Weingarten calculus on computers and its application to

random quantum Gaussian states.

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# 1 Weingarten Calculus on computers

"RTNI - A symbolic integrator for Haar-random tensor networks", arXiv:1902.08539 [quant-ph] with Nechita and Koenig.

# 1.1 Computer packages for average over Haar unitary matrices

#### Preceding packages:

- IntU, a Mathematica package, [Puchała and Miszczak (2017)].
- IntHaar, a Maple package, [Ginory and Kim (2016)]. (includes Haar orthogonal and symplectic cases.)

### Our package:

RTNI (Random Tensor Network Integrator), Mathematica and Python packages.

#### Differences:

- Others calculate averages of monomials in the entries of a random unitary matrices.
- Our package calculates averages symbolically, which allows tensor structures easily.

# 1.2 Polynomial version of Weingarten calculus

 $\mathcal{U}(n)$ : group of  $n \times n$  unitary matrices, and dU: normalized Haar measure.

Fix  $p \in \mathbb{N}$  and let  $i = (i_1, \ldots, i_p)$ ,  $i' = (i'_1, \ldots, i'_p)$ ,  $j = (j_1, \ldots, j_p)$ ,  $j' = (j'_1, \ldots, j'_p)$  be p-tuples of positive integers from  $\{1, 2, \ldots, n\}$ . [Collins and Sniady (2006)]

$$\int_{\mathcal{U}(n)} U_{i_1 j_1} \cdots U_{i_p j_p} \bar{U}_{i'_1 j'_1} \cdots \bar{U}_{i'_p j'_p} dU$$

$$= \sum_{\alpha, \beta \in S_n} \delta_{i_1 i'_{\alpha(1)}} \dots \delta_{i_p i'_{\alpha(p)}} \delta_{j_1 j'_{\beta(1)}} \dots \delta_{j_p j'_{\beta(p)}} \operatorname{Wg}_n(\alpha^{-1} \beta) .$$

Here,  $\mathcal{S}_p$  is the permutation group and  $Wg_n(\cdot)$  is called the *unitary Weingarten function*. Also, note that

$$\int_{\mathcal{U}(p)} U_{i_1 j_1} \cdots U_{i_p j_p} \bar{U}_{i'_1 j'_1} \cdots \bar{U}_{i'_p, j'_{p'}} dU = 0 \quad \text{for} \quad p \neq p'.$$

For integer  $n \geq p$ , the Weingarten functions can be written as

$$Wg_n(\sigma) = \frac{1}{(p!)^2} \sum_{\lambda \vdash n} \frac{(\chi^{\lambda}(e))^2}{s_{\lambda,n}(1)} \chi^{\lambda}(\sigma) .$$

- $\lambda \vdash p$  means that  $\lambda$  is a partition of the integer p.
- $\chi^{\lambda}$  is the character of the irreducible representation of the symmetric group  $S_p$  specified by  $\lambda$ .
- $s_{\lambda,n}(1)$  is the Schur polynomial evaluated at the identity:

$$s_{\lambda,n}(1) = \prod_{1 \le i \le n} \frac{\lambda_i - \lambda_j + j - i}{i - j}$$
.

Use Murnaghan–Nakayama rule to generate the character tables of  $\mathcal{S}_p.$ 

- Non-recursive method: Young tableau.
- Recursive method: Young diagrams.

#### Examples.

$$\operatorname{Wg}_n((1)) = \frac{1}{n}, \quad \operatorname{Wg}_n((1,1)) = \frac{1}{n^2 - 1}, \quad \operatorname{Wg}_n((2)) = \frac{-1}{n(n^2 - 1)}.$$

$$\int_{\mathcal{U}(n)} U_{ij} \bar{U}_{i'j'} dU = \sum_{\alpha, \beta \in S_1} \delta_{i_1 i'_{\alpha(1)}} \delta_{j_1 j'_{\beta(1)}} \operatorname{Wg}_n(\alpha^{-1}\beta) = \frac{1}{n} \delta_{ii'} \delta_{jj'}.$$

$$\mathbb{E}[(UAU^*)_{k\ell}] = \sum_{r,s=1}^n \mathbb{E}[U_{kr}A_{rs}U_{s\ell}^*] = \sum_{r,s=1}^n \mathbb{E}[U_{kr}A_{rs}\bar{U}_{\ell s}]$$
$$= \sum_{r=1}^n \frac{1}{n}\delta_{k\ell}\,\delta_{rs}\,A_{rs}.$$

$$\mathbb{E}[(UA\bar{U})_{k\ell}] = \sum_{r,s=1}^{n} \mathbb{E}[U_{kr}A_{rs}\bar{U}_{s\ell}] = \sum_{r,s=1}^{n} \frac{1}{n} \delta_{ks} \, \delta_{r\ell} \, A_{rs} .$$

## 1.3 Graphical version

Examples - continued.

$$\mathbb{E}[(UAU^*)_{k\ell}] = \sum_{r,s=1}^{n} \frac{1}{n} \delta_{k\ell} \, \delta_{rs} \, A_{rs} = \frac{1}{n} \delta_{k\ell} \sum_{r=1}^{n} A_{rr} .$$

This means that

$$\mathbb{E}[UAU^*] = \frac{\text{Tr}[A]}{n} I_n .$$

$$\mathbb{E}[(UA\bar{U})_{k\ell}] = \sum_{l=1}^{n} \frac{1}{n} \delta_{ks} \, \delta_{r\ell} \, A_{rs} = \sum_{l=1}^{n} \frac{1}{n} A_{lk} .$$

This means that

$$\mathbb{E}[UAU^*] = \frac{1}{4}A^T.$$

## Graphical calculus [Collins and Nechita (2011)]

$$\mathbb{E}[UAU^*] = \frac{\text{Tr}[A]}{n} I_n \qquad \qquad \mathbb{E}[UAU^*] = \frac{1}{n} A^T$$

Original contractions:

Original contractions:

$$\begin{split} & [\mathsf{L}\text{-end}, (U, \mathsf{out})], [(U, \mathsf{in}), (A, \mathsf{out})], \ \ [\mathsf{L}\text{-end}, (U, \mathsf{out})], [(U, \mathsf{in}), (A, \mathsf{out})], \\ & [(A, \mathsf{in}), (U^*, \mathsf{out})], [(U^*, \mathsf{in}), \mathsf{R}\text{-end}] \ \ [(A, \mathsf{in}), (U^*, \mathsf{in})], [(U^*, \mathsf{out}), \mathsf{R}\text{-end}] \end{split}$$

After: After:

 $[\mathsf{L}\text{-end},\mathsf{R}\text{-end}],[[A,\mathsf{out}],[A,\mathsf{in}]] \qquad [\mathsf{L}\text{-end},[A,\mathsf{in}]],[[A,\mathsf{out}],\mathsf{R}\text{-end}]$ 

## 1.4 "Computer version"

 $[id \otimes Tr](UAU^*)$ :

```
In[1]: from IHU_source import *
In[2]: e1 = [["A", 1, "out", 1], ["U", 1, "in", 1]]
In[3]: e2 = [["A", 1, "out", 2], ["U", 1, "in", 2]]
In[4]: e3 = [["U*", 1, "out", 1], ["A", 1, "in", 1]]
In[5]: e4 = [["U*", 1, "out", 2], ["A", 1, "in", 2]]
In[6]: e5 = [["U", 1, "out", 2], ["U*", 1, "in", 2]]
In[7]: g = [e1, e2, e3, e4, e5]
In [8]: gw = [g,1]
In [9]: visualizeTN (gw)
                                         - lin1:out1][in2:out2]-
```

\_ [out1:in1][out2:in2]

```
\mathbb{E}[[\mathrm{id} \otimes \mathrm{Tr}](UAU^*)]:
ln[10]: k, n = symbols('k n')
In[11]: rm = ["U", [n,k], [n,k], n*k]
In[12]: Eg = integrateHaarUnitary(gw,rm)
In[13]: print(Eg)
In [14]: visualizeTN(Eg)
Out[1]: [[[[['@U*', 1, 'in', 1], ['@U', 1, 'out', 1]],
[['A', 1, 'out', 1], ['A', 1, 'in', 1]],
[['A', 1, 'out', 2], ['A', 1, 'in', 2]]], 1/n]]
```

This figure is followed by the new weight 1/n in the program.

# 1.5 Demo (had been done.)

To get RTNI, go to https://github.com/MotohisaFukuda/RTNI

To get a user-friendly but limited version, called RTNI\_light, go to https://github.com/MotohisaFukuda/RTNI\_light

To try a web version of RTNI\_light, go to https://motohisafukuda.pythonanywhere.com

## 2 Random Gaussian states

"Typical entanglement for Gaussian states", arXiv:1903.04126 [quant-ph] with Koenig.

## 2.1 Motivation and past research

Typical entanglement entropy of random bipartite quantum states.

- Lubkin('78); Lloyd & Pagels('88); Page('93); Foong & Kanno('94): average of entanglement entropy.
- Hayden, Leung and Winter ('04): concentration of entanglement entropy.

- In the setting of random Gaussian states
   Serafini, Dahlsten, Gross & Plenio ('07):
  - micto-canonical measure and canonical measure.
  - inverse squared purity for subsystem of one mode is calculated.
  - Us:
    - direct access to symplectic eigenvalues
    - the number of modes of subsystem can grow as  $n^{\kappa}$  with  $\kappa < 1$  with the total mode is n, in the regime  $n \to \infty$ .
  - Others:

## 2.2 Passive Gaussian unitary operations

The set of Gaussian unitary maps has one-to-one correspondence to the real symplectic group  $\operatorname{Sp}(2n)=\{S\in M_{2n\times 2n}(\mathbb{R})\mid SJS^T=J\}$ ; a Gaussian unitary evolution of a Gaussian state means, for  $S\in\operatorname{Sp}(2n)$ 

$$M \mapsto SMS^T$$

where M is the covariance matrix of the state.

Replacing the real symplectic group  $\mathrm{Sp}(2n)$  by the orthogonal symplectic group  $\mathrm{K}(2n)=\mathrm{Sp}(2n)\cap\mathrm{O}(2n)$ , which is compact. The map

$$\mathsf{U}(n) \to \mathsf{K}(2n)$$

$$U \mapsto \begin{pmatrix} \mathsf{Re}(U) & \mathsf{Im}(U) \\ -\mathsf{Im}(U) & \mathsf{Re}(U) \end{pmatrix} \equiv \frac{1}{2} \begin{pmatrix} I_n & iI_n \\ iI_n & I_n \end{pmatrix} \begin{pmatrix} U & 0 \\ 0 & \overline{U} \end{pmatrix} \begin{pmatrix} I_n & -iI_n \\ -iI_n & I_n \end{pmatrix}$$

induces the measure on K(2n).

# 2.3 Partial trace as Gaussian operation

Suppose we have a Gaussian state of n-mode and want to get the covariance matrix of the reduced k-mode state (with redundant spaces).

For the covariance  $(2n \times 2n)$  matrix  $M_n$  of the n-mode, calculate

$$M_k = \hat{\Pi}_{n,k} M_n \hat{\Pi}_{n,k}$$

with the projection

$$\hat{\Pi}_{n,k} = \begin{pmatrix} \Pi_{n,k} & 0_n \\ 0_n & \Pi_{n,k} \end{pmatrix},$$

where  $\Pi_{n,k} = \operatorname{diag}(\underbrace{1,\ldots,1}_{k \text{ times}},\underbrace{0,\ldots,0}_{n-k \text{ times}})$  is a projection matrix of rank k.

## 2.4 Symplectic eigenvalues

A valid covariance matrix  $(M+iJ\geq 0)$  has Williamson normal form: there is  $S\in \operatorname{Sp}(2n)=\{S\in M_{2n\times 2n}(\mathbb{R})\mid SJS^T=J\}$  such that

$$SMS^T = diag(\lambda_1, \lambda_1, \lambda_2, \lambda_2, \dots, \lambda_n, \lambda_n)$$
.

where  $\lambda_j$  satisfy  $\lambda_j \geq 1$  and are called the *symplectic eigenvalues* .

They can be obtained by computing the spectrum of the matrix  ${\cal J}M$ , which consists of complex conjugate pairs as follows:

$$\operatorname{spec}(JM) = \bigcup_{j=1}^{n} \{\pm i\lambda_j\} .$$

Note that 
$$J = \begin{pmatrix} 0_n & -I_n \\ I_n & 0_n \end{pmatrix}$$
.

# 2.5 Random Gaussian pure states

The covariance matrix M of a pure n-mode Gaussian state can be diagonalized by Euler/Bloch-Messiah decomposition;

$$M = S \begin{pmatrix} Z_n & 0 \\ 0 & Z_n^{-1} \end{pmatrix} S^T$$
 for some  $S \in \mathsf{K}(2n)$ .

Here,  $Z_n = \operatorname{diag}(z_1, \ldots, z_n)$  with  $z_i \geq 1$ .

The induced measure on  $\mathsf{K}(2n)$  generates random pure Gaussian states.

We call  $\{z_j\}_{j=1}^n$  the *squeezing parameters*. Observe that such a state has energy:

$$\operatorname{Tr}\left[Z_n+Z_n^{-1}\right]=\sum_{j=1}^n E_j$$
 where  $E_j=rac{1}{2}(z_j+1/z_j)$  .

We can understand  $E_j$  as the energy in the j-th mode after a suitable Gaussian unitary rotation.

# 2.6 Symplectic eigenvalues of reduced states

Let  ${\cal M}_k$  be the covariant matrix of k-mode, reduced from the random state of n mode. Then,

$$JM_k = \frac{1}{2} \begin{pmatrix} i\Pi & \Pi \\ -\Pi & -i\Pi \end{pmatrix} \begin{pmatrix} UAU^T & -iUBU^* \\ -i\bar{U}BU^T & -\bar{U}AU^* \end{pmatrix} \begin{pmatrix} \Pi & i\Pi \\ i\Pi & \Pi \end{pmatrix}$$

where  $A = (Z_n - Z_n^{-1})/2$  and  $B = (Z_n + Z_n^{-1})/2$ .

Define

$$f(U) = \operatorname{Tr}\left[\left((JM_k)^2 + \lambda^2 I_{2k}\right)^2\right] = 2\sum^k \left(\lambda_j^2 - \lambda^2\right)^2.$$

Here,  $\{\lambda_j\}_{j=1}^k$  are the symplectic eigenvalues of  $M_k$  and  $\lambda=\mathrm{Tr}[B]/n$  is the averaged energy.

The calculation of  $\mathbb{E}[f(U)]$  involves Weingarten calculus with  $\mathcal{S}_4$ .

#### 2.7 Concentration

#### Concentration of symplectic eigenvalues:

For fixed k, suppose the energy of the initial pure Gaussian state is universally bounded;  $\sup_n \|Z_n\|_{\infty} < \infty$  ( $n^z$  with z < 1/8 is possible). Then

$$\Pr\{f(U) > \epsilon\} < \exp(-c\epsilon n)$$

for some universal constant c > 0.

#### Concentration of entanglement entropy:

In addition, if  $\lambda = \lambda_n$  is bounded below by  $\mu > 1$ , then

$$\Pr\{|S(M_k) - S(\lambda I)| > \epsilon\} < \exp\left(-c\frac{\epsilon^4}{\beta(\mu)^4}n\right)$$

for some universal constant c>0;  $S(\cdot)$  is von Neumann entropy, and

$$\beta(\mu) = \log \frac{\mu + 1}{\mu - 1} \ .$$

is called the inverse temperature.

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