How Much More is Different? Large Quantum Numbers and the Modern Correspondence Principle

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I'm excited for the opportunity to give this colloquium on the theme of quantum field theory and the emergence of classical physics from it under certain circumstances.

My subject is the simplification of otherwise-strongly-coupled quantum systems in the limit of large quantum number, which I'll refer to generically as "*J*".

By "otherwise strongly coupled" I'll mean outside of any simplifying limit where the theory becomes semiclassical for other reasons or possibly in a simplifying limit but with the quantum number taken so large that the system behaves differently than you might have expected despite being weakly coupled.

The primary question in such a talk is, is this even a subject?



The answer is, yes, and in some sense it's an old one; many examples have appeared in the literature going far back into the past. Recently there have been a number of groups focusing on systematizing this point of view and applying it more broadly.

Pre-history:

- Atomic hypothesis
- Quantum theory and the correspondence principle



I think a lot of you are from different fields and I'd like as much as possible to give an idea of the important things that are known, so you can understand the context and why this phenomenon is interesting and useful.

This is after all the Institute for the Physics and Mathematics of the UNIVERSE, so I'd like to start by mentioning the fundamental laws that govern the Universe. They are "gravity plus quantum field theory".

The fundamental laws that govern the world – in the absence of gravity – are called the "Standard Model" of particle physics, and they describe a numberof particle types and forces and interactions between particles.

When one is dealing with a few particles at a time, it's generally possible to compute their interactions from first principles, often to incredible accuracy.

For instance, the quantum theory of electrons and their electromagnetic interactions -called "QED" – predicts a tiny quantum correction to the semi-classical magnetic moment of the spinning electron.

Instead of "2" in natural units, quantum corrections computed in QED put the magnetic moment of the electron at $2 \times [1.00115965218073(28)].$

This number is the "anomalous magnetic moment of the electron", whose experimental measurement agrees with the theoretical prediction of QED to 14 digits of precision.

The two digits in parentheses at the end are an experimental , not a theoretical uncertainty.

This level of accuracy – and its less-precise but still comparable extension to other particles and interactions represents a triumph of the basic paradigm of reductionism that many physicists have followed since the early 20th century particularly, when the 2000-year-old atomic hypothesis was stunningly verified in Albert Einstein's least famous paper of 1905.



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But you might want to ask, wait, if the basic building blocks of matter – atoms, then electrons and nuclei, then the more fundamental particles of the Standard Model – were only discovered starting in 1905, how did anything ever get done before then? How was there physics at all?

Well, the natural laws as they were understood prior to the 20th century, are now understood as emergent effective laws resulting from the presence of a large number of particles. Statistical mechanics is a particularly famous example, but there are many other such emergent laws that were studied in the 19th century and much earlier.

When natural laws are "emergent", one recurring theme is that the details of the "fundamental" laws are usually not very important. Maybe one or two or some finite number of details matter for the determination of the natural laws, but most of them are irrelevant. All but a finite number of them should be irrelevant in fact. Otherwise you couldn't have any confidence that any particular law would emerge!

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In relativistic quantum field theory there is a well-developed technology to study emergence, called the "renormalization group", that was developed by many people but most dominantly by Kenneth Wilson. The "renormalization group" is a deceptively, falsely modest name for an extremely radical set of ideas.



The renormalization group is a quantitative theory of how the effective dynamics of a statistical or quantum theory evolve as a function of the distance scale at which the theory is viewed.





In relativistic theories, the speed of light in vacuum is an absolute constant called "*c*", the same for all observers and in all circumstances.

It is part of the definition of the structural relationship between space and time.

Given the absolute speed of light, every distance scale is automatically associated with a time scale, namely the amount of time it takes a light signal takes to travel that particular distance.



This idea is the core idea of special relativity, which was first put forward by Albert Einstein in his third least famous paper of 1905.





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Zur Elektrodynamik bewegter Körper; von A. Einstein.

Daß die Elektrodynamik Maxwells -- wie dieselbe gegenwärtig aufgefaßt zu werden pflegt - in ihrer Anwendung auf bewegte Körper zu Asymmetrien führt, welche den Phänomenen nicht anzuhaften scheinen, ist bekannt. Man denke z. B. an die elektrodynamische Wechselwirkung zwischen einem Magneten und einem Leiter. Das beobachtbare Phänomen hängt hier nur ab von der Reintivbewegung von Leiter und Magnet, während nach der üblichen Auffassung die beiden Fälle, daß der eine oder der andere dieser Körper der bewegte sei, streng voneinander zu trennen sind. Bewegt sich nämlich der Magnet und ruht der Leiter, so entsteht in der Umgebung des Magneten ain elektrisches Feld von gewissem Energiewerte, welches an den Orten, wo sich Teile des Leiters befinden, einen Strom erzeugt. Ruht aber der Magnet und bewegt sich der Leiter, so entsteht in der Umgebung des Magneten kein elektrisches Feld, dagegen im Leiter eine elektromotorische Kraft, welcher an sich keine Energie entspricht, die aber - Gleichheit der Relativbewegung bei den beiden ins Auge gefaßten Fällen vorausgesetzt - zu elektrischen Strömen von derselben Größe und demselben Verlaufo Veranlassung gibt, wie im ersten Falle die elektrischen Kräfte.

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So in relativistic physical theories, the renormalization group describes the way physical theories change when viewed on increasingly longer time scales as well as distance scales.



To explain why this makes the modern renormalization group a particularly powerful idea, I need to mention a third major ingredient, that of quantum mechanics.



Quantum mechanics started out as the idea that a physical system can only contain excitations of energy E in discrete integer multiples of the vibrational frequency ω of the degree of freedom carrying the energy.



This idea was proposed by Max Planck in 1900 as a rather ad hoc and abstract rule to explain the universal spectrum of light emitted from a heated body, which otherwise had no sensible explanation in classical statistical mechanics or thermodynamics.





The rule for the size of the discrete energies is stated as

 $E = n \hbar \omega$

The quantity \hbar is an empirically measured quantity chosen so that the intensity spectrum of emitted light as a function of the frequency of light, fits the measured curve under the assumption that the electromagnetic field only has energies in these "quantized units" $\hbar\omega$.





As radical as it was, Planck's hypothesis was soon vindicated in spectacular form when Einstein used it to explain the photoelectric effect, in the only paper he managed to write in 1905 that was actually adequate enough to earn a Nobel Prize.



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The number *n* expressing the number of minimal units of energy is called the "quantum number" of some particular oscillation.

The quantum number will be a central idea in what I'm going to try to express to you today.

Most all of you know what I've told you very well already, but I'm reviewing it to emphasize that the quantization of energy has profound implications for the notion of emergent laws.



I've so far expressed the quantization of energy as Planck did, in terms of oscillations, but really, when you think about it, everything is an oscillation...





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By Fourier transforming the way in which any particle or field changes in time we can always decompose any notion of dynamics into a linear superposition of oscillators.




Those oscillators have then got to obey the laws of quantization of energy according to Planck's hypothesis, which we now know to be a universal law of nature which is the antecedant to modern quantum theory.





So we needn't think of every degree of freedom in the universe as being literally a harmonic oscillator of definite frequency in order to understand the significance of the quantization of energy...





...we can simply use the law of Fourier analysis, as embodied by the oscilloscope shown here, to understand that all motion is a superposition of oscillations, and the faster the rate of change of a physical quantity, the more high frequency oscillations involved in the frequency decomposition of its motion





...and the greater the amount of energy that motion must carry, since you can never have less than one quantum of energy $\hbar\omega$ involved in an oscillation if it is excited at all .





We can think of this principle as defining a "time-energy complementarity" or a "time-energy uncertainty relation "

A motion or signal of any kind that is very localized in time , necessarily has a Fourier transform that is very spread out in frequency space, and so it necessarily has components of very high energy:





$$< E > \simeq \hbar < \omega > \simeq \hbar \times \frac{1}{\text{[thing]}} \times \frac{d[\text{thing]}}{dt}$$

This rule is completely robust, so long as "[thing]" is any kind of observable physical quantity .

This is actually logically equivalent to saying that if we have a limited amount of energy E to work with, we can never view any physical process with a time resolution that is sharper than $\Delta t \sim \frac{\hbar}{E}$.





All this is very familiar to most/all of you, but the implications for our view of theory space are profound, because energy is a resource – the amount of available energy is always limited.

It tells us that the time scale on which we study our theory is determined by the energy budget with which we do our experiments ... or if we are theorists, our thought experiments .





For ordinary quantum mechanics with a fixed, finite set of degrees of freedom, this does not necessarily lead to very interesting consequences.

But once relativity is added to the conceptual ensemble, the situation changes a lot .





In a relativistic theory, the speed of light provides a universal speed limit and separated objects can never interact instantaneously.

But Hamiltonian dynamics, whether classical or quantum , is always micro-causal , that is, it describes instantaneous time evolution so it must be formulated in terms of local degrees of freedom interacting only with their infinitesimal neighbors .





In other words, relativistic interactions must be formulated in terms of local fields rather than particles with some sort of inter-particle potential :





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$$H = \sum_{i \neq j} \frac{e^2}{|\vec{x}_i - \vec{x}_j|} \leftarrow \text{NO!}$$
$$= \int d^3 \vec{x} \frac{\vec{\mathsf{E}}^2 + \vec{\mathsf{B}}^2}{2} - e A_\mu J^\mu \leftarrow \text{OK}$$

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Another major consequence of relativity and quantum mechanics put together is just that time and space are on an equal footing, just as energy and momentum are on an equal footing.



This doesn't even require special relativity, even just Galilean relativity of inerertial frames is enough to tell you that.



Energy is not invariant under a change of intertial frame... so in order to have a time-energy uncertainty relation hold in every intertial frame, one needs to incorporate a position-momentum uncertainty relation so that the uncertainty relations will transform covariantly :

 $(\Delta E)(\Delta t) \ge \hbar + [\text{relativity}] \Rightarrow (\Delta p)(\Delta x) \ge \hbar$





The relation $(\Delta p)(\Delta x) \ge \hbar$ is implemented by de Broglie in modern quantum theory by the precise replacement





$$p \rightarrow -i\hbar \frac{\partial}{\partial x}$$

and the the time-energy uncertainty relation $(\Delta E)(\Delta t) \ge \hbar$ is implemented as the Schrödinger equation

$$H = +i\hbar \frac{\partial}{\partial t} = \frac{p^2}{2m} + V(x) = -\frac{\hbar^2}{2m} \frac{\partial^2}{\partial x^2} + V(x)$$

acting on wavefunctions $\psi(x,t)$



This leads to all sorts of well-known and fascinating effects like matter particles becoming waves and interfering with themselves.





All this part holds even without special relativity, since as I said, even Galilean relativity suffices to derive the energy-momentum uncertainty principle.



The promotion of matter to waves looks much more "natural" in special relativity, when we are forced to turn particles into fields in order to for the interactions to respect locality and causality.







The photon γ was already understood as the quantum of the electromagnetic field $\gamma \rightarrow A_{\mu} = (-\phi, \vec{A})$ as argued by Einstein by way of the photoelectric effect...





So it was inevitable that the electron would be promoted to a quantum of the electron field $e^- \rightarrow \psi_{\alpha}(x, t)$ as eventually understood by Dirac, Pauli and others, and other matter fields followed.

From these ideas quantum field theory (QFT) was invented.



 QFT is a theoretical structure that incorporates all the theoretical priors of:

- Causality
- Quantum mechanics, and
- Lorentz invariance.



From these it automatically follows that all processes have a probability amplitude for particle creation if allowed by conservation of energy, because $E = mc^2$.





The rule for "doing" QFT is easy to state, thanks to Feynman's formulation.



Take any classical Lagrangian density that you would use to define a classical field theory by the principle of least action .

For instance, if you want a quantum theory of Maxwell's equations, you would start with the classical Lagrangian for Maxwell's equations, $\mathcal{L} = -\frac{1}{4}F_{\mu\nu}F^{\mu\nu} + eA_{\mu}J^{\nu}$.

Then integrate it over spacetime and you get the action:

$$S=\int d^D x \mathcal{L}$$
.

The key input in the theory is the local Lagrangian density.

Then if you wanted to do classical field theory you would do differential functional calculus and take its functional derivative to get the Euler-Lagrange equations which give you Maxwell's equation:

$$\partial_{\mu}F^{\mu
u} = -e\,J^{
u}$$

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OK so that's classical field theory.

If you want to do quantum field theory then you just do integral functional calculus instead.

That is, instead of differentiating with respect to every possible direction in field space and setting the functional derivative to zero -

Instead you integrate over every possible direction in field space, with the weight given by the exponential of $\frac{i}{\hbar}$ times the action :

[probability amplitude for anything] = $\int \mathcal{D}[\frac{\text{field}}{\text{configurations}}] \exp\left(\frac{i}{\hbar}S\right)$.

And that's it! In Feynman's formulation that's literally (in principle) all you need to know, about the rules of quantum mechanics or quantum field theory.

In practice there are various ambigutities:

- You don't know what the infinite-dimensional measure *D*[field configurations] is;
- The path integral has all kinds of divergences at high energies and you don't know how to cut them off or what the cutoff means ; and
- > You dont know what the local Lagrangian density actually is !

Fortunately these ambiguities are all the same thing .

The measure $\mathcal{D}[\text{field configurations}]$ is defined by cutting off the effects of the short distance/time degrees of freedom, for instance by discretizing space and time or something with the same effect as that.

Since short distance/time equals high momentum/energy in quantum theory, this cuts off the high energy divergences at a cut-off energy Λ . (These are called "ultraviolet divergences".)

The how of the cut-off doesn't matter , because any (sufficiently "local") measure $\mathcal{D}[{\rm fields}]$ for field configurations is equivalent to any other , up to equivalent local terms in the lagrangian density.

Quantum field theory

That is, any change of cutoff procedure

 $[{\rm one \ kind \ of \ cut} - {\rm off \ at \ energy} \ \sim \Lambda] \Rightarrow$

[another kind of cut – off at energy $\sim \Lambda$]

is equivalent to a change of measure

 $d[\text{field configurations}] \rightarrow d[\text{field configurations}]$

where the change of measure can be compensated by a change of the Lagrangian density by local terms with their coefficients given by dimensional analysis in terms of powers of the cutoff :

 $e^{\frac{i}{\hbar}S} d$ [field configurations] = $e^{\frac{i}{\hbar}S} d$ [field configurations],

 $\mathcal{L} \Rightarrow \widetilde{\mathcal{L}}$

 $= \mathcal{L} + [\text{sum of local terms with coefficients as powers of } \Lambda]$

These compensating terms are called counterterms and they just encode the common-sense fact that you never actually knew what your theory actually was at unlimitedly short distances anyway, due to your limited resources of energy.

Quantum field theory

That's renormalization in a nutshell, and it's just not the big deal people used to think it was.

If there is a prescription for defining the theory with the energy cutoff Λ taken all the way up to infinity, the theory is called "renormalizable" .

Otherwise it is known as an "effective [quantum] field theory" or EFT , with a finite energy cutoff Λ .

What makes QFT forbidding is not the metaphysical issues of renormalization, it's the practical issue that QFT, unlike quantum mechanics, has an operator algebra and Hilbert space generated by arbitrarily many degrees of freedom as you increase your energy budget.

Much modern research in quantum field theory is devoted not to "taming ultraviolet divergences" but to parametrizing and understanding this huge complex jungle of theories and behaviors of theories.

In short, we want to map out the large-scale structure of theory space and the gross structure of behaviors within each theory, and hopefully organize theories into families with helpfully strong family resemblances of some kind.

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The most helpful organizing tools are symmetries .

Quantum field theory

- In classical mechanics, symmetries are merely descriptive.
- But in quantum mechanics, every symmetry describing the laws of nature is associated with a conservation law.
- ► For instance, the fact that the laws of Nature look the same everywhere , implies the law of conservation of momentum .
- Similarly, the fact that the laws of nature look the same no matter what direction you are facing, is associated with conservation of angular momentum.
- Conserved quantities in quantum mechanics –
- typically come in integer multiples of some minimum amount .
- For instance, angular momentum in quantum mechanics is famously quantized in units of (half of) Planck's constant ħ.

- The relation between symmetries and conservation laws actually goes back to classical mechanics in the Hamiltonian formulation.
- There, the relation between observables and operations is reflected in the structure of the Poisson bracket { , }, where every observable on phase space implements an infinitesimal operation on the system –
- just Heisenberg's famous commutator does in quantum mechanics.

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Symmetries in Quantum Theory

- The key difference from quantum theory lies in the word "infinitesimal" when discussing the Poisson bracket.
- In contrast to the Poisson bracket, extracting information via an observation in quantum mechanics disturbs the system by a minimum finite amount.
- ▶ When the observation corresponds to a symmetry operation -
- such as rotating the system or moving it in some direction
- one always disturbs the system by a minimum finite amount, due to Heisenberg's uncertainty principle.
- ▶ This is expressed by the replacement of the Poisson bracket by the algebraic commutator $\{ \ , \ \} \rightarrow [\ , \]$ when the system is quantized .

- You might wonder what happens to the quantization of angular momentum in everyday life.
- Well, the typical amount of angular momentum is so huge that the quantization is invisible:

 $J=N\,\hbar$

where N is Avogadro's number or something.

Nowadays physicists sometimes refer to this as the recovery of classical physics in the macroscopic limit.

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To show the power of symmetries to help cut through the complexity of QFT, let us focus on a theoretical structure that incorporates all the theoretical priors of quantum field theory,

Causality

- Quantum mechanics
- Lorentz invariance, and,

plus one more:

Scale invariance

Such a structure is called a conformal field theory (CFT).
Scale invariance is not a symmetry of our world!

Scale invariance is not a symmetry of our world!





Scale invariance is not a symmetry of our world!



Nonetheless it plays many important roles in our understanding of theoretical physics.



Most importantly , CFT is a consistency condition for general QFT at short distance



as illustrated by scaling behavior in QCD at high energies.

Conformal field theory is a major focus of my own research!



Given the significance of conformal field theory, we should know more about the space of possibilities.



Conformal field theories can contain arbitrarily complicated phenomena, including chaos...



...thermalization...



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including phenomena that are exactly mathematically equivalent to those of black holes.



The best-studied CFT – the exactly solveable CFT – contain none of these phenomena, and comprise an infinitesimal and non-representative subclass of CFT

No chaos or thermalization!

We would like to extract information about the generic case.

If we want to use the rules of CFT, we ought to explain what a $\ensuremath{\mathsf{CFT}}$ is.

A CFT is an object defined by a set of local operators $\mathcal{O}_i(z)$ and an operator product expansion.

$$\mathcal{O}_i(z_1)\cdot\mathcal{O}_j(z_2)=\sum_k f_{ij}{}^k(z_1,z_2)\mathcal{O}_k(z_2) ,$$

including the identity $\mathcal{O}_0 = 1$.

These local operators define a set of expectation values such that the OPE is satisfied inside the expectation value.

$$\langle \mathcal{O}_{i_1}(z_1)\mathcal{O}_{i_2}(z_2) \cdot (\text{other ops}) \rangle = \sum_j f_{i_1i_2}^{\ j}(z_1, z_2) \langle \mathcal{O}_j(z_3) \cdot (\text{other ops}) \rangle$$

This product is taken to be associative, and the expansion is convergent, for z_1 sufficiently close to z_2 .

One of these operators is taken to be the stress tensor T_{ab} .

Furthermore the theory is taken to be defined on an arbitrary^{*} manifold M with an arbitrary ^{*} background geometry g_{ab} .

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CFT basics

Finally the stress tensor is given by the variation of the theory with resepect to g_{ab} :

$$\left\langle T^{ab}(z) \cdot (\text{operators}) \right\rangle = \frac{\delta}{\delta g_{ab}} \left\langle (\text{operators}) \right\rangle$$

For a theory depending only on the conformal structure, and not on the local scale, the stress tensor must be traceless: $T_a{}^a = 0$.

In particular, our expectation values depend only on the intrinsic geometry and topology and not on the coordinate system.

The invariance under infinitesimal coordinate transformations is equivalent to the condition that the stress tensor is conserved, $\nabla^b T_{ab} = 0$, and the invariance under coordinate transformations not connected to the identity is referred to as modular invariance.

The basic rules of CFT are actually very constraining

The associativity of the operator algebra alone, together with conformal invariance and unitarity, turns out to impose a huge number of consistency conditions on a correlation function.

In a (very) few cases these constraints are so severe as to allow a highly precise numerical solution for the amplitude .

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This enterprise is known as the conformal bootstrap.

- ► The goals of the large quantum number expansion are largely to answer the same questions as the conformal bootstrap:
- Learn to systematically and efficiently analyze QFT (in practice usually CFT) that have no exact solution in terms of explicit functions.

- We'd all like to know "what does theory space look like": Generic theories, generic amplitudes.
- This is a very consequential question for field theory, mathematics, quantum gravity, and cosmology.
- Most theories are not integrable, and we need to learn how to attack them in general circumstances.
- "Direct" numerical bootstrap methods are remarkably efficient, power-law in number of operators exchanged in the amplitude.

BUT...

Critique of Pure Bootstrap

- Since number of operators grows exponentially with dimension / central charge / other quantum number, direct numerical attack is still intractable in extreme limits.
- Fortunately, known "extreme limits" appear to have simplifying behaviors in many (all known?) known circumstances. This is broadly a generalization of the notion of "duality".
- In the case of large spin in a single plane, the limit has been analyzed within the bootstrap itself.
- The relative ease of this is related to the fact that the spacetime coordinates themselves carry the quantum number.
- For other quantum numbers, this is not the case. For instance, there is no known analytic bootstrap method to attack the case of large spin in multiple planes in D ≥ 4.

- In many cases such limits are accessible to some new kinds of EFT in regions where bootstrap methods slow down.
- As we'll see, there's also a excellent agreement where the two methods overlap.
- Where does this leave us? What do we hope to accomplish ?

- (*) Most modestly: Translate EFT behavior into bootstrap terms, say what it means for CFT data. Operator dimensions and OPE coefficients.
- (***) Most grandiosely: Derive EFT behavior from bootstrap equations, and use it to solve everything in every limit where direct numerical methods break down.
- (**) Intermediate: Use some small subset of EFT inputs, and obtain some subset of CFT data not directly numerically accessible.
- Grandiose goal (***) appears out of reach for now. (I tried!)

So now I'll tell you about some of our progress on modest goal (*).

Large charge J in the O(2) model

- Simplest example: The conformal Wilson-Fisher O(2) model at large O(2) charge J.
- ► This is a complex scalar field ϕ in D = 3 with potential $V(\phi) = g^2 |\phi|^4 m^2 |\phi|^2$ with g being taken to infinity but m being tuned so that the scalar field stays massless in spite of quantum effects.
- ► Canonical question: What is the dimension Δ_J of the lowest operator O_J at large J?
- Translated via radial quantization: Energy of lowest state of charge J on unit S²?
- ► Renormalization-group analysis reveals the low-lying large-charge sector is described by an EFT of a single compact scalar χ , which can be thought of as the phase variable of the complex scalar $\phi = |\phi|e^{i\chi}$.

Large charge J in the O(2) model

The leading-order Lagrangian of the EFT is remarkably simple:

$$\mathcal{L}_{ ext{leading-order}} = b |\partial \chi|^3$$

- The coefficient b is not something we know how to compute analytically; nonetheless the simple structure of this EFT has sharp and unexpected consequences.
- The immediate consequence of the structure of the EFT is that the lowest operator is a scalar, of dimension

$$\Delta_J \simeq c_{rac{3}{2}} J^{rac{3}{2}}$$

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where $c_{\frac{3}{2}}$ has a simple expression in terms of *b*.

- ► The leading-order EFT predicts more than just the leading power law, because quantum loop effects in the EFT are suppressed at large J, so the EFT can be quantized as a weakly-coupled effective action with effective loop-counting parameter J^{-3/2}.
- For instance we can compute the entire spectrum of low-lying excited primaries.
- ► The dimensions, spins, and degeneracies of the excited primaries, are those of a Fock space of oscillators of spin *l*, with *l* ≥ 2.

Large charge J in the O(2) model

- The propagation speed of the χ-field is equal to ¹/_{√2} times the speed of light.
- So the frequencies of the oscillators are

$$\omega_\ell = rac{1}{\sqrt{2}} \sqrt{\ell(\ell+1)} \;, \qquad \qquad \ell \geq 1 \;.$$

- The ℓ = 1 oscillator is also present, but exciting it only gives descendants; the leading-order condition for a state to be a primary is that there be no ℓ = 1 oscillators excited.
- ► So for instance, the first excited primary of charge J always has spin $\ell = 2$ and dimension $\Delta_J^{(1)} = \Delta_J + \sqrt{3}$.

- Subleading terms can be computed as well.
- ► These depend on higher-derivative terms in the effective action with powers of $|\partial \chi|$ in the denominator .
- These counterterms have a natural hierarchical organization in J:

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- At any given order in derivatives, there are only a finite number of such terms.
- As a result, at a given order in the large-J expansion, only a finite number of these terms contribute.
- Since there are far more observables than effective terms, there are an infinite number of theory-independent relations among terms in the asymptotic expansions of various observables.

► Our gradient-cubed term is the only term allowed by the symmetries at order J³/₂, and there is only one other term contributing with a nonnegative power of J, namely

$$\mathcal{L}_{j^{+\frac{1}{2}}} = b_{\frac{1}{2}} \left[\left| \partial \chi \right| \texttt{Ric}_3 + 2 \frac{(\partial \left| \partial \chi \right|)^2}{\left| \partial \chi \right|} \right]$$

▶ In particular, there are no terms in the EFT of order J^0 , with the result that the J^0 term in the expansion of Δ_J is calculable, independent of the unknown coefficients in the effective lagrangian.

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$$\Delta_J = c_{\frac{3}{2}} J^{+\frac{3}{2}} + c_{\frac{1}{2}} J^{+\frac{1}{2}} -0.0937256\cdots$$

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$$\Delta_J = c_{\frac{3}{2}} J^{+\frac{3}{2}} + c_{\frac{1}{2}} J^{+\frac{1}{2}}$$

- This universal term and the other universal large-J relations in the O(2) model don't have any fudge factors or adjustable parameters;
- Given the identification of the universality class, these values and relations are universal and absolute;

Similar predictions have been made for OPE coefficients
• You might think that there is something "weird" or "inconsistent" or "uncontrolled" about a Lagrangian like $\mathcal{L} = |\partial \chi|^3$.

So, let me anticipate some frequently asked questions:

- Q: Isn't this Lagrangian singular?? It is a nonanalytic functional of the fields, so when you expand it around χ = 0, you will get ill-defined amplitudes.
- A: Yes, but you aren't supposed to use the Lagrangian there. It is only meant to be expanded around the large charge vacuum, which at large J is the classical solution

 $\chi = \mu t$,

with

$$\mu = O(\sqrt{\rho}) = O(J^{\frac{1}{2}}) \; .$$

► The expansion into vev and fluctuations carries a suppression of µ⁻¹ or more for each fluctuation.

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 (parenthetical comment:) There are already many well-known effective actions of this kind, including the Nambu-Goto action.



- Q: Isn't this effective theory ultraviolet-divergent ? That means that loop corrections are incalculable and observables are meainingless beyond leading order.
- A: No. The EFT is quantized in a limit where loop corrections are small. Our UV cutoff Λ for the EFT is taken to satisfy

$$E_{\rm IR} = R_{\rm S^2}^{-1} \quad \ll \quad \Lambda \quad \ll E_{\rm UV} = \sqrt{\rho} \propto J^{+rac{1}{2}} R_{\rm S^2}^{-1}$$

► Loop divergences go as powers of ³/ρ³/₂ ≪ 1, and are proportional to nonconformal local terms which are to be subtracted off to maintain conformal invariance of the EFT.

- Q: OK but then don't the counterterms ruin everything? Don't they render the theory incalculable?
- A: No. As usual in EFT the counterterm ambiguities of subtraction correspond one-to-one with terms in the original action allowed by symmetries;
- As we've mentioned there are only a finite and small number of those contributing at any given order in the expansion, and at some orders there are no ambiguities at all.

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- Q: You're saying that every CFT with a conserved global charge has this exact same asymptotic expansion. But here's a counterexample! (describes theory SH didn't say anything about)
- A: I didn't make any claim that broad. Our RG analysis applies to many but not all CFT with a conserved global charge. More generally, CFT can be organized into large-charge universality classes.
- ► For instance, free complex fermions as well as free complex scalars in D = 3 are in different large-J universality classes.
- The large-J universality class of the O(2) model contains many other interesting theories, such as
 - The $\mathbb{CIP}(n)$ models at large topological charge ;
 - The D = 3, N = 2 superconformal fixed point for a chiral superfield with W = Φ³ superpotential, at large *R*-charge;
 - Probably others o o o

Other large-J universality classes

- Many other interesting universality classes in D = 3:
- ► Large Noether charge in the higher Wilson-Fisher *O*(*N*) [Alvarez-Gaumé, Loukas, Reffert, Orlando 2016] and *U*(*N*) models;
- ► Also the CIP(n) [de la Fuente] and higher Grassmanian models real and complex ; [Loukas, Reffert, Orlando 2017]
- Large baryon charge in the SU(N) Chern-Simons-matter theories;
- Large monopole charge in the U(N) Chern-Simons-matter theories;
- Of course these last two are dual to one another and would be interesting to investigate.

- I didn't say anything yet about supersymmetry but you can think of it as just a very nice, constraining type of symmetry that relates fermions and bosons.
- For conformal supersymmetric theories in D = 4 there is always at least a continuous global symmetry commuting nontrivially with the super-generators, called an R -symmetry.
- Often theories with SUSY have non-unique ground states, even non-unique up to symmetry rotation.
- These are said to have moduli spaces of vacua or vacuum manifolds.
- Among the most tractable universality classes are large R-charge in extended superconformal theories with moduli spaces of supersymmetric vacua.

► Simplest case is the N = 2, D = 3 superconformal fixed point of three chiral superfields with superpotential W = XYZ.

- Its vaccum manifold has three one-complex-dimensional branches: X, Y, Z ≠ 0.
- ► WLOG consider the X-branch.

- The X-branch has coordinate ring spanned by X^J , $J \ge 0$.
- These BPS scalar chiral primary operators are the (X-branch part of the) chiral ring of the theory.
- ► The dimension of X^J is exactly equal to its R-charge J and protected from all quantum corrections: In this case the formula for the dimension Δ_J is boring :

$$\Delta_J = 1 \cdot J \quad \Leftarrow \quad \text{BORING!}$$

► The formula for the dimension of the second-lowest primary of J_R = J_X = J is also boring; it lies an a protected scalar semishort representation with only 12 Poincaré superpartners:

 $\Delta_J^{(+1)} = 1 \cdot J + 1 \qquad \qquad \Leftarrow \quad \text{also boring!}$

Nonetheless we would like to see this explicitly in a large-J expansion, and also be able to compute non-protected large-J quantities such as third-lowest operator dimensions and also OPE coefficients.

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- ► The effective theory describing the lowest state of $J_X = J_R = J$, is simply the moduli space effective action , appearing in the same role as the gradient-cubed theory for the O(2) model.
- Unlike the O(2) model EFT, here the leading effective action is simply free :

$$\mathcal{L} = \int d^2 \theta \, d^2 \overline{\theta} \, \Phi^{\dagger} \Phi \;, \qquad \qquad \Phi = (\text{const.}) \times X^{\frac{3}{4}} + \cdots \;,$$

where the · · · are higher-derivative D-terms .

- ► To compute operator dimensions, quantize the theory around the lowest classical solution with given large J on an S² spatial slice:
- Here, the classical solution is

 $\phi = \mathbf{v} \exp\left(i\mu t\right) \;,$

$$\mu = rac{1}{2R} \; , \qquad \qquad \mathbf{v} = \sqrt{rac{J}{2\pi R}} \; .$$

Note here the frequency of the solution (chemical potential) is determined by supersymmetry (the BPS bound on operator dimensions) rather than the unknown coefficients in the Lagrangian.

The results of the direct diagrammatic quantization are as follows, for the lowest and second-lowest states:

$$\Delta_J = J$$

 $+0 \times J^{0} + 0 \times J^{-1} + 0 \times J^{-2} + 0 \times J^{-3}$ $+O(J^{-4}) \qquad \Leftarrow \text{ three loops!}$

 $\Delta_J^{(+1)} = J + 1 \times J^0$ +0 × J⁻¹ + 0 × J⁻² + 0 × J⁻³ +O(J⁻⁴) \Leftarrow two loops!,

confirming the predictions of supersymmetry to the order we can calculate .

The third-lowest primary is a non-BPS scalar, with dimension

$$\Delta_J^{(+2)} = J + 2 \cdot J^0$$

 $+0 \times J^{-1} + 0 \times J^{-2}$

 $-\kappa \times 192 \,\pi^2 \times J^{-3}$

 $+O(J^{-4}) \quad \leftarrow \text{ one loop! },$

where κ the coefficient of the leading interaction term in the *EFT*.

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The form of the leading interaction term is a D-term, consisting of a four-derivative bosonic component

$$\mathcal{L}_{-1} \equiv +4 \kappa_{\mathrm{FTP}} \, \frac{|\partial \phi|^4}{|\phi|^6} \; ,$$

plus conformally and superconformally completing terms worked out by many authors $% \left({{{\left({{{{{\bf{n}}}} \right)}_{i}}}_{i}}} \right)$.

We don't know the value of κ for the XYZ model, but we do know its sign :

 $\kappa > 0$ (superluminality constraint)

So the first nonprotected operator dimension gets a contribution of order J⁻³ with a negative coefficient of unknown magnitude .

- It is more fun to compute quantities which are both nontrivial in the large-J expansion and checkable in principle by exact supersymmetric methods.
- One nice example is the two-point functions of chiral primary operators in 8-supercharge theories.
- ► The technically simplest class of examples are the chiral primaries spanning the Coulomb branch chiral ring in D = 4, N = 2 theories, in the special case the gauge group has rank one.

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Examples include

- $\mathcal{N} = 4$ SYM with G = SU(2),
- $\mathcal{N} = 2$ SQCD with $N_c = 2$, $N_f = 4$,
- Many rank-one nonlagrangian Argyres-Douglas theories with one-dimensional Coulomb branch,
- including the recently discovered $\mathcal{N} = 3$ examples.
- Some of these are Lagrangian theories with marginal coupling, and some of them are non-Lagrangian theories with more abstract descriptions, but we can treat them all on an equal footing.

The Coulomb branch chiral ring in a rank-one theory is spanned by

$$\mathcal{O}_{\mathcal{J}} \equiv \mathcal{O}^n_{\Delta} \;, \qquad \qquad \mathcal{J} = n\Delta \;,$$

where the^(*) generator \mathcal{O}_{Δ} of the chiral ring has $U(1)_R$ -charge $J_R = \Delta$.

- (*) This assumes the chiral ring is freely generated; there are no known counterexamples, but see recent work for counterexamples in higher rank.
- At large charge in radial quantization these correspond to classical solutions on the sphere where the Coulomb branch scalar \hat{a} gets a vev proportional to $\sqrt{J/R}$.

- ► For Lagrangian theories the generator \mathcal{O} is $\operatorname{tr}(\hat{\phi}^2)$ and $\Delta = 2$.
- ► For non-Lagrangian theories the dimension △ of the generator can take certain other values.
- ► These are constrained to some extent and recently it was proven that △ is always rational
- We can write the large-J effective action in terms of an effective field φ ≡ (O_Δ)^{1/Δ}. The singularity in the change of variables is invisible in large-J perturbation theory because the quantum state field is supported far away from φ = 0.

- ► The leading-order action is again the free action for φ, and the leading interaction term is the anomaly term compensating the difference in Weyl a- anomaly and U(1)_R-anomalies between the underlying interacting SCFT and the free vector multiplet.
- The leading interaction term is

$$\mathcal{L}_{\mathrm{anom}} \equiv \pmb{lpha} \, \int \, d^4 \theta \, d^4 ar{ heta} \log(\phi) \log(ar{\phi})$$

+(curvature and $U(1)_{\rm R}$ connection terms),

where the coefficient α is proportional to the Weyl-anomaly mismatch:

$$\alpha = +2 (a_{\rm CFT} - a_{\rm EFT})^{\rm [AEFGJ units]}$$

- Some comments on this interaction term:
- It was first written down by as the unique four-derivative term in the Coulomb branch EFT of an N = 2 gauge theory;
- ► It is formally an N = 2 D- term, i.e. a full-superspace integrand ···
- ▶ ... but only formally, since it is non-single-valued; its single-valued version can be obtained as an F -term, *i.e.* an integral over only the θ 's and not the $\overline{\theta}$'s.
- Its bosonic content comprises the famous Wess-Zumino term for the Weyl a-anomaly that was used to prove the a-theorem in four dimensions.
- This is why its coefficient α is proportional to the a-anomaly mismatch.

- One other remarkable fact about rank-one theories, is that the anomaly term is that it is unique as a (quasi-)*F*-term on conformally flat space.
- That is, there are an infinite number of higher-derivative D-terms, but there are no higher-derivative *F*-terms one can construct out of a single vector multiplet in a superconformal *N* = 2 theory.
- The simple explanation: An N = 2 superconformal theory is super-Weyl invariant, with the super-Weyl transformation parametrized by a chiral superfield Ω:

 $\phi \to \exp\left(\Omega\right) \cdot \phi$.

► In the regime of the validity of the effective theory, φ has a nonzero vev, and in flat space we can super-Weyl transform the vector multiplet to 1.

The EFT is therefore^(*)

 $\mathcal{L} = \mathcal{L}_{\rm free} + \mathcal{L}_{\rm anomaly} + \mathcal{L}_{\rm higher \ D-term}$

- For quantities insensitive to D-terms, this simple, two-term effective action, can be quantized meaningfully, and gives unambiguous answers to all orders in ¹/₇ perturbation theory.
- Note that the dimension △ of the generator of the chiral ring does not enter into the EFT at all, nor does the marginal coupling *τ* or any other parameter.
- ▶ In other words, any purely F-term-dependent observable has a large-J expansion that is uniquely determined by the anomaly coefficient α and nothing else, for a one-dimensional Coulomb branch of an $\mathcal{N} = 2$ gauge theory.

 One set of such observables are the Coulomb branch correlation functions

$$\exp(q_n) \equiv Z_n \equiv Z_{S^4} \times |x-y|^{2\mathcal{J}} \left\langle (\mathcal{O}(x)_{\Delta})^n (\overline{\mathcal{O}}(y)_{\Delta})^n \right\rangle_{S^4}$$

► The insertions φ^J(x) and φ^J(y) can be taken into the exponent as

$$S_{
m sources} \equiv -\mathcal{J}\log\left[\phi(x)
ight] - \mathcal{J}\log\left[ar{\phi}(y)
ight]$$

► This quantity $Z_n = \exp(q_n)$ is partition function of the EFT with sources:

$$Z_n = \int \mathcal{D}\Phi \,\mathcal{D}\Phi^{\dagger} \exp\left(-S_{\rm EFT} - S_{\rm sources}\right)$$

► This quantity is scheme-dependent, and dependent on the normalization of O_Δ, but these dependences cancel out in the double difference observables

$$\frac{Z_{n+1}Z_{n-1}}{Z_n^2} = \exp\left(q_{n+1} - 2q_n + q_{n-1}\right) \;.$$

These can now in principle be evaluated straightforwardly as functions of *J* and *α* using Ferynman diagrams, with no further input from the underlying CFT, as long as we are in large-*J* perturbation theory.

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The form of the expansion is

$$q_n = \mathbf{A} n + \mathbf{B} + \mathcal{J} \log(\mathcal{J}) + \left(\frac{\alpha}{2} + \frac{1}{2} \right) \log(\mathcal{J}) + \sum_{m \geq 1} \frac{\hat{K}_m(\alpha)}{\mathcal{J}^m}.$$

- The first two terms are the scheme and normalization ambiguities, the third term is the classical value of the source term, one loop free term, and classical anomaly term contributions.
- The last is the series of power-law corrections coming from loop diagrams with interaction vertices coming from the source term and the anomaly term, with the anomaly term vertices carrying powers of α .
- ► The structure of the EFT makes the polynomials $\hat{K}_m(\alpha)$ a polynomial in α of order m + 1:

$$\hat{\mathcal{K}}_m(lpha) = \sum_{\ell=0}^{m+1} \, \hat{\mathcal{K}}_{m,\ell} \, lpha^\ell \; .$$



Table 1 – Diagrams appearing at order 1/3.

- Of course, actually directly evaluating multiloop diagrams in an EFT is hard;
- To evaluate the power-law corrections, my collaborators and I used a combination of
 - Direct evaluation of some low-order diagrams;
 - \blacktriangleright Use of known data for some theories such as the free vector multiplet and $\mathcal{N}=4$ SYM ;
 - Supersymmetric recursion relations [Papadodimas 2009];
 - Embedding of the Coulomb-branch EFT into nonunitary UV completions invoving ghost hypermultiplets to apply the recursion relations to arbitrary values of α .

▶ With this combination of tricks, we were able to solve all the power-law corrections for any value of α , with the result:

$$q_n = \mathbf{A} n + \mathbf{B} + \log \left[\Gamma \left(\mathcal{J} + \alpha + 1 \right) \right]$$

+smaller than any power of ${\mathcal J}$.

I'll comment on those exponentially small corrections in a moment.

Confirmation of the large- \mathcal{J} expansion

- But first, let me talk about some evidence for this picture of large-J self-perturbatization of strongly coupled theories.
- Starting with our predictions for the O(2) model, where we predicted a formula

$$\Delta_J = \Delta_J = c_{\frac{3}{2}} J^{+\frac{3}{2}} + c_{\frac{1}{2}} J^{+\frac{1}{2}} -0.0937256\cdots$$

- It would be good to compare with bootstrap calculations in the O(2) model; at the moment bootstrap methods can only reach J ≤ 2 with any precision. [Kos, Poland, Simmons-Duffin 2013].
- It would be good if bootstrap methods could be developed to the point of being able to confirm our results, or add something substantial to them.
- But at the moment that hasn't happened, so let's move on to other avenues of confirmation.

► The first really nontrivial confirmation came from a Monte Carlo analysis up to J = 15 in the O(2) model, independently computing charged operator dimensions and estimating the leading Lagrangian coefficient b from the energies of charged ground states on the torus.

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 These results are from a PRL by [Banerjee, Orlando, Chandrasakhran 2017].

Monte Carlo numerics [Banerjee, Chandrasekharan, Orlando 2017]



Figure: Operator dimensions with the $c_{3/2}$, $c_{1/2}$ coefficients in the EFT prediction fit to data, giving $c_{3/2} = 1.195/\sqrt{4\pi}$ and $c_{1/2} = 0.075\sqrt{4\pi}$.

Monte Carlo numerics



Figure: Note the coefficients are fit with high-J data for operator dimensions and torus energies, and yet the leading-order prediction extrapolates extremely well down to J = 2.

Confirmation of the large- \mathcal{J} expansion

Though precise bootstrap results only exist up to J = 2, note that the values of the EFT parameters calculated from Monte Carlo calculation give

 $\Delta_{J=2} = 1.236(1)$

which one can compare to the bootstrap result

 $\Delta_{J=2} = 1.236(3)$

 There are other high-precision agreements between large-J theory and MC simulation in [Banerjee, Chandrasekharan, Orlando 2017].

.

 WARNING: The bootstrap result may have improved recently.

Confirmation of the large- \mathcal{J} expansion

- Moving beyond the O(2) case to other models in the same large- J universality class, one can look at dimensions of operators carrying topological charge J in the CIP(n) models.
- This analysis was done by , using a combination of large-N methods and numerical methods, with the result

$$\Delta_J^{\mathbb{CP}(n)} = c_{\frac{3}{2}}(n) J^{\frac{3}{2}} + c_{\frac{1}{2}}(n) J^{\frac{1}{2}} + c_0 + O(J^{-\frac{1}{2}}) ,$$

where the first two coefficients depend on the *n* of the model, but the J^0 term does not; in particular he finds

 $c_0 = -0.0935 \pm 0.0003$,

as compared to the EFT prediction

 $c_0 = -0.0937 \cdots$.

So the error bars are less than one percent , and the EFT prediction sits inside of them.
Confirmation of the large- \mathcal{J} expansion

- Now let's move on to our predictions for D = 4, N = 2 superconformal theories with one-dimensional Coulomb branch.
- ▶ For the case of free Abelian gauge theory and $\mathcal{N} = 4$ SYM with G = SU(2) our all-orders-in-J formula agrees with the exact expression:

 $Z_n^{(\text{EFT})} = Z_n^{(\text{CFT})} = n!$, free vector multiplet,

$$Z_n^{(\text{EFT})} = Z_n^{(\text{CFT})} = (2n+1)!$$
, $\mathcal{N} = 4$ SYM.

In these cases, there are no exponentially small corrections to the formula.

Confirmation of the large- \mathcal{J} expansion

- For other cases, the correlation functions are D-term independent and can be evaluated by exact supersymmetric methods involving localization and supersymmetric recursion relations,
- • • though at present these methods are limited to theories with a marginal coupling.
- Even using these methods, the recursion relations grow more challenging in application to compute corelators of higher J owing to the complication of the sphere partition function as a function of the coupling.
- ▶ Nonetheless we have been able to carry the recursion relations to $J \sim 76$ in the case of $\mathcal{N} = 2$ SQCD with $N_c = 2$, $N_f = 4$.

Numerics (Localization)



Figure 4.1 – Second difference in n for $\triangle_n^2 q_n^{(nc)}$ (dots) and for $\triangle_n^2 q_n^{EFT}$ (continuous lines) as function of $Im\tau$ at fixed values of n. The numerical results quickly reach a τ -independent value that is well approximated by the asymptotic formula when n is larger than $n \gtrsim 5$.

Confirmation of the large- \mathcal{J} expansion

- It is interesting to try to understand the disagreement between the all-orders-¹/₁ formula and the exact localization results.
- Our framework for large-J analysis dictates that any disagreement must be smaller than any power of J and associated with a breakdown of the Coulomb-branch EFT.
- ► The natural candidate for such an effect would be propagation of a massive particle over the infrared scale R = |x y|.
- Therefore we would expect the leading difference between the localization result and the EFT prediction, to be of the form

$$q_n^{(\mathrm{loc})} - q_n^{(\mathrm{EFT})}$$

 $\sim {
m const.} imes \exp\left(-M_{
m BPS \ particle} imes R
ight)$

$$= \text{const.} \times \exp\left(-(\text{const.})\sqrt{\frac{\mathcal{J}}{\text{Im}(\tau)}}\right).$$

- We compared the difference between EFT and exact results in the scaling limit of , where J is taken large with this exponent held fixed and fit it to this virtual-BPS-dyon ansatz for the exponentially small correction .
- We found the difference $q_n^{(loc)} q_n^{(EFT)}$ fits very well to

$$q_n^{(\mathrm{loc})} - q_n^{(\mathrm{EFT})} \simeq 1.6 \, e^{-rac{1}{2}\sqrt{\pi\,\lambda}}$$

 $\lambda \equiv 2\pi \mathcal{J}/ \texttt{Im}(au)$.

Numerics (Localization)



Figure 6.1 – Second difference in π for the discrepancy between localization and EFT results $\Delta_n^2(q_n^{(loc)}-q_n^{EFT})$ (dots) compared to $\Delta_n^2(1.6~e^{-\sqrt{\pi\lambda}/2})$ (continuous lines) as functions of Im τ at fixed values of $n/Im~\tau=\lambda/(4\pi)$. The agreement is quite good already for $\lambda=3$.

Summary so far

- In this first part of this colloquium, we have seen:
 - The large-quantum number expansion gives an asymptotic expansion for various observables that is complementary to ordinary perturbation theory and seemingly completentary to the conventional conformal bootstrap;
 - These methods are applicable to large global charge in generic critical points with global symmetries as well as large R-charge in superconformal fixed points at strong coupling;
 - ► The large-quantum-number limit gives a controlled expansion of many quantities with some universal and some nonuniversal terms in the series, with the nonuniversal terms always corresponding to unknown Wilson coefficients in the action of the large-charge EFT.
 - The form of the large-charge EFT can be quite distinct from any underlying Lagrangian realization of the full CFT, if such a realization even exists.

Numerics (Localization)



Figure 6.1 – Second difference in π for the discrepancy between localization and EFT results $\Delta_n^2(q_n^{(loc)}-q_n^{EFT})$ (dots) compared to $\Delta_n^2(1.6~e^{-\sqrt{\pi\lambda}/2})$ (continuous lines) as functions of Im τ at fixed values of $n/Im~\tau=\lambda/(4\pi)$. The agreement is quite good already for $\lambda=3$.

- So this is a rather interesting situation.
- Due to the magic of supersymmetry, not only can we compute all power-law corrections exactly modulo the scheme-dependent coefficients, we are actually able to compare to exact results to a precision where we can see the qualitative breakdown of the effective theory that we used to generate the all orders approximation.
- Seeing this, one is naturally tempted to try and go further and compare the exponentially small correction with physcial expectations at a precision level as well.

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- In order to do this, one really has to take on the "non-(super-)universal"^(*) coefficients A and B.
- The sum rules are fine for checking power law corrections, where all three adjacent terms in the sum rule have the same order of magnitude,
- but when checking exponentially small corrections which are rapidly decreasing as a function of *n*, the sum rule tends to introduce large relative errors and one would like to do better by deriving the actual value of the coefficients *A* and *B*.

The exponentially small correction

- The main challenge in doing this, is that the A and B coefficients are not only dependent on the marginal parameter τ, they are also scheme dependent.
- Often in the literature, including in the literature on supersymmetric localization, a "scheme dependent" coefficient is often treated as synonymous with an "inherently ambiguous" coefficient.
- This point of view is often used as a rationale for not doing certain kinds of computations, but it is simply wrong.
- Having a scheme-dependent coefficient in a microscopic or effective lagrangian, just means that you have to be careful about how operationally you are defining your renormalized lagrangian paramters relative to the UV completion or renormalization procedure being used.

The exponentially small correction

- For generic theories with marginal parameters this is often a bit involved; but
- for theories with extended supersymmetry the scheme dependence can often be reduced to an ambiguity by a holomorphic function of the complex coupling constant; and
- for theories such as SQCD which have an S-duality symmetry, even the holomorphic ambiguity can be reduced to a finite parameter, which

- can then be eliminated altogether by matching with perturbation theory .
- So, that is the course we are going to take here.

The holomorphic reparametrization scheme-dependence

- ► The first scheme dependence to discuss is the one that affects the *A* coefficient.
- It is a kind of "classical" scheme dependence having to do with the parametrization of the holomorphic gauge coupling.
- ► The Coulomb-branch chiral primary \$\mathcal{O} \equiv \$\mathcal{O}_2\$ \equiv \$\Tr(\hat{\phi}^2)\$ is uniquely defined up to an overall normalization, characterized by its supersymmetry properties and by its dimension and R-charge .
- However the overall normalization is exactly what matters so we have to specify it.

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The holomorphic reparametrization scheme-dependence

- ► In the literature the way mostly used to normalize *O* is by its relation to a marginal operator.
- ► After all, \mathcal{O} can be thought of as the $\mathcal{N} = 2$ F-term superspace integrand over all four positively R-charged Grassman coordinates θ_+ to generate the holomorphic half of the marginal operator that adjusts the gauge coupling $\tau \equiv \frac{4\pi i}{g_{YM}^2} + \frac{\theta}{2\pi}$:

 $\int d^4\theta_+ \mathcal{O}_2 = [\text{theory} - \text{independent constant}] \times \text{Tr}(F_+^2) + \cdots,$

where F_+ is the self-dual piece of the Yang-Mills field strength and the \cdots are the kinetic terms for the scalars and fermions.

So the normalization of O is related to the normalization of the dimension-two chiral primary operator O is naturally linked to the normalization of the marginal operator that is a superconformal descendant in the same multiplet.

The holomorphic reparametrization scheme-dependence

- However this doesn't resolve the question because a marginal operator doesn't have a universal natural normalization either.
- Rather, a (chiral half of a complex) marginal operator transforms under reparametrizations of the coupling constant as a section of the holomorphic cotangent bundle of theory space.
- That is, it transforms as

$$\mathcal{M}_{[\tau]} = rac{d au'}{d au}\,\mathcal{M}_{[au']}\;, \qquad \qquad \mathcal{M} \equiv {
m Tr}(F_+^2) + \cdots$$

and the chiral primary \mathcal{O} has the same transformation, since its normalization is canonically related to the normalization of \mathcal{M} :

$$\mathcal{O}_{[\tau]} = rac{d au'}{d au} \, \mathcal{O}_{[au']} \; ,$$

under a holomorphic reparametrization $\tau' = f(\tau)$.

 Under this coupling reparametrization scheme transformation, the exponentiated A-coefficient transforms as the norm-squared of the chiral primary itself

$$\exp\left(\mathsf{A}_{[au]}
ight) = \left|rac{d au'}{d au}
ight|^2 \exp\left(\mathsf{A}_{[au']}
ight)$$

We will exploit this transformation law to solve for A in a particularly simple holomorphic coordinate and then write the transformation law in any other holomorphic coordinate including the natural Lagrangian parameter τ.

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- There is a second, less obvious scheme ambiguity related to the Euler-density counterterm E_4 .
- First of all it is very non-obvious why this counterterm should even be relevant at all for the computation of two-point functions!
- But some elementary deduction shows that it is.
- After all, two-point functions on flat space are conformally equivalent to two-point functions on the four-sphere, and

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the four-sphere has a nonzero Euler number .

- So the sphere partition function transforms multiplicatively under an additive shift of the coefficient of the Euler counterterm.
- Supersymmetry does allow the Euler counterterm to appear in the action.
- However this term is in some sense an N = 2 F-term, so it can only appear with a (holomorphic) + (antiholomorphic) dependence on the holomorphic gauge coupling.
- ► Since the $Z_n = e^{q_n}$ are unnormalized partition functions with sources, they are affected by the same counterterm ambiguity as the sphere partition function without sources.

Euler-counterterm ambiguity

► The B coefficient is the n⁰ term in the large-n expansion of the q_n, so e^B transforms the same way under the Euler-counterterm ambiguity as does the sphere partition function :

 $\mathcal{L} \to \mathcal{L} - \operatorname{Re}[\operatorname{Log}[P(\tau)]] E_4$,

 $Z_{S^4} o |P(\tau)|^2 Z_{S^4} \;, \qquad \qquad e^B o |P(\tau)|^2 e^B \;.$

This transformation law means we must assign B a scheme label as well:

$$\exp\left(B_{\text{scheme 2}}\right) = \frac{Z_{\text{scheme 2}}}{Z_{\text{scheme 1}}} \exp\left(B_{\text{scheme 1}}\right)$$

S-duality

- Fixing the scheme-ambiguities is greatly simplified in a theory with an S-duality.
- In terms of the exponentiated gauge coupling

 $q \equiv e^{2\pi i \tau} \; ,$

the S-duality symmetry acts as:

- $S: \qquad q o 1-q \;, \qquad \qquad T: \qquad q o rac{q}{q-1} \;.$
- This is not quite the familiar fractional linear transformation by which the S-duality acts in N = 4 super-Yang-Mills.

S-duality

The infrared effective Abelian gauge coupling σ is the one that transforms in the familiar way by fractional linear transformations,

$$\sigma \to \frac{a\sigma + b}{c\sigma + d}, \qquad \qquad \begin{pmatrix} a & b \\ c & d \end{pmatrix} \in SL(2, \mathbb{Z})$$

with the generators acting by

$$S: \qquad \sigma \to -\frac{1}{\sigma} ,$$

 $T: \qquad \sigma o \sigma + 1$.

The relationship between the two couplings is given by the modular Lambda function

$$q = e^{2\pi i\tau} = \lambda(\sigma) ,$$

$$\sigma = 2\tau + \frac{4i}{\pi} \log[2] - \frac{i}{\pi} \left[\frac{q}{2} + \frac{13}{64} q^2 + \frac{23}{192} q^3 + \frac{2,701}{32,768} q^4 + \cdots \right]$$

- Given our transformation law for coupling reparametrizations we can take modular transformations as a special case.
- It follows that the chiral marginal operator *M*_[σ] and the chiral primary *O*_[σ] in the σ−frame, transform as holomorphic modular forms of weight 2.
- ► From there we can see that the A- coefficient transforms as a nonholomorphic modular form of weights (2, 2).

- ► The next ingredient is the recursion relations discovered by as a generalization of the tt^* equations to D = 4.
- These relations say that

$$\partial_{\sigma}\partial_{\bar{\sigma}} q_n = e^{q_{n+1}-q_n} - e^{q_n-q_{n-1}}$$

▶ When applied to the power law corrections they uniquely fix the form of q_n to be the Γ -function $\Gamma(2n + \frac{5}{2})$ up to the terms An + B.

Recursion relations and their duality-covariant solution

- They also give equations for the coupling dependence of the A- and B- terms.
- For the A- function they give

 $\partial_{\sigma}\partial_{\bar{\sigma}}A_{[\sigma]} = 8 e^{A_{[\sigma]}}$

For the B- function they give

 $\partial_{\sigma}\partial_{\bar{\sigma}}(B-A)=0$.

Note that these equations are covariant under both the holomorphic reparametrization scheme-dependence, and under the Euler counterterm scheme dependence, both of which shift A and/or B by a holomorphic plus antiholomorphic function of the complex coupling.

- That means that we can solve these equations in any scheme we like and transform it to whatever other scheme we like.
- It is simplest to solve in the σ -coordinate.
- In the σ-coordinate, the Liouville equation, the modular property, and the correct match with tree-level double-scaled perturbation theory uniquely fix the result to be

$$e^{\mathcal{A}_{[\sigma]}} = rac{16}{[\mathrm{Im}[\sigma]]^2}$$
 .

Recursion relations and their duality-covariant solution

➤ To specify the Euler counterterm scheme choice, we will compare with the scheme in which the sphere partition function was originally calculated by Pestun using the U(2) instanton partition function computed by Nekrasov.

The partition function as computed in this scheme has a derivable duality transformation given by:

$$q
ightarrow 1-q: \qquad \exp\left(B_{ ext{Pestun-}}\left[1-q
ight]
ight) = rac{|q|^2}{|1-q|^2}\exp\left(B_{ ext{Pestun-}}\left[q
ight]
ight) \;,$$

$$q
ightarrow rac{1}{q}: \qquad \exp\left(egin{smallmatrix} B_{ ext{Pestun-}}\left[rac{1}{q}
ight]
ight) = |q|^{-4}\exp\left(egin{smallmatrix} B_{ ext{Pestun-}}\left[q
ight]
ight)$$

- I say "derivable" rather than "derived" because the transformation does not appear AFAIK in the literature.
- In order to find it, it was essential to relate the Pestun-Nekrasov scheme to a slightly different scheme used by in which the duality transfomation law is more manifest by its relation to the crossing-symmetry transformation of a four-point function in two-dimesional Liouville theory under the well-known AGT correspondence.

► With this transformation law for the *B*-coefficient in the Pestun-Nekrasov scheme, and the general constraint from the recursion relations

 $\exp(B) = |\text{some holomorphic function}|^2 \times \exp(A_{[\sigma]})$,

we have the solution

$$\exp\left(B_{\frac{\text{Pestun-}}{\text{Nekrasov}}}\right) = [\text{const.}] \times \frac{|\lambda(\sigma)|^{+\frac{2}{3}} |1 - \lambda(\sigma)|^{+\frac{8}{3}}}{|\eta(\sigma)|^8 [\text{Im}(\sigma)]^2}$$

Recursion relations and their duality-covariant solution

- Again we have an ambiguity by a coupling-independent constant which we can fix again by matching with double-scaled perturbation theory.
- This time it is simpler to match at strong double-scaled coupling λ.
- We are able to do this by making use of the exact solution to the one-loop double-scaled coupling dependence found by .
- The result is

$$\exp\left(B_{\text{Pestun-}\atop\text{Nekrasov}}\right) = \gamma_{\text{G}}^{+12} e^{-1} 2^{-\frac{9}{2}} \pi^{-\frac{3}{2}} \frac{|\lambda(\sigma)|^{+\frac{2}{3}} |1-\lambda(\sigma)|^{+\frac{8}{3}}}{|\eta(\sigma)|^8 [\text{Im}(\sigma)]^2}$$

- Having in hand the explicit expressions for the A- and B-coefficients, we can now compare the full EFT approximation at large R-charge with expectations from localization, 000
- • • without taking double differences;
- this improves the numerical accuracy of the match, at basically all values of τ, to the point where the error in making the EFT approximation, compared to the exact result, has to be exaggerated on a plot in order to be visible at all.

Localization Compared With Full EFT Solution – Fixed n as a function of τ



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Localization Compared With Full EFT Solution – Fixed τ as a function of n



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So now we have solved for the full EFT approximation to the correlation functions:

$$Z_n^{(\mathrm{exact})} = Z_n^{(\mathrm{eft})} \times Z_n^{(\mathrm{mmp})}$$
,

$$Z_n^{(\mathrm{eft})} = e^{q_n^{(\mathrm{eft})}} = e^{An+B} imes \Gamma(2n+rac{5}{2}) \; ,$$

where the factor $Z_n^{(\text{mmp})} = e^{q_n^{(\text{mmp})}}$ is the set of exponentially small corrections describing massive macroscopic propagation of virtual BPS particles .

• We can get a handle on these too by the same strategy.

Exponentially small corrections

- Here's how we do it .
- ► We use the fact that the connected MMP term q_n^(mmp) is exactly what is left over when we take the full connected partititon function with sources q_n = Log[Z_n] and subtract the connected EFT contribution q_n^(eft) for which we now have an exact formula :

$$q_n^{(\text{mmp})} \equiv q_n - \mathbf{A} n - \mathbf{B} - \text{Log}\left[\Gamma(2n + \frac{5}{2})\right]$$

Using this identity we can rewrite the recursion relation for the full connected amplitude with sources q_n as a recursion relation for the macroscopic massive propagation contribution q_n^(mmp).

Exponentially small corrections

The resulting equation of variation for q_n^(mmp) takes the form
 [LHS] = [RHS]

with

 $[LHS] \equiv 16 \, \mathrm{Im}[\sigma]^2 \, \partial_{\sigma} \bar{\partial}_{\bar{\sigma}} q_n^{(\mathrm{MMP})}$

• • • • and

$$[\text{RHS}] \equiv (2n + \frac{7}{2})(2n + \frac{5}{2}) \left[\frac{Z_{n+1}^{(\text{MMP})}}{Z_n^{(\text{MMP})}} - 1 \right] - (2n + \frac{3}{2})(2n + \frac{1}{2}) \left[\frac{Z_n^{(\text{MMP})}}{Z_{n-1}^{(\text{MMP})}} - 1 \right]$$

where $Z_n^{(mmp)} \equiv e^{q_n^{(mmp)}}$

- The recursion relation is one input.
- The next input is the structure of the asymptotic expansion as dictated by effective field theoretic considerations.

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- To understand the structure of the expansion, we have to think carefully about what it is the MMP corrections represent.
- Mathematically they are the sum of all connected diagrams minus the EFT contributions ····
- with the latter representing propagation of virtual massless particles together with microscopic propagation of massive particles, which are absorbed into effective couplings of the EFT.

After these are subtracted we are left with connected diagrams which each have at least one macroscopic propagator for a massive particle.

- The lightest massive particle is the doublet hypermultiplet .
- In terms of the R-charge and gauge coupling its mass is given by

$$M_{\mathrm{hyper}} = rac{1}{R} \sqrt{rac{\mathcal{J}}{\pi \, \mathrm{Im}[\sigma]}} = rac{1}{R} \sqrt{rac{2n}{\pi \, \mathrm{Im}[\sigma]}}$$

where R is the radius of the three-sphere in radial quantization .

Remember we are always working in the limit

$$E_{
m IR} = R_{
m sphere}^{-1} \ll \Lambda \ll E_{
m UV} = \sqrt{\rho} \propto rac{\sqrt{\mathcal{J}}}{R} \propto rac{\sqrt{n}}{R} \propto M_{
m hyper} \; ,$$

so at fixed coupling and large R-charge \mathcal{J} the mass of the hyper is parametrically above the cutoff Λ .

- There is nothing inconsistent about including heavy particles above the cutoff in an effective field theory
- • • so long as we do it consistently!
- Actually such treatments of heavy supercutoff objects in EFT are well-understood and familiar in many contexts where the heavy state is stable or approximately stable.
- Examples include heavy quark effective theory effective string theory , the D-brane action , gapped goldstones , , and other examples .
- These examples can all be described in terms of a second quantized Hilbert space coupled to a first quantized dynamics of motions of the heavy particle.

- As we have seen, large R-charge is a semi-classical limit \cdots
- so we expect the leading contribution of the virtual massive particle at large R-charge to come from a classical configuration of the action for a massive BPS partcle coupled to masless vector multiplet.
- The heavy particle is conformal and gets its mass strictly from the magnitude of the vev of the vector multiplet, which is consant in the conformal frame of the cylinder.
- So we have to look for finite action classical trajectories of a particle of constant mass in on the cylinder.

- This narrows it down a lot because there aren't very many finite action trajectories for a massive hyper on the cylinder.
- ► In fact the only such trajectories are great circles of the spatial S³ at a fixed value of the radial time coordinate.



Figure: The leading contribution to the connected MMP function.

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► This means that the MMP function q_n^(mmp) has the asymptotic expansion at fixed coupling and large charge that is of the form q_n = e^{-𝔐} with

 $\mathfrak{W} = [$ worldline instanton action $S_{WLI}] + [$ parametrically smaller in n]

 Using the BPS mass formula we find that the worldline instanton action is

$$S_{\mathrm{WLI}} = \sqrt{rac{8\pi n}{\mathrm{Im}[\sigma]}}$$

► The first subleading correction is given by the quantum fluctuation determinant of the geometric worldline action about the classical trajectory. It contributes to 𝔐 proportional to Log[n].

- There are also contributions from the classical and quantum back-reaction of the massive hyper on the degrees of freedom of the massless abelian vector multiplet.
- ► These contribute to \mathfrak{W} with an *n*-dependence of at most $n^{-\frac{1}{2}}$.
- Incorporating additional higher-order geometric fluctuations of the massive trajectory and additional loops of the massless vector multiplet gives contributions which are suppressed by further powers of n^{-1/2}.
- So we have an asymptotic expansion of the form

$$-\mathrm{Log}[q_n] \equiv \mathfrak{W} = \sqrt{\frac{8\pi n}{\mathrm{Im}[\sigma]}} + \gamma[\sigma] \mathrm{Log}[n]$$

- At this point we could in principle just calculate all these terms directly in the effective theory of a geometric fluctuations of a massive worldline coupled to massless fields in the bulk about a nontrivial classical solution.
- But it turns out we have to do very little caculation.
- The recursion relations give PDE s for the σ -dependence of the functions $\gamma[\sigma]$, $w_p[\sigma]$ at each order , and we have enough information about boundary conditions to find the physically correct solution to each PDE .

For instance, the recursion relation at order Log[n] gives

 $(\partial_{\sigma} - \partial_{\bar{\sigma}})\gamma[\sigma] = 0$.

- ► This means γ can depend only on the real part of σ which is proportional to the infrared θ -angle θ_{IR} .
- But the dynamics must be independent of θ_{IR} at weak coupling, so γ must be independent of σ identically:

$$\gamma[\sigma] = (\sigma - \text{independent}) = \gamma$$
.

- To find the actual value of γ we must match with double-scaled perturbation theory again.
- Taking the double-scaling limit of 22 and then taking the strong coupling expansion of that double-scaling limit we find

$$\gamma = \lim_{\lambda \text{ fixed} \atop \lambda \text{ fixed}} \mathfrak{W} \bigg|_{\lambda \text{ term}} = \mathcal{F}^{(\text{inst})}[\lambda] \bigg|_{\lambda \text{ term}} = -\frac{1}{4} .$$

- ► Here the quantity F^(inst)[λ] is 's "worldline instanton partition function " which sums up all the terms scaling as n⁰ in the double scaling limit of the MMP function, without any terms of order n⁻¹ or smaller, and also without any EFT contributions.
- ► The function F^(inst)[λ] can be thought of as the sum over massive macroscopic worldlines and the first-quantized quantum fluctuations of their worldlines about the classical trajectory while discarding all quantum fluctuations of the massless fields.

- The functions $w_p[\sigma]$ can be found similarly.
- At each *p* the recursion relation gives a first-order PDE for σ ;
- ► The boundary condition at weak coupling forces the correct solution to depend on $s \equiv \text{Im}[\sigma]$ only;
- ► This determines the solution up to a single σ independent constant ····
- ► ··· which can be fixed by taking the double scaling limit and matching with the function F^(inst)[λ] of .

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The first few terms are:

$$w_{1} = \frac{1}{48(s/2\pi)^{3/2}} + \frac{1}{\sqrt{(s/2\pi)}} - \frac{11\sqrt{(s/2\pi)}}{16}$$
$$w_{2} = -\frac{1}{4} - \frac{1}{64(s/2\pi)} + \frac{19(s/2\pi)}{64}$$
$$w_{3} = -\frac{1}{5120(s/2\pi)^{5/2}} - \frac{1}{96(s/2\pi)^{3/2}}$$
$$-\frac{119}{512\sqrt{(s/2\pi)}} + \frac{11\sqrt{(s/2\pi)}}{32} - \frac{527(s/2\pi)^{3/2}}{3072}$$

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► $w_4 = \frac{119}{1024} + \frac{1}{2048(s/2\pi)^2}$ $+\frac{1}{64(s/2\pi)}-\frac{19(s/2\pi)}{64}+\frac{235(s/2\pi)^2}{2048}$ $w_5 = \frac{1}{229,376(s/2\pi)^{7/2}} + \frac{3}{10,240(s/2\pi)^{5/2}}$ $+\frac{737}{98,304(s/2\pi)^{3/2}}+\frac{101}{1024\sqrt{(s/2\pi)}}$ $-\frac{8,155\sqrt{(s/2\pi)}}{32,768}+\frac{527(s/2\pi)^{3/2}}{2048}$ 14,083 $(s/2\pi)^{5/2}$ 163.840



Figure: Plot giving the accuracy of the fixed-coupling large-charge estimates of the MMP function through N⁶LO, plotted as the number of digits of accuracy of each of the estimates, as a function of *n*. The quantity being plotted is $-\frac{1}{\text{Log}[10]}$ the logarithm of the relative error in the estimate of the MMP function. The horizontal axis is *n*, and the vertical axis is $-\frac{1}{\text{Log}[10]}$ Log $\left| \frac{q_n^{(\text{MMP})} - (q_n^{(\text{MMP})})_{\text{estimate}}}{q_n^{(\text{MMP})}} \right|$. The LO, NLO, N²LO, N³LO, N⁴LO, N⁵LO and N⁶LO estimates are given by the blue, yellow, green, red, and purple, brown, and light blue curves respectively, which are in ascending order on the chart for $n \gtrsim 65$.



Figure: Plot of the giving the accuracy of the double-scaled large-charge estimates of the MMP function through N⁵LO. The quantity being plotted is $-\frac{1}{\text{Log}[10]}$ the logarithm of the relative error in the estimate of the MMP function. The horizontal axis is *n*, and the vertical axis is $-\frac{1}{\text{Log}[10]} \text{Log} \left[\frac{|q_n^{(\text{MMP})} - (q_n^{(\text{MMP})})_{\text{estimate}}|}{|q_n^{(\text{MMP})}|} \right]$ The LO, NLO, N²LO, N³LO, N⁴LO and N⁵LO double-scaled estimates are given by the blue, yellow, green, red, and purple, and brown dots respectively.

Some data points for the MMP function

	e ^{-S} WLI w/	estimate w/	estimate w/	estimate w/	estimate w/	estimate_w/	exact
	prefactor	$O(n^{-\frac{1}{2}})$	$O(n^{-1})$	$O(n^{-\frac{3}{2}})$	O(n-2)	$O(n^{-\frac{5}{2}})$	$q_n^{(MMP)}$
1	0.3306547971	0.5494056108	0.3208548662	0.4473688062	0.3855992091	0.4826779760	0.4263073863
2	0.2369155213	0.3392528251	0.2592574487	0.2915876823	0.2809550113	0.2923319506	0.2924432054
3	0.1777356756	0.2382818085	0.1991725121	0.2123297908	0.2088531261	0.2118834217	0.2140116919
4	0.1376160664	0.1773896477	0.1550715601	0.1616504023	0.1601561764	0.1612839808	0.1629019007
5	0.1090165541	0.1368079122	0.1228554207	0.1265627818	0.1258128005	0.1263191935	0.1274138410
6	0.08788478715	0.1081284329	0.09885739579	0.1011186866	0.1007021939	0.1009589550	0.1016922546
7	0.07184428016	0.08704428588	0.08060668448	0.08206644241	0.08181796572	0.08195980455	0.08245642599
8	0.05940855833	0.07109089681	0.06646850876	0.06745213130	0.06729571354	0.06737924445	0.06772080050
9	0.04960141702	0.05874895841	0.05534091587	0.05602642358	0.05592374401	0.05597544585	0.05621405675
10	0.04175689585	0.04903003507	0.04646262665	0.04695358198	0.04688386790	0.04691717164	0.04708632558

Table: The successive refined estimates for the massive macroscopic propagation function $q_n^{(\text{MMP})} \equiv q_n - q_n^{(\text{eft})}$, at the coupling $\tau = \frac{25i}{\pi}$.

Some data points for the MMP function

	e ^{-S} WLI w/ prefactor	estimate w/ $O(n^{-\frac{1}{2}})$	estimate w/ $O(n^{-1})$	estimate w/ $O(n^{-\frac{3}{2}})$	estimate w/ $O(n^{-2})$	estimate w/ $O(n^{-\frac{5}{2}})$	$\frac{\text{exact}}{q_n^{(\text{MMP})}}$
20	0.01000372820	0.01120653279	0.01090917838	0.01094979540	0.01094572872	0.01094710280	0.01095666332
30	0.003237876660	0.003552394945	0.003489274164	0.003496339719	0.003495762541	0.003495921786	0.003496908245
40	0.001234172590	0.001337343268	0.001319481435	0.001321216247	0.001321093558	0.001321122874	0.001321264040
50	0.0005236492943	0.0005626347012	0.0005566148966	0.0005571384461	0.0005571053342	0.0005571124109	0.0005571373807
60	0.0002400576708	0.0002563211089	0.0002540336796	0.0002542154292	0.0002542049370	0.0002542069841	0.0002542121164
70	0.0001167774246	0.0001240840233	0.0001231342717	0.0001232041768	0.0001232004409	0.0001232011157	0.0001232022950
80	0.00005956426027	0.00006304351857	0.00006262109222	0.00006265018869	0.00006264873419	0.00006264897995	0.00006264927477
90	0.00003159033412	0.00003332720441	0.00003312863197	0.00003314153162	0.00003314092369	0.00003314102053	0.00003314109898
100	0.00001731365088	0.00001821547422	0.00001811776564	0.00001812378889	0.00001812351960	0.00001812356029	0.00001812358207
110	9.760096897×10 ⁻⁶	0.00001024423944	0.00001019427220	0.00001019720974	0.00001019708452	0.00001019710256	0.00001019710872
1120	5.638450512×10 ⁻⁶	5.905956503×10 ⁻⁶	5.879544842×10 ⁻⁶	5.881031737×10 ⁻⁶	5.880971055×10 ⁻⁶	5.880979426×10 ⁻⁶	5.880981146×10 ⁻⁶

Table: The successive refined estimates for the massive macroscopic propagation function $q_n^{(\text{MMP})} \equiv q_n - q_n^{(\text{eft})}$, at the coupling $\tau = \frac{25i}{\pi}$.

More Recent History:

- N = 4 SYM at large R-charge [Bernstein, Maldacena, Nastase]
- and large spin [Belistsky, Basso, Korchemsky, Mueller], [Alday, Maldacena]
- Large-spin expansion in general CFT from light-cone bootstrap [Komargodski-Zhiboedov], [Fitzpatrick, Kaplan, Poland, Simmons-Duffin], [Alday 2016]
- Large-spin expansion in hadrons [SH, Swanson], [SH, Maeda, Maltz, Swanson], [Caron-Huot, Komargodski, Sever, Zhiboedov], [Sever, Zhiboedov]

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Other stuff I didn't get to mention

Modern:

- Large-charge expansion in generic systems with abelian global symmetries: [SH, Orlando, Reffert, Watanabe 2015], [Monin 2016], [Monin, Pirtskhalava, Rattazzi, Seibold 2016], [Loukas 2016]
- Nonabelian symmetries: [Alvarez-Gaume, Loukas, Orlando, Reffert 2016], [Loukas, Orlando, Reffert 2016], [SH, Kobayashi, Maeda, Watanabe 2017], [Loukas 2017], [SH, Kobayashi, Maeda, Watanabe 2018]
- Charge AND spin: [Cuomo, de la Fuente, Monin, Pirtskhalava, Rattazzi 2017]
- Topological charge: [Pufu, Sachdev 2013] [Dyer, Mezei, Pufu, Sachdev 2015], [de la Fuente 2018]
- EFT connection with bootstrap: [Jafferis, Mukhametzhanov, Zhiboedov 2017]
- Large charge limit in gravity: [Nakayama, Nomura 2016], [Loukas, Orlando, Reffert, Sarkar 2018]

Vacuum manifolds \Leftrightarrow chiral rings at large-R-charge:

- $D = 3, \ \mathcal{N} \ge 2$ theories : [SH, Maeda, Watanabe 2016]
- ▶ D = 4, N ≥ 2 theories : [SH, Maeda 2017], [SH, Maeda, Orlando, Reffert, Watanabe 2018], [SH, Maeda, Orlando, Reffert, Watanabe 2020], [SH, Orlando 2021], [SH 2021]
- ▶ Double-scaling limit in lagrangian N ≥ 2 theories: [Bourget, Rodriguez-Gomez, Russo 2018], [Grassi, Komargodski, Tizzano 2019]

Other stuff I didn't get to mention

- In addition, there has been a great deal of fascinating work in this area in the past few years that I don't have the space to do justice to in the references here.
- A sampling includes: [Favrod, Orlando, Reffert 2018] [Loukas, Orlando, Reffert, Sarkar 2018] [Kravec, Pal 2018] [Bourget, Rodriguez-Gomez, Russo 2018] [Badel, Cuomo, Monin, Rattazzi 2019] [Alvarez-Gaume, Orlando, Reffert 2019] [Arias-Tamargo, Rodriguez-Gomez, Russo 2019] [Badel, Cuomo, Monin, Rattazzi 2020] [Delacretaz 2020] [Cuomo, Esposito, Gendy, Khmelnitsky, Monin, Rattazzi 2020] [Cuomo 2020] [Orlando, Reffert, Sannino 2020] [Antipin, Bersini, Sannino, Wang, Zhang 2020] [Komargodski, Mezei, Pal, Raviv-Moshe 2021] [Cuomo, Delacretaz, Mehta 2021] [Orlando, Pellizzani, Reffert 2021] [Dondi, Kalogerakis, Orlando, Reffert 2021] [Cassani, Komargodski 2021]

Conclusions

- The large-J expansion gives an analytically controlled way to compute CFT data outside of any other sort of simplifying limit, particularly illuminating simple behavior in regimes where numerical bootstrap methods cannot currently access, despite formal similarity of the expansions.
- ► The large- J predictions in cases such as the O(2) model and various D = 4, N = 2 superconformal theories with one-dimensional Coulomb branch, agree extremely well even at low J with Monte Carlo, bootstrap, and exact supersymmetric methods.
- These results have greatly improved our quantitative control and conceptual understanding of even the simplest strongly-coupled CFT.
- Analysis of more examples is sure to yield further interesting surprises about the large-scale structure of theory space.

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- Thank you.
- LOOK AT THIS PHOTOGRAPH.