

Sofie Martins, University of Graz, 22.1.2026

**FWF** Österreichischer  
Wissenschaftsfonds

# Computing cross-sections from the lattice

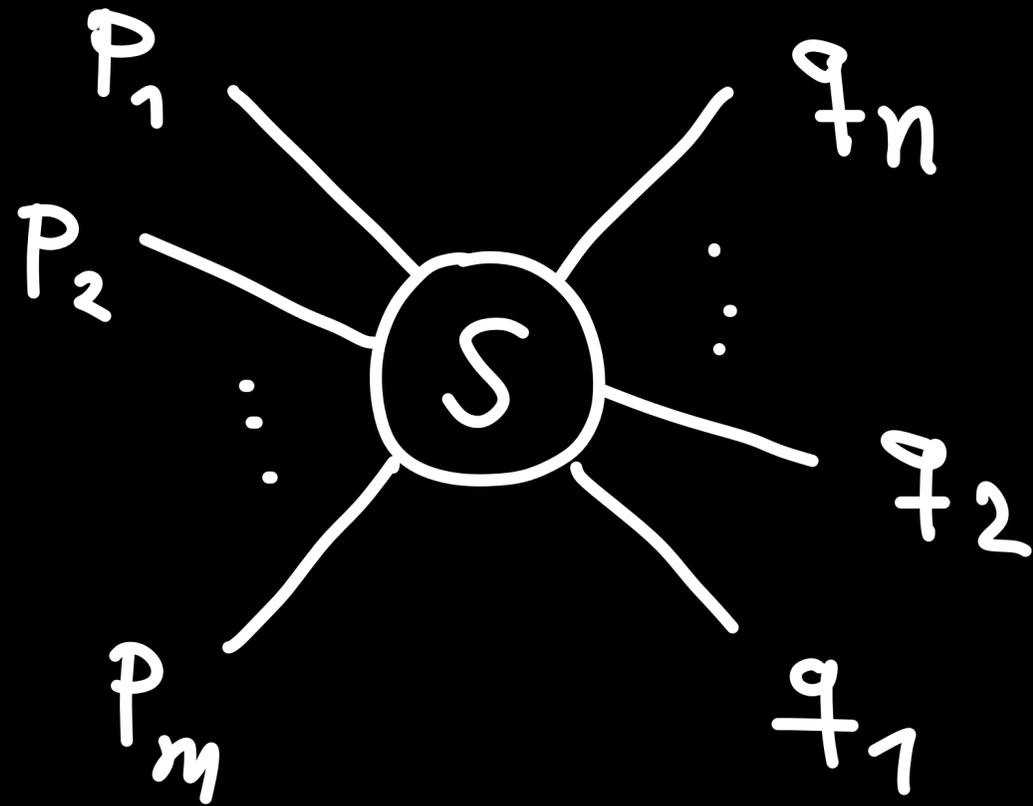
A recent shift in lattice thinking paradigms?

# Outline

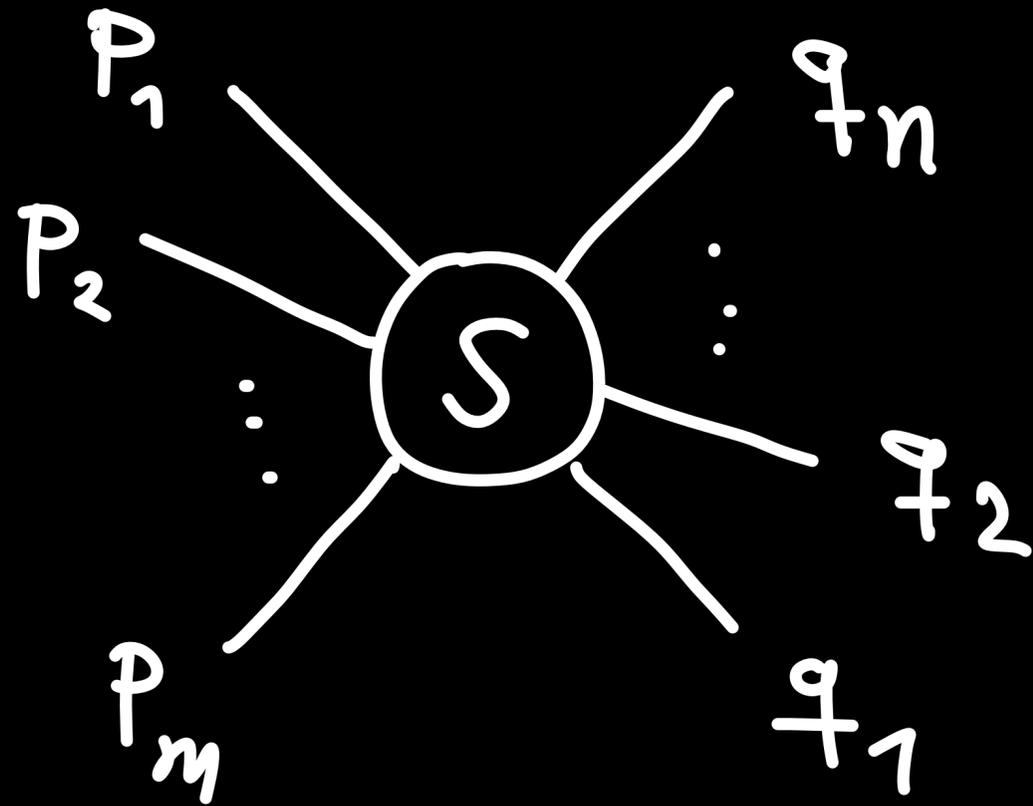
- What are scattering amplitudes?
- Our test case
- Computing spectral densities from lattice data
- Method by Hansen and Bulava
- Method by Patella and Tantalo
- Conclusions and Outlook

**What are scattering amplitudes?**

# Computing a transition probability

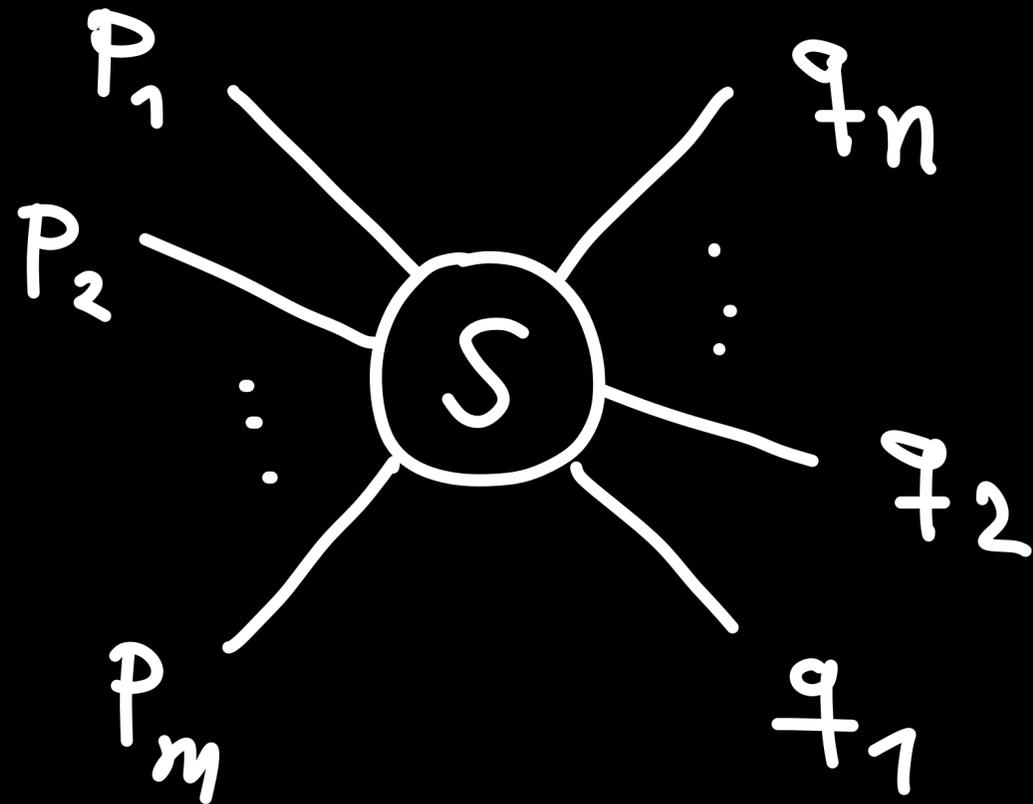


# Computing a transition probability



$$\leadsto P = |\langle \text{out} | \{P_f\} | \{q_i\} \rangle_{in}|^2$$

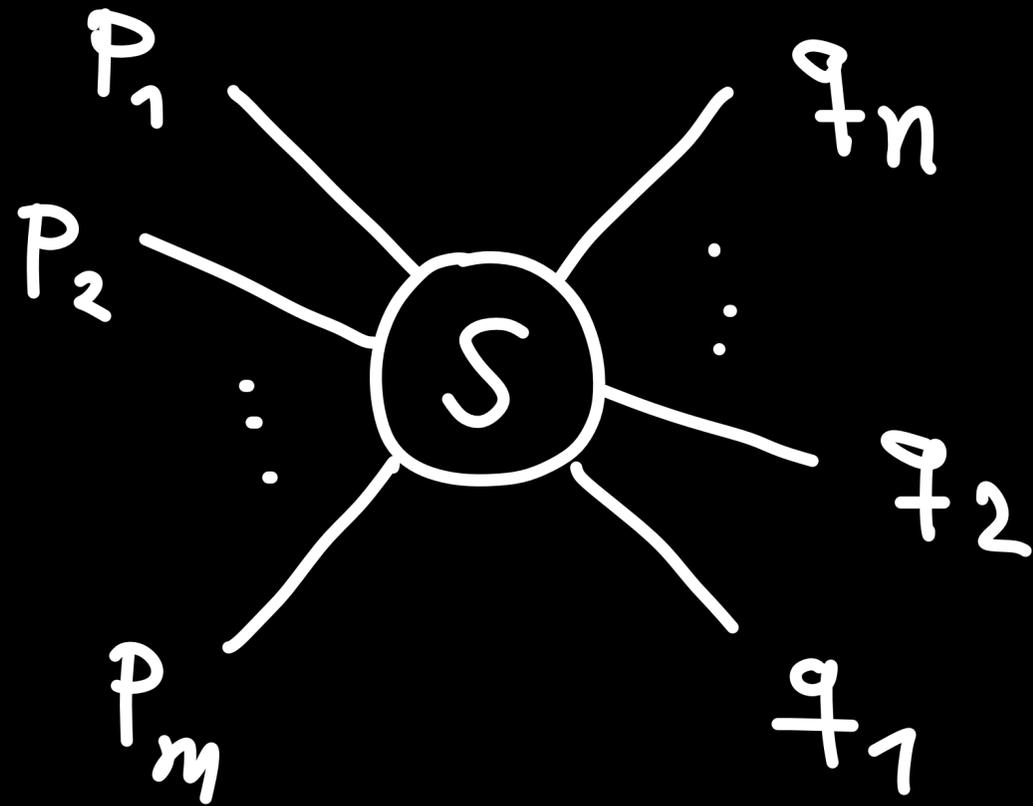
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With definite momenta  $p_f$  and  $q_i$

# Computing a transition probability



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With definite momenta  $p_f$  and  $q_i$

Physical theory encoded in S-Matrix

# Cross sections conceptually

$$d\sigma(\dots) = \int |_{out} \langle \{p_f\} | \{q_i\} \rangle_{in}|^2 \times \text{KINEMATICS}$$

# Cross sections conceptually

known



$$d\sigma(\dots) = \int |_{out} \langle \{p_f\} | \{q_i\} \rangle_{in}|^2 \times \text{KINEMATICS}$$

known



# Cross sections conceptually

possibly non-perturbative  
physics

$$d\sigma(\dots) = \int \left| \langle \{p_f\} | \{q_i\} \rangle_{in} \right|^2 \times \text{KINEMATICS}$$

# Perturbative approach LSZ reduction

Connect matrix element to time-ordered correlation functions

$${}_{out} \langle \{p_f\} | \{q_i\} \rangle_{in} \sim \langle 0 | T \{ \phi(x_1) \cdots \phi(x_n) \phi(y_1) \cdots \phi(y_m) \} | 0 \rangle$$

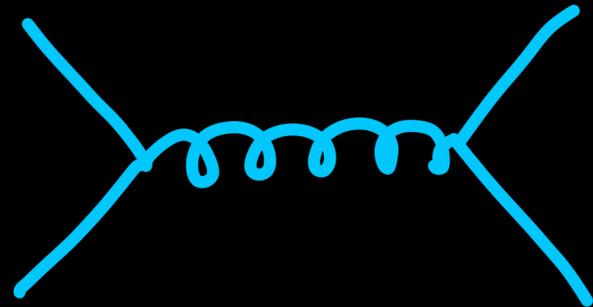
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$$\sim \sum \text{Feynman Diagrams}$$

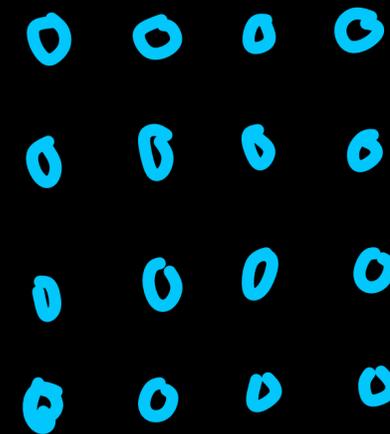
# Perturbation theory vs the Lattice



What we compute in  
perturbation theory

$$\langle 0 | T \{ \phi(x_1) \cdots \phi(x_n) \phi(y_1) \cdots \phi(y_m) \} | 0 \rangle$$

Time-ordered correlation functions



What we compute on the  
lattice

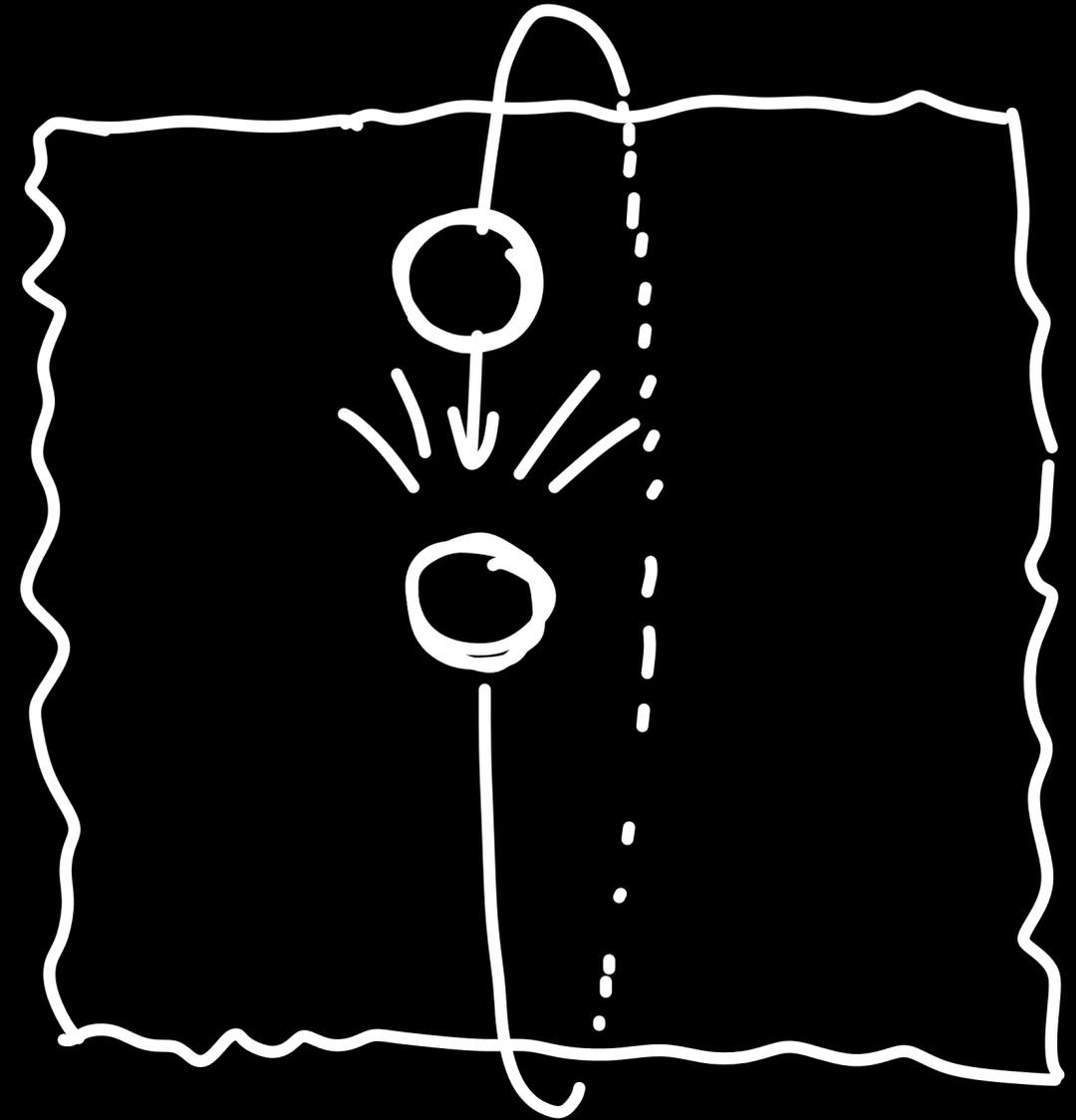
$$\langle 0 | \phi(x_1) \cdots \phi(x_n) \phi(y_1) \cdots \phi(y_m) | 0 \rangle$$

No time ordering

# Conversion: Lüscher Method

[Lüscher, 1986]

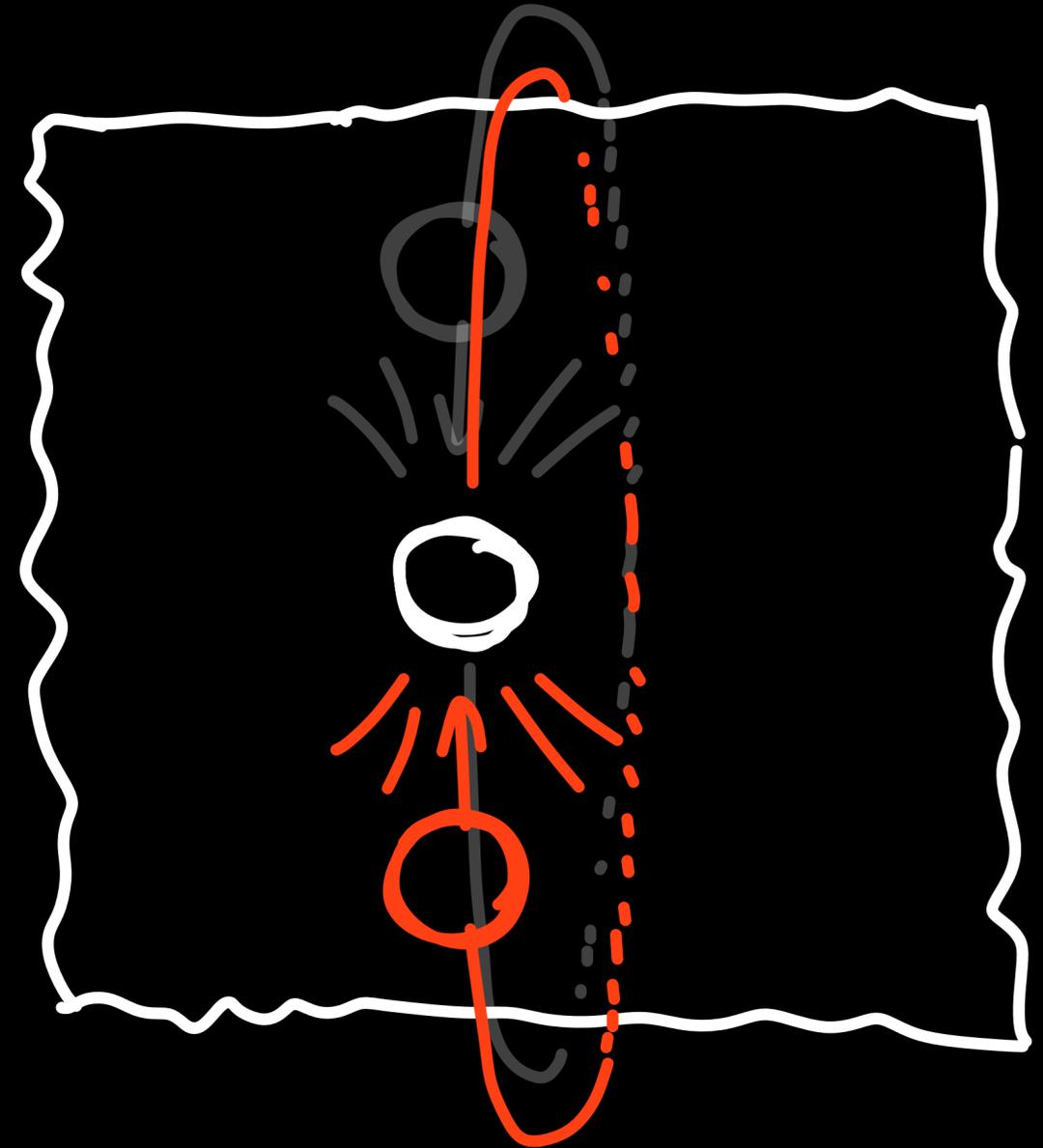
- Due to the finite volume any particle interacts with itself over the boundary



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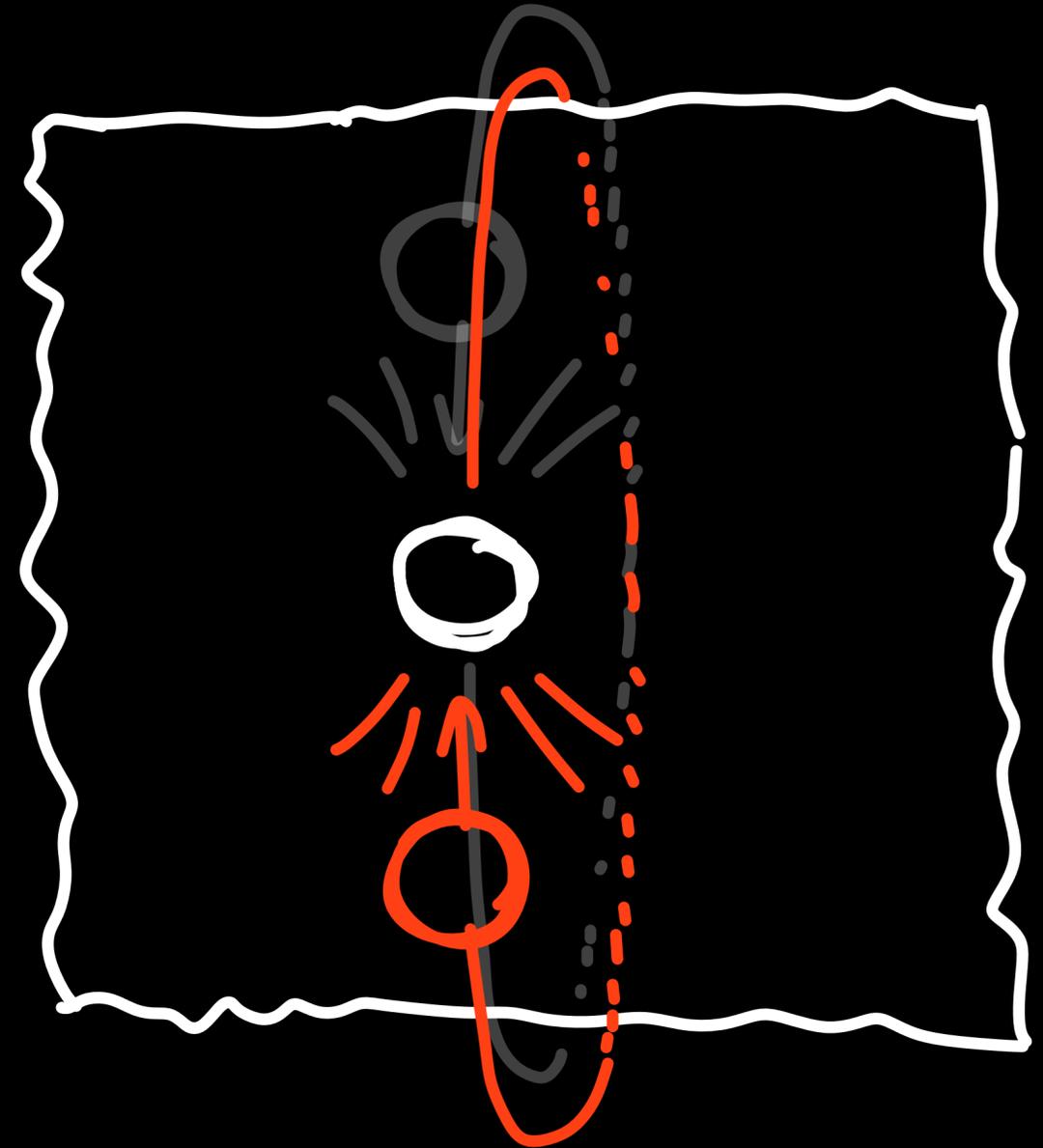
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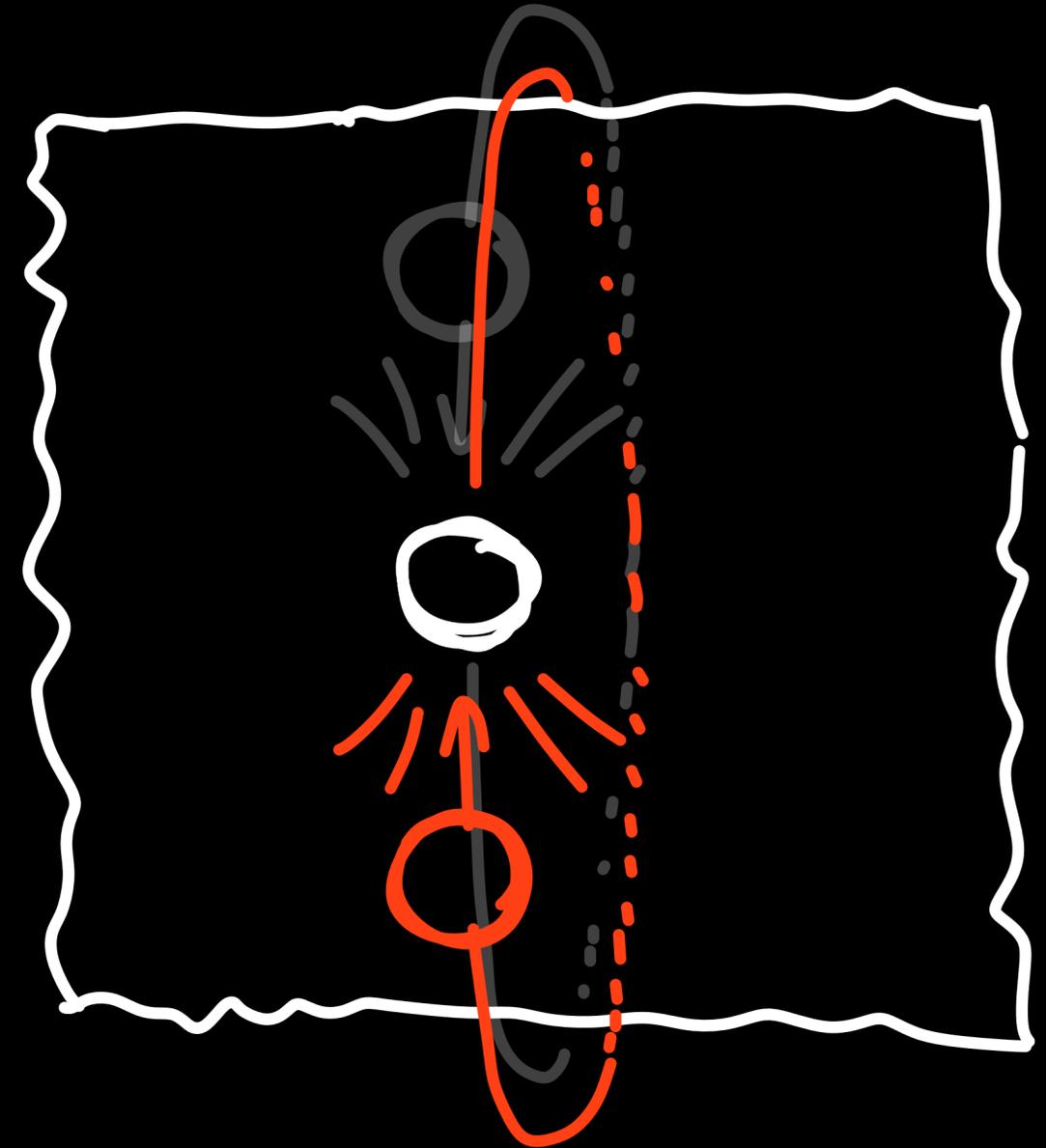
- Due to the finite volume any particle interacts with itself over the boundary
- This interaction is forward and back in time
- Cancellation reconstructs time ordering



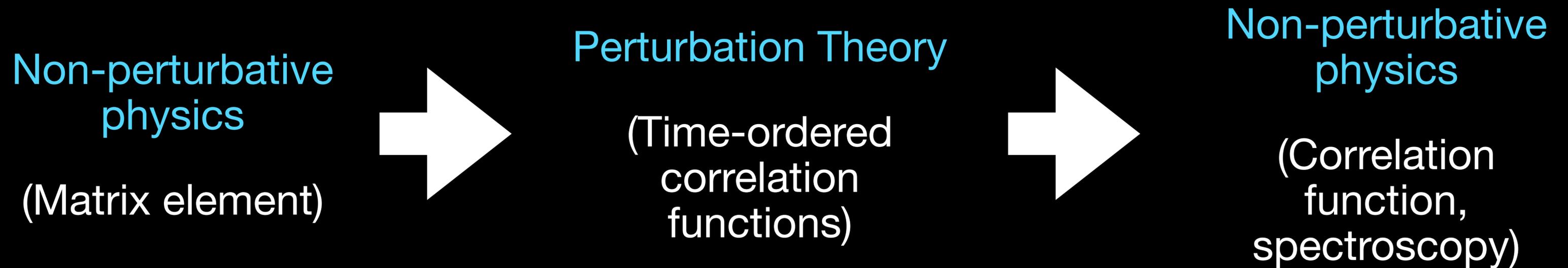
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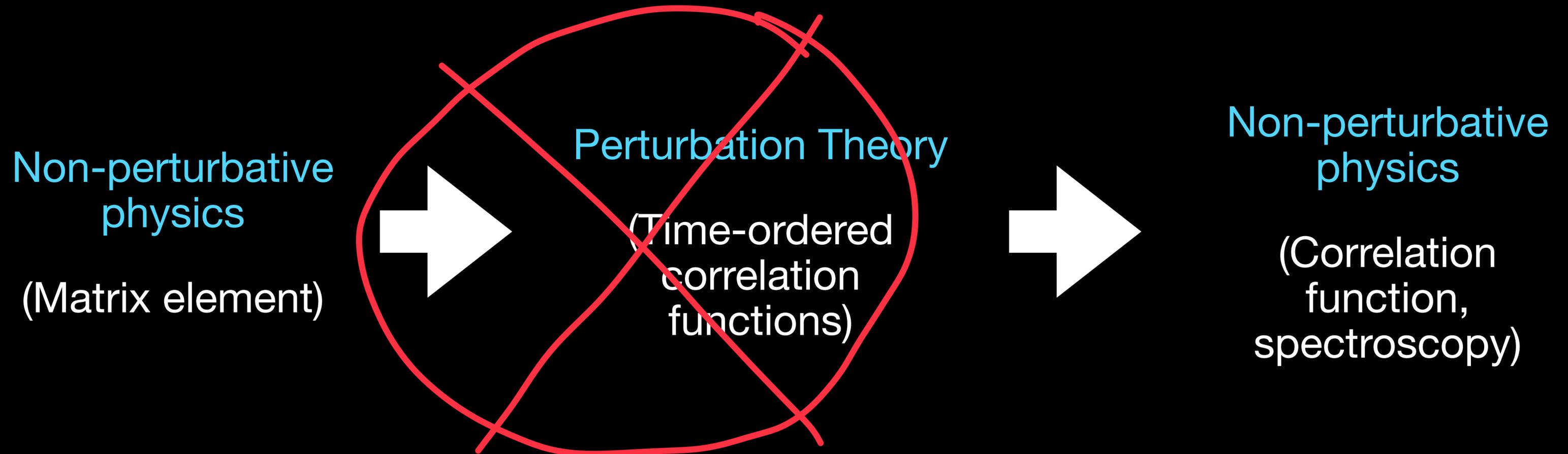
- Due to the finite volume any particle interacts with itself over the boundary
- This interaction is forward and back in time
- Cancellation reconstructs time ordering
- Relation between finite volume energy levels evaluated in Euclidean space to infinite-volume Minkowski space Scattering Amplitudes



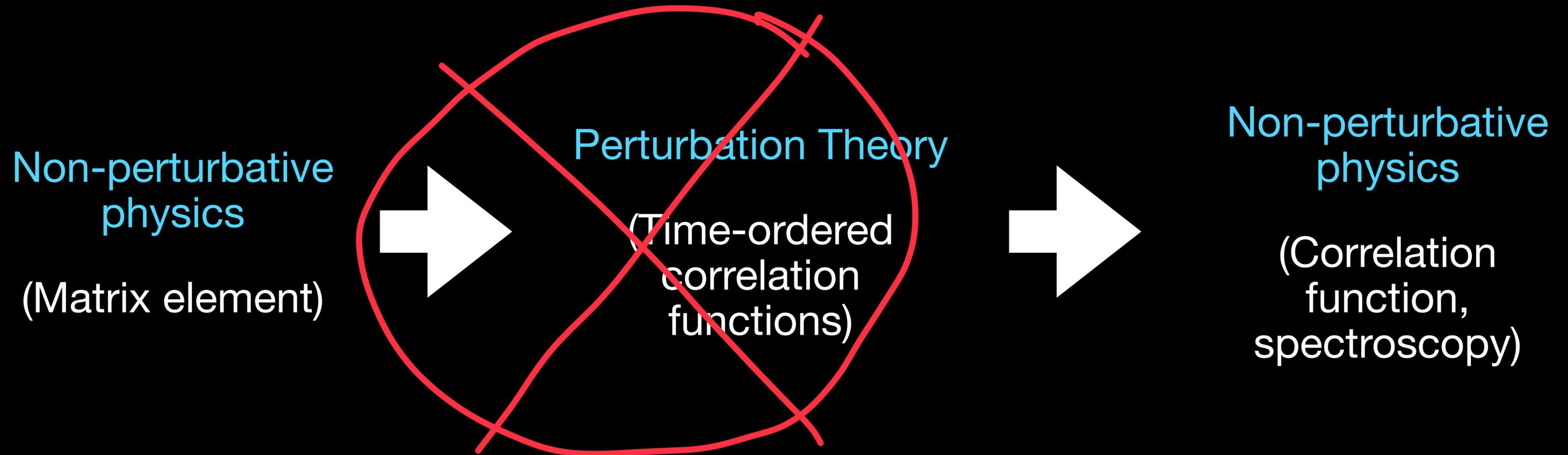
# A comment



# Rethinking the lattice extraction method



# Rethinking the lattice extraction method



Use spectral densities instead of correlation functions

**Apply LSZ again**

$$\langle 0 | \phi^\dagger(x) \phi(y) | 0 \rangle$$

Apply LSZ again

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# Apply LSZ again

$$\langle 0 | \underbrace{\phi^\dagger(x) \phi(y)}_{\mathbb{1}} | 0 \rangle = \sum_{\mathcal{h}} \langle 0 | \phi^\dagger(x) | \mathcal{h} \rangle \langle \mathcal{h} | \phi(y) | 0 \rangle$$

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$$\langle 0 | \underbrace{\phi^\dagger(x) \phi(y)}_{\mathbb{1}} | 0 \rangle = \sum_n \langle 0 | \phi^\dagger(x) | n \rangle \langle n | \phi(y) | 0 \rangle$$
$$= \int \frac{d^4 p}{(2\pi)^4} e^{-ip(x-y)} \delta(p^2)$$

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Extract this from the lattice

# Apply LSZ again

$$\langle 0 | \underbrace{\phi^\dagger(x) \phi(y)}_{\mathbb{1}} | 0 \rangle = \sum_n \langle 0 | \phi^\dagger(x) | n \rangle \langle n | \phi(y) | 0 \rangle$$
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Valid for general energies

# Apply LSZ again

$$\langle 0 | \phi^\dagger(x) \phi(y) | 0 \rangle = \sum_n \langle 0 | \phi^\dagger(x) | n \rangle \langle n | \phi(y) | 0 \rangle$$

$\uparrow$   
 $\mathbb{1}$

$$= \int \frac{d^4 p}{(2\pi)^4} e^{-ip(x-y)} \delta(p^2)$$



Valid for general energies



LSZ requires a mass gap, so not applicable for QCD+QED

# Our testing theory

# Testing scattering for weak physics

QCD

Strong coupling

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- Fine lattice spacing required

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In practice: Master Field Simulations

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Electroweak

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In practice: Master Field Simulations

[Francis et. al., 2020]

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- No continuum extrapolation

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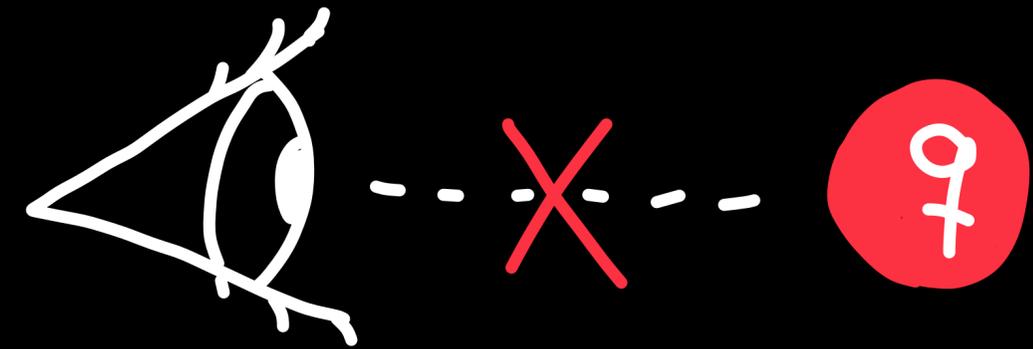
- No continuum extrapolation
- No LOCP interpolation
- Easily controllable FV effects

↳ how to simulate massive gauge bosons?  
Spontaneous symmetry breaking?

# FMS Framework

[Maas, 2019] & references therein

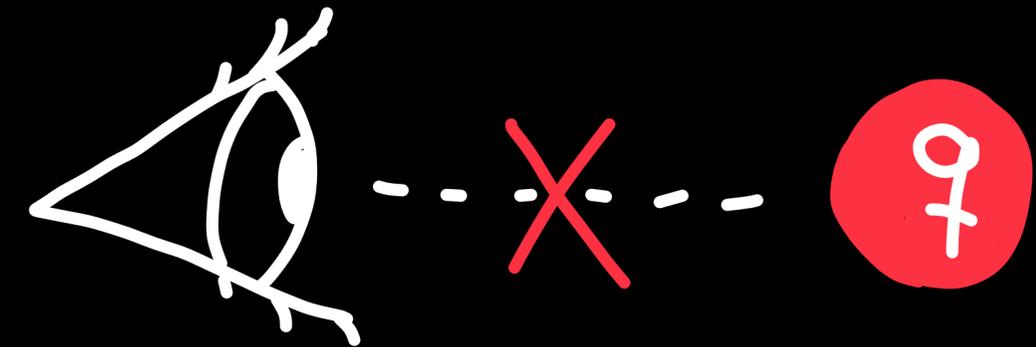
1. Observers are color blind



# FMS Framework

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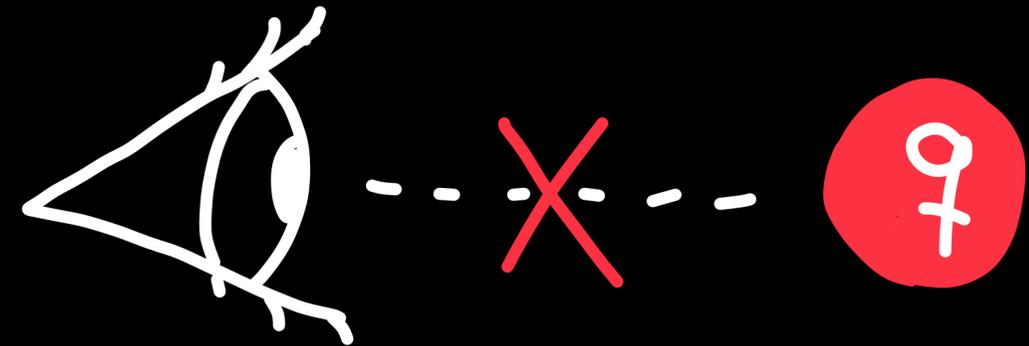


2. Physical objects are color singlets

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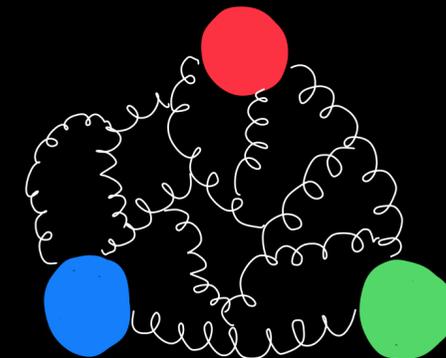
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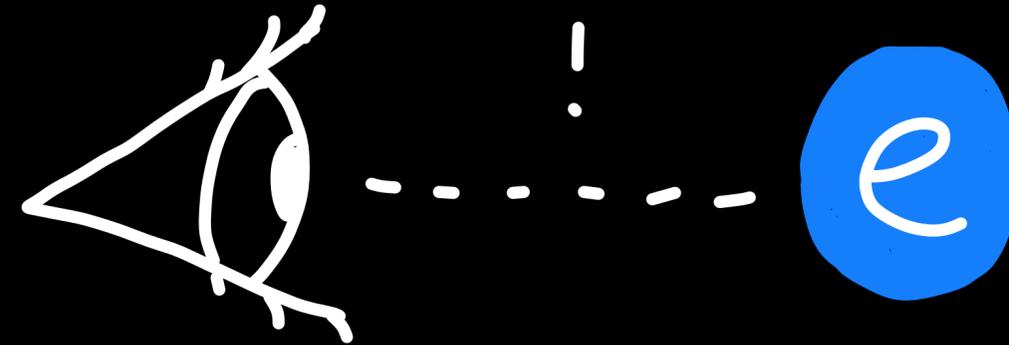
3. Predictions require non-perturbative techniques



# FMS Framework

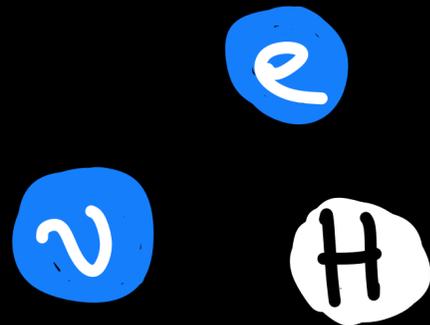
1. Observers are **not** weak isospin blind

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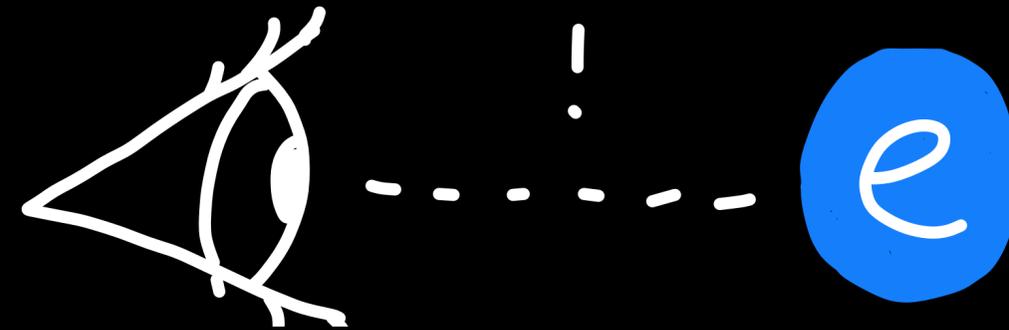


# FMS Framework

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[Maas, 2019] & references therein



2. Leptons and neutrinos are individually observable

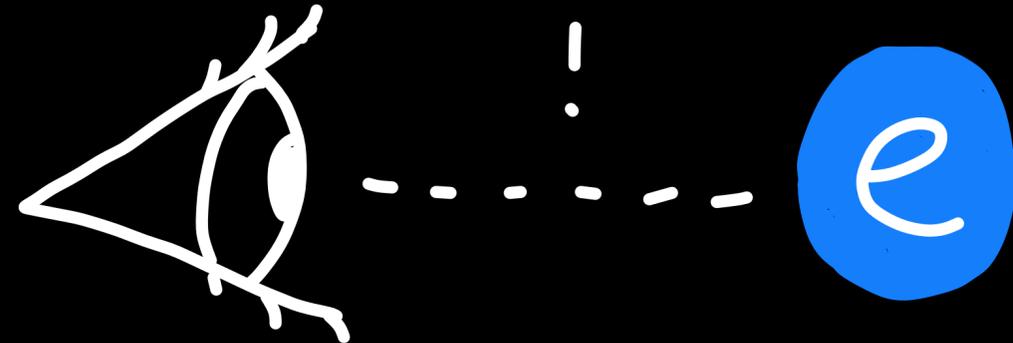
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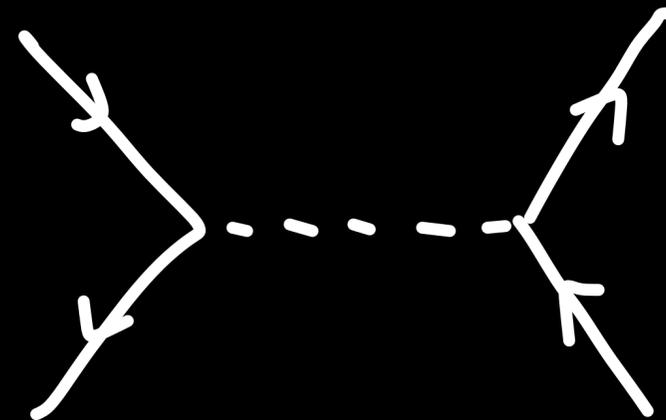


3. Perturbative treatment possible using the BRST construction

[Maas, 2019] & references therein

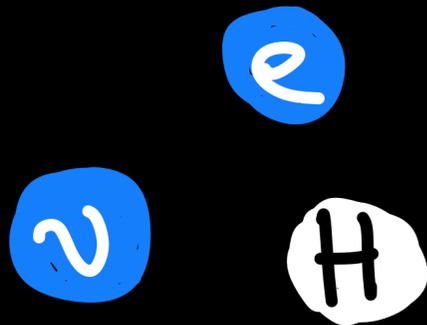


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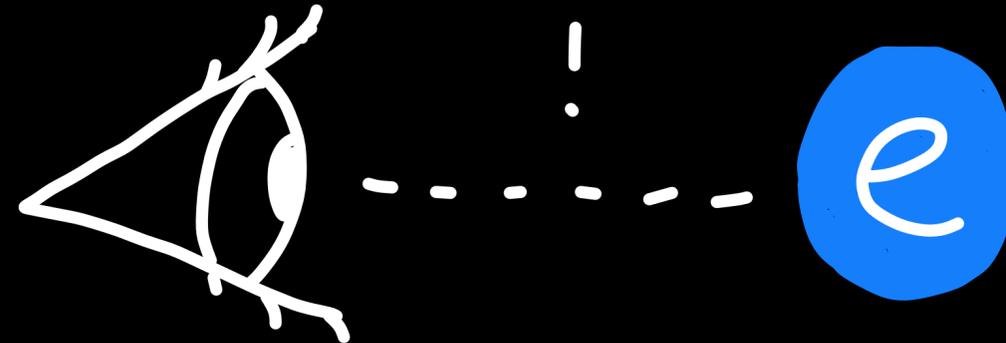
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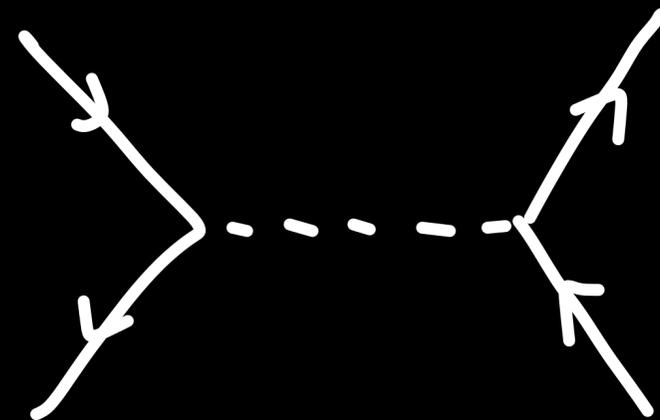


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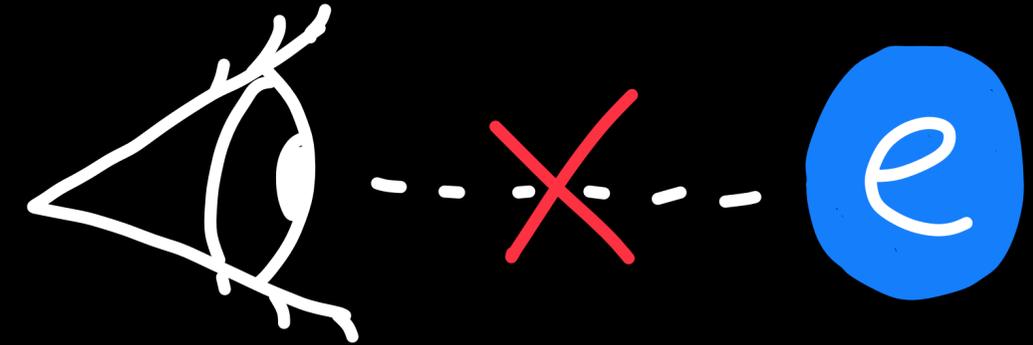


**Elitzur's theorem: Gauge symmetries cannot be spontaneously broken!**

# FMS Framework

[Maas, 2019] & references therein

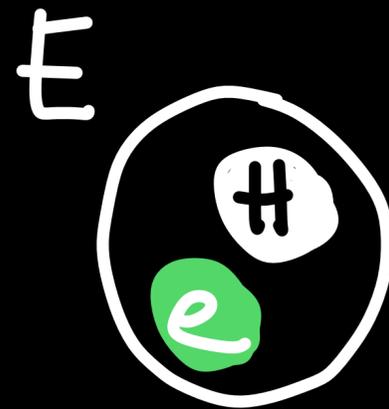
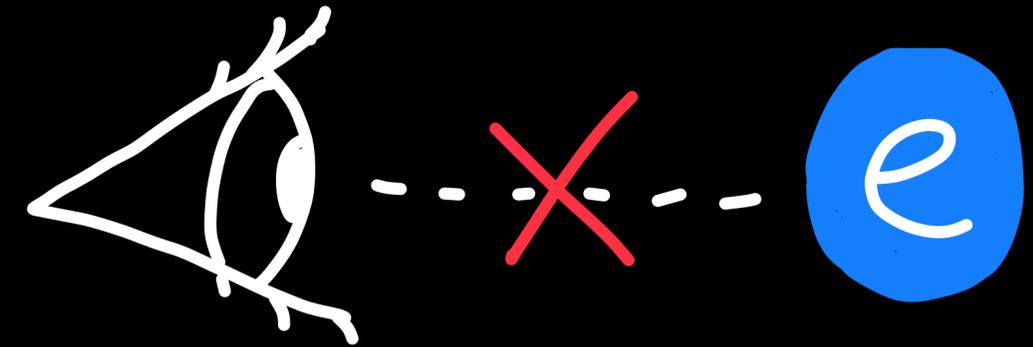
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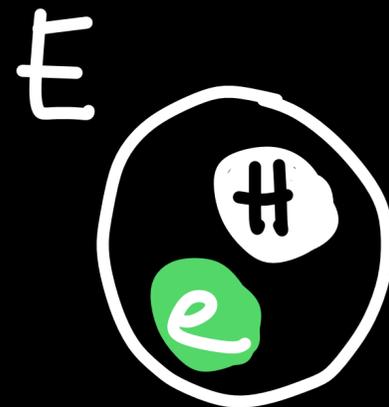
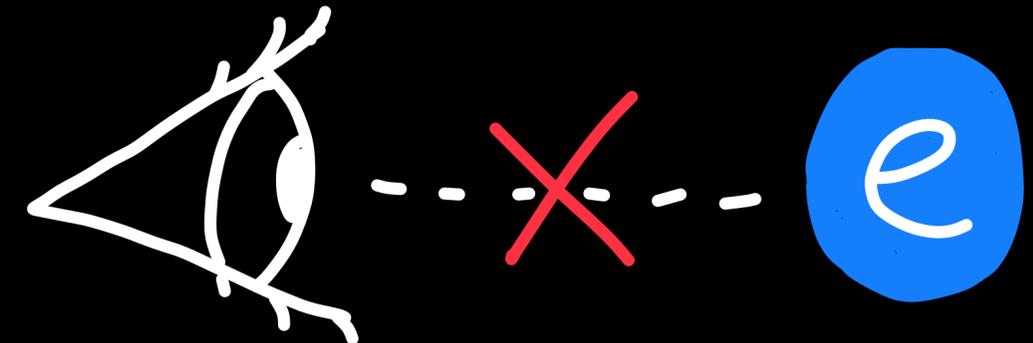


2. Physical fermions are composite

# FMS Framework

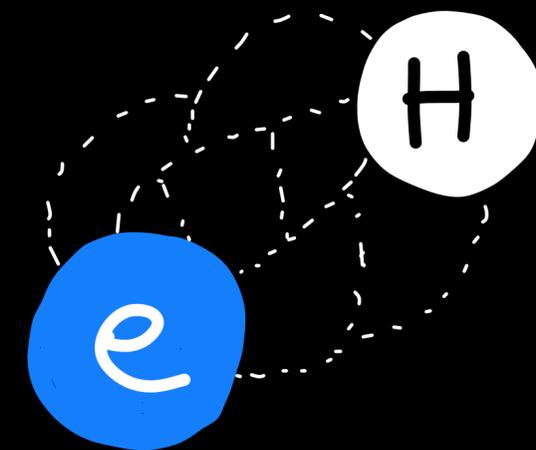
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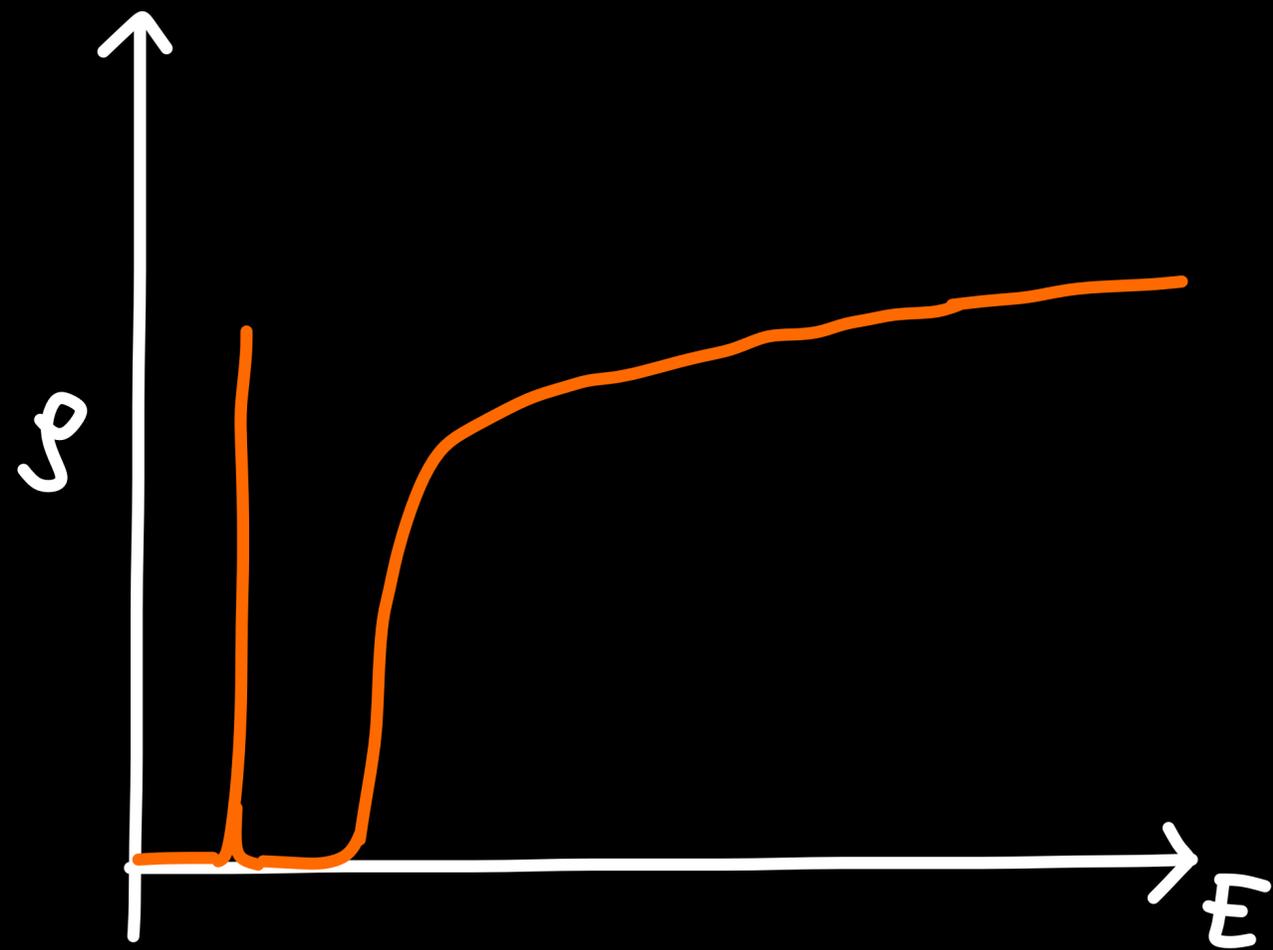
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# Spectral Densities from the Lattice

# Backus Gilbert Approach [Backus, Gilbert, 1968 & 1970]

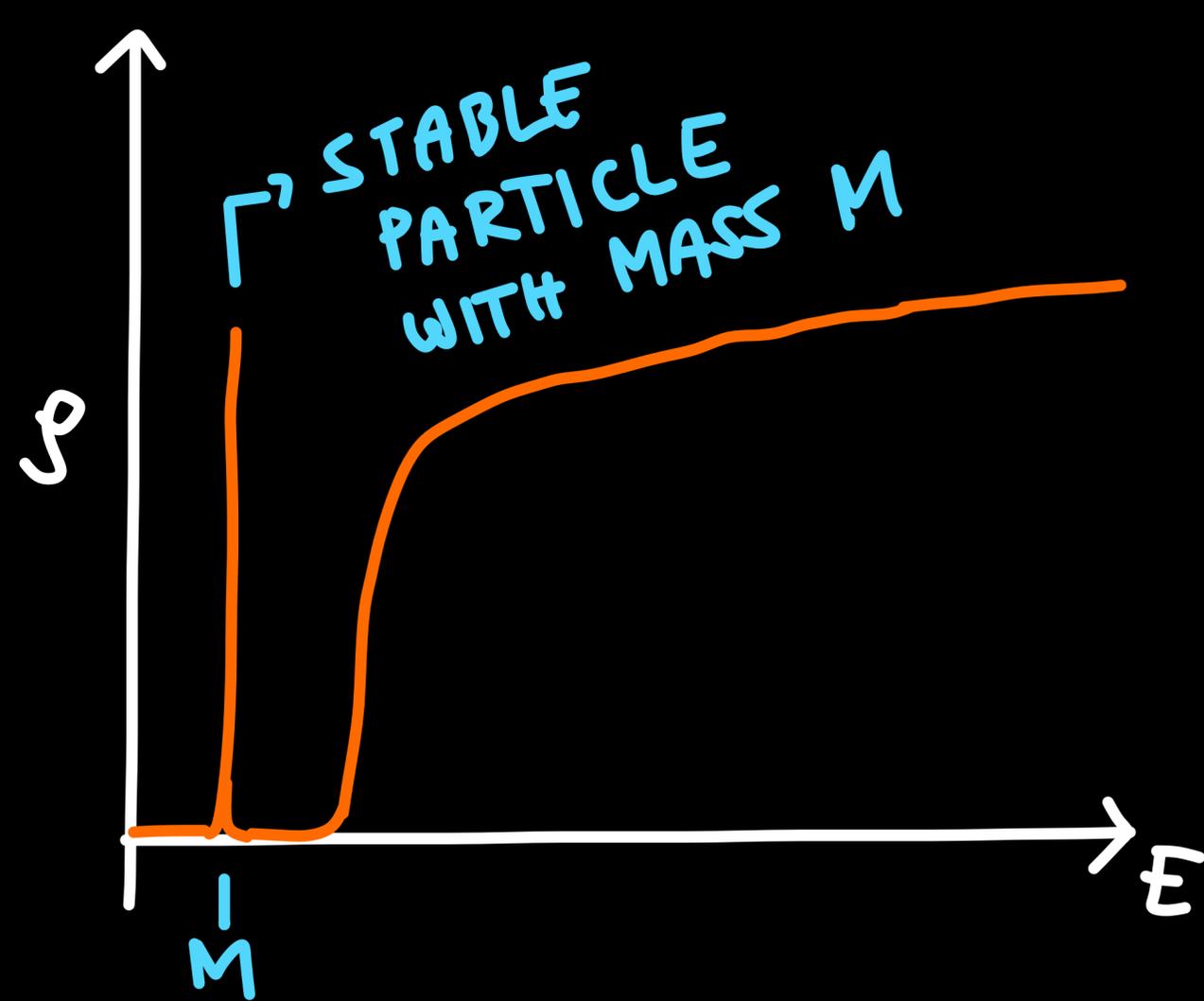


Continuum



Lattice / Finite Volume

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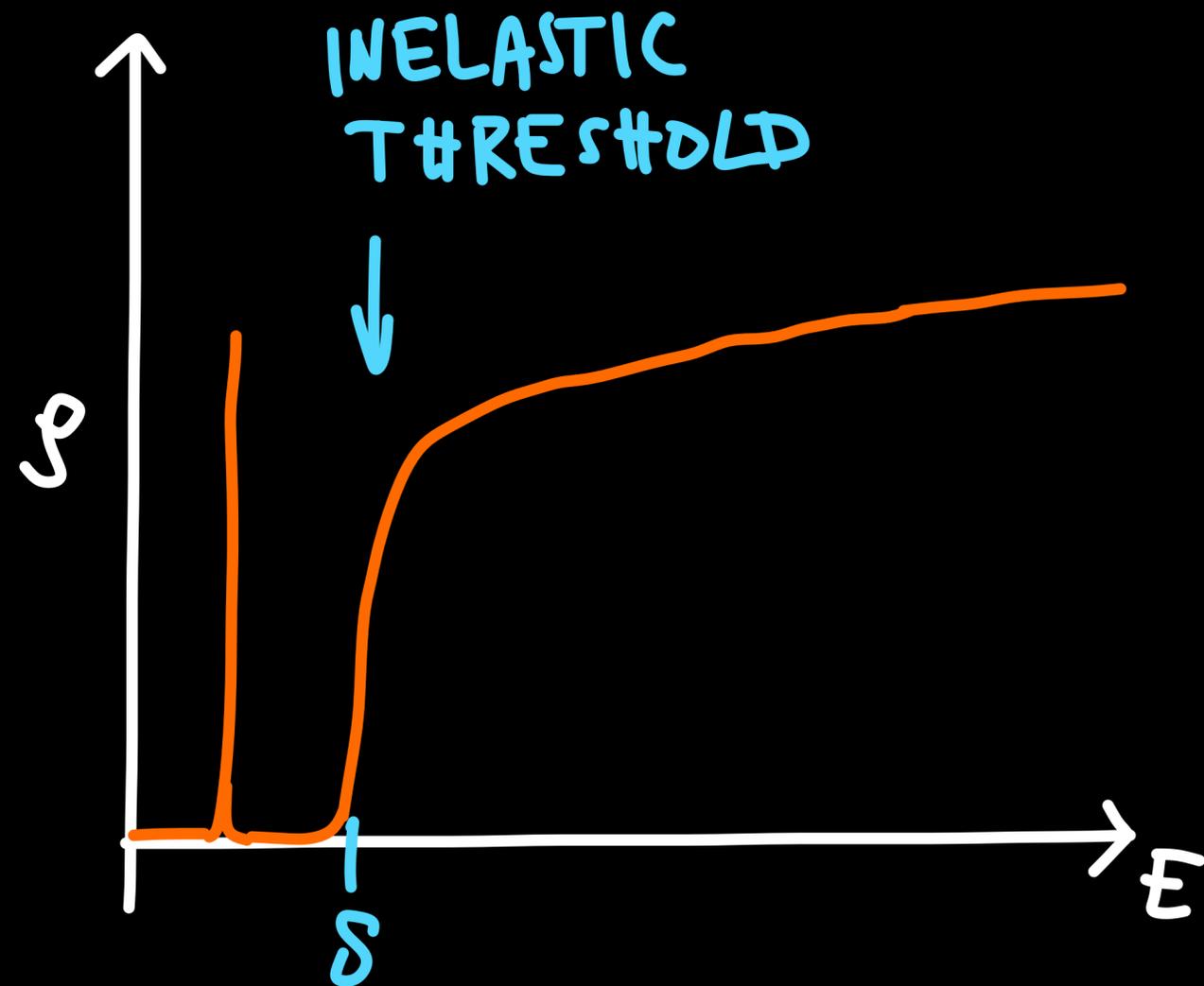


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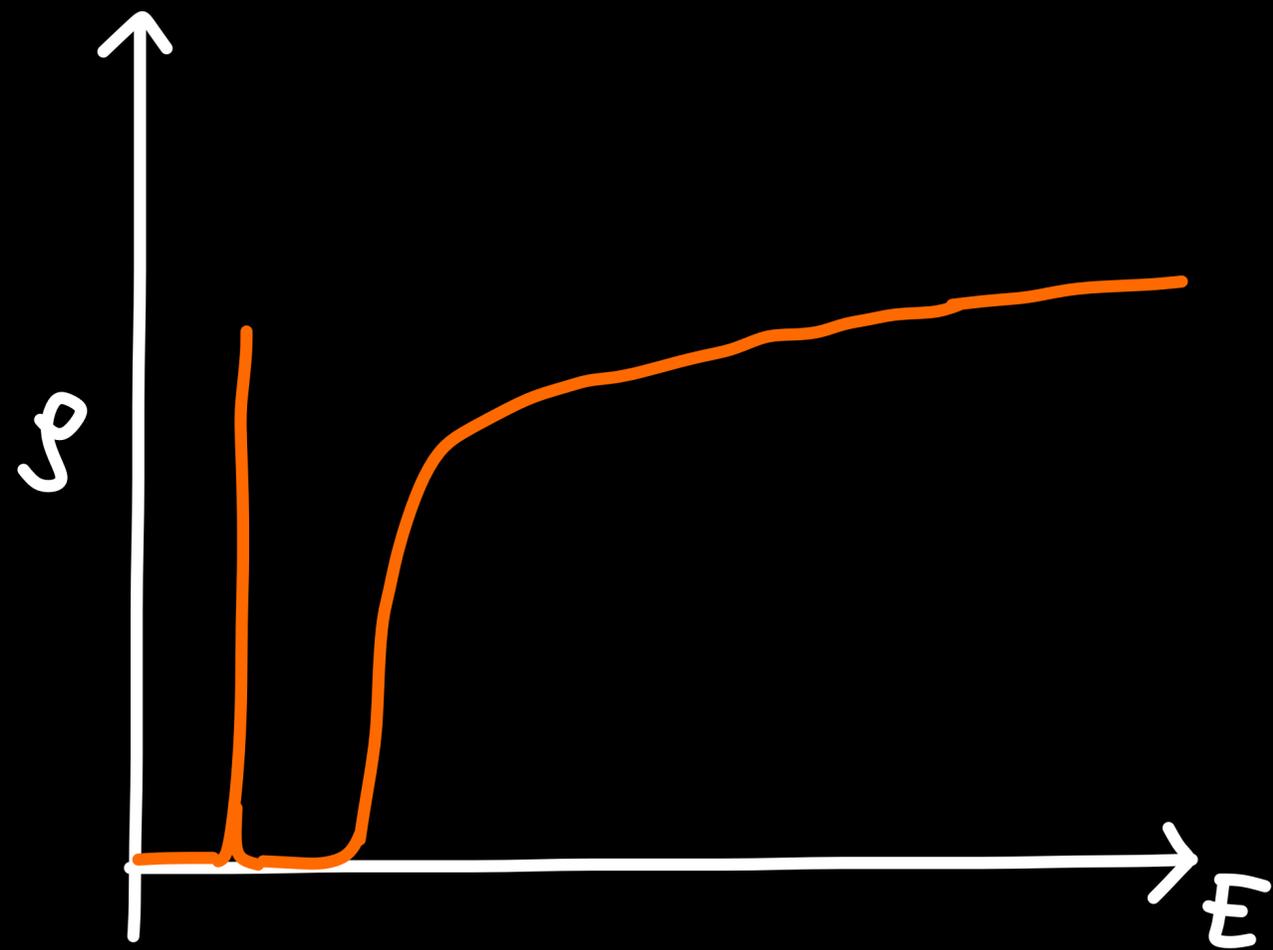


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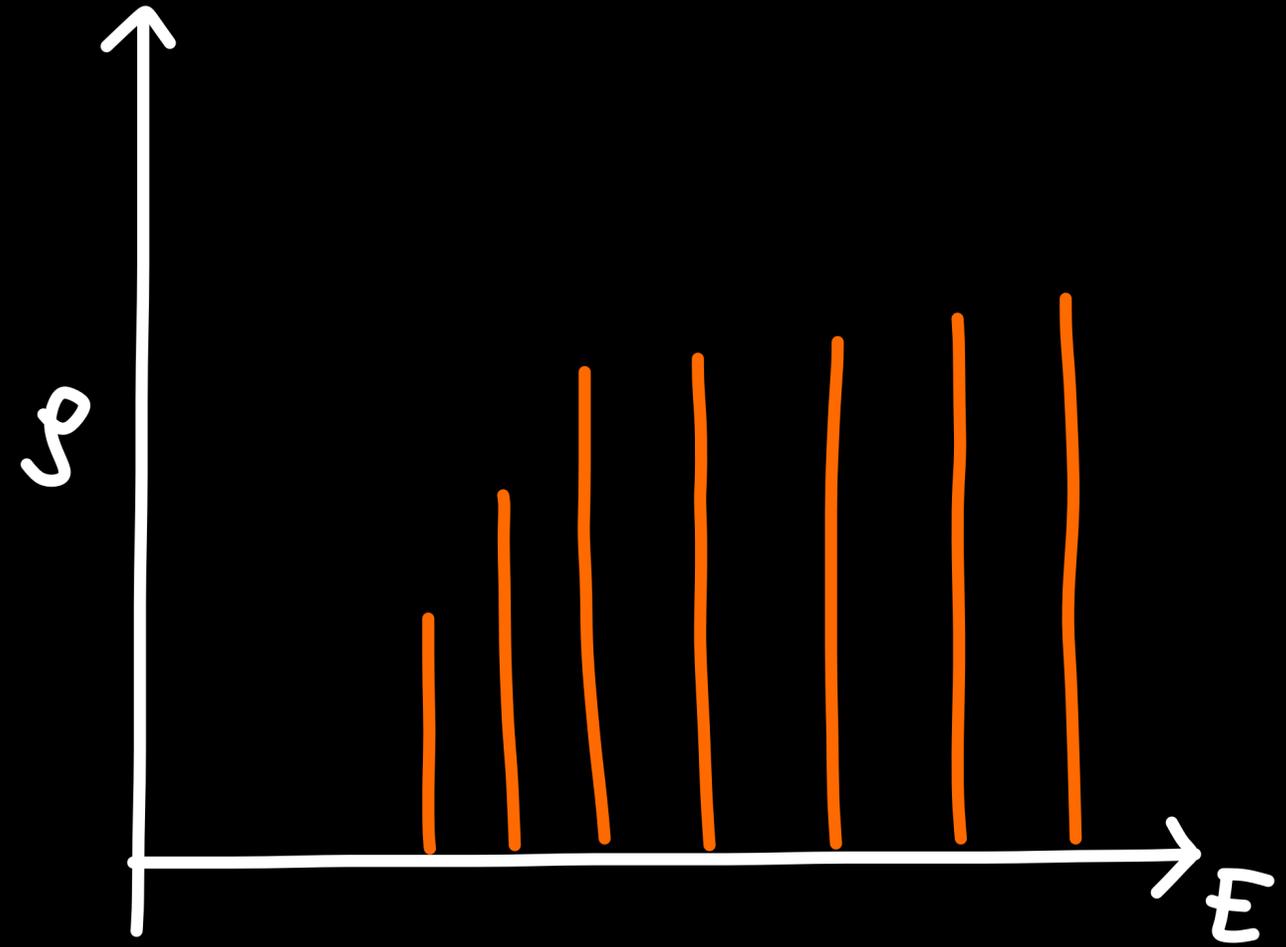


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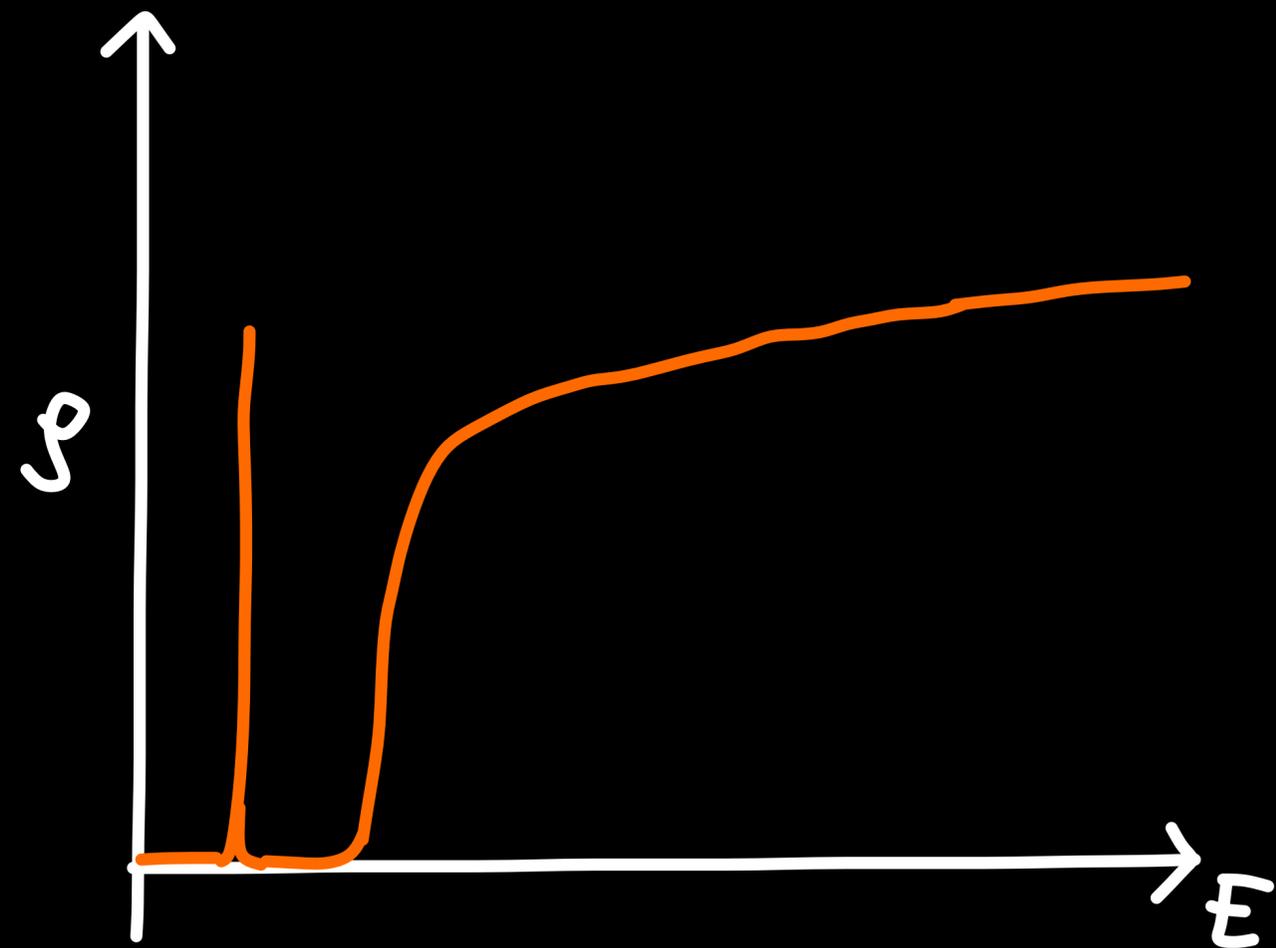
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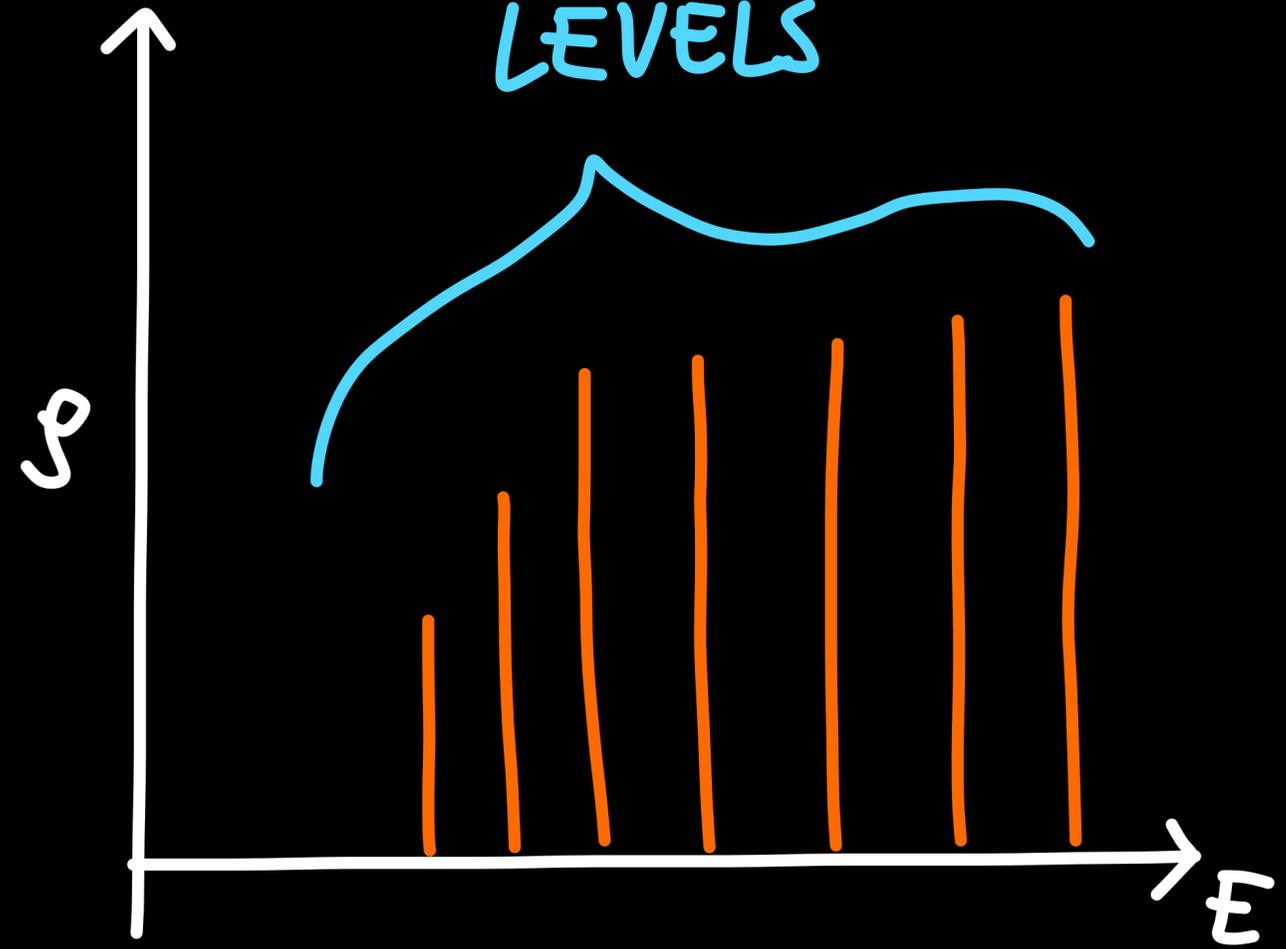
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Continuum

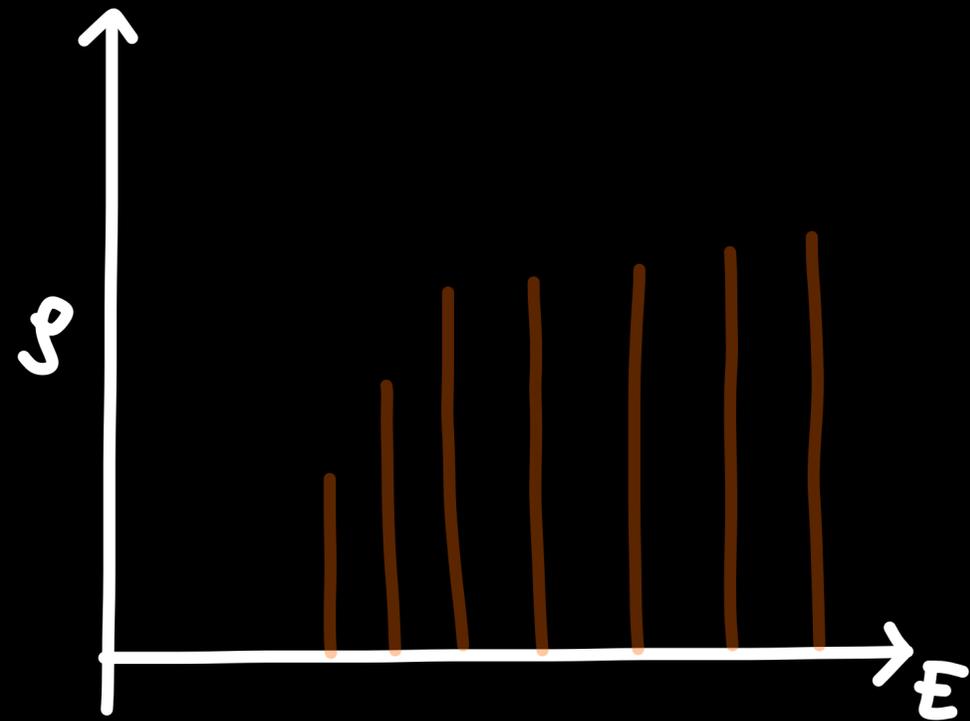
DISCRETE ENERGY LEVELS



Lattice / Finite Volume

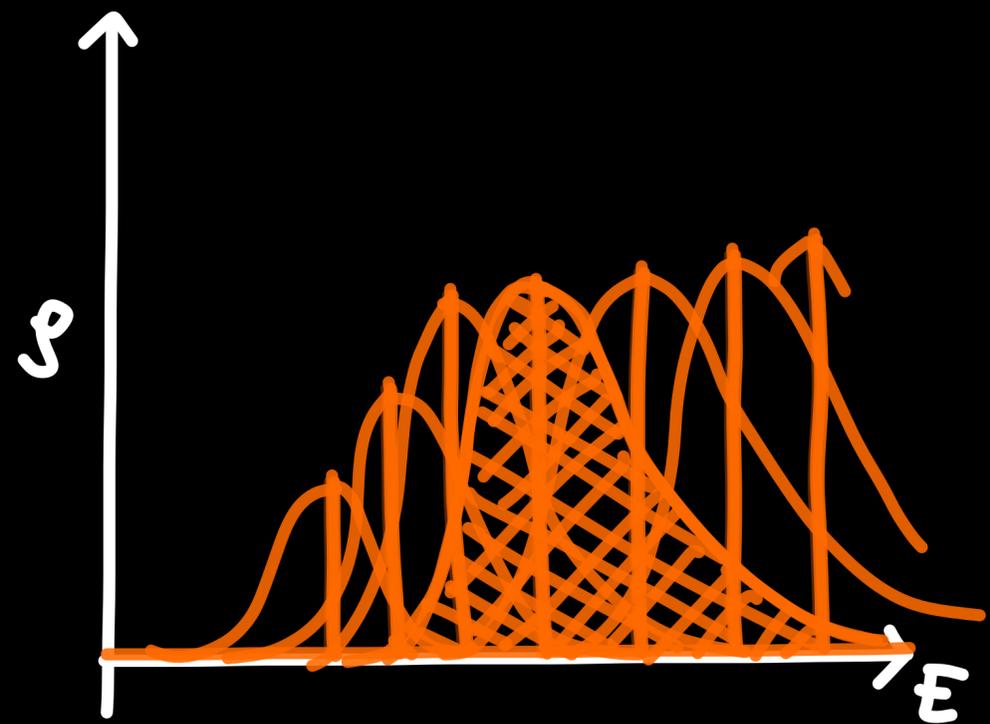
# Backus Gilbert Approach [Backus, Gilbert, 1968 & 1970]

## Smearing



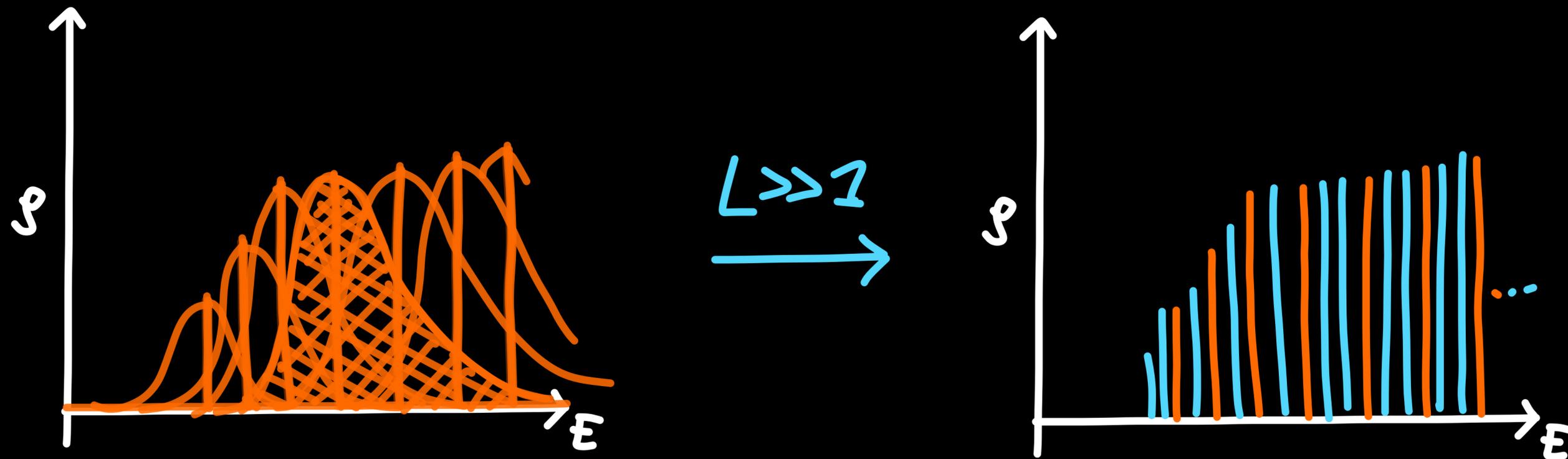
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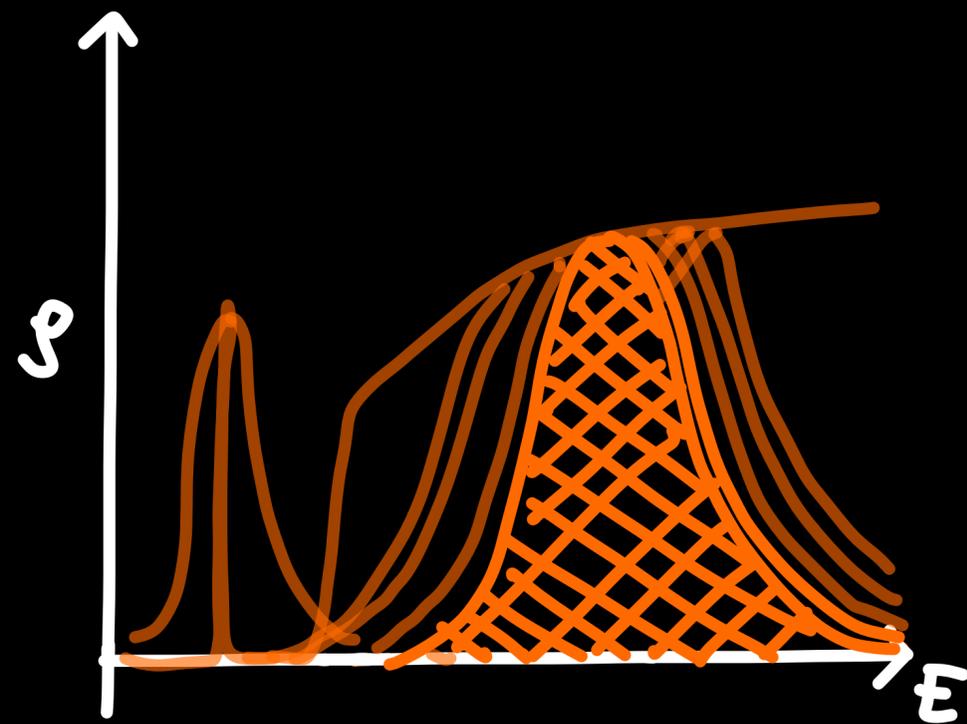
## Smearing



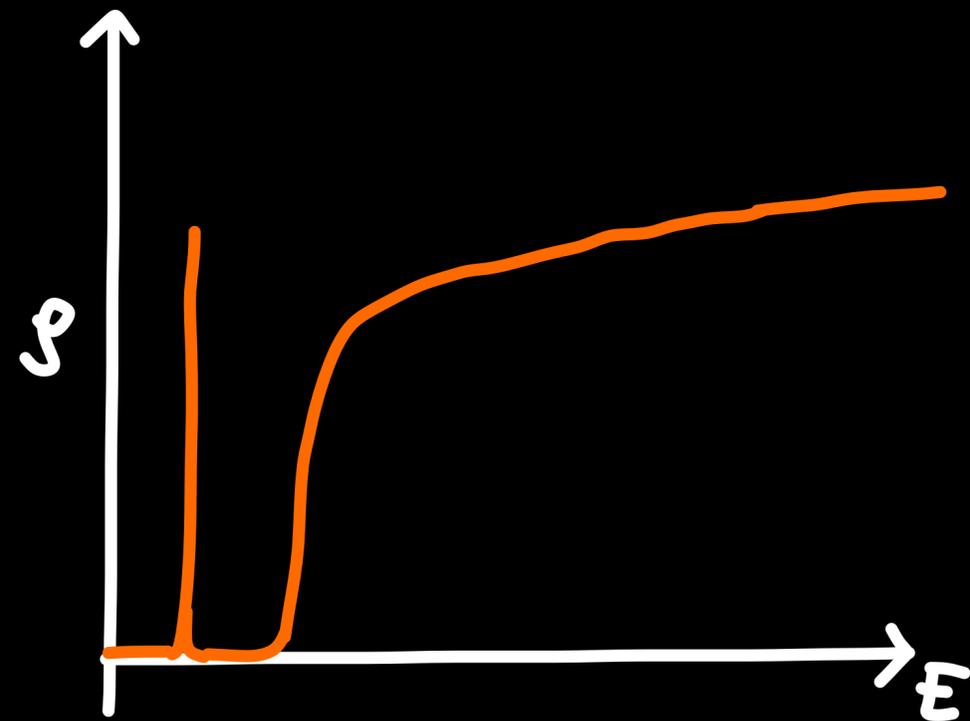
$$\hat{S}_L(\sigma, E_*) = \int_0^{\infty} dE \Delta_{\sigma}(E_*, E) \rho_L(E)$$

# Backus Gilbert Approach [Backus, Gilbert, 1968 & 1970]

## Smearing



$\sigma \rightarrow 0$   
 $\longrightarrow$

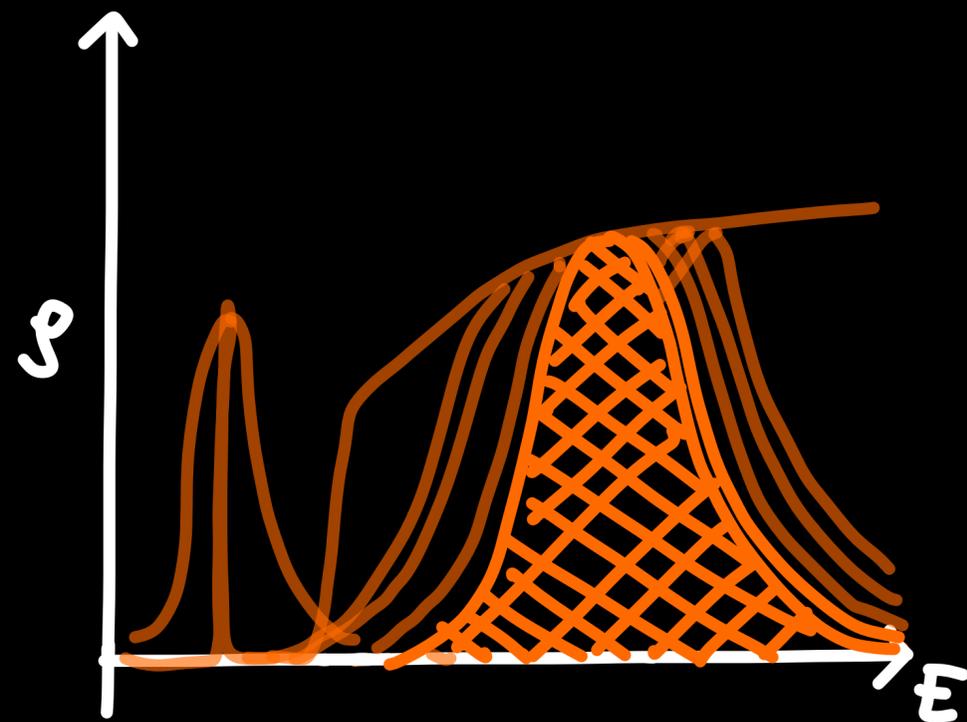


$$\lim_{L \rightarrow \infty} \hat{g}_L(\sigma, E_*)$$

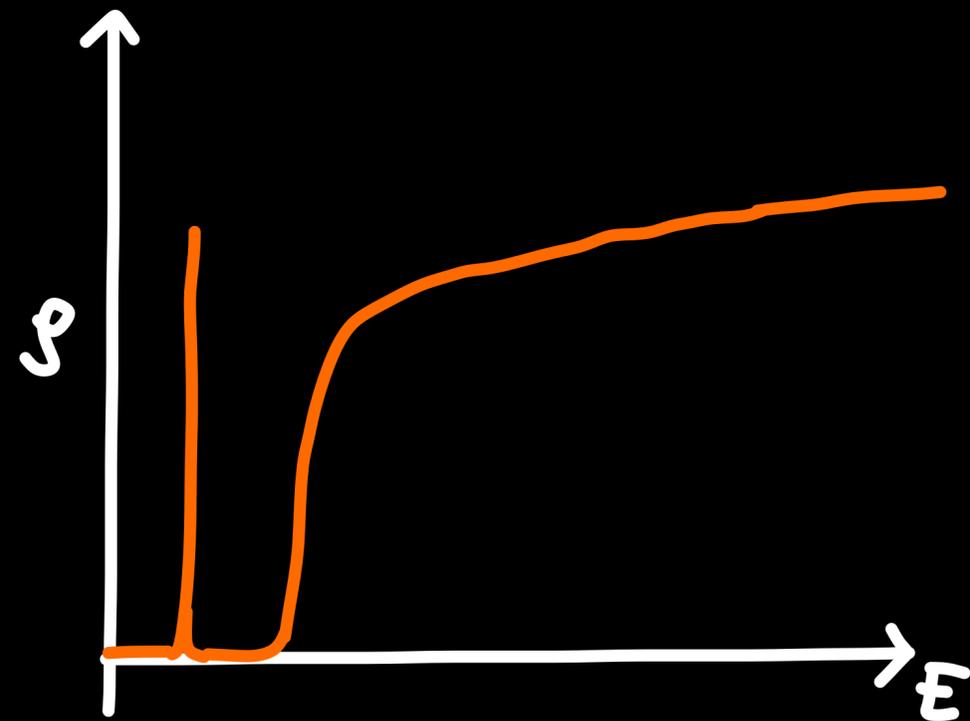
$$g(E_*) = \lim_{\sigma \rightarrow 0} \lim_{L \rightarrow \infty} \hat{g}_L(\sigma, E_*)$$

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$$\lim_{L \rightarrow \infty} \hat{g}_L(\sigma, E_*)$$

$$g(E_*) = \lim_{\sigma \rightarrow 0} \lim_{L \rightarrow \infty} \hat{g}_L(\sigma, E_*)$$

**!! IN THIS ORDER !!**

[Hansen et. al, 2019]

# Hansen-Lupo-Tantalo Method

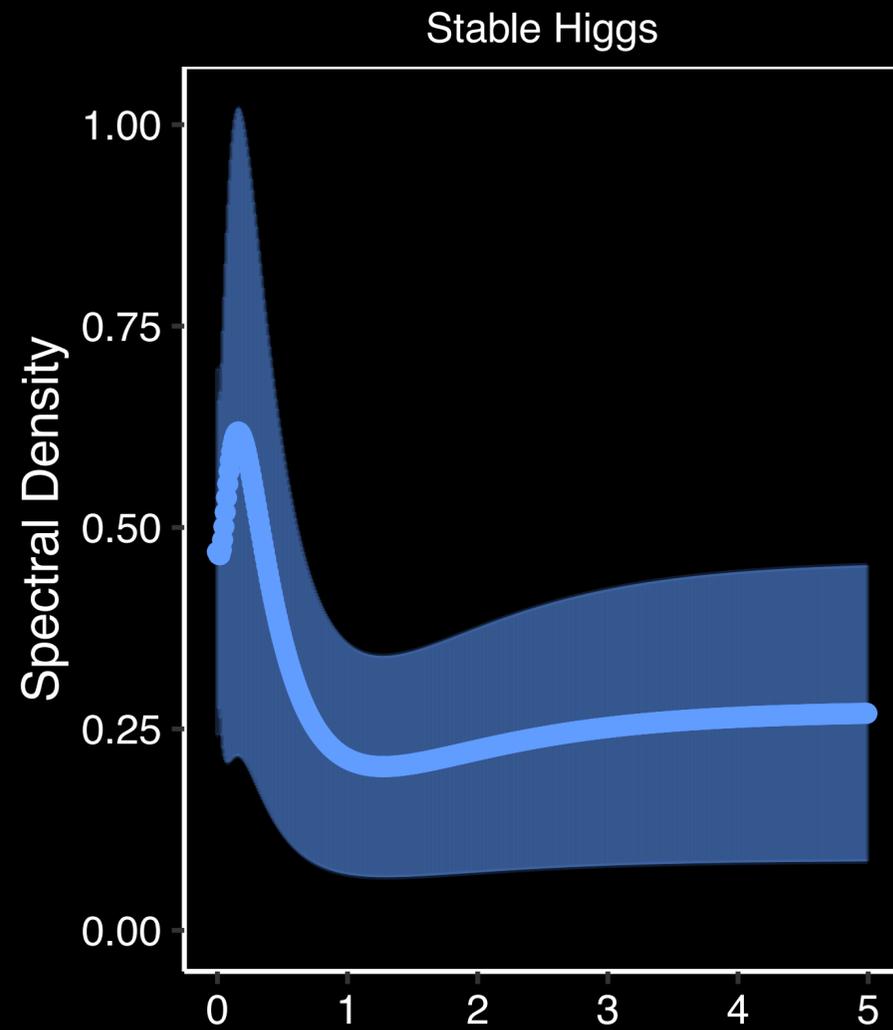
Expand kernel in functional basis preserving backpropagation

$$\Delta(E_*, E) = \sum_{t=0}^T g_t(E_*) \left[ e^{-tE} + e^{-(T-t)E} \right]$$

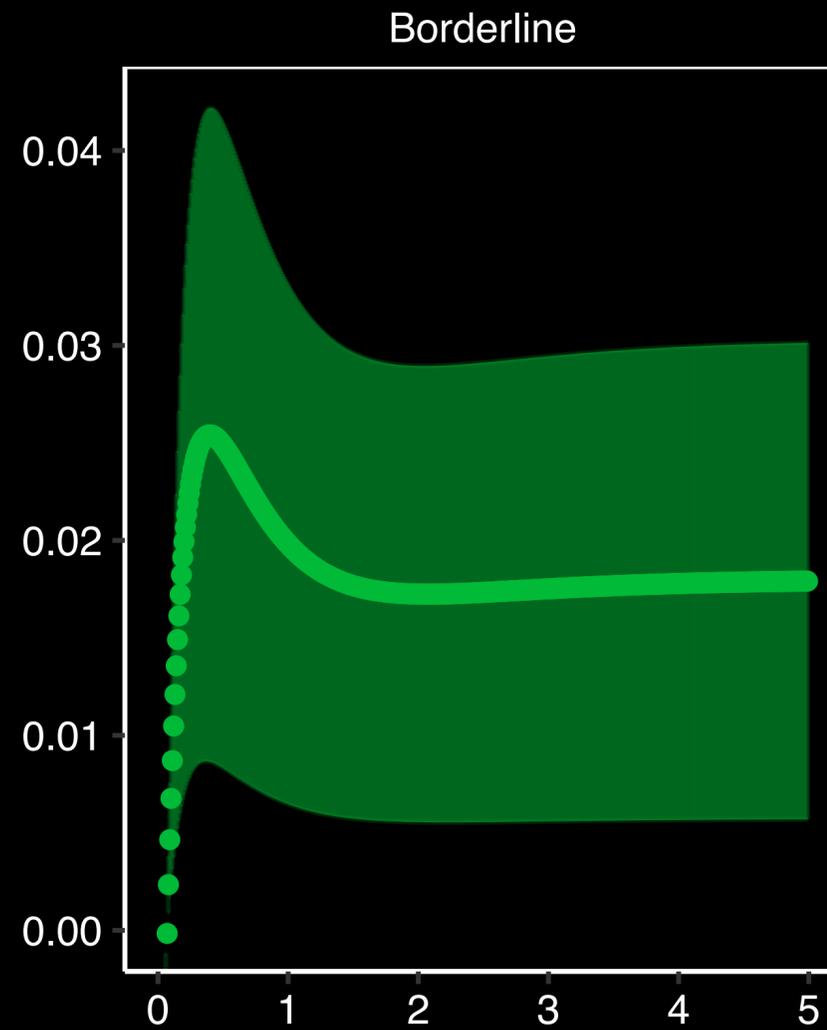
Balance **systematic** and **statistical** uncertainties

$$A[g] = \int_{E_0}^{\infty} dE \left| \overline{\Delta}_{\sigma}(E_*, E) - \Delta_{\sigma}(E_*, E) \right|^2 \quad \frac{\text{Var}(c(t))}{c^2(0)}$$

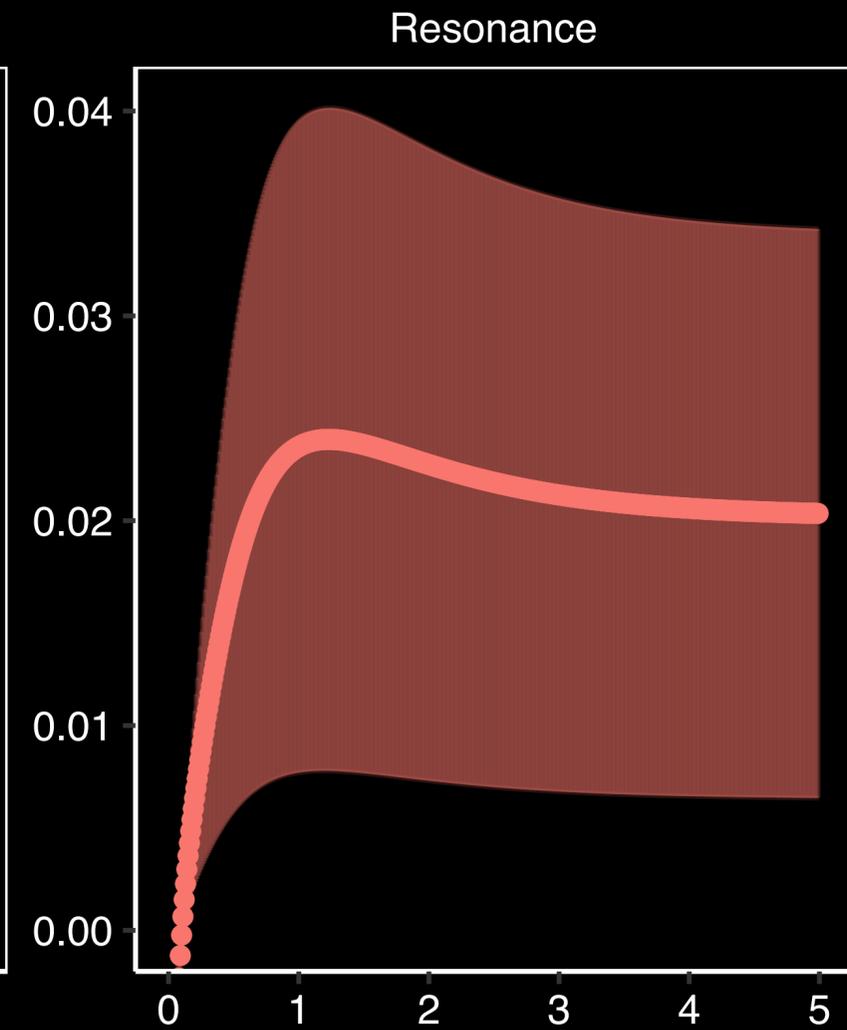
# Higgs stability



$m_W \approx 0.2$   
 $m_H \approx 0.3$   
 $H \not\rightarrow WW$



$m_W \approx 0.3$   
 $m_H \approx 0.6$   
 $H \overset{?}{\rightarrow} WW$



$m_W \approx 0.3$   
 $m_H \approx 1.2$   
 $H \rightarrow WW \checkmark$

Lattice size

- 16
- 20
- 28

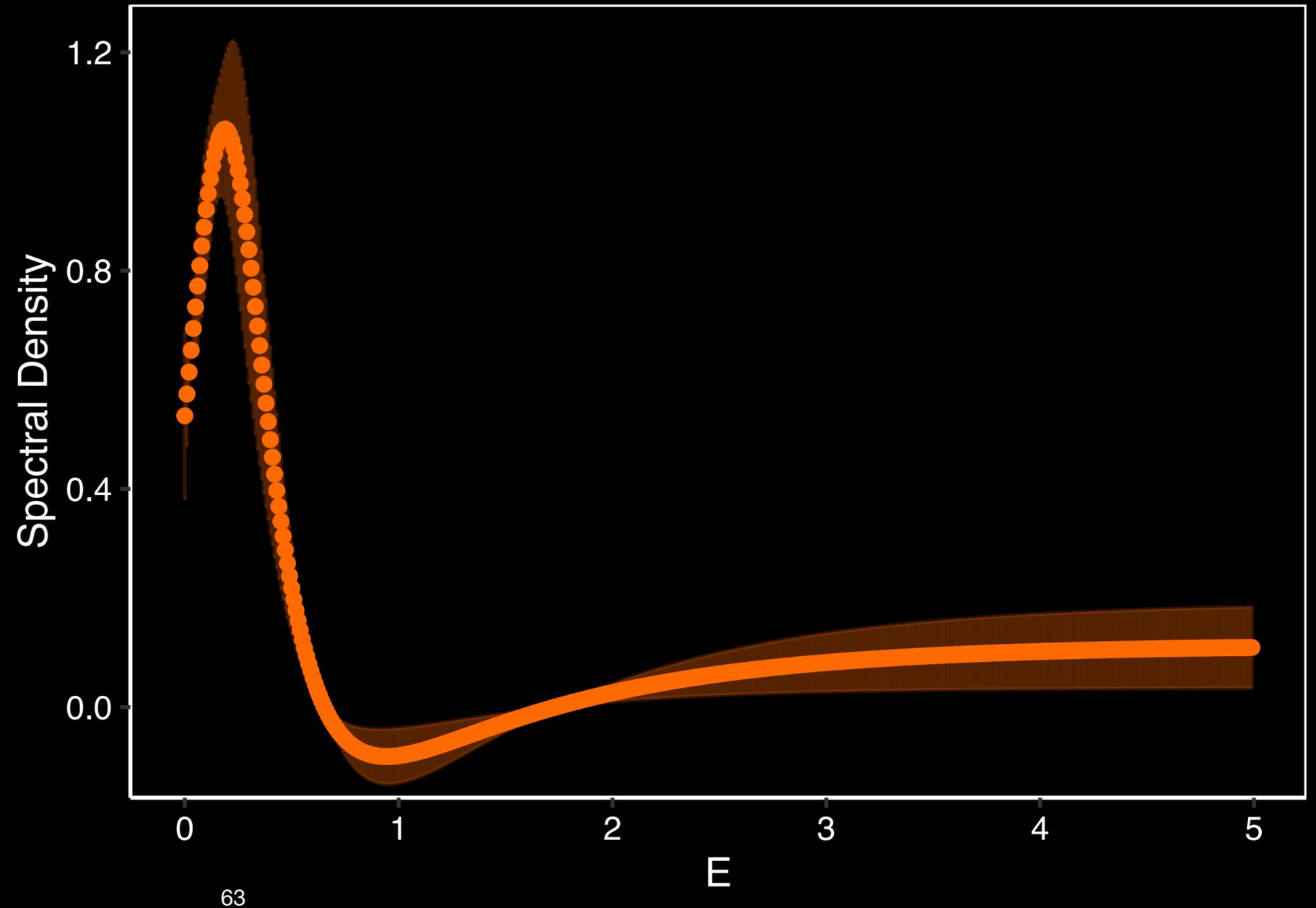
# Example ensemble

The vector channel is much more precise

$$m_w \approx 0.2$$

$$m_H \approx 0.3$$

$$T \times L^3 = 32 \times 32^3$$



# Summary

## Spectral densities

- We can extend scattering amplitudes above the **inelastic threshold** by defining them in terms of **spectral densities**

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# Summary

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- Backus & Gilbert suggested to **smear** this discrete data into continuous data and perform limits to infinite volume and vanishing smearing

# Summary

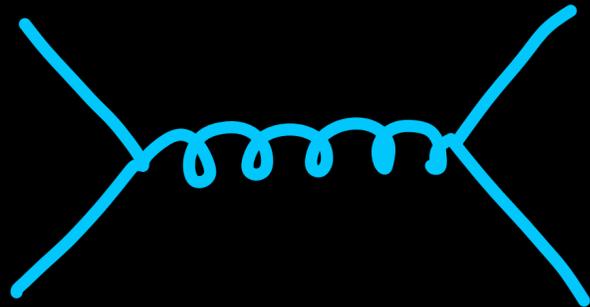
## Spectral densities

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- Computation of spectral densities on the lattice is an **ill-posed problem** because the finite volume enforces discrete energy levels
- Backus & Gilbert suggested to **smear** this discrete data into continuous data and perform limits to infinite volume and vanishing smearing
- HLT improved this by **optimization** of the functional basis w.r.t. **systematic** and **statistical** errors

# Method by Hansen and Bulava

[Bulava, Hansen, 2019]

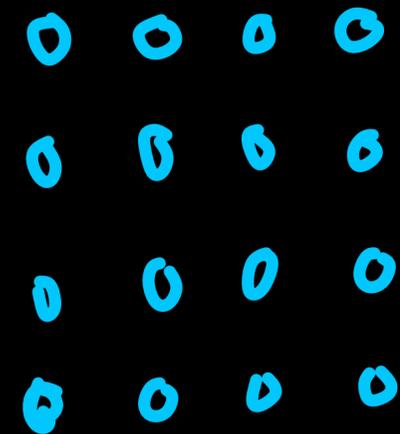
# Idea



What we compute in  
perturbation theory

$$\langle 0 | T \{ \phi(x_1) \cdots \phi(x_n) \phi(y_1) \cdots \phi(y_m) \} | 0 \rangle$$

Time-ordered correlation functions

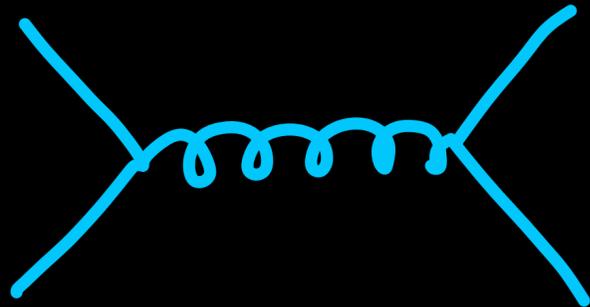


What we compute on the  
lattice

$$\langle 0 | \phi(x_1) \cdots \phi(x_n) \phi(y_1) \cdots \phi(y_m) | 0 \rangle$$

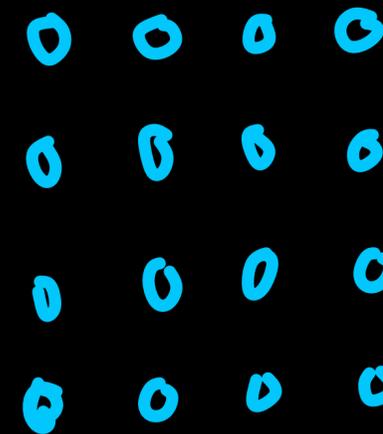
No time ordering

# Idea



What we compute in  
perturbation theory

$$\langle 0 | T \{ \phi(x_1) \cdots \phi(x_n) \phi(y_1) \cdots \phi(y_m) \} | 0 \rangle$$



What we compute on the  
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$$\langle 0 | \phi(x_1) \cdots \phi(x_n) \phi(y_1) \cdots \phi(y_m) | 0 \rangle$$



Identical on-shell pole structure

**Express amplitude as spectral density**

*$m$  particles out,  $n$  particles in*

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$$i\mathcal{M}_c(\{P_f\}|\{q_i\}) \propto S_{P_m q_1}(\{P_f\}\backslash P_m, \{q_i\}\backslash q_1)$$

- Compute this spectral function by solving inverse problem

$$C_{P_m q_1}(\{P_f\}\backslash P_m, \{q_i\}\backslash q_1) = \int \frac{d^r E}{\pi^r} e^{-\Sigma E \Delta t} S_{P_m q_1}(\{P_f\}\backslash P_m, \{q_i\}\backslash q_1)$$

# Compute spectral density

$m$  particles out,  $n$  particles in

- The previously mentioned correlation function is constructed by "amputating" an  $m+n$ -point correlation function

$$C_{P_m q_1}(\{\mathcal{P}_f\} \setminus p_m, \{q_i\} \setminus q_1) = 2 \sqrt{E(p_m) E(q_1)} L^3 z^{\frac{1}{2}}(p_m) z^{\frac{1}{2}}(q_1)$$

$$\lim_{\Delta t_i \rightarrow \infty} \frac{C_{m+n}(\{\mathcal{P}_f\}, \{q_i\}, \{\Delta t_i\})}{C_{2pt}(p_m, \Delta t_{m+4}) C_{2pt}(q_1, \Delta t_1)}$$

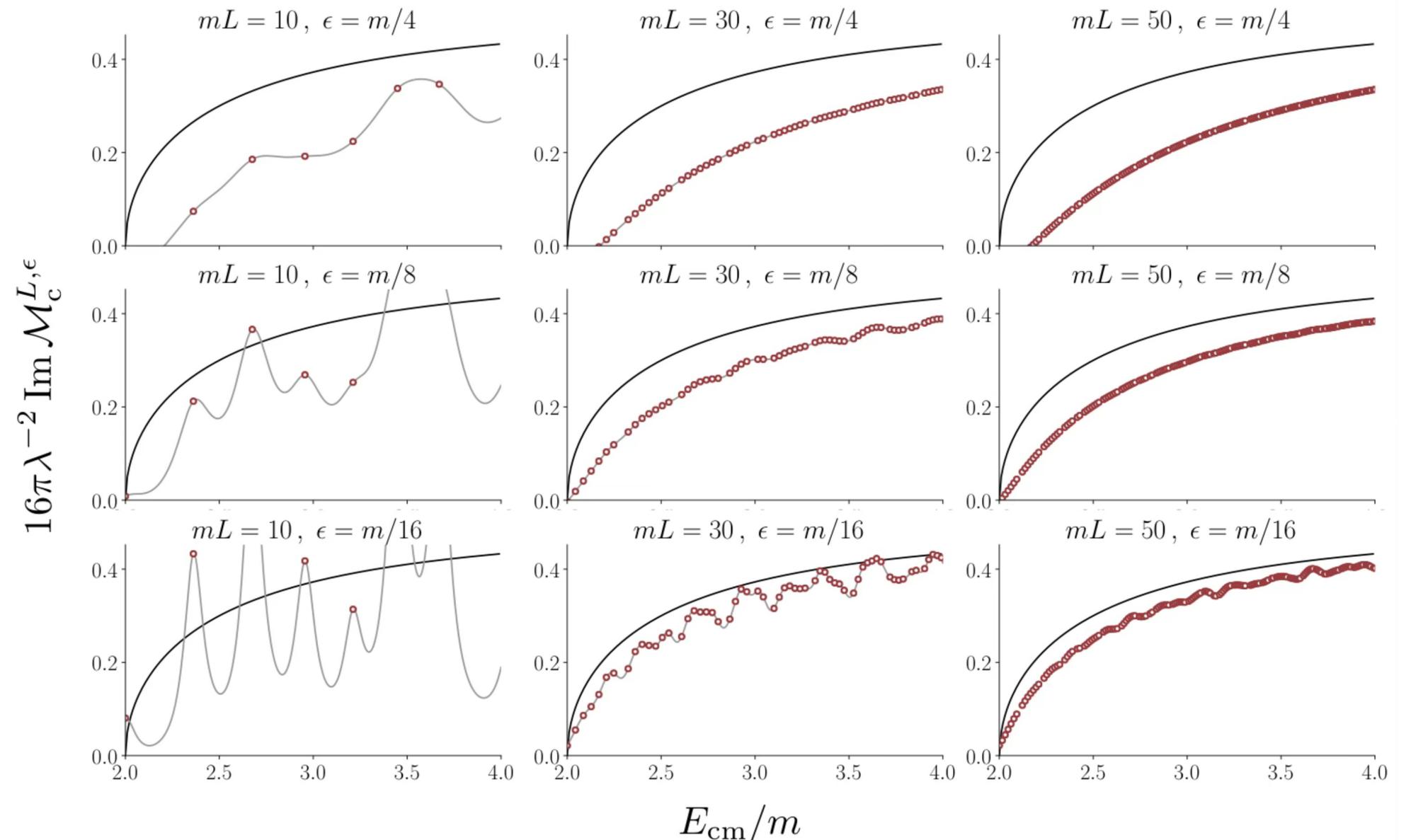
**Examples for  $2 \rightarrow 2$  exclusive  
amplitudes**

# Test of the method

[Bulava, Hansen, 2019]

$\phi^4$ -theory NLO

- Reconstructing next-to-leading order  $2 \rightarrow 2$  process
- Analytical result available (solid line)
- Large finite volume effects

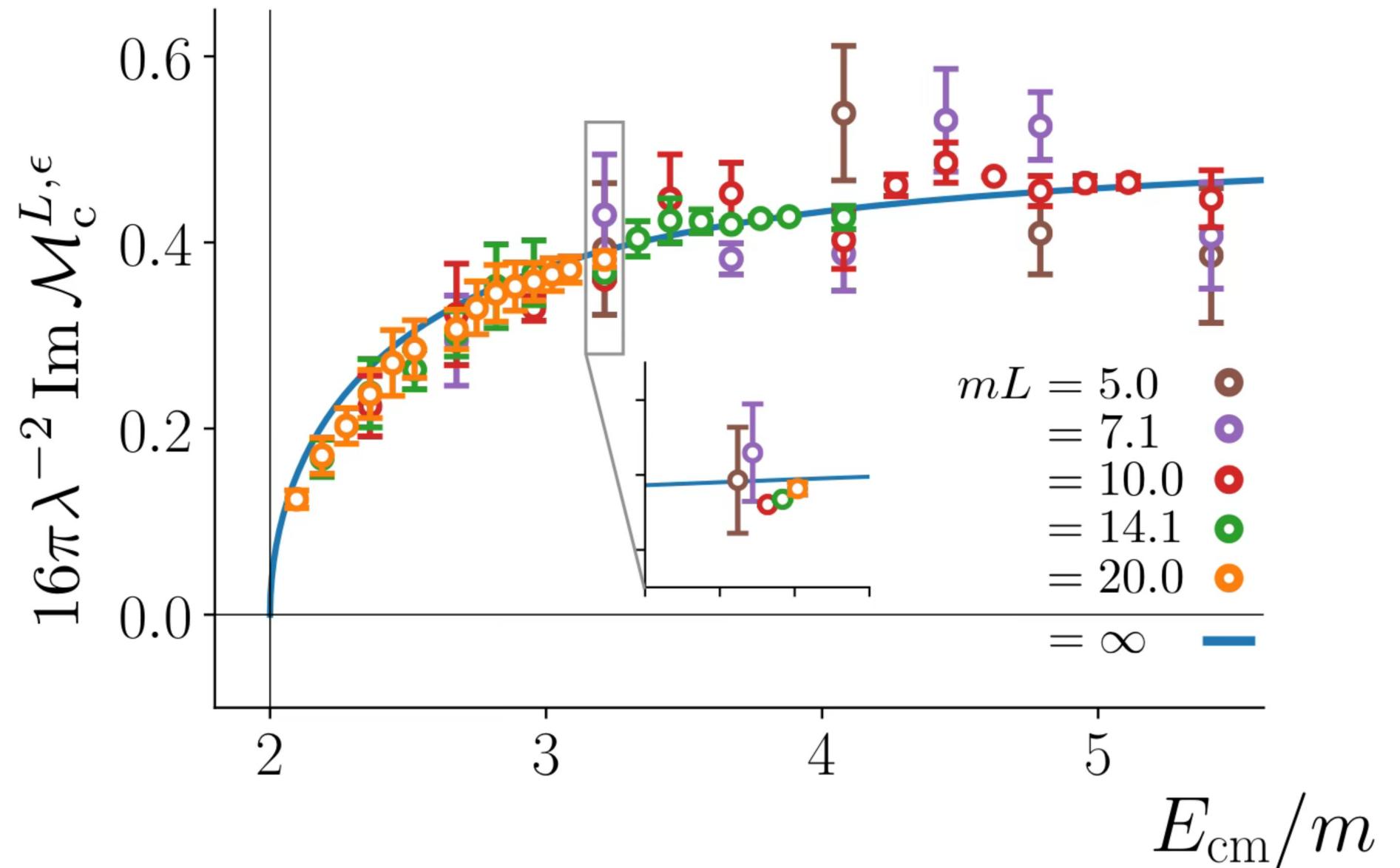


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$\phi^4$ -theory NLO

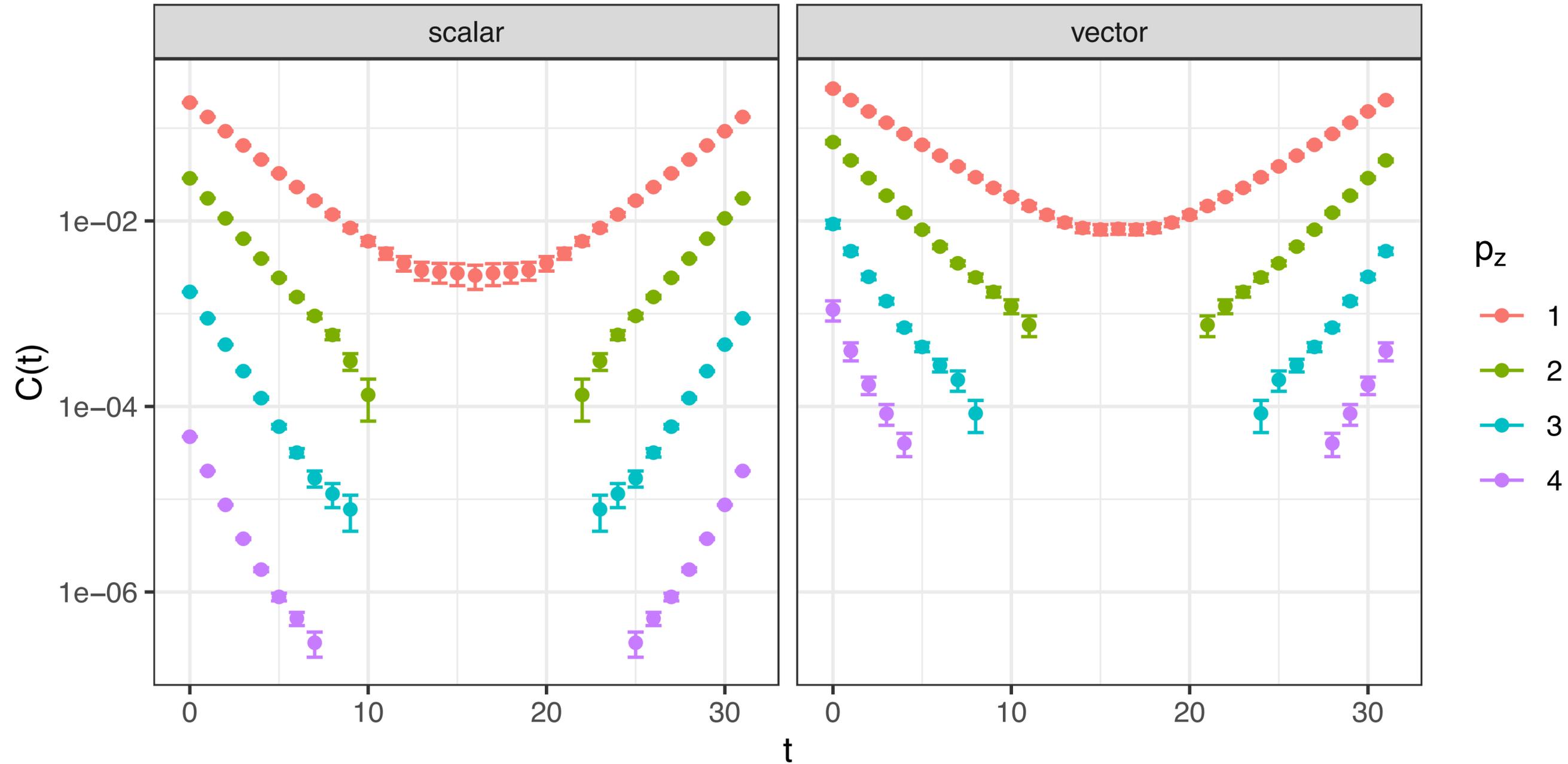
[Bulava, Hansen, 2019]

- Extrapolation to vanishing smearing width



# Electroweak theory

## Boosted correlation functions

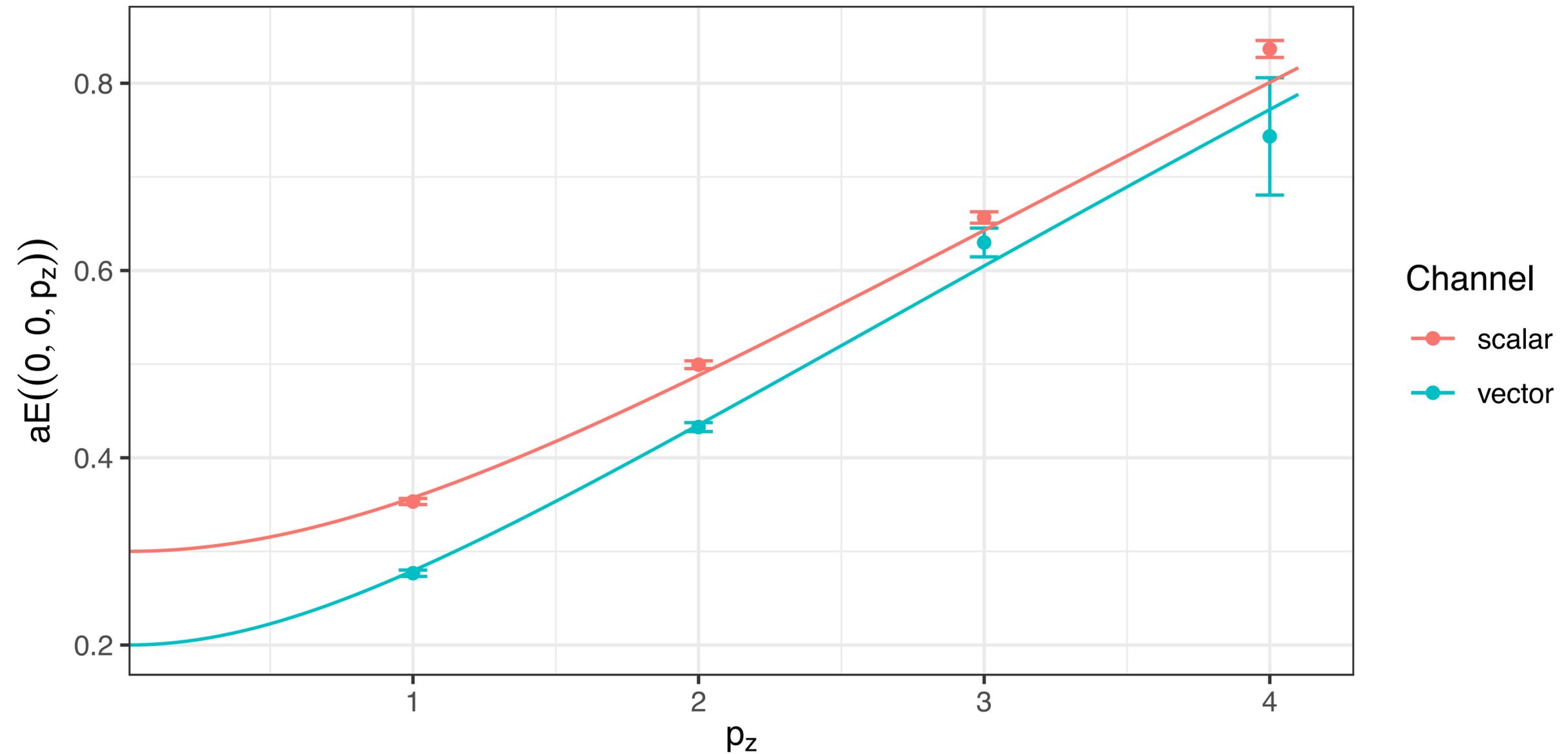


# Electroweak theory

## Naive energies through dispersion relation

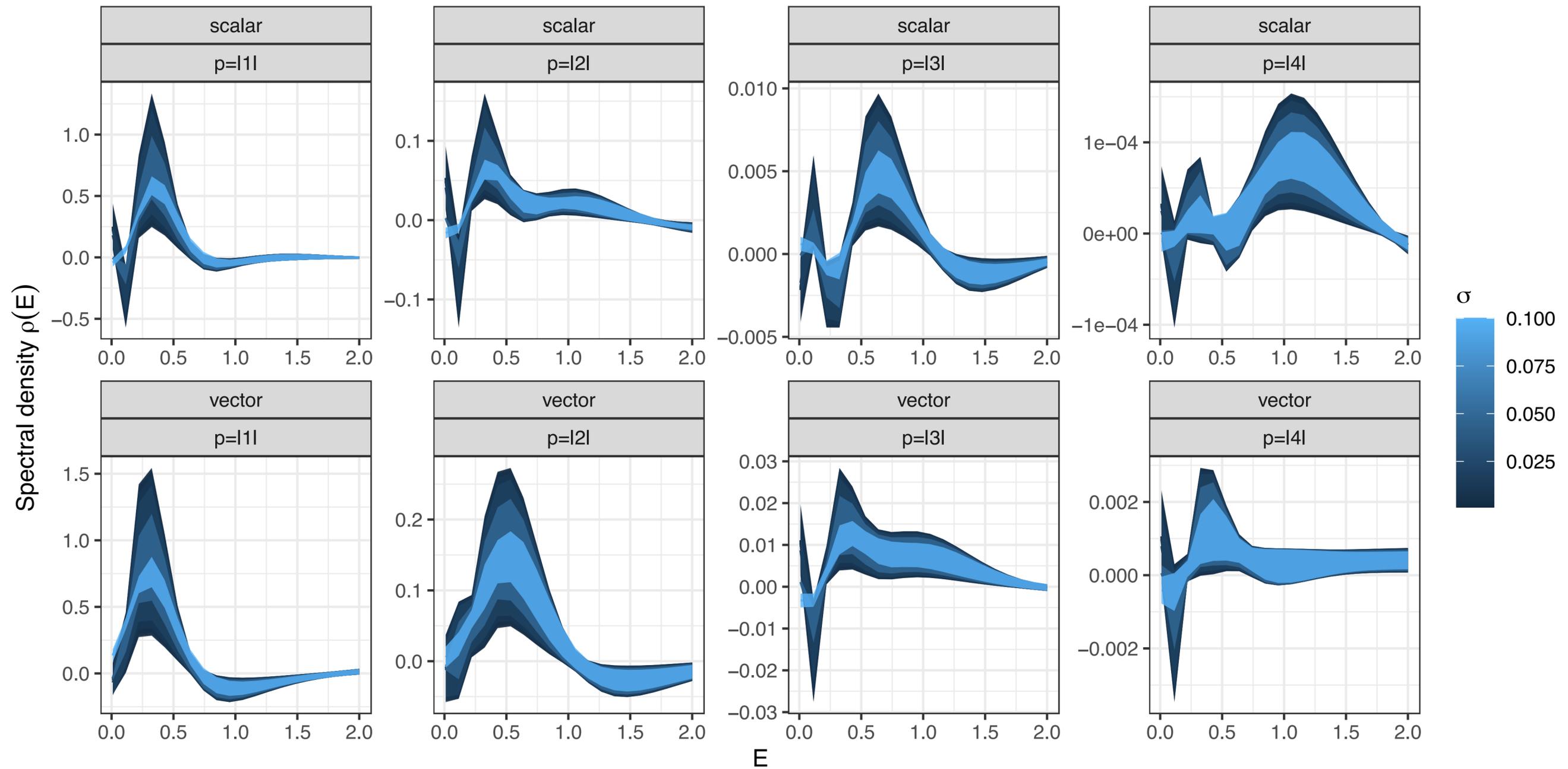
Momentum dependence of mass vs. expected analytical behavior

$$aE(p) = a\cosh(\cosh(am) + 1 - \cos(2 * \pi p_z/L))$$



# Electroweak theory

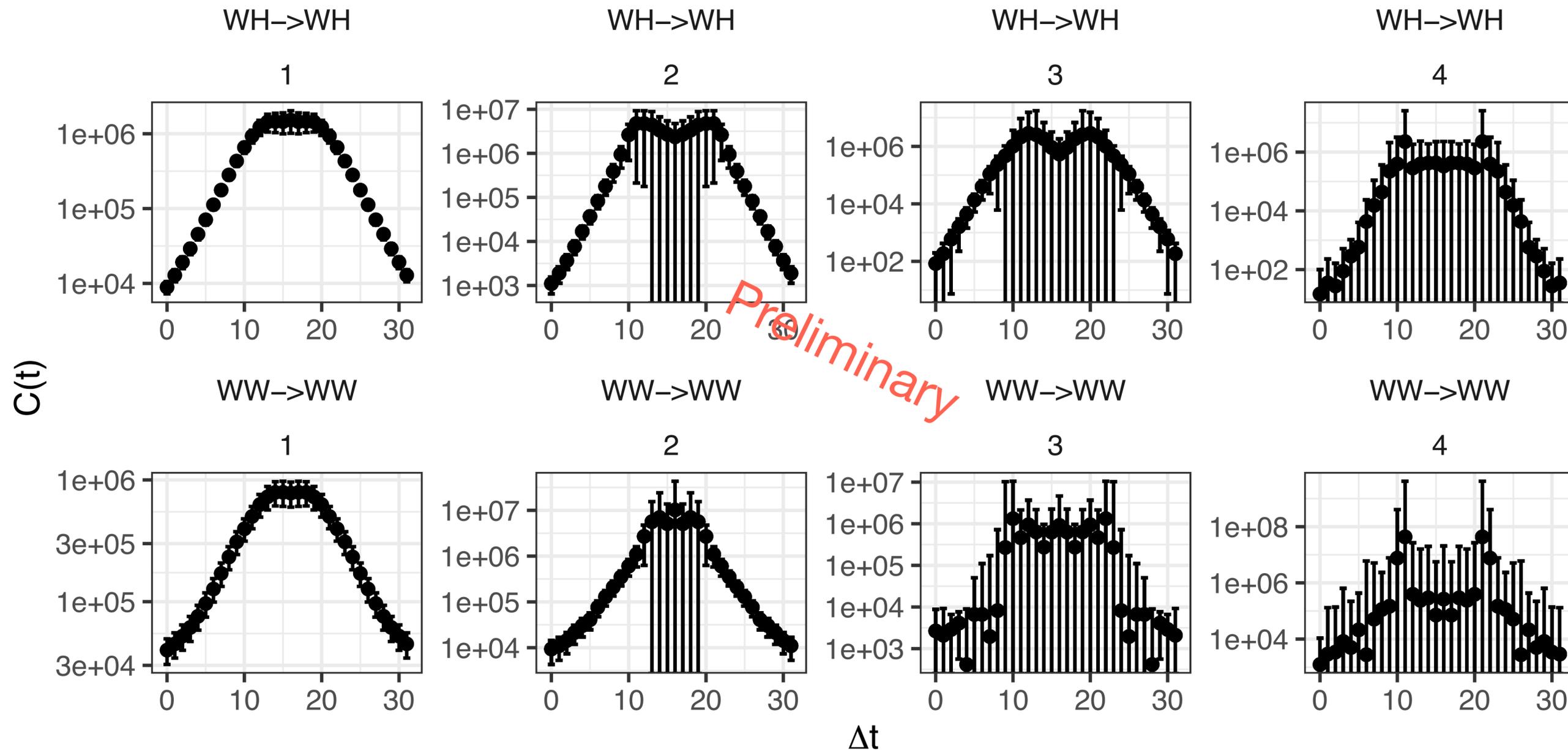
## Boosted spectral densities (two-point functions)



# Electroweak theory

## Ratio functions

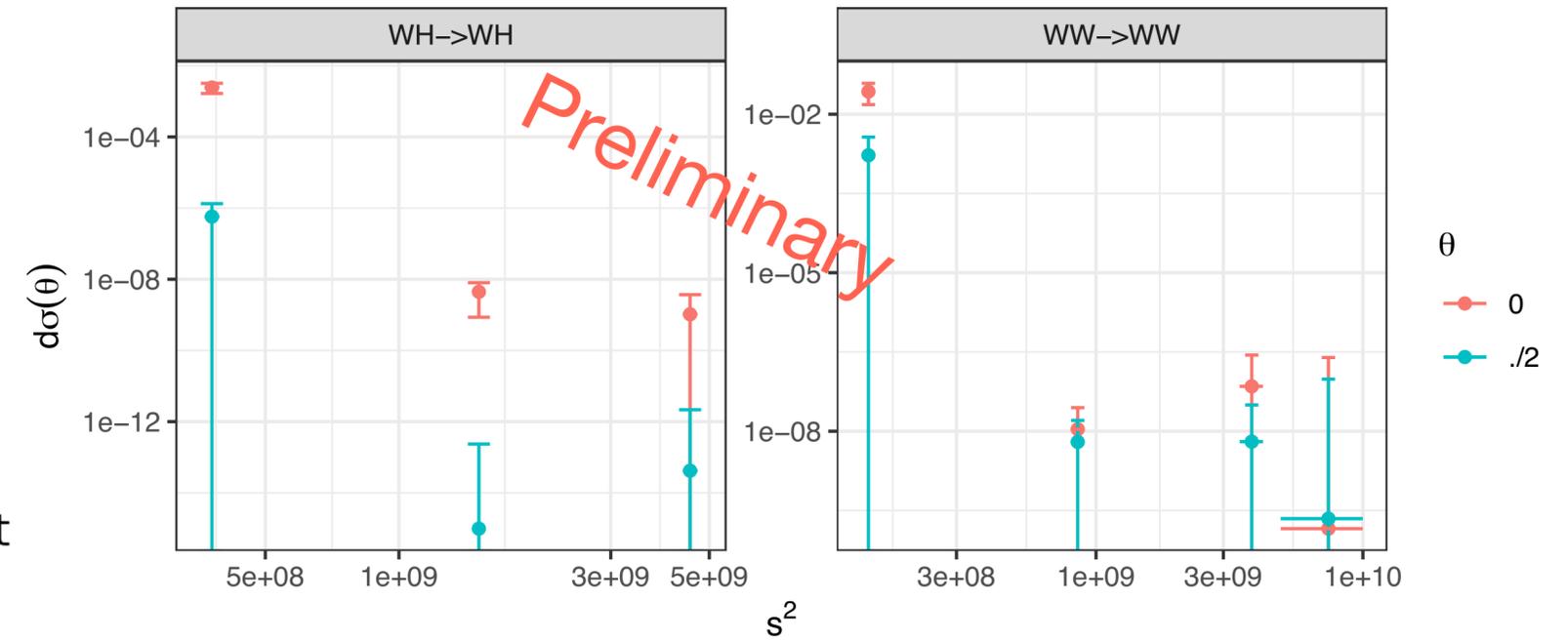
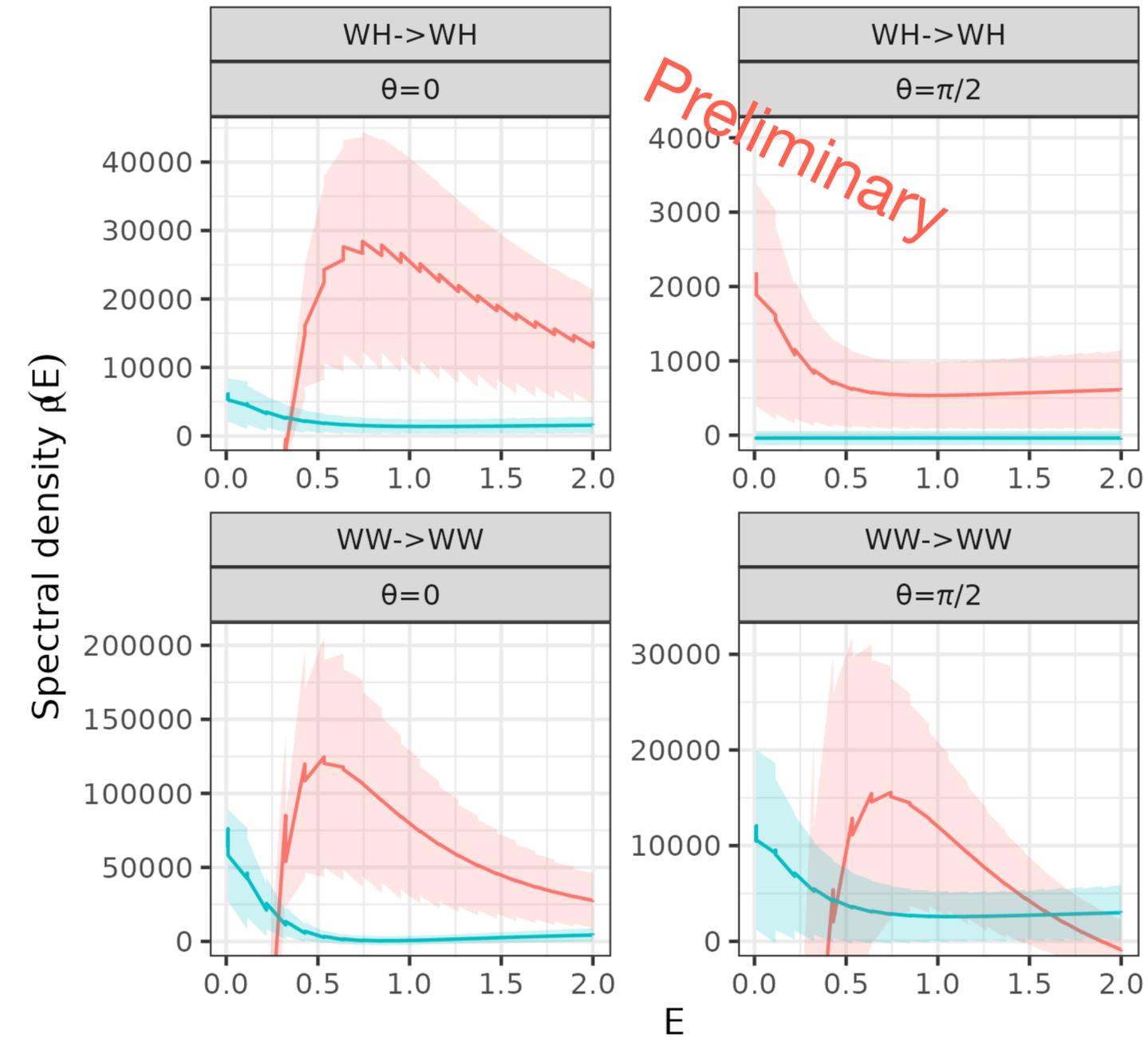
$$C_{P_{in} q_1}(\{P_f\} \setminus P_{in}, \{q_i\} \setminus q_1, \Delta t)$$



# Electroweak theory

## Spectral density of ratio

$$\mathcal{S}_{P_m q_1}(\{P_f\} \setminus P_m, \{q_i\} \setminus q_1)$$



# Patella-Tantalo Approach

[Patella, Tantalo, 2024]

# Skipping the spectral density

Expressing the scattering amplitudes as linear combinations of correlation functions

$$S_c(\sigma, \epsilon) = \sum_{n_1, \dots, n_{n+l-1} \geq 1} \sum_{b \geq 0} \omega_{n,b}^{\sigma, \epsilon}$$

$$\times \int \left[ \prod_{A=1}^{n+l} \frac{d^3 p_A}{(2\pi)^3} f_A^{(*)}(p_A) \right] \tilde{h}(\Delta(p)) \Upsilon(\tau n; P) [\Delta(p)]^b \hat{C}_c(\tau n; P)$$

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Expressing the scattering amplitudes as linear combinations of correlation functions

NUMERICAL  
OPTIMIZATION

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$$\times \int \left[ \prod_{A=1}^{m+l} \frac{d^3 p_A}{(2\pi)^3} f_A^{(*)}(p_A) \right] \tilde{h}(\Delta(p)) \Upsilon(\tau u; p) [\Delta(p)]^b \hat{C}_c(\tau u; p)$$

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KNOWN FUNCTIONS /

KERNELS / CONSERVATION

# Skipping the spectral density

Expressing the scattering amplitudes as linear combinations of correlation functions

Smearing  
Precision

$$S_c(\sigma, \epsilon) = \sum_{n_1, \dots, n_{m+l} - 1 \geq 1} \sum_{b \geq 0} \omega_{n, b}^{\sigma, \epsilon}$$

$$\times \int \left[ \prod_{A=1}^{m+l} \frac{d^3 p_A}{(2\pi)^3} f_A^{(*)}(p_A) \right] \tilde{h}(\Delta(p)) \Upsilon(\tau u; p) [\Delta(p)]^b \hat{C}_c(\tau u; p)$$

No numerical results available (yet)

# Summary and Outlook

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- Hansen-Bulava Method allows computation of general scattering amplitudes from spectral densities: The bottleneck is the precision that the data can achieve in the  $n+m$  point correlation functions

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- Hansen-Bulava Method allows computation of general scattering amplitudes from spectral densities: The bottleneck is the precision that the data can achieve in the  $n+m$  point correlation functions
- Patella-Tantalo Method could skip the ill-posed inversion and directly use the correlation functions. However, no numerical results exist at this point.

# Outlook

- Improve precision of spectral density computation

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↳ Nevanlinna-Pick-Interpolation [Fields, Christ, 2025]

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↳ Add results for resonance Higgs

- Scan physical parameter space

**Thank you!**