formalism. The super vertex operator $V_\alpha(\omega) = e^{\alpha\Phi}(\omega)$ has holomorphic normalization [35]

$$\langle V_{\alpha_1}(\omega_1) V_{\alpha_2}(\omega_2) \rangle = (w_1 - w_2 - \theta_1\theta_2)^{-\alpha_1\alpha_2},$$

where Φ is the superfield and $\omega = (w, \theta)$ is the holomorphic super-coordinate.

We consider super irregular operator $W^{(q)}(\omega)$ of rank q

$$W^{(q)}(\omega) = \exp \left(\sum_{k=0}^{2q} \frac{\gamma_k}{\hbar} D_\omega^k \Phi(w, \theta) \right),$$

where $D_\omega = \theta \partial_w + \partial_\theta$ is the super-derivative. γ_k is a commuting (anti-commuting) number when k is even (odd). The irregular state lies in NS sector if p is an integer and in R -sector if p is a half-odd integer. The adjoint operator is defined as

$$\hat{W}^{(p)}(\hat{\omega}) = R^{(p)} \exp \left(\sum_{k=0}^{2p} \frac{\hat{\gamma}_k}{\hbar} D_\zeta^k \Phi(1/\zeta, \hat{\theta}) \right),$$

where $\hat{\omega}$ is another super coordinate $(\zeta, \hat{\theta})$ and $R^{(p)}$ is the scale factor

$$R^{(p)} = \exp \left(\sum_{\ell=0}^{2p} \frac{(\hat{\gamma}_0 - \epsilon) \hat{\gamma}_\ell}{\hbar^2} D_\zeta^\ell \log(\zeta) \right).$$

Using the irregular operator we define the super irregular state and its adjoint

$$|W^{(q)}\rangle = \lim_{w \rightarrow 0} W^{(q)}(w, \theta)|0\rangle, \quad \langle W^{(p)}| = \lim_{\zeta \rightarrow 0} \langle 0|\hat{W}^{(p)}(\zeta, \hat{\theta}).$$

Inner product has the explicit form

$$\begin{aligned} \langle W^{(p)}|W^{(q)}\rangle &= \lim_{\zeta, w \rightarrow 0} \langle 0|\hat{W}^{(p)}(\zeta)W^{(q)}(w)|0\rangle = \exp \left(\frac{K^{(p;q)}}{\hbar^2} \right), \\ K^{(p;q)} &= \sum_{k \geq 1} \left(\frac{\hat{c}_k c_k}{k} + \hat{\xi}_k \xi_{k-1} \right), \end{aligned}$$

where $\hat{\gamma}_0 + \gamma_0 = 0$ and no background charge ($\epsilon = 0$) is assumed. $c_k = (\gamma_{2k} + \gamma_{2k-1}\theta)k!$ and $\hat{c}_k = (\hat{\gamma}_{2k} + \hat{\gamma}_{2k-1}\hat{\theta})k!$ are commuting. $\xi_k = (\gamma_{2k}\theta + \gamma_{2k+1})k!$ and $\hat{\xi}_k = (\hat{\gamma}_{2k}\hat{\theta} + \hat{\gamma}_{2k+1})k!$ are anti-commuting. Here we use the super derivative identities $Dw^l = l(\theta w^{l-1})$, $D(\theta w^l) = (w^l)$, $D^2 w^l = l(w^{l-1})$, and $D^2(\theta w^l) = l(w^{l-1}\theta)$, $D^3(\theta w^l) = l(w^{l-1})$, $D^3(w^l) = l(l-1)(\theta w^{l-2})$. These identities provide non-vanishing quantities as $w \rightarrow 0$: $D^{2n} w^n = n!$, $D^{2n}(\theta w^n) = n!\theta$, $D^{2n-1}(w^n) = n!\theta$ and $D^{2n+1}(\theta w^n) = n!$.

In the presence of background charge we need screening operators $\oint d\zeta V_b(\zeta)$ so that neutrality condition $\hat{\gamma}_0 + \gamma_0 + \hbar N b = \epsilon$ holds. Then, the inner product has the form

$$\begin{aligned} \langle W^{(p)}|W^{(q)}\rangle_N &= \langle W_0^{(p)}| \left(\oint d\zeta V_b(\zeta) \right)^N |W_0^{(q)}\rangle = e^{K^{(p;q)}/\hbar^2} Y_N^{(p;q)}, \\ Y_N^{(p;q)} &= \int \left[\prod_{I=1}^N dz_I d\theta_I \right] \prod_{I < J} (z_{IJ} - \theta_I \theta_J)^\beta e^{\frac{\sqrt{\beta}}{g_s} \sum_I V^{(p;q)}(\zeta_I)}. \end{aligned} \quad (4.3)$$

Here we use the notation $g_s = i\hbar$. The partition function $Y_N^{(p;q)}$ will be called the irregular super matrix model. The potential $V(\zeta_I)$ of the form $V(\zeta_I) = V_B(z_I) + \theta_I V_F(z_I)$: $V_B(z)$ and $V_F(z)$ are bosonic and fermionic part of super-potential

$$V_B(z_I) = c_0 \ln(z_I) - \sum_{k \geq 1}^q \frac{c_k}{k z_I^k} - \sum_{k \geq 1}^p \frac{\hat{c}_k z_I^k}{k}, \quad V_F(z_I) = - \sum_{k \geq 0}^q \frac{\xi_k}{z_I^{k+1}} + \sum_{k \geq 1}^p \hat{\xi}_k z_I^{k-1},$$

where $c_0 = \gamma_0$ is used. The potential can be put into more symmetric way if we put $\hat{c}_k = -c_{-k}$ and $\hat{\xi}_k = -\xi_{-k}$,

$$V_B(z_I) = c_0 \ln(z_I) - \sum'_{k=-p}^q \frac{c_k}{k z_I^k}, \quad V_F(z_I) = - \sum'_{k=-p}^q \frac{\xi_k}{z_I^{k+1}}, \quad (4.4)$$

where the prime in the summation denotes that no sum on $k = 0$. It is noted that the super irregular matrix model (4.3) is also obtained from the colliding limit of the superconformal system with central charge $c = 3/2(1 + 2Q^2)$.

The loop equation becomes the super-spectral curve [13, 24] for the super-conformal transformation $z_I \rightarrow z_I + \frac{\delta}{z-z_I}$ and $\theta_I \rightarrow \theta_I(1 + \frac{\delta}{2(z-z_I)^2})$:

$$\begin{aligned} x_B(z)x_F(z) + \epsilon x'_F(z) &= F_F(z), \\ x_B(z)^2 + \epsilon x'_B(z) + x_F(z)V'_F(z) - x'_F(z)V_F(z) &= 2F_B(z), \end{aligned} \quad (4.5)$$

where $x_F(z)$ ($x_B(z)$) is anti-commuting (commuting) one-point resolvent $\omega_F(z)$ ($\omega_B(z)$) shifted by potential term, $x_F(z) = \omega_F(z) - V_F(z)$ ($x_B(z) = \omega_B(z) + V'_B(z)$)

$$\omega_B(z) = \epsilon \left\langle \sum_I \frac{1}{z - z_I} \right\rangle, \quad \omega_F(z) = \epsilon \left\langle \sum_I \frac{\theta_I}{z - z_I} \right\rangle.$$

F_F (F_B) is also anti-commuting (commuting) holomorphic function

$$F_F(z) = f_F(z) - V'_B(z)V_F(z) - \epsilon V'_F(z), \quad F_B(z) = f_B(z) + \frac{1}{2}V_B'^2 + \epsilon V'_B(z)$$

and represent spin 3/2 supercurrent (spin 2 Virasoro) symmetry of the partition function

$$\begin{aligned} f_F(z) &\equiv \epsilon \left\langle \frac{(V'_B(z) - V'_B(z_I))\theta_I - (V_F(z) - V_F(z_I))}{z - z_I} \right\rangle, \\ f_B(z) &= \epsilon \left\langle \sum_I \frac{(V'_B(z) - V'_B(z_I)) + \theta_I(V'_F(z) - V'_F(z_I))}{z - z_I} + \frac{1}{2} \frac{\theta_I(V_F(z) - V_F(z_I))}{(z - z_I)^2} \right\rangle. \end{aligned}$$

Suppose the potential is given in (4.4). Then, we have F_F (F_B) of the form

$$F_F(z) = \sum_{r=1/2-2p}^{2q-1/2} \frac{\Omega_r}{z^{3/2+r}} + \sum_{r=1/2-p}^{q-1/2} \frac{\eta_r}{z^{3/2+r}}, \quad F_B(z) = \sum_{t=-2p}^{2q} \frac{\Lambda_t}{z^{2+t}} + \sum_{t=1-p}^{q-1} \frac{d_t}{z^{2+t}}, \quad (4.6)$$

where $\Omega_r = \sum_{k+\ell=r-1/2} c_k \xi_\ell + \epsilon(r+1/2)\xi_{r-1/2}$ is anti-commuting and $\Omega_{r \geq q+1/2}$ are the eigenvalues of spin 1/2 positive modes. $\Lambda_t = \sum_{k+\ell=t} c_k c_\ell / 2 - \epsilon(l+1)c_t / 2$ is commuting and $\Lambda_{t \geq q}$ are the eigenvalues of the Virasoro positive modes.

The highest mode eigenvalue Λ_{2q} and $\Omega_{2q-1/2}$ are not vanishing for NS sector. This is because the coefficients in the potential can be made of even mode γ_k (k even) only. The commuting number $c_q = \gamma_{2q}\theta q!$ contains the commuting mode γ_{2q} . The anti-commuting mode $\xi_{q-1} = \gamma_{2q-2}\theta(q-1)!$ also contains the commuting mode γ_{2q-2} . As a result, $\Lambda_{2q} = c_q^2/2$ and $\Omega_{2q-1/2} = c_m \xi_{q-1}$ are not vanishing.

However, this is not the case for the Ramond sector where odd modes $\gamma_{k=\text{odd}}$ are essential ingredients. Especially, $c_m = \gamma_{2q-1}\theta q!$ and $\xi_{q-1} = (\gamma_{2q-1})(q-1)!$ contain the odd mode γ_{2q-1} only. As a result the highest positive mode eigenvalues $\Lambda_{2q} = c_q^2/2$ and $\Omega_{2q-1/2} = c_q \xi_{q-1}$ should vanish. This is the crucial difference between the NS sector and R sector.

η_r and d_t are expectation values of supercurrent representations g_r and ℓ_t (corresponding to right action). η_r is anti-commuting

$$\eta_r = g_r(-\hbar^2 \ln Y), \quad g_r = \sum_{\ell-k=r} \left(k \xi_{\ell-1/2} \frac{\partial}{\partial c_k} - c_\ell \frac{\partial}{\partial \xi_{k-1/2}} \right). \quad (4.7)$$

The relation is obtained if one notices that

$$\frac{\sqrt{\beta}}{g} \left\langle \frac{1}{z_I^{k+1}} \right\rangle = k \frac{\partial}{\partial c_k} \ln Y, \quad \frac{\sqrt{\beta}}{g} \left\langle \frac{\theta_I}{z_I^k} \right\rangle = \frac{\partial}{\partial \xi_k} \ln Y.$$

d_t is commuting

$$d_t = \ell_t(-\hbar^2 \ln Y), \quad \ell_t = \sum_{\ell-k=t} \left(k c_\ell \frac{\partial}{\partial c_k} + \left(\frac{k+\ell}{2} \right) \xi_{\ell-1/2} \frac{\partial}{\partial \xi_{k-1/2}} \right). \quad (4.8)$$

The super-flow equation (4.7) and (4.8) provides a way to find the partition function directly from the spectral curve (4.5). The currents g_r in (4.7) and ℓ_ℓ in (4.8) satisfy the commutation relation of right action of the super algebra

$$[l_\ell, g_r] = \left(r - \frac{\ell}{2} \right) g_{r+\ell}, \quad \{g_r, g_s\} = -2l_{r+s}, \quad [l_k, l_\ell] = -(k-\ell)l_{k+\ell}.$$

This commutation relation provide the consistency condition for the partition function. One may better the derivatives $\partial/\partial c_k$ and $\partial/\partial \xi_{\ell-1/2}$ in terms of ℓ_k and g_r . Denoting the derivative elements as $\partial = (\{\partial_m\}) = (\{\frac{\partial}{\partial c_k}\}, \{\frac{\partial}{\partial \xi_{\ell-1/2}}\})$ and $\zeta = (\{\zeta_m\}) = (\{\ell_k\}, \{g_r\})$ one has $\partial = \mathcal{U}\zeta$, where \mathcal{U} is a $2q \times 2q$ super-matrix. The partition function is given as

$$d(-\hbar^2 \log Y) = d\zeta \mathcal{U} \tau,$$

where $\tau = (\{\tau_n\}) = (\{d_k\}, \{\eta_\ell\})$ in (4.6).

If one considers the degenerate primary superfield $V_{-1/b}$ which has the null state at level $3/2$, then the normalized expectation value $\Psi^{(p,q)}(z, \theta)$ of $V_{-1/b}$ is written as following

$$\begin{aligned} \Psi^{(p,q)}(z, \theta) &= \exp\left(\frac{V^{(p,q)}}{\epsilon}\right) \left\langle \prod (z - \lambda_I - \theta \theta_I) \right\rangle = \Psi_B^{(p,q)} \Psi_F^{(p,q)}, \\ \Psi_B^{(p,q)}(z, \theta) &= \exp\left(\frac{1}{\epsilon} \int^z x_B(z') dz'\right) = P(z) \exp\left(\frac{V^{(p,q)}}{\epsilon}\right), \\ \Psi_F^{(p,q)}(z, \theta) &= \exp\left(-\frac{\theta}{\epsilon} x_F(z)\right), \end{aligned}$$

where $P(z) = \langle \prod (z - \lambda_I) \rangle$ is the N -th order monic polynomial. As in the Virasoro case, the bosonic resolvent is given in terms of the polynomial at the classical/NS limit, $\omega_B(z) = \epsilon P'(z)/P(z)$. The super spectral curve (4.5) reduces to the second-order differential equation of the polynomial $P(z)$ and therefore, the bosonic contribution $\Psi_B^{(p,q)}(z, \theta)$ is given in terms of the polynomial solution. Then, the super-spectral curve (4.5) is rewritten as

$$\begin{aligned} \epsilon \frac{\partial}{\partial z} (\Psi_B^{(p,q)} x_F) - \Psi_B^{(p,q)} F_F &= 0, \\ \epsilon^2 \frac{\partial^2}{\partial z^2} \Psi_B^{(p,q)} + \left(x_F \frac{\partial V_F}{\partial z} - \frac{\partial x_F}{\partial z} V_F - 2F_B \right) \Psi_B^{(p,q)} &= 0. \end{aligned} \quad (4.9)$$

One thing to note is that the pole structure of bosonic resolvent ω_B shares with that of the fermionic one ω_F . This can be seen from (4.9). One has the solution of the form $\epsilon x_F(z) = \tau_F(z)/\Psi_B^{(p,q)}(z)$, where $\tau_F(z) = \int^z \Psi_B^{(p,q)}(z') F_F(z') dz'$. Since $\tau_F(z_\alpha)$ is not zero in general, the obvious conclusion is that the zero z_α of $P(z)$ in $\Psi_B^{(p,q)}(z)$ becomes the pole position of $\omega_F(z)$.

5 Conclusion

Free field formalism is a convenient way to construct the irregular states. We provide a few examples; Virasoro states using one boson field, W states using multiple boson fields, supersymmetric states with super-fields. Screening operators provide more degrees of freedom and the role of screening operators are given in terms of irregular matrix model. Therefore, the free field formalism together with the screening operators demonstrates that the irregular conformal block can be constructed without using the colliding limit.

The partition function is evaluated with the help of the (super) conformal symmetry. The loop equation, or (deformed) spectral curve contains the (super) conformal symmetry. This leads to the flow equations of the partition function on the parameter space of the irregular states. In evaluating the partition function, the analytic structure of the spectral curve plays the crucial role.

The spectral curve also provides a way of connecting the irregular matrix model with $N = 2$ super gauge theory. The spectral curve is the (deformed) Seiberg–Witten curve of the Hitchin system and the filling fraction (number of screening operator) is identified with the Coulomb branch parameter of the gauge theory. Therefore, the flow equation obtained from the spectral curve effectively determines the partition function of the Argyres–Douglas gauge theory.

The inner product is the two-point correlation. One may find multi-point correlations by considering the expectation value of product of regular and/or irregular vertex operators. The expectation value is the product of the free field contribution and irregular matrix model which is due to the screening operators insertion. The irregular matrix model is simply modified by adding the potential terms at each position of the inserted regular and/or irregular vertex operators. Therefore, the spectral curve looks the same, only changing the analytic property due to the conformal current for the same type of conformal symmetry: Virasoro symmetry has the quadratic form in x , and $W^{(s+1)}$ symmetry has the s -th powered form in x . The difference of the irregular singularity from the regular one lies in ξ_s 's which represents the conformal symmetry representation. Depending on the type of vertex operators one has the different type of analytic structure in ξ_s , regular or irregular poles and zeros. As a result, the flow equation changes accordingly and so does the partition function. Depending on the symmetry representation, one has different analytic structure of the partition function. The understanding of the mathematical structure will be a challenging problem.

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