

Bootstrability with Improved Truncation Methods

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based on 2108.08859, 2108.09330, 2209.02801, **2306.15730**

with **G. Kantor, C. Papageorgakis, P. Richmond, A.G. Stapleton, M. Woolley** (QMUL)



This talk is about:

alternative strategies in the (numerical) conformal bootstrap

1. **Physics (QFT)** part:

alternative **non-perturbative** solution methods in CFTs

[CFT problem]

The standard **linear functional method** vs other options

2. **Computational** part:

non-convex, large-scale optimization problems

Is **Machine Learning (ML) / Artificial Intelligence (AI)** a useful tool?

[numerical, Computer Science problem]

Example: Bootstrability [Cavaglia et al. '22]



in 1d CFT on 1/2-BPS Wilson line in planar N=4 SYM theory

Sample of results for 3 squared (unprotected) OPE-coefficients

SDPB (~1000 digit precision)

C_1^2	C_2^2	C_3^2
$0.294014873 \pm 4.88 \cdot 10^{-8}$	$0.039788 \pm 4.10 \cdot 10^{-4}$	$0.146757 \pm 5.82 \cdot 10^{-4}$
$0.294014228 \pm 6.77 \cdot 10^{-7}$	$0.041832 \pm 1.86 \cdot 10^{-3}$	$0.144100 \pm 2.39 \cdot 10^{-3}$

$$\lambda = (4\pi)^2 \simeq 157.91$$

improved truncation (16 digit precision)

- For the **first time** a non-rigorous method, **not relying on positivity**, **competes so directly** with the so-far standard rigorous tools in numerical conformal bootstrap!

That opens up many possibilities...

- **AI** algos could play an interesting new role in theoretical problems

I will comment on a comparison between ML and non-ML algorithms in a specific example

Some motivation

- Why [Conformal Field Theories \(CFTs\)](#)?
 - UV/IR behaviour of QFTs
 - Phase transitions
 - Quantum Gravity via the AdS/CFT correspondence
& via the worldsheet description of strings...

- Why [bootstrap](#)?

The non-perturbative structure of QFT is rich, but poses a **hard** conceptual and computational problem

- Real-world physical systems with strong interactions
- QFT-QFT (strong-weak) dualities
- QFT-gravity dualities (holography)
- New exotic **non-Lagrangian** QFTs from String Theory...
(kills' Lattice)

We need a (new) powerful framework for all these cases...

- The **conformal bootstrap programme**

[Ferrara-Grillo-Gatto '73, Polyakov '74] + [Rattazzi-Rychkov-Tonni-Vichi '08]

aims to solve CFTs non-perturbatively leveraging general principles of symmetry (*without using the path-integral*)

— Caution: Symmetry alone cannot not be enough!

Combine: analytical, numerical, exact, perturbative...

- Recent progress in many fronts:

analytical/numerical conformal bootstrap, new exact methods in SUSY gauge theories (e.g. integrability, localization) ...

at the center of this talk

How symmetry helps in CFTs

Local CFT data: operators $\mathcal{O}(x)$ with some quantum numbers under global symmetries [scaling dimension Δ , spin s , charges Q ...]

2-point correlation functions:
$$\langle \mathcal{O}_{\Delta}^{(1)}(x_1) \mathcal{O}_{\Delta}^{(2)}(x_2) \rangle = \frac{G_{12}}{|x_1 - x_2|^{2\Delta}}$$

3-point functions:
$$\langle \mathcal{O}_{\Delta_1}(x_1) \mathcal{O}_{\Delta_2}(x_2) \mathcal{O}_{\Delta_3}(x_3) \rangle = \frac{C_{123}}{|x_{12}|^{\Delta_{12,3}} |x_{23}|^{\Delta_{23,1}} |x_{13}|^{\Delta_{13,2}}}$$

$$\Delta_{ij,k} := \Delta_i + \Delta_j - \Delta_k$$

Example: 4-point functions (identical operators)

$$\langle \mathcal{O}(x_1)\mathcal{O}(x_2)\mathcal{O}(x_3)\mathcal{O}(x_4) \rangle = \frac{g(u, v)}{|x_{12}|^{2\Delta} |x_{34}|^{2\Delta}}$$

$$u = \frac{x_{12}^2 x_{34}^2}{x_{13}^2 x_{24}^2}, \quad v = \frac{x_{23}^2 x_{14}^2}{x_{13}^2 x_{24}^2} \quad (\text{cross ratios})$$

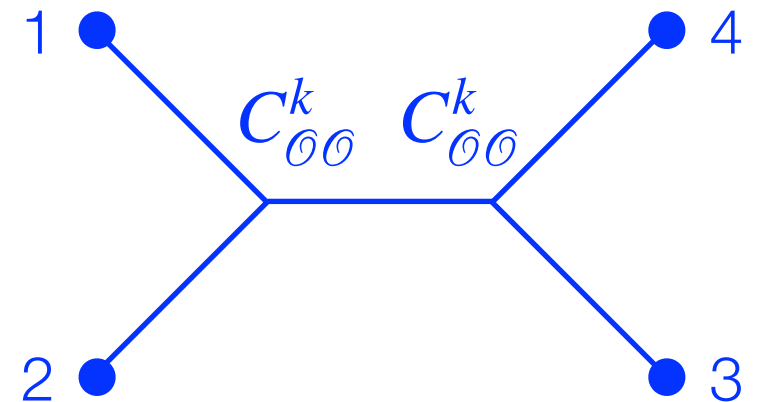
$$u := z\bar{z}$$

$$v := (1-z)(1-\bar{z})$$

Using the (12) and (34) OPEs we get an expansion of the form:

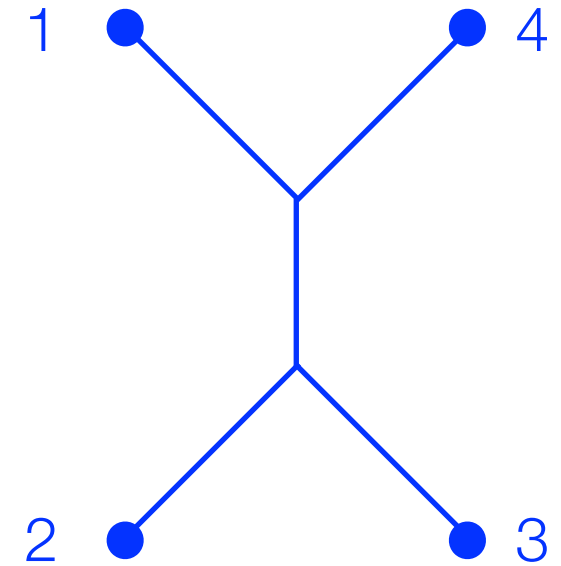
$$g(u, v) = \sum_k \left(C_{\mathcal{O}\mathcal{O}}^k \right)^2 \boxed{g_{\Delta_k, \ell_k}(u, v)} \longrightarrow$$

conformal blocks
known functions



The same function $g(u, v)$ can be obtained with the (14)-(23) OPE

$$g(u, v) = \left(\frac{u}{v}\right)^\Delta \sum_k (C_{\mathcal{O}\mathcal{O}}^k)^2 g_{\Delta_k, \ell_k}(v, u)$$



Yields a crossing equation of the form

$$\sum_k (C_{\mathcal{O}\mathcal{O}}^k)^2 F_{\Delta_k}(u, v) = 0 \quad \text{to be solved for the CFT data } (\Delta_k, C_{\mathcal{O}\mathcal{O}}^k)$$

Converted to an optimization problem:

- linear functional method (standard, convex optimization)
- we will discuss alternatives (non-convex opt) ➔ ML/AI comes in

Solving the crossing equations

$$\mathfrak{C}_0 := (C_{000}^k)^2$$

$$\sum_{\mathcal{O}} \mathfrak{C}_{\mathcal{O}_{\Delta,S}} F_{\mathcal{O}_{\Delta,S}}(z, \bar{z}) = 0$$

linear in \mathfrak{C} *non-linear in Δ*

- functional dependence on z, \bar{z}
- infinite unknowns \mathfrak{C}, Δ
- infinite number of 4-point functions (& similar sum-rules)

Linear functional method

[Rattazi-Rychkov-Tonni-Vichi, '08]

- Act on crossing eq. with linear functionals α

$$\sum_{\mathcal{O}} C_{\mathcal{O}_{\Delta,S}} \alpha(F_{\mathcal{O}_{\Delta,S}}) = 0 \quad [\alpha(F_{\mathcal{O}_{0,0}}) = 1]$$

≥ 0 ≥ 0

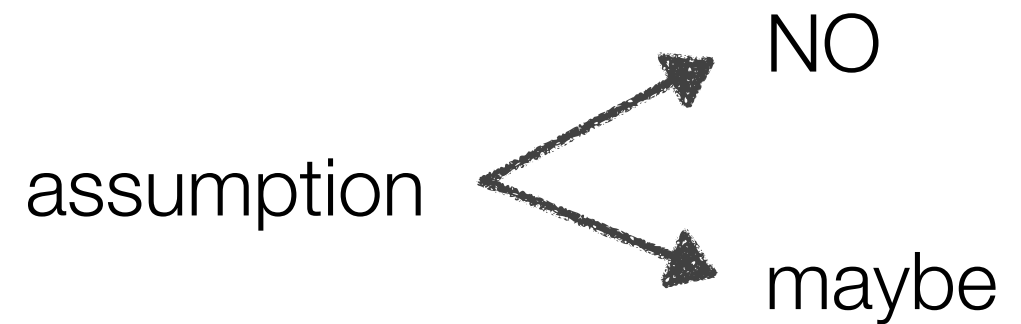
- Oracle assumptions, e.g. $\Delta_s \geq \Delta_{\min}$ for some operator

- Search for α :

$$\alpha[F_{\mathcal{O}_{\Delta,S}}] \geq 0 \quad \forall \Delta \neq 0$$

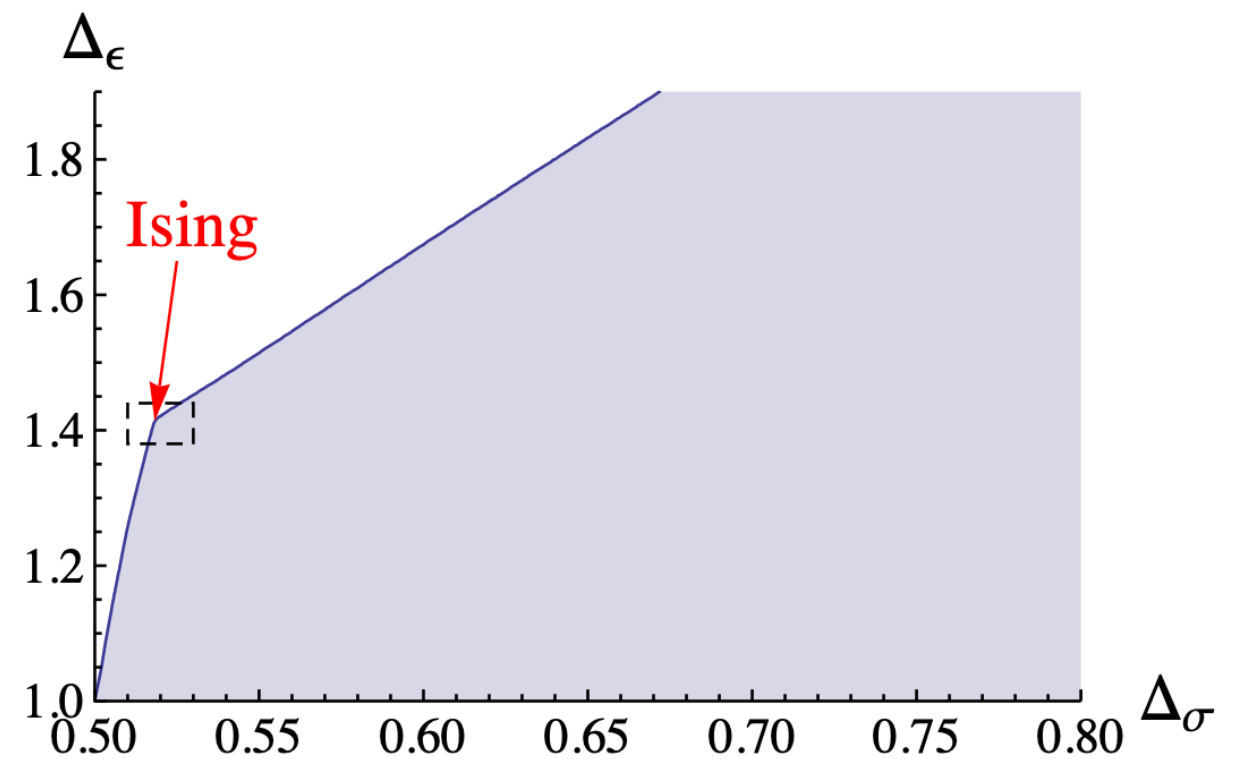
- If α exists, the assumption is eliminated

This is a **classification algorithm**:



Pros:

- Rigorous
- Generic statements for CFTs
- Sometimes cusps/islands
- Powerful Semi-Definite Programming algorithm **SDPB** [Simmons-Duffin '15] (convex optimization)



El-Showk et al, '12

Cons:

- Typically, the assumptions are blind. Difficult to:
 - Scan high-dimensional parameter spaces
 - Explore
 - *“Solve my theory”*
- Positivity $\mathfrak{C}_k \geq 0$ is not available in many contexts of interest
 - higher-point bootstrap
 - boundary CFT bootstrap
 - non-unitary theories

The diagram illustrates an equality between two expressions. On the left, a sum over k (indicated by a green \sum_k) is applied to a vertex diagram. This diagram consists of a horizontal line at the bottom with a label b_k below it. A vertical line extends upwards from the center of this line to a vertex. From this vertex, two lines branch out upwards and outwards to two external legs, each labeled with a circle containing the letter \mathcal{O} . The vertex itself is labeled with \mathcal{O}_k to its right. The lines connecting the vertex to the external legs are labeled with C_{12}^k in green. On the right side of the equation, there is an equals sign followed by a sum over k' (indicated by a green $\sum_{k'}$). This sum is applied to two separate diagrams. Each diagram has a horizontal line at the bottom with a label $b_{k'}$ below it. A single vertical line extends upwards from the center of each horizontal line to a vertex labeled with a circle containing the letter \mathcal{O} .

Cons:

- SDPB is powerful but expensive
(100s of digits precision required)
- Multiple-correlator bootstrap is a challenge



desirable to supplement the linear functional method with other methods

Truncation methods offer an alternative but are notoriously messy...

Let's examine the issues. Our exact crossing eq. reads:

$$\sum_n C_n F_n(z, \bar{z}) + r(z, \bar{z}) = 0 \quad (*)$$

known functions of Δ_n

unknown C_n, Δ_n

includes exactly known contributions

Step 1: discretize the z, \bar{z} dependence

e.g., evaluation on a lattice or z, \bar{z} -derivatives on $z = \bar{z} = \frac{1}{2}$

Yields a finite system of equations (still exact)

$$\sum_n \bigoplus_n \mathbb{T}_n + \mathbb{r}_b = 0$$

Step 2: Identify a subset \mathcal{S} of “most significant” operators in \sum_n

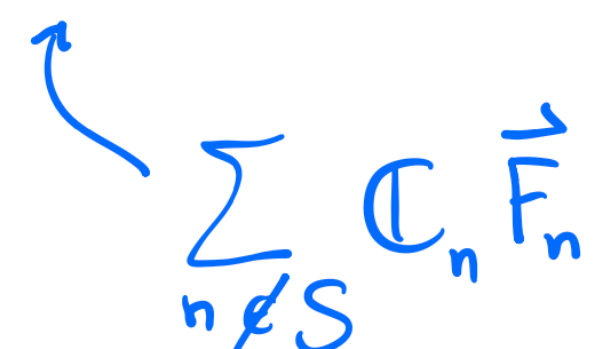
spin-partition: consider spins up to some max spin & for each spin s

assume N_s operators

(don't need to know their Δ 's)

Then (*) becomes (still exact)

$$\sum_{n \in S} \mathbb{C}_n \vec{F}_n + \vec{T} + \vec{r} = 0$$



(Common) Step 3: Set $\vec{T} = 0$ for the 'tail' and solve!

[Gliozzi '13] Pretend $\sum_{n \in S} \mathbb{C}_n \vec{F}_n + \vec{r} = 0$ is an exact over-constrained

systems and try to solve it.

[Technicalities: drastic truncation, typically $O(10)$ operators]

- [Li '17] Formulate a cost function \mathcal{L} (e.g., L₂-norm of \vec{E})

$$\text{Cost} \left[\left\{ \Delta_n, \mathbf{C}_n \right\}_{n \in \mathcal{S}} \right] := \mathcal{L}(\vec{E})$$

and minimize it

$$\vec{E} \equiv \sum_{n \in \mathcal{S}} \mathbf{C}_n \vec{F}_n + \vec{r}$$

This is now a highly non-trivial **non-convex** optimization problem

In [2108.08859, 2108.09330] we started thinking about this as a

large scale optimization problem (thinking that the larger \mathcal{S} the better!)

This framework is still not good enough!

[1] It is not algorithmic

Q: How do you select the spin-partition?

[2] Q: What is the effect of the infinite number of operators in the dropped
`tail' contribution \overrightarrow{T} ? (across parameter spaces?)

[3] Near-degeneracies can grow very quickly with scaling dimension
(and spectrum can be chaotic)

Q: How do you track this complexity?

[4] It is not a straightforward optimization problem!

- Multi-nodal high-dimensional landscape of \mathcal{L}
- Multiple basins of minima, complicated microstructure
(& many configurations with comparable cost value)

You may not be interested in the global minimum!

The fastest optimizer may not be the right one!

Efficiency in 'local guided search'
Markov-chain algorithms
may be more suitable! Precisely
where AI/ML comes in for us...

[5] Related to the complexities of [3]+[4]:

adding more operators does not necessarily improve the accuracy

A viewpoint shift and proposed improvements

We now focus on families of CFTs: $\text{CFT}[\lambda]$

- Consider adiabatic deformations of a CFT solved at some λ^*
[improves [1], consistent with the “solve my CFT” perspective]

- **DO NOT drop** the tail \vec{T} , approximate it!

[improves [2]]

Static approximation: At the known solution point λ^*

$$\sum_n \mathbb{C}_n^* \vec{F}_n^* + \vec{T}^* + \vec{r}^* = 0 \quad (\text{exact})$$

Allows us to determine the λ^* -tail from a finite number of CFT data:

$$\vec{T}^* = - \sum_{n \in S} C_n^* \vec{F}_n - \vec{r}^*$$

Assume that $\vec{T}(\lambda) \simeq \vec{T}(\lambda^*)$ (at least in some regime of λ)

Then,

$$\sum_{n \in S} C_n \vec{F}_n + \vec{r} - \sum_{n \in S} C_n^* \vec{F}_n^* - \vec{r}^* \simeq 0$$

$$\vec{E} - \vec{E}^* \simeq 0$$

to be solved by minimizing a new cost function $\mathcal{L}(\vec{E} - \vec{E}^*)$

- The contribution of **nearly-degenerate** operators can be approximated by effective operators [addresses [3]]

$$\sum_{\Delta \epsilon \text{ band } C S} C_n^{(\text{exact})} F_n^{(\text{exact})} \approx \sum_{n_{\text{eff}}} C_{n_{\text{eff}}} F_{n_{\text{eff}}}$$

reduced # of ops

- Effective high- Δ operators can also be used to [address [5]]
 - “dynamical soft tail”

Different optimization algorithms may treat the soft tail differently!!

THE COMPUTATIONAL PROBLEM

Determine (local) minima of $\mathcal{L}(\vec{E} - \vec{E}^*)$.

This is a non-convex optimization problem for the CFT data
(actual+effective) $\{\Delta_n, \mathfrak{C}_n\} \in \mathcal{S}$ (order 100-1000 unknowns)



BootSTOP (Bootstrap STOchastic OPTimiser)

From [BootSTOP repository README](#)

- [Contents](#)
 - [Overview](#)
 - [Installation](#)
 - [Running the code](#)
 - [References](#)

PyGMO (Python Parallel Global Multiobjective Optimizer)
Izzo-Biscani — European Space Agency

SAC (Soft-Actor-Critic)
developed for task control in robotics

Overview

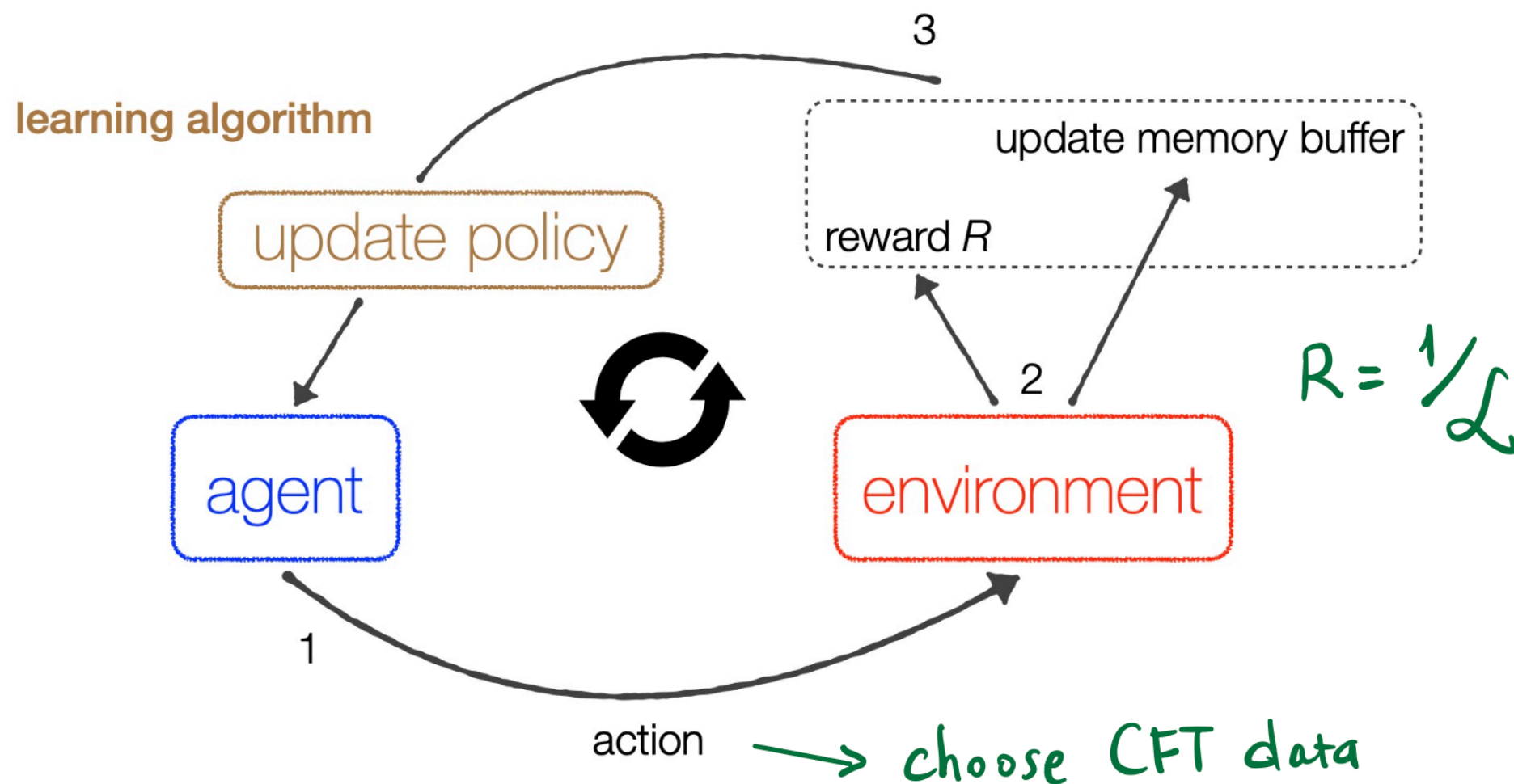
BootSTOP is a Python package for determining CFT data (OPE-coefficients squared and scaling dimensions) which minimise a theory's truncated crossing equation. To do this the code can apply either a custom PyTorch implementation of the Soft-Actor-Critic algorithm or one of the algorithms within the PyGMO package (information about PyGMO can be found [on the PyGMO website](#)).

At present the crossing equation for each of the following CFTs is coded within BootSTOP: 1D defect CFT (see [4]),

Soft Actor-Critic (SAC) algorithm

[Harnooja, Zhou, Abbeel, Levine '18]

- Stochastic optimization as a **Markov Decision Process**
- Handles continuous actions and state spaces



Technical features of the 1d CFT application that follows

- 2 sets of runs on [QMUL HPC Apocrita](#)
 - $(4 \times 10^8 \text{ agents, } \sim 15 \text{ mins})$
 - IPOPT C PyGMO
 - SAC
 - $(200 \text{ parallel agents, } \sim 12 \text{ hrs})$
- 124 CFT data
- Crossing eqs with up to 260 or 700 derivatives (preloaded in BootSTOP)
- Machine precision (16 digits)

Application: 1d defect CFT on Wilson lines

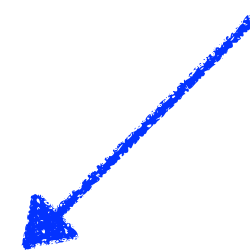
Infinite straight (1/2-BPS) Wilson line in 4d N=4 SYM

$$\mathcal{W} = \text{Tr} P \exp \int_{-\infty}^{\infty} dt \left(iA_t + \Phi_{\parallel} \right)$$

Local operators $\mathcal{O}(t)$ inserted on Wilson line

captured by a 1d CFT

Interested in correlation functions



$$\langle\langle \mathcal{O}_1(t_1) \cdots \mathcal{O}_n(t_n) \rangle\rangle := \langle \text{Tr} W_{-\infty}^{t_1} \mathcal{O}_1(t_1) W_{t_1}^{t_2} \mathcal{O}_2(t_2) \cdots \mathcal{O}_n(t_n) W_{t_n}^{+\infty} \rangle$$

Focus on $\langle\langle\Phi_{\perp}^1(x_1)\Phi_{\perp}^1(x_2)\Phi_{\perp}^1(x_3)\Phi_{\perp}^1(x_4)\rangle\rangle$

- Φ_{\perp}^1 is one of the 5 transverse scalars of N=4 SYM
- Planar limit, dependence on the 't Hooft coupling $\lambda = g_{YM}^2 N$

Definition: $g := \frac{\sqrt{\lambda}}{4\pi}$

- Dependence of the 4-point function on the single cross-ratio

$$\chi = \frac{x_{12}x_{34}}{x_{13}x_{24}}$$

- Operators are organized in **superconformal multiplets**

— protected \mathcal{B}_k , $k = 1, 2, \dots$ with $\Delta_k = k$ ($\Phi_{\perp}^1 \in \mathcal{B}_1$)

— long \mathcal{L}_{Δ}

- OPE:

$$\mathcal{B}_1 \times \mathcal{B}_1 = 1 + \mathcal{B}_2 + \sum_n \mathcal{L}_{\Delta_n}$$

- Crossing equation ...

$$\sum_n C_n^2 G_{\Delta_n}(\chi) = H(\chi)$$

known, depends on λ

unknown

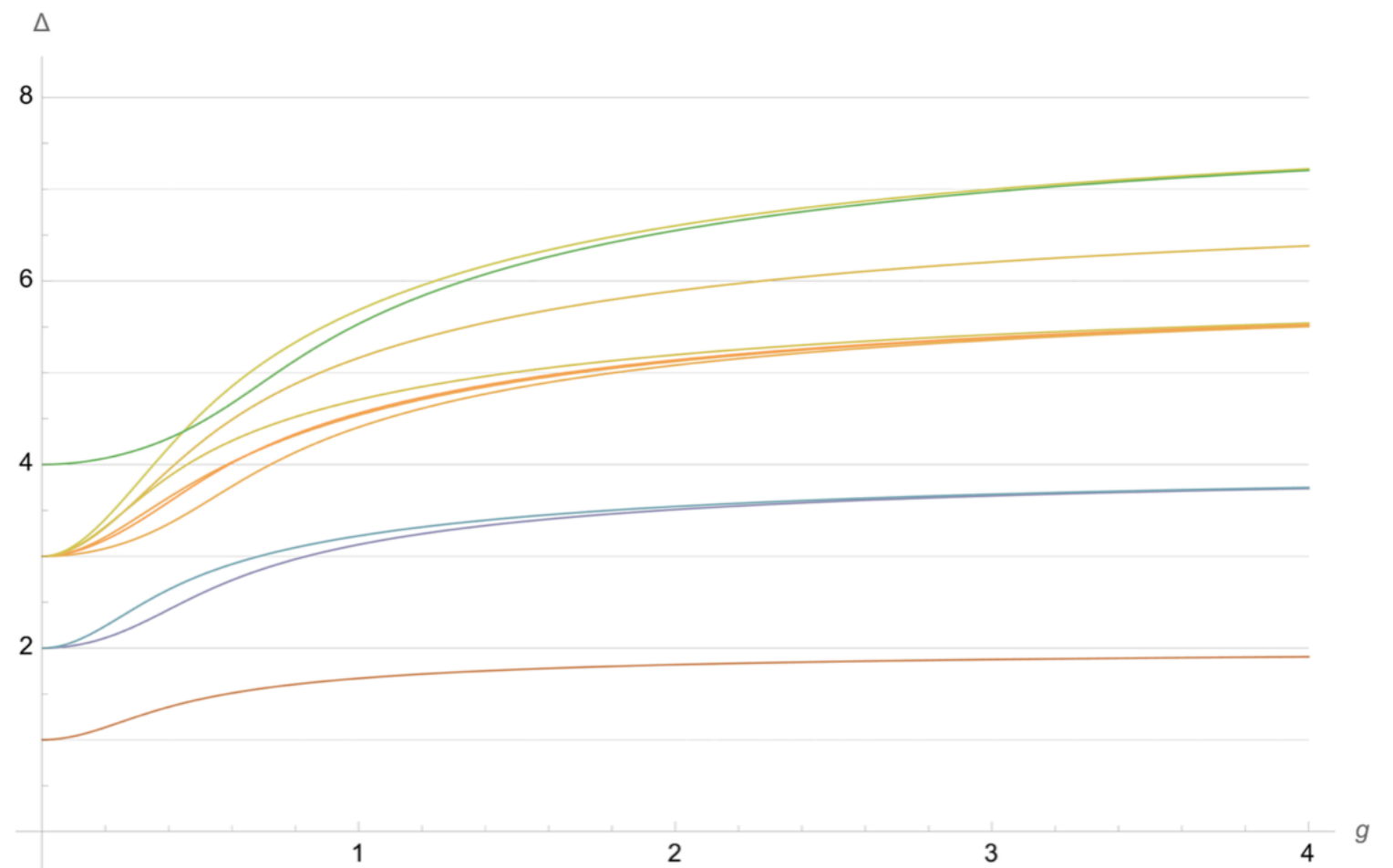
Bootstrability [Cavaglia-Gromov-Julius-Preti, '21]

- Use integrability (QSC) to fix (some of the) planar Δ_n

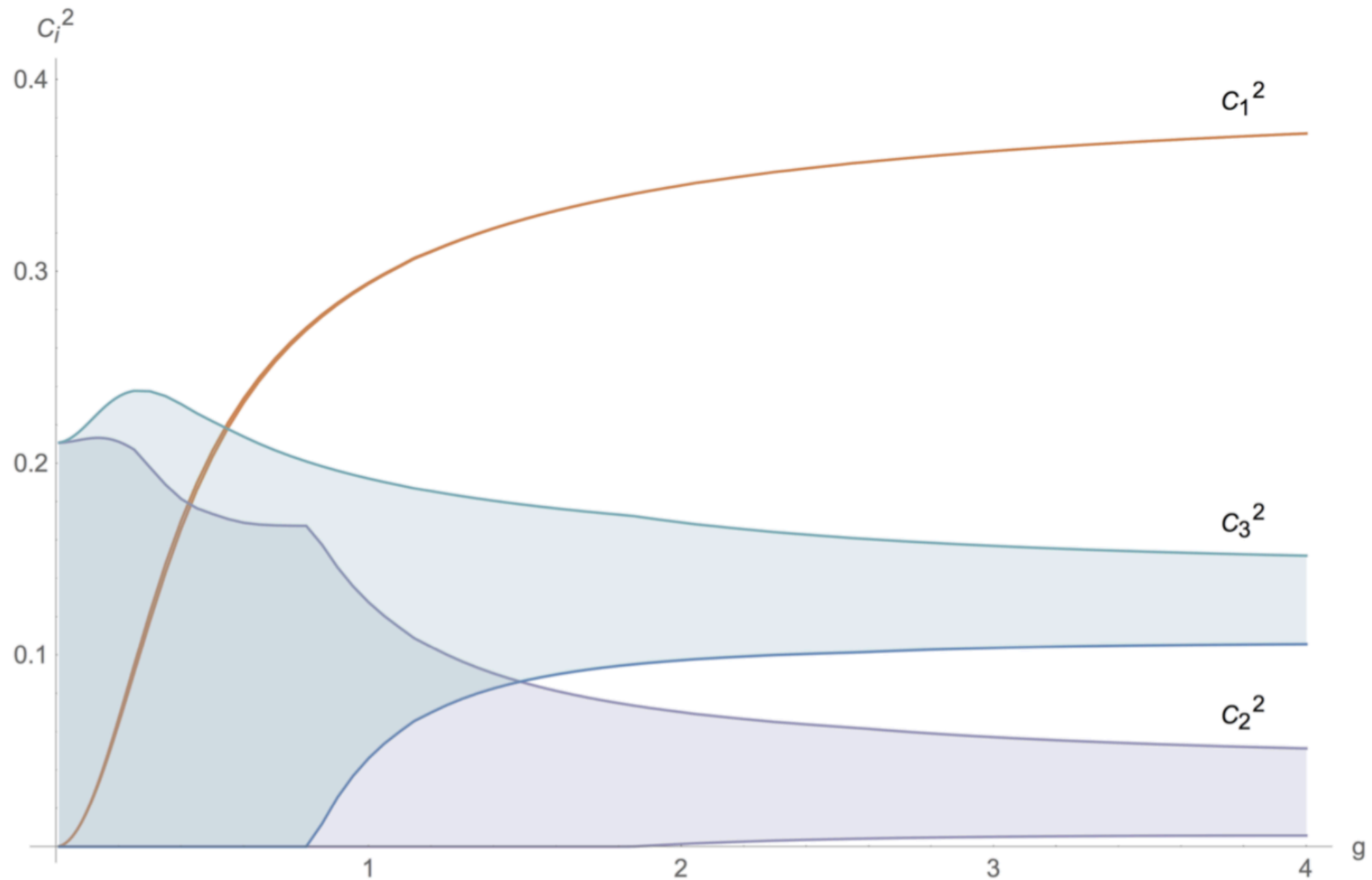
- Use bootstrap to determine C_n^2 ($C_n := C_{\mathcal{B}_1 \mathcal{B}_1 \mathcal{L}_{\Delta_n}}$)

Fix 10 long Δ_n 's and

determine C_1^2, C_2^2, C_3^2



Strategy 1: Linear functional method [Cavaglia et al, '21]



Strategy 2: Traditional truncation

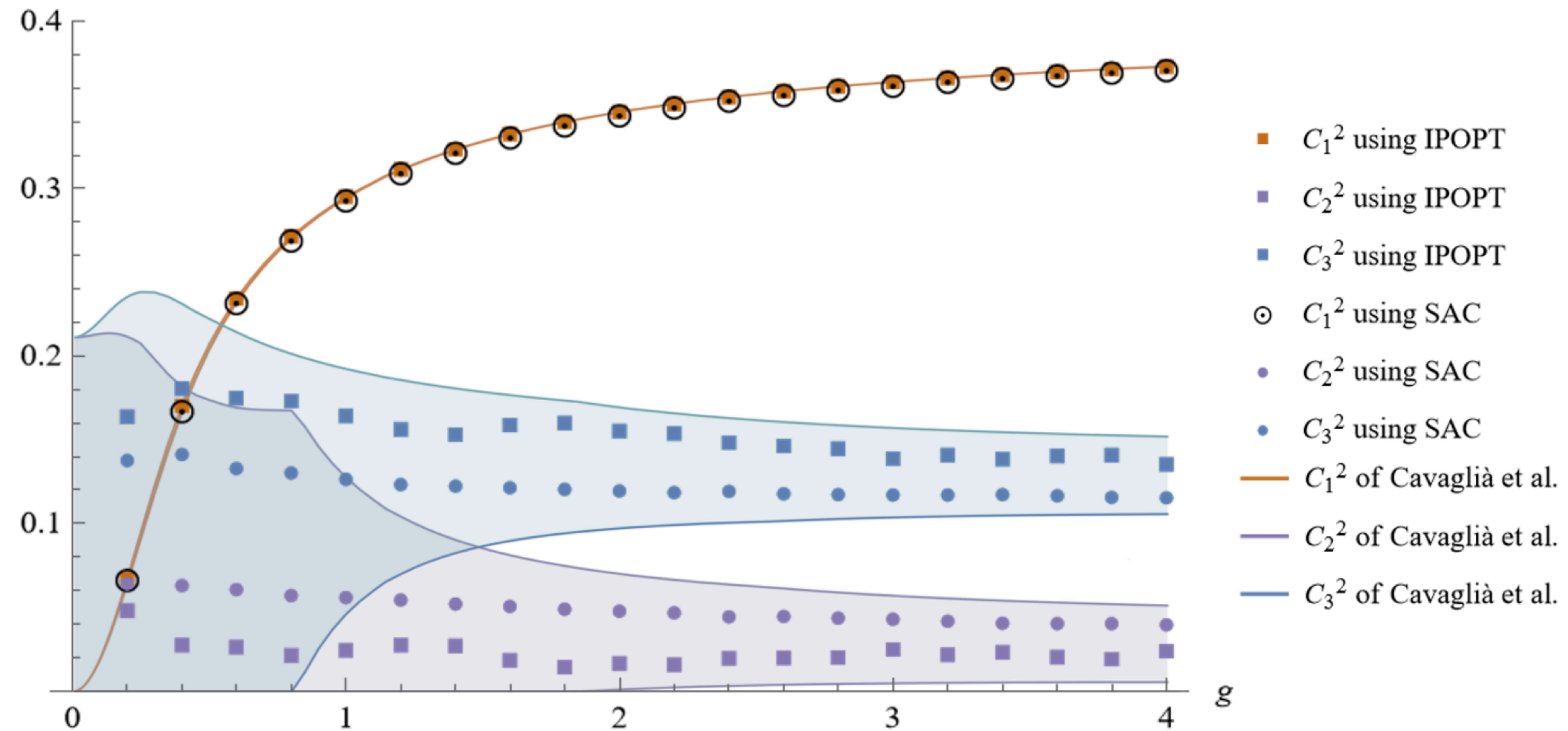
- Truncate to 10 operators with fixed scaling dimensions
- Minimize the root mean square cost function

With linear dependence on the unknowns $\mathfrak{C}_n := C_n^2$ this is a
linear-regression problem

The results are **not** impressive

Strategy 3: Improved Truncation


[MN-Papageorgakis-Richmond-Stapleton-Woolley, '23]




- C_1^2 results reproduced quite well

e.g., at $g = 0.2$

SDPB	0.0663 $\pm 1.9 \times 10^{-3}$
IPOPT	0.066073 $\pm 4.2 \times 10^{-5}$
SAC	0.067339 $\pm 1.26 \times 10^{-3}$

 rigorous

 statistical

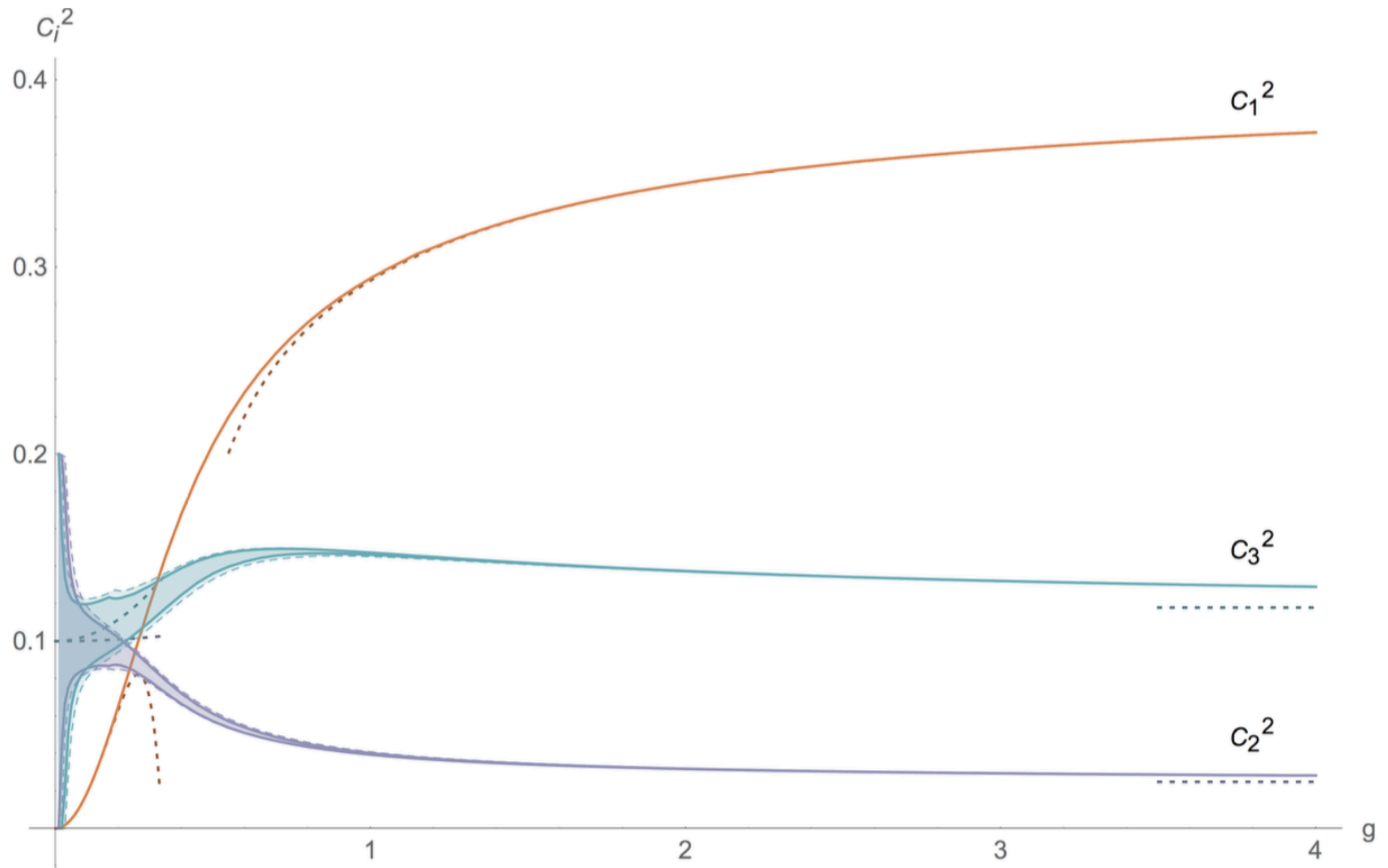
- IPOPT & SAC treat the effective spectrum differently. Spread between IPOPT-SAC shows correlation with size of SDPB allowed regions
- SAC mean is surprisingly accurate !!

Bootstrability with integrated constraints [Cavaglia et al, '22]

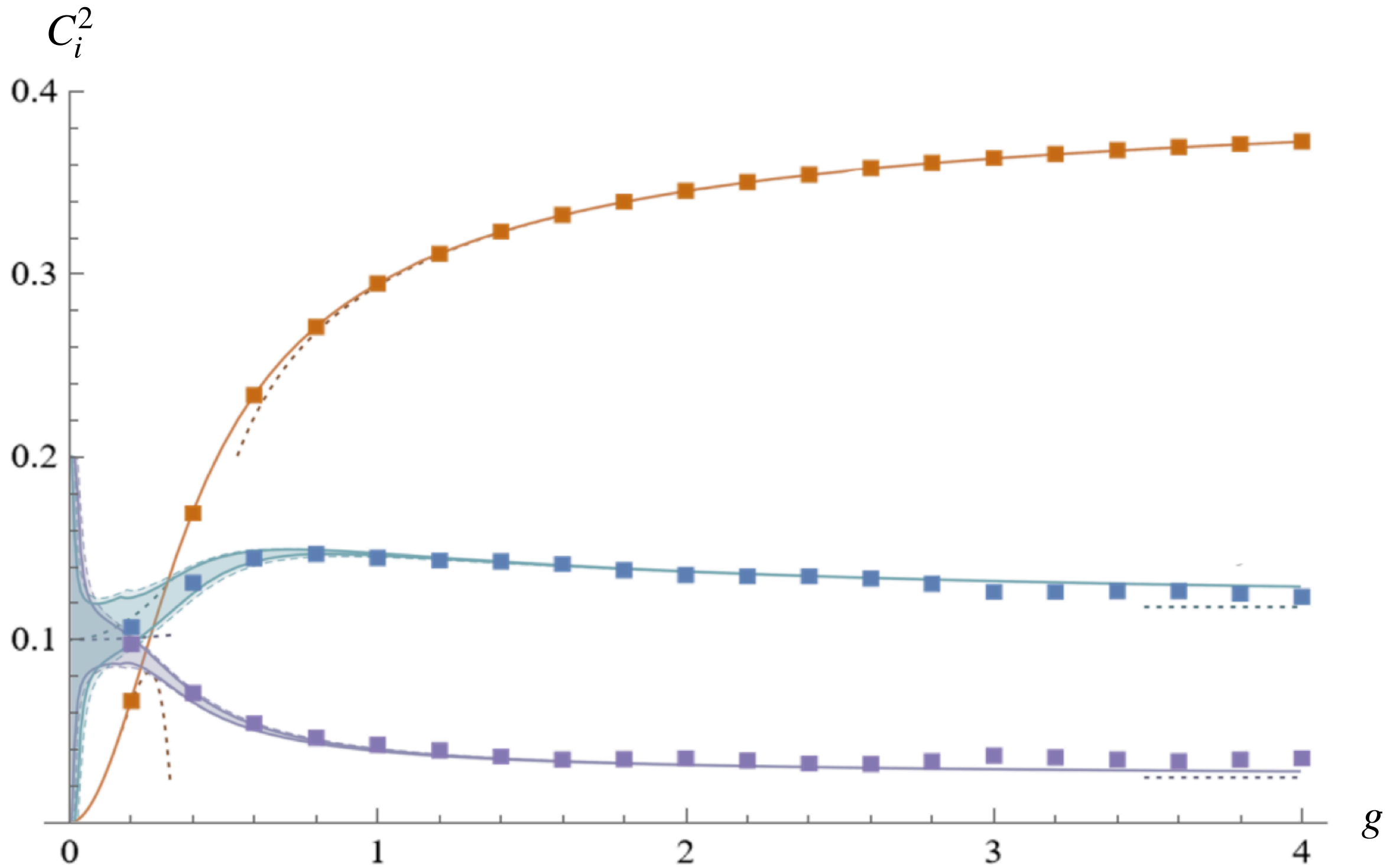
Integrated correlators [Drukker et al, '22] [Cavaglia et al, '22] yield 2 extra sum rules for the defect CFT data **that depend on g** .

Use them to tighten the bootstrap bounds !

Linear functional method [Cavaglia et al, '22]



Improved Truncation [\[MN-Papageorgakis-Richmond-Stapleton-Woolley, '23\]](#)



- Improved truncation performs as well (sometimes even better!) with much less demand on numerical precision (**16 vs 1000 digits**)

Method	g	C_1^2	C_2^2	C_3^2
[2]	0.2	$0.065679029 \pm 6.95 \cdot 10^{-7}$	$0.09452 \pm 7.25 \cdot 10^{-3}$	$0.1101 \pm 1.27 \cdot 10^{-2}$
IPOPT	0.2	$0.06567873 \pm 1.55 \cdot 10^{-7}$	$0.09683 \pm 1.41 \cdot 10^{-3}$	$0.1063 \pm 2.42 \cdot 10^{-3}$
[2]	0.4	$0.16838882 \pm 1.29 \cdot 10^{-6}$	$0.06925 \pm 2.80 \cdot 10^{-3}$	$0.13196 \pm 7.16 \cdot 10^{-3}$
IPOPT	0.4	$0.16838814 \pm 6.13 \cdot 10^{-7}$	$0.07010 \pm 1.06 \cdot 10^{-3}$	$0.13026 \pm 2.58 \cdot 10^{-3}$
[2]	0.6	$0.233041731 \pm 4.49 \cdot 10^{-7}$	$0.05246 \pm 1.47 \cdot 10^{-3}$	$0.14546 \pm 2.99 \cdot 10^{-3}$
IPOPT	0.6	$0.233041064 \pm 8.18 \cdot 10^{-7}$	$0.05347 \pm 1.30 \cdot 10^{-3}$	$0.14376 \pm 2.37 \cdot 10^{-3}$

- The soft tail works well everywhere from weak to strong coupling!
- New predictions for higher states are possible
(improvements expected with more QSC input)
- Multiple-correlator bootstrap is cheap and technically possible

Take home messages

- Improved truncation methods are flexible, cheap & can be efficient and accurate!
- We can use them to attack physically interesting situations where positivity is absent
- AI-powered methods may play a useful role
 - SAC mean is impressively accurate considering the numbers we ran (200 SAC vs 4×10^8 IPOPT agents)
 - Better RL implementations? Collaborative AI, e.g. MARL?

Some (preliminary results) on the 6d $N=(2,0)$ superconformal bootstrap

(without improvements in truncation method)

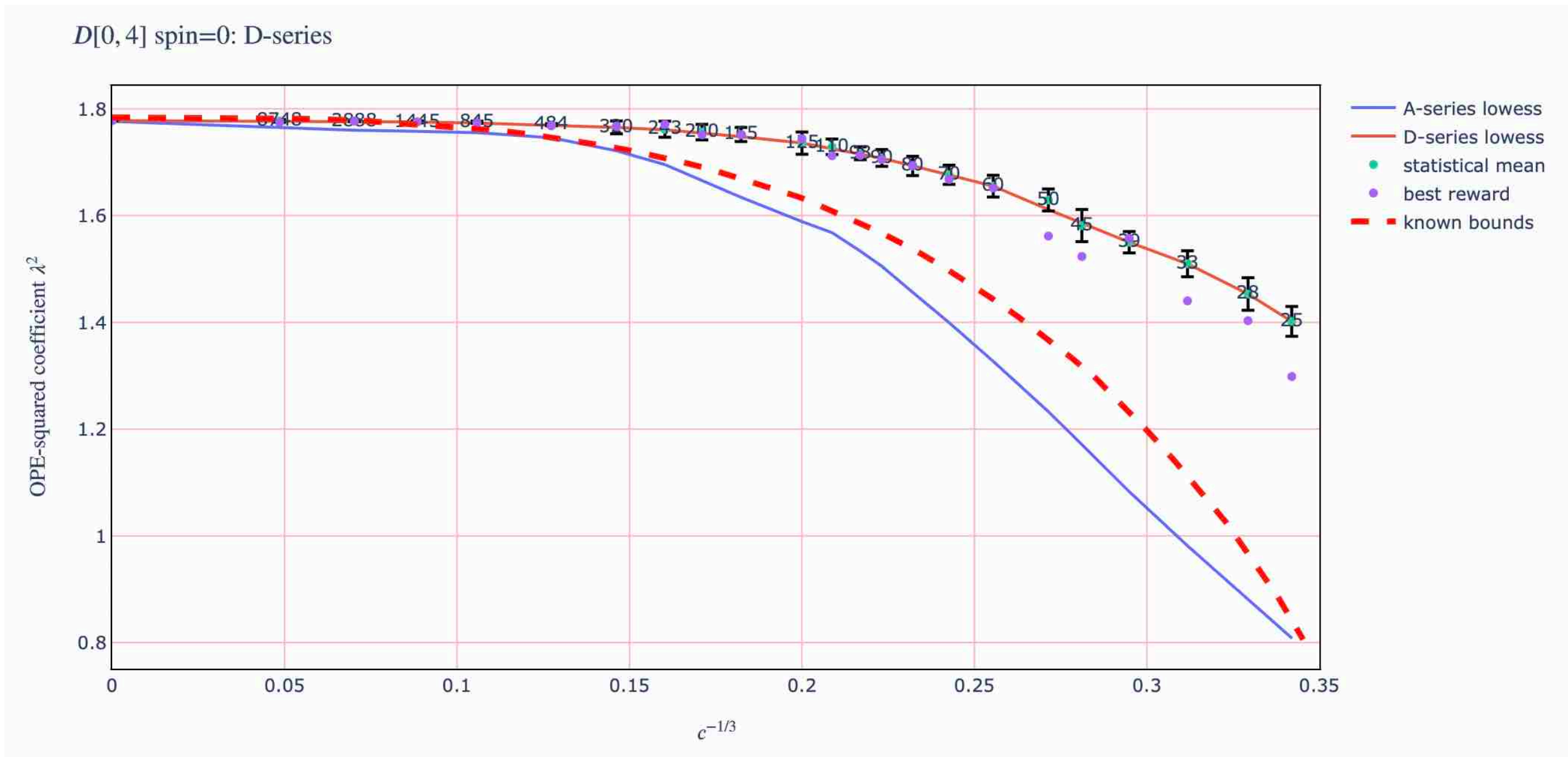
[Kantor-VN-Papageorgakis-Richmond]

- No Lagrangian — only AdS/CFT at $c \rightarrow \infty$
- A- and D-series theories. Can bootstrap distinguish them?

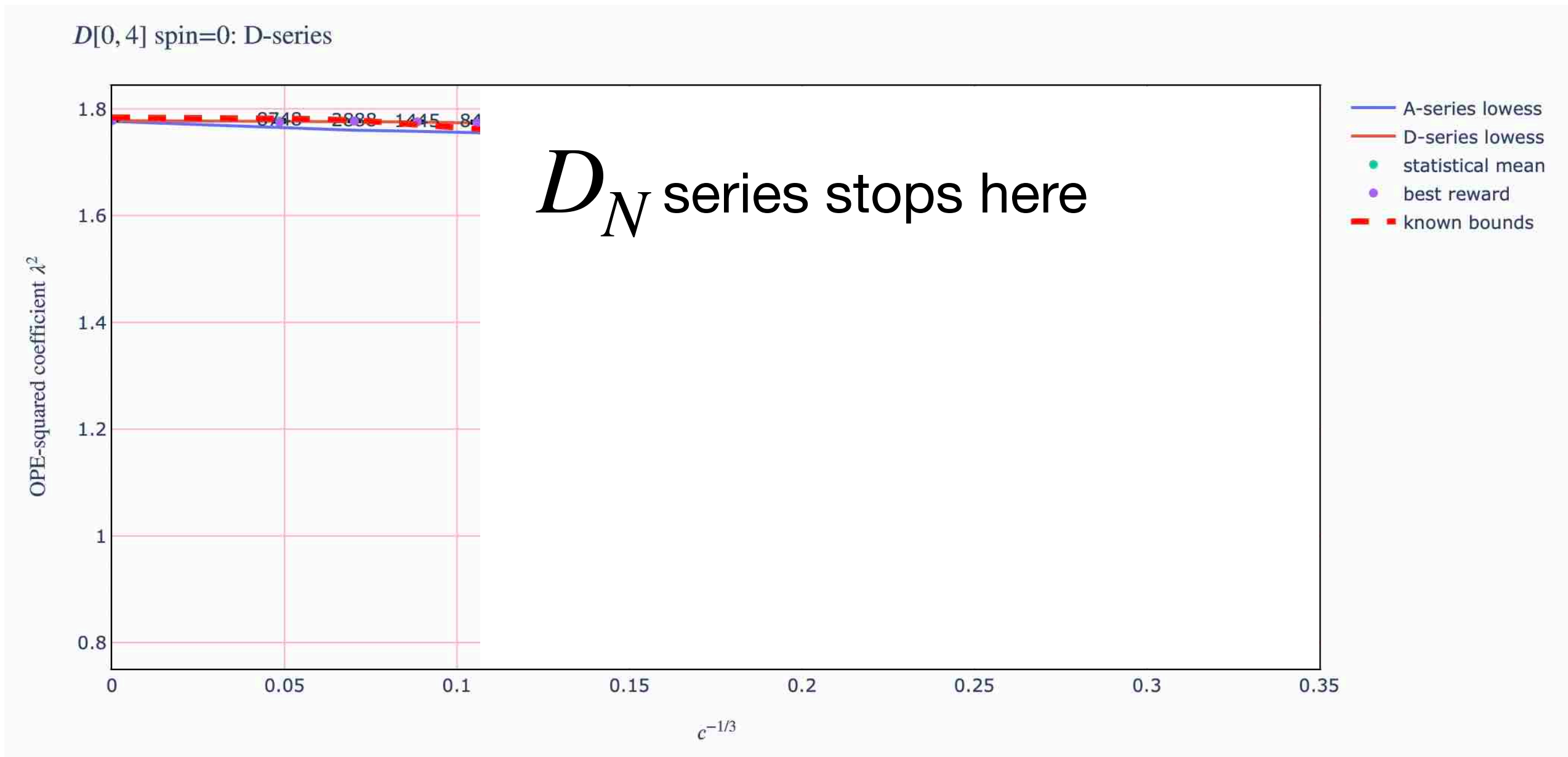
A-series: $c = 25$ to $c = \infty$

D-series: $c = 676$ to $c = \infty$

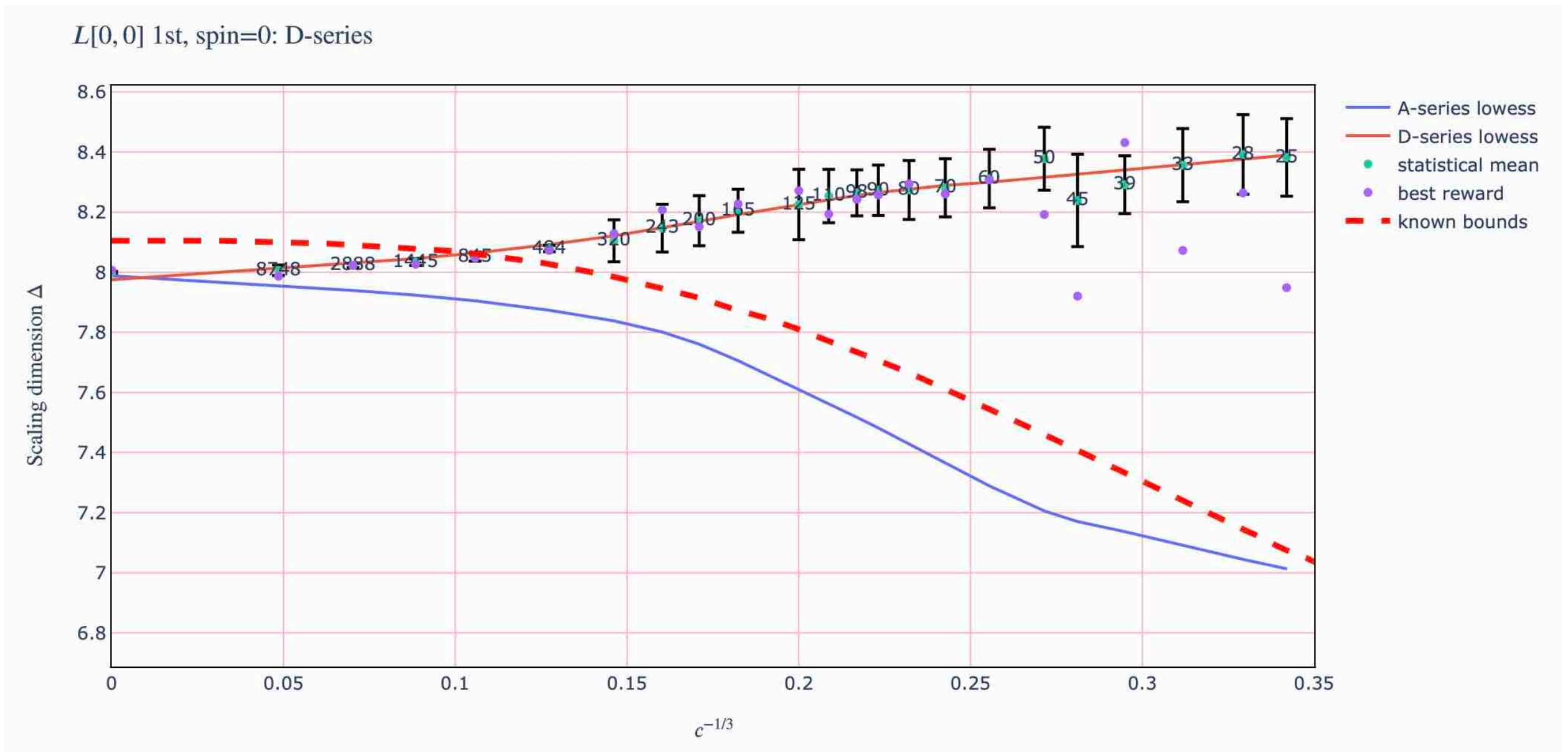
Results (lowest protected multiplet “*D* series”)



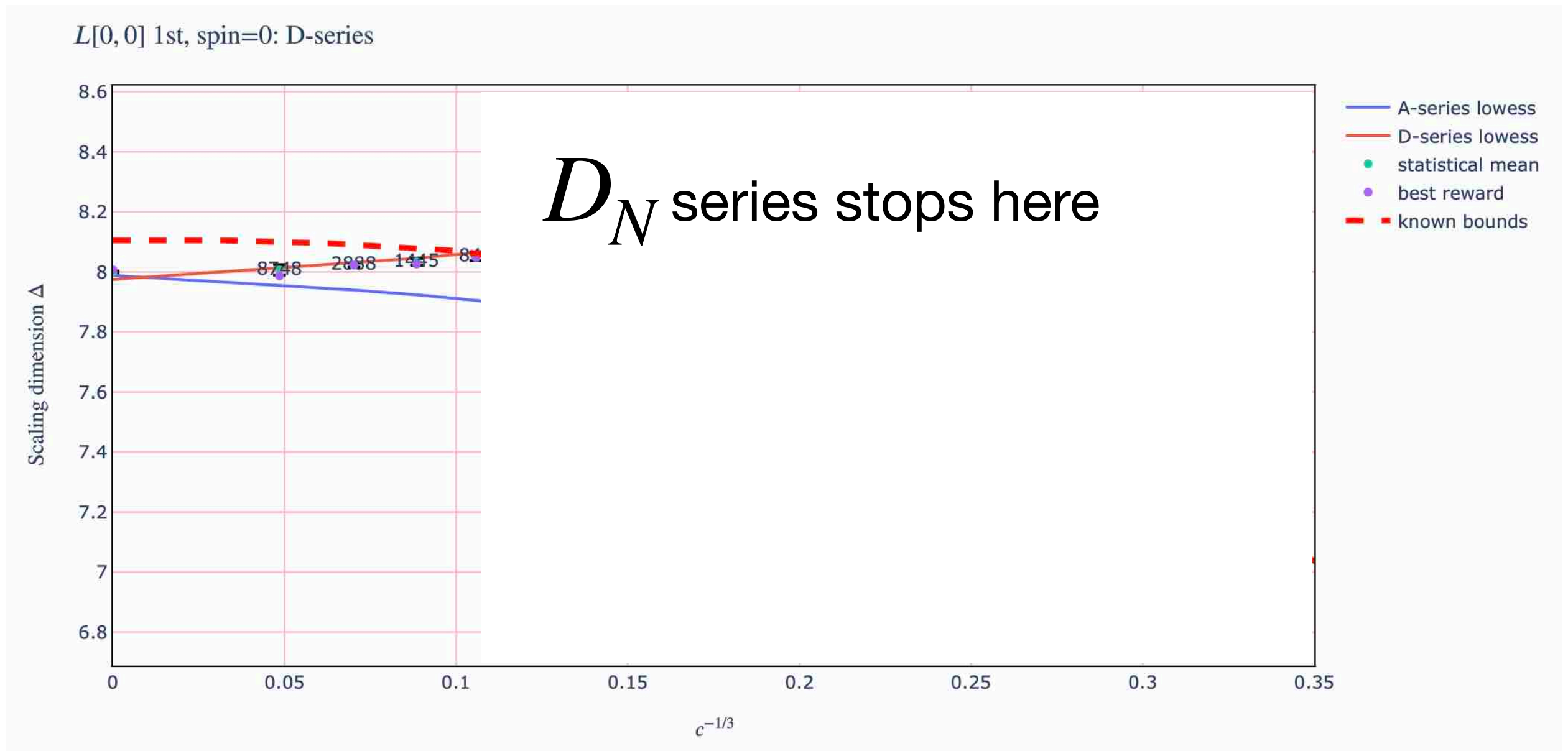
Results (lowest protected multiplet “ D series”)



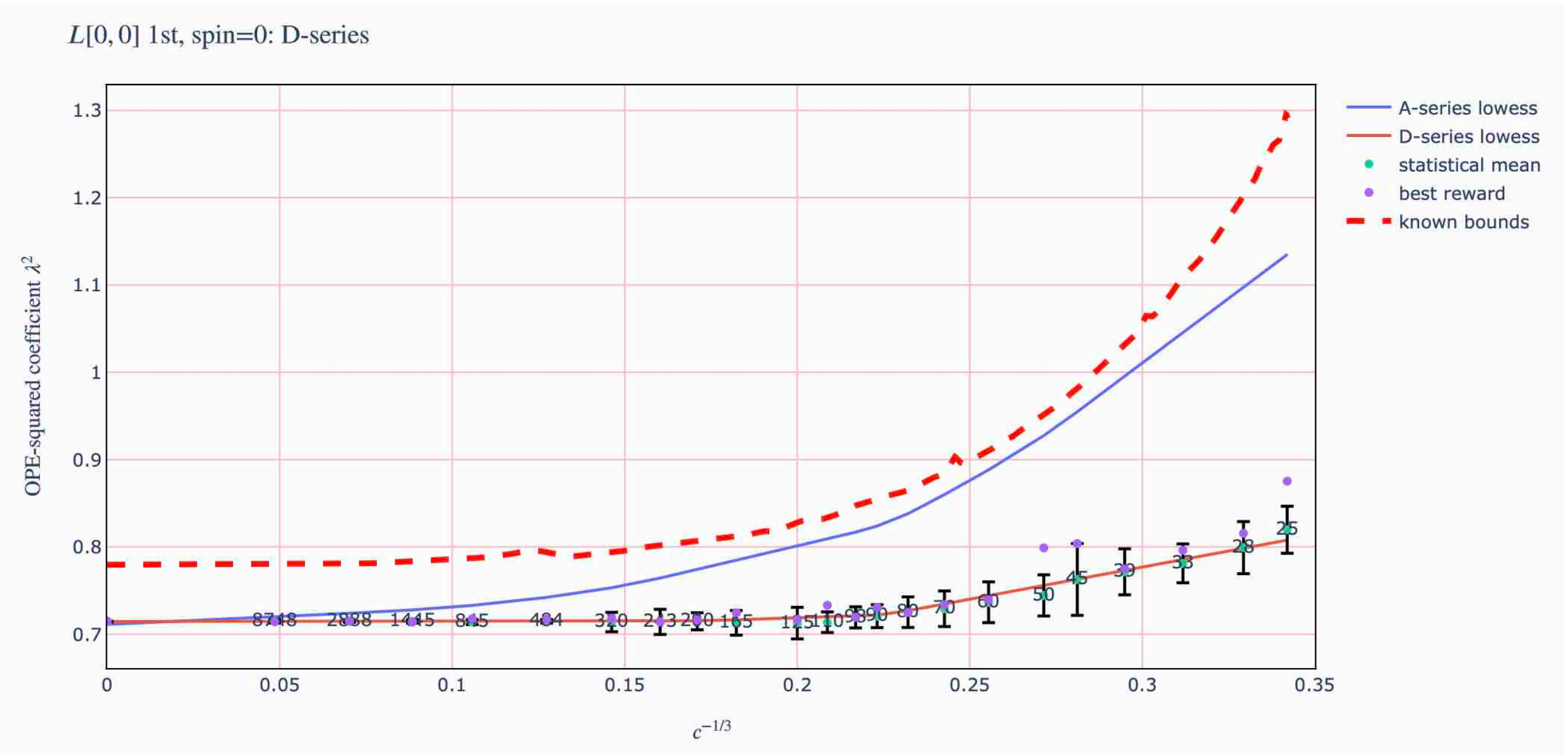
Results (lowest unprotected multiplet “*D* series”)



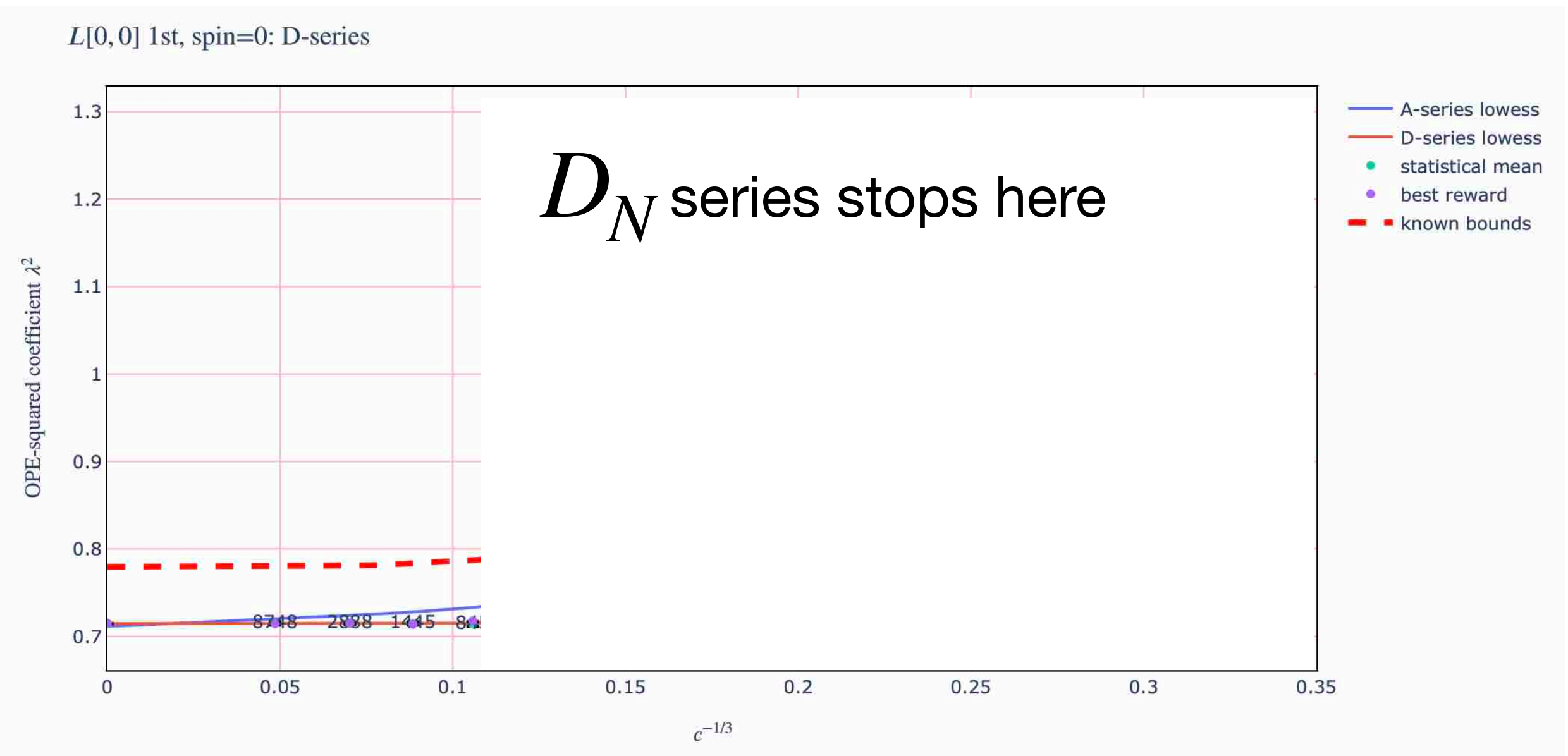
Results (lowest unprotected multiplet “ D series”)



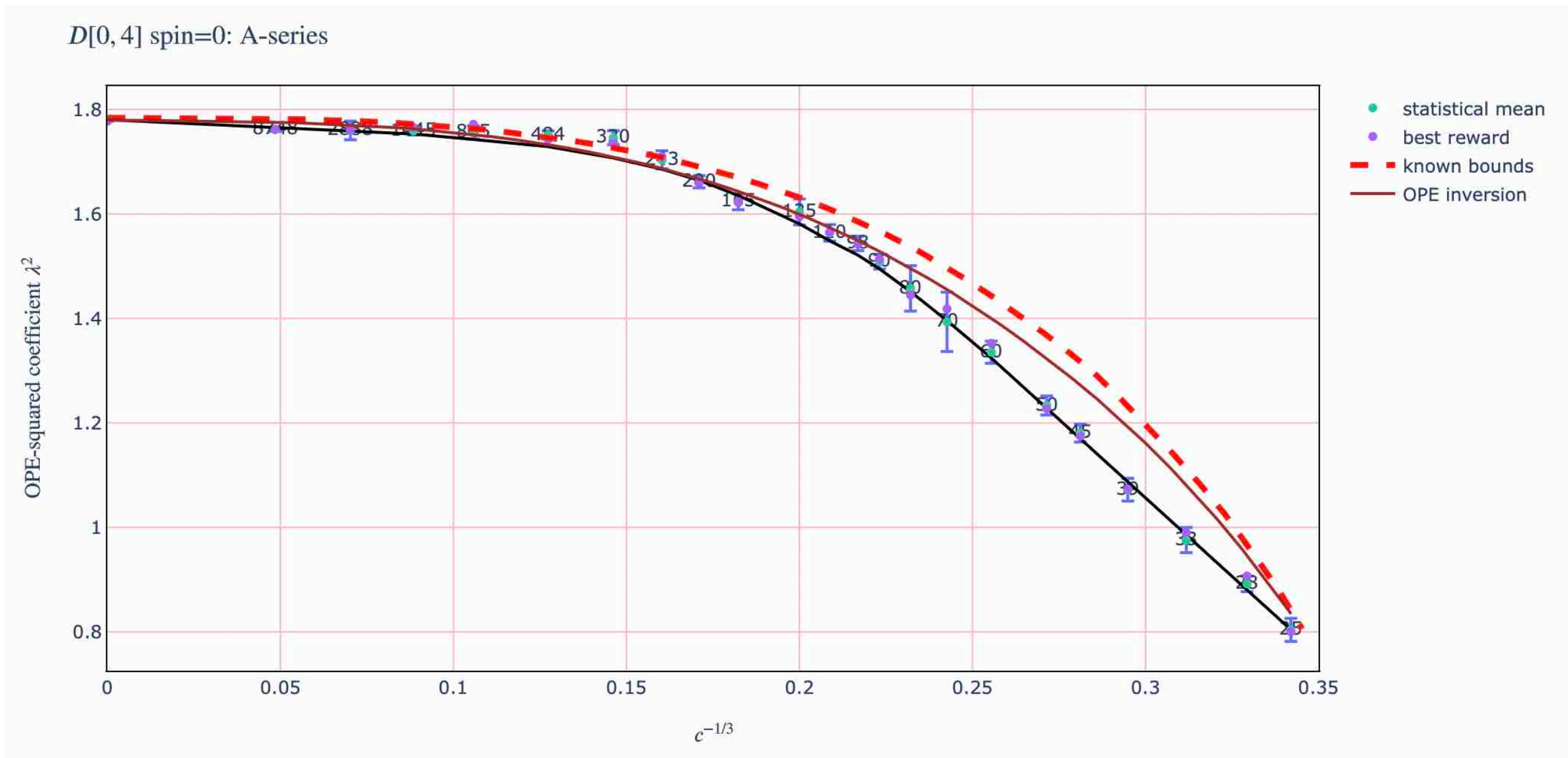
Results (lowest unprotected multiplet “D series”)



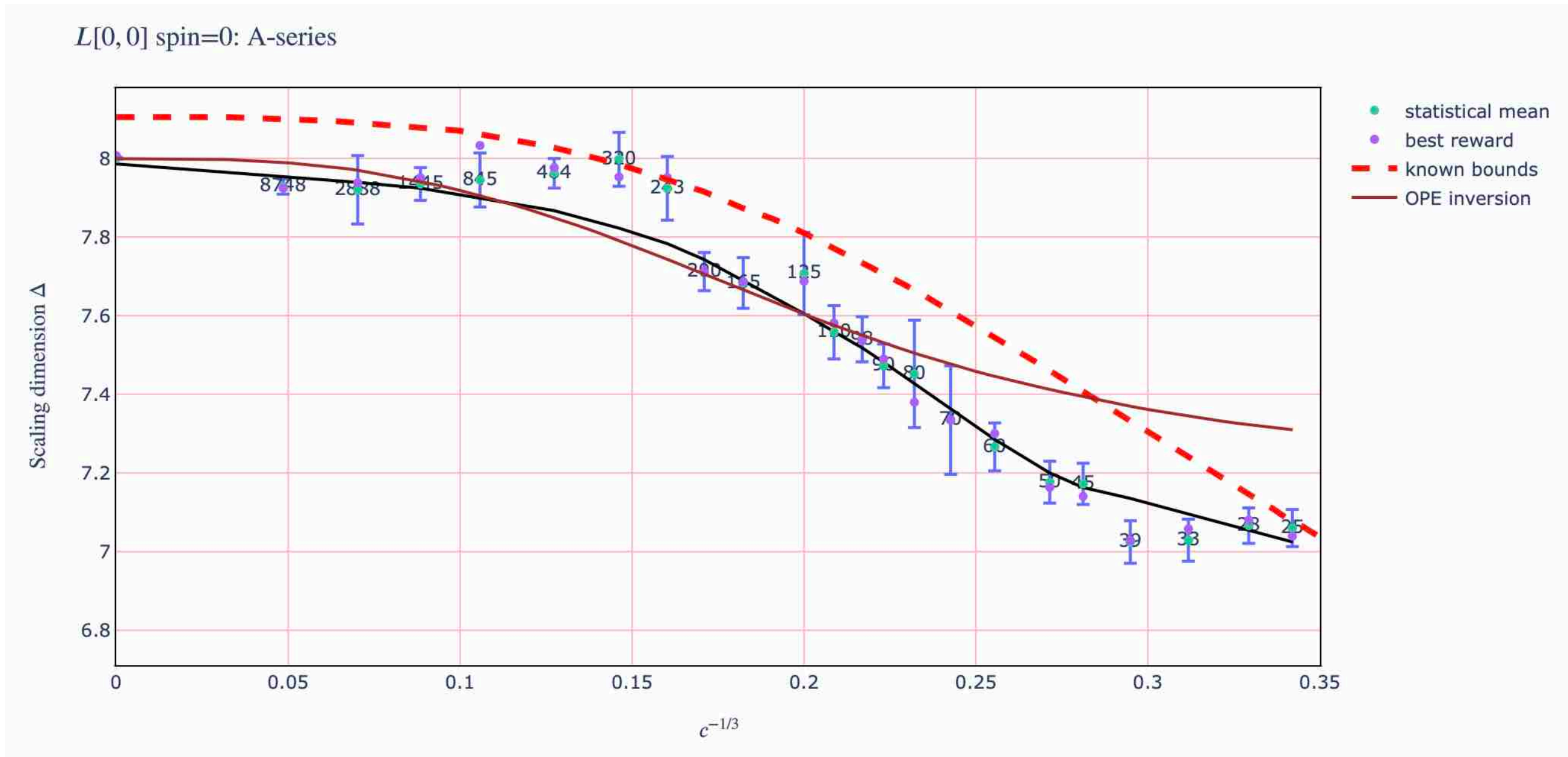
Results (lowest unprotected multiplet “ D series”)



Results (lowest protected multiplet “A series”)



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