

Recent developments in **integral transform** approaches

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Integral Transforms as an alternative method for dealing with the continuum

Reactions to continuum

perturbative (electro-weak)



non-perturbative (hadronic)



*a, b, c, d... are single nucleons or bound nuclear systems
These reactions involve the dynamics of A nucleons
A-BODY PROBLEM IN THE CONTINUUM!*

Reactions to continuum

Framework:

- Energies in the non-relativistic regime
→ Non-Relativistic Quantum Mechanics
(including *Translation*, *Galileian*, *Rotational* ... invariances)
 $[H, \mathbf{P}_{cm}] = 0$ $[H, \mathbf{R}_{cm}] = 0$ $[H, \mathbf{J}] = 0$...
- Degrees of freedom: all A nucleons
("microscopic" model)

- $H = T + V$ $V = \sum_{ij} v_{ij} + \sum_{ijk} v_{ijk} + \dots$

Reactions to continuum

perturbative (electro-weak)

$$\gamma^{(*)} + b \text{ ----> } c + d + \dots$$

Reactions to continuum

perturbative (electro-weak)

$$\gamma^{(*)} + b \longrightarrow c + d + \dots$$

- **First order** perturbation theory
(*Fermi-Golden Rule*)
- **Linear** Response theory

Reactions to continuum

perturbative (electro-weak)



- **First order** perturbation theory (*Fermi-Golden Rule*)
- **Linear** Response theory

$$\sigma(\omega) \sim \left| \langle \mathbf{n} | \Theta | \mathbf{0} \rangle \right|^2 \delta(\omega - E_n + E_0)$$

$$H | \mathbf{n} \rangle = E_n | \mathbf{n} \rangle$$

Reactions to continuum

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Energy transferred by
the perturbative probe

perturbative (electro-weak)



Reactions to continuum

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$$\sigma(\omega) \sim \left| \langle n | \Theta | 0 \rangle \right|^2 \delta(\omega - E_n + E_0)$$

Ground state of the target
A-body bound state!

perturbative (electro-weak)



Reactions to continuum

- **First order** perturbation theory
(*Fermi-Golden Rule*)
- **Linear** Response theory

$$\sigma(\omega) \sim \left| \langle \mathbf{n} | \Theta | \mathbf{0} \rangle \right|^2 \delta(\omega - \mathbf{E}_n + E_0)$$

Fragmented target
A-body continuum state!

perturbative (electro-weak)



Reactions to continuum

- **First order** perturbation theory
(*Fermi-Golden Rule*)
- **Linear** Response theory

$$\sigma(\omega) \sim |\langle \mathbf{n} | \Theta | \mathbf{0} \rangle|^2 \delta(\omega - E_n + E_0)$$

Property of the target responsible
of the interaction with the perturbative probe
e.g. charge-current density

perturbative (electro-weak)



Reactions to continuum

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(*Fermi-Golden Rule*)
- **Linear** Response theory

$$\sigma(\omega) \sim \sum_n |\langle n | \Theta | 0 \rangle|^2 \delta(\omega - E_n + E_0)$$

perturbative (electro-weak) **INCLUSIVE**

$$\gamma^{(*)} + b \longrightarrow b^* \quad (b^* = c+d+\dots \text{ or } e+f+\dots \text{ or } \dots)$$

Reactions to continuum

- **First order** perturbation theory
(*Fermi-Golden Rule*)
- **Linear** Response theory

$$\sigma(\omega) \sim \sum_n |\langle \mathbf{n} | \Theta | \mathbf{0} \rangle|^2 \delta(\omega - E_n + E_0)$$

$$\sum_n | \mathbf{n} \rangle \langle \mathbf{n} | = \mathbf{I}$$

$$\mathbf{H} | \mathbf{n} \rangle = E_n | \mathbf{n} \rangle$$

Reactions to continuum

PERTURBATIVE INCLUSIVE

$$S(\omega) = \sum_n |\langle \mathbf{n} | \Theta | \mathbf{0} \rangle|^2 \delta(\omega - E_n + E_0)$$

$S(\omega)$ represents the crucial quantity
Requires the solution of both
the **bound** and **continuum** A-body problem

We see next that in case of non perturbative reactions the crucial quantity for calculating the cross section has a very similar form

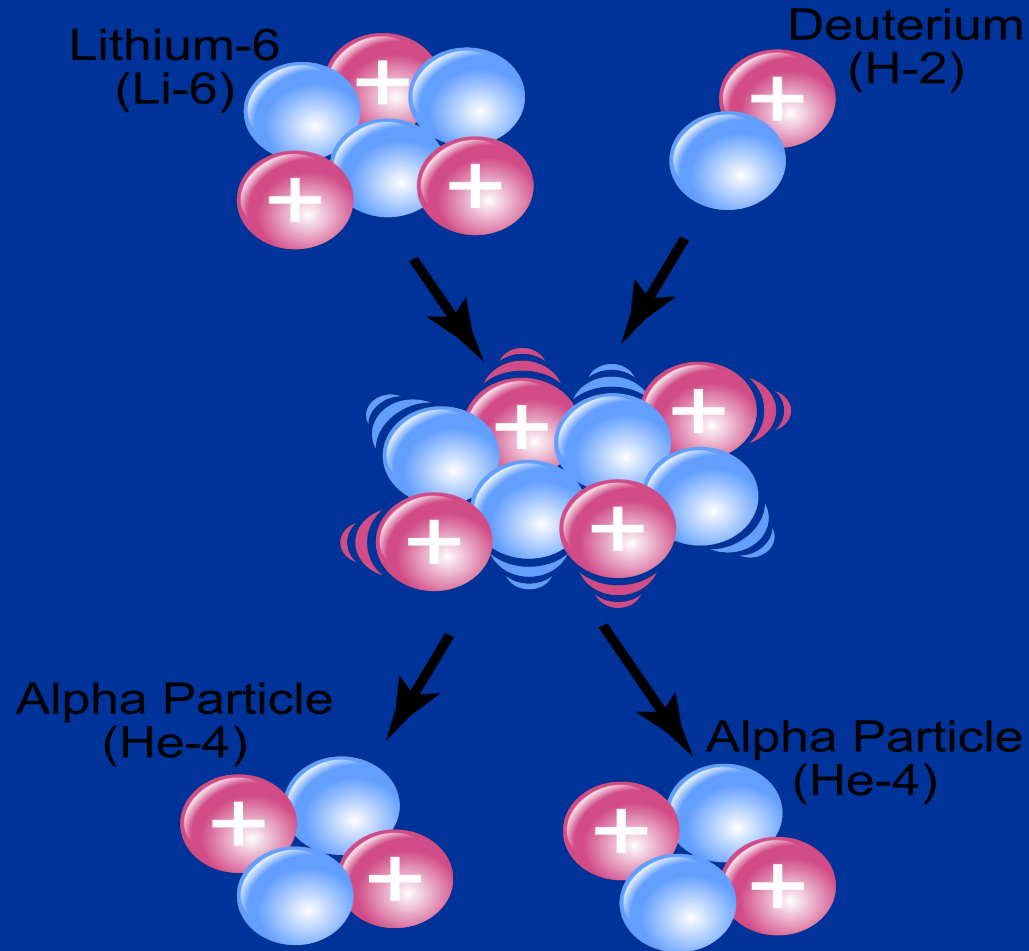
Reactions to continuum

non-perturbative (hadronic)

$a + b \rightarrow c + d + \dots$

$$\sigma(\omega) \sim |T_{\beta\alpha}(E)|^2$$

H is the Hamiltonian of the 8-body system



Lithium-6 – Deuterium Reaction

Reactions to continuum

non-perturbative (hadronic)

$a + b \rightarrow c + d + \dots$

$$\sigma(\omega) \sim |T_{\beta\alpha}(E)|^2$$

Reactions to continuum

non-perturbative (hadronic)



$$\sigma(\omega) \sim |T_{\beta\alpha}(E)|^2$$

General form of T-matrix

(notation as Goldberger-Watson Collision Theory)

$$T_{\beta\alpha}(E) = \langle \chi_{\beta} \mathcal{V}_{\alpha} \chi_{\alpha} \rangle + \langle \chi_{\beta} \mathcal{V}_{\beta} (E - H + i\eta)^{-1} \mathcal{V}_{\alpha} \chi_{\alpha} \rangle$$

Reactions to continuum

non-perturbative (hadronic)



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A-body continuum energy

Reactions to continuum

non-perturbative (hadronic)



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General form of T-matrix

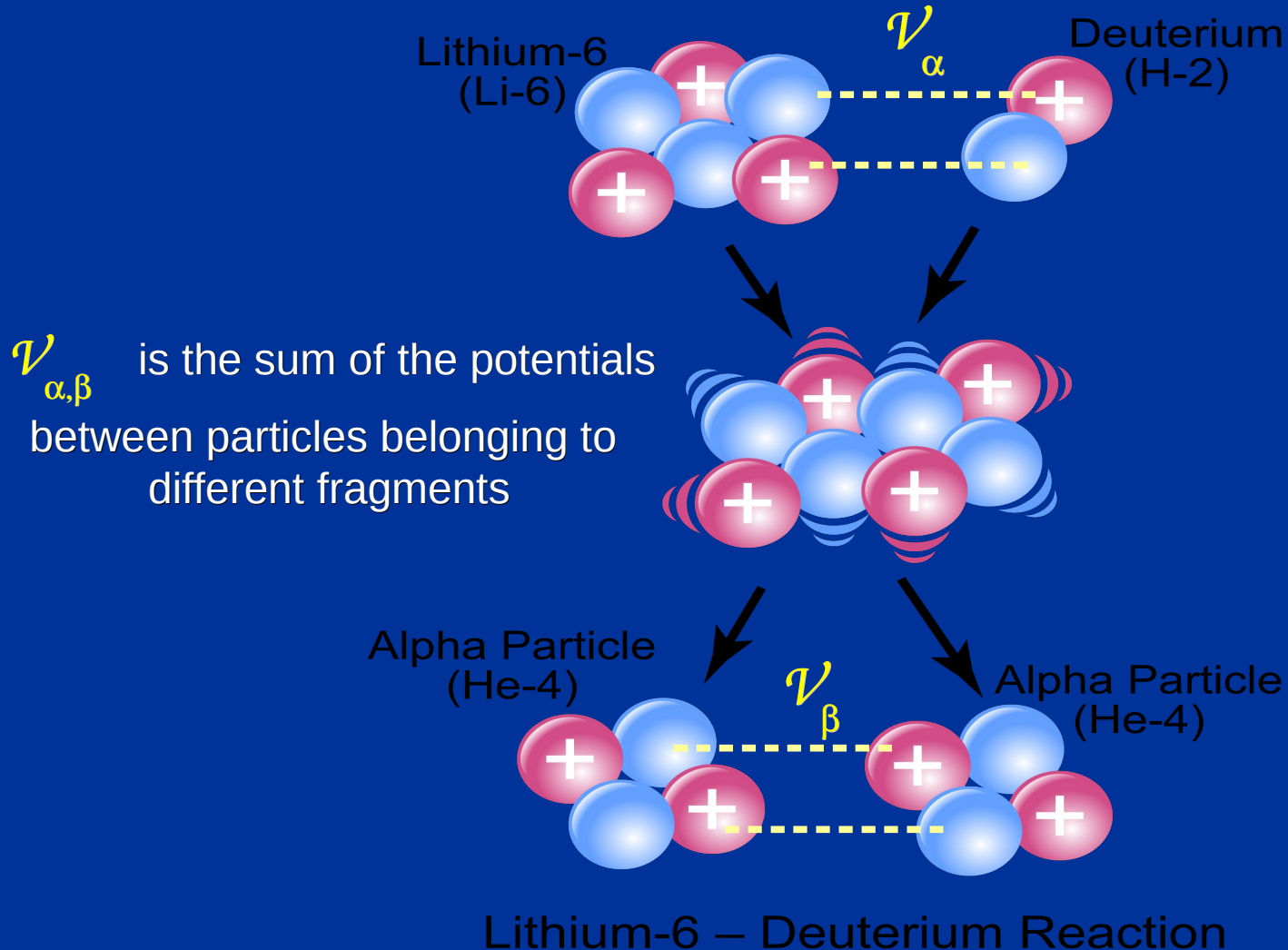
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χ_{β} and χ_{α} are the “channel functions” (with proper antisymmetrization), namely products of the **bound states** of **a** and **b**, times a relative **Plane Wave**

$$|\chi_{\alpha}\rangle = \mathcal{A} |a\rangle |b\rangle |PW\rangle$$

H is the Hamiltonian of the 8-body system



General form of T-matrix

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“easier” part



Very difficult part



General form of T-matrix

$$T_{\beta\alpha}(E) = \langle \chi_{\beta} \mathcal{V}_{\alpha} \chi_{\alpha} \rangle + \langle \chi_{\beta} \mathcal{V}_{\beta} (E - H + i\eta)^{-1} \mathcal{V}_{\alpha} \chi_{\alpha} \rangle$$

“easier” part

Very difficult part

If we denote $\mathcal{V}_{\alpha,\beta} \chi_{\alpha,\beta} = \phi_{\alpha,\beta}$

$\mathcal{V}_{\alpha,\beta}$ is the sum of the potentials between particles belonging to different fragments

General form of T-matrix

$$T_{\beta\alpha}(E) = \langle \chi_\beta | \mathcal{V}_\alpha | \chi_\alpha \rangle + \langle \chi_\beta | \mathcal{V}_\beta (E - H + i\eta)^{-1} \mathcal{V}_\alpha | \chi_\alpha \rangle$$

easier part

Very difficult part

$$\langle \phi_\beta | (E - H + i\eta)^{-1} | \phi_\alpha \rangle$$

One can manipulate the non trivial part:

$$\langle \phi_\beta | (\mathbf{E} - \mathbf{H} + i \eta)^{-1} | \phi_\alpha \rangle =$$

Step 1) Insert **completeness of eigenstates** $|n\rangle$ of H: $\sum_n |n\rangle\langle n| = I$

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$$= \int d\omega \sum_n \delta(\omega - \mathbf{E}_n) (\mathbf{E} - \omega + i\eta)^{-1} \langle \phi_\beta | n \rangle \langle n | \phi_\alpha \rangle =$$

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$$= \int d\omega (\mathbf{E} - \omega + i\eta)^{-1} \mathbf{F}_{\alpha\beta}(\omega) =$$

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the problem reduces to calculate the function $\mathbf{F}_{\alpha\beta}(\omega)$

$$\mathbf{F}_{\alpha\beta}(\omega) = \sum_n \delta(\omega - \mathbf{E}_n) \langle \phi_\beta | n \rangle \langle n | \phi_\alpha \rangle$$

Similar expressions!

Non-Pert.

$$F_{\alpha\beta}(\omega) = \sum_{\mathbf{n}} \langle \phi_{\beta} | \mathbf{n} \rangle \langle \mathbf{n} | \phi_{\alpha} \rangle \delta(\omega - E_{\mathbf{n}})$$

Pert.

$$S(\omega) = \sum_{\mathbf{n}} \langle 0 | \Theta^+ | \mathbf{n} \rangle \langle \mathbf{n} | \Theta | 0 \rangle \delta(\omega - E_{\mathbf{n}} + E_0)$$

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$|0\rangle, |\phi_\alpha\rangle, |\phi_\beta\rangle$ Needs only to be able to calculate bound states!

(Remember: $\phi_\alpha = \mathcal{V}_\alpha \chi_\alpha = \mathcal{V}_\alpha \mathcal{A} |a\rangle |b\rangle |PW\rangle$)

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$$F_{\alpha\beta}(\omega) = \sum_n \langle \phi_\beta | \mathbf{n} \rangle \langle \mathbf{n} | \phi_\alpha \rangle \delta(\omega - E_n)$$

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$|\mathbf{n}\rangle$

*All eigenstates of the Hamiltonian,
Bound and continuum*

**At present there are
only few methods for
reactions and they are
limited to $A=3,4$.**

Why?

scattering many-body problem!

**In configuration space
(Schroedinger)**

**Difficulties in dealing with
asymptotic conditions in the
solution of the many coupled
differential equations**

scattering many-body problem

**In momentum space
(Faddeev-Yakubowski)**

**Difficulties in coping with
complicated poles in the many
coupled integral equations**

The integral transform approach may help

Integral transforms:

$$\Phi(\sigma) = \int d\omega K(\omega, \sigma) F(\omega)$$

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Integral transforms:



$$\Phi(\sigma) = \int d\omega K(\omega, \sigma) F(\omega)$$



Sometimes problematic!

In fact:

$$\Phi(\sigma) = \int d\omega \, K(\omega, \sigma) F(\omega)$$

Suppose to have a different $F'(\omega)$



$$[F(\omega) + A \sin(\nu\omega)]$$

$$\Phi(\sigma) = \int d\omega K(\omega, \sigma) F(\omega)$$

Suppose to have a different $F'(\omega)$

$$\Phi(\sigma) + A \Delta \Phi(\nu) = \int d\omega K(\omega, \sigma) [F(\omega) + A \sin(\nu\omega)]$$

$$\Phi(\sigma) = \int d\omega K(\omega, \sigma) F(\omega)$$

Error in the transform

Suppose an error

$$\Phi(\sigma) + A \Delta \Phi(\nu) = \int d\omega K(\omega, \sigma) [F(\omega) + A \sin(\nu\omega)]$$

for very large ν

0

independently on the
amplitude A of the error!

More remarks later on this aspect

Remember the similar expressions of the crucial quantities for perturbative and Non-perturbative reactions:

Non-Pert.

$$F_{\alpha\beta}(\omega) = \sum_n \langle \phi_\beta | \mathbf{n} \rangle \langle \mathbf{n} | \phi_\alpha \rangle \delta(E_n - \omega)$$

Pert.

$$S(\omega) = \sum_n \langle 0 | \Theta^+ | \mathbf{n} \rangle \langle \mathbf{n} | \Theta | 0 \rangle \delta(\omega - E_n + E_0)$$

$|\mathbf{n}\rangle$

*All eigenstates of the Hamiltonian,
Bound and continuum*

Let us consider e.g. $S(\omega)$

(similar for $F_{\alpha\beta}(\omega)$ with $\Theta | 0 \rangle$ replaced by $|\phi_\alpha\rangle$

and $\langle 0 | \Theta^\dagger$ replaced by $\langle \phi_\beta |$)

$$S(\omega) = \sum_n |\langle n | \Theta | 0 \rangle|^2 \delta(\omega - E_n + E_0)$$

Scattering states

Energies in the continuum

$$S(\omega) = \sum_n |\langle n | \Theta | 0 \rangle|^2 \delta(\omega - E_n + E_0)$$

$$\Phi(\sigma) = \int S(\omega) K(\omega, \sigma) d\omega =$$

1) exchange \int and \sum

$$\Phi(\sigma) = \sum_n \int d\omega \delta(\omega - E_n + E_0) K(\omega, \sigma) |\langle 0 | \Theta | n \rangle|^2$$

2) integrate the δ -function

$$\begin{aligned} \Phi(\sigma) &= \sum_n K(E_n - E_0, \sigma) \langle 0 | \Theta^\dagger | n \rangle \langle n | \Theta | 0 \rangle = \\ &= \sum_n \langle 0 | \Theta^\dagger K(H - E_0, \sigma) | n \rangle \langle n | \Theta | 0 \rangle \end{aligned}$$

3) Use completeness of H eigenstates

$$\sum_n |n\rangle \langle n| = I$$

$$\begin{aligned}\Phi(\sigma) &= \sum_n \langle 0 | \Theta^\dagger K(H-E_0, \sigma) | n \rangle \langle n | \Theta | 0 \rangle \\ &= \langle 0 | \Theta^\dagger K(H-E_0, \sigma) \Theta | 0 \rangle\end{aligned}$$

$$\Phi(\sigma) = \int S(\omega) K(\omega, \sigma) d\omega =$$



$$\langle 0 | \Theta^\dagger K(H - E_0, \sigma) \Theta | 0 \rangle$$

The calculation of **ANY** transform seems to require, **in principle**, only the knowledge of the ground state!

However,

$K(H-E_0, \sigma)$ can be quite a complicate operator.

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The calculation of **ANY** transform seems to require, **in principle**, only the knowledge of the ground state!

However,

$K(H-E_0, \sigma)$ can be quite a complicated operator.

So, which kernel is suitable for calculation of this?


$$\Phi(\sigma) = \langle 0 | \Theta^\dagger K(H-E_0, \sigma) \Theta | 0 \rangle$$

One familiar example: sum rules!

Sum rules are transforms in terms of “*Moments*”

$$K(\omega, \sigma) = \omega^\sigma \text{ with } \sigma \text{ integer}$$

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One familiar example: sum rules!

Sum rules are transforms in terms of “*Moments*”

$$K(\omega, \sigma) = \omega^\sigma \text{ with } \sigma \text{ integer}$$

*To obtain $S(\omega)$ the inversion of the transform
is equivalent to the reconstruction of $S(\omega)$
in terms of its moments (theory of moments)*

One familiar example: sum rules!

Sum rules are transforms in terms of “*Moments*”

$$K(\omega, \sigma) = \omega^\sigma \text{ with } \sigma \text{ integer}$$

however,

$$\Phi(\sigma) = \int d\omega \omega^\sigma S(\omega)$$

may be *infinite* for some $\sigma > \bar{\sigma}$!!!

Another common example:

The Laplace Kernel:

$$\Phi(\sigma) = \int e^{-\omega\sigma} S(\omega) d\omega = \langle 0 | \Theta^+ e^{i(H-E_0)\sigma} \Theta | 0 \rangle$$

The Laplace Kernel:

$$\Phi(\sigma) = \int e^{-\omega\sigma} S(\omega) d\omega = \langle 0 | \Theta^+ e^{i(H-E_0)(i\sigma)} \Theta | 0 \rangle$$

In Condensed Matter Physics:

In QCD

In Nuclear Physics:

The Laplace Kernel:

$$\Phi(\sigma) = \int e^{-\omega\sigma} S(\omega) d\omega = \langle 0 | \Theta^+ e^{-(H-E_0)\tau} \Theta | 0 \rangle$$

In Condensed Matter Physics:

In Nuclear Physics:

In QCD

$\sigma = \tau = it$ imaginary time!

$\Phi(\tau)$ is calculated with Monte Carlo Methods

The Laplace Kernel:

$$\Phi(\sigma) = \int e^{-\omega\sigma} S(\omega) d\omega = \langle 0 | \Theta^+ e^{-(H-E_0)\tau} \Theta | 0 \rangle$$

In Condensed Matter Physics:

In Nuclear Physics:

In QCD

$\sigma = \tau = it$ imaginary time!

$\Phi(\tau)$ is calculated with **Monte Carlo Methods**
and then inverted with **methods**
based on **Bayesian theorem (MEM)**

**Are there other kernels
suitable for diagonalization
methods on
*finite norm basis functions***

???

$$\Phi(\sigma) = \int S(\omega) K(\omega, \sigma) d\omega =$$



$$\langle 0 | \Theta^\dagger K(H - E_0, \sigma) \Theta | 0 \rangle$$

Notice !!!

$$S(\omega) = \sum_n |\langle n | \Theta | 0 \rangle|^2 \delta(\omega - E_n + E_0)$$

$$S(\omega) =$$

$$= -1/\pi \operatorname{Im} \left[\sum_n \langle 0 | \Theta^\dagger | n \rangle \langle n | \Theta | 0 \rangle \right] (\omega - E_n + E_0 + i\varepsilon)^{-1} \Big]$$

$$= 1/\pi \operatorname{Im} \left[\sum_n \langle 0 | \Theta^\dagger (\mathbf{H} - \omega - E_0 + i\varepsilon)^{-1} | n \rangle \langle n | \Theta | 0 \rangle \right]$$

$$= -1/\pi \operatorname{Im} \left[\langle 0 | \Theta^\dagger (\mathbf{H} - \omega - E_0 + i\varepsilon)^{-1} \Theta | 0 \rangle \right]$$

If we had to deal with a “**confined**” system one could represent H on **bound states eigenfunctions** $|v\rangle$

$$\langle 0 | \Theta^+ K(H - E_0, \sigma) \Theta | 0 \rangle = \Phi(\sigma) =$$

$$\sum_{\mu\nu} \langle 0 | \Theta^+ |\mu\rangle \langle \mu | K(H_{\mu\nu} - E_0, \sigma) |\nu\rangle \langle \nu | \Theta | 0 \rangle$$

If we had to deal with a “**confined**” system one could represent H on **bound states eigenfunctions** $|v\rangle$

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$$\sum_{\mu\nu} \langle 0 | \Theta^+ |\mu\rangle \langle \mu | K(H_{\mu\nu} - E_0, \sigma) | \nu \rangle \langle \nu | \Theta | 0 \rangle$$

After diagonalizing $H_{\mu\nu}$ the transform would be simply

$$\sum_{\lambda} K(\epsilon_{\lambda} - E_0, \sigma) |\langle \lambda | \Theta | 0 \rangle|^2 = \Phi(\sigma)$$

However, a nucleus is NOT “**confined**”!
The nuclear **H** has positive energy eigenstates
and therefore, in general, CANNOT be represented
on **b.s. eigenfunctions** $|\nu\rangle$
(Continuum discretization approximation)

THE GOOD NEWS:

The representation of H on **b.s. eigenfunctions** $|v\rangle$ and therefore the calculation of the transform via

$$\Phi(\sigma) = \sum_{\lambda} K(\varepsilon_{\lambda} - E_0, \sigma) |\langle \lambda | \Theta | 0 \rangle|^2$$

is **allowed** for **specific kernels** $K(\omega, \sigma)$!

**No approximation!**

Conditions required:

$$1) \quad \Phi(\sigma) = \int S(\omega) K(\omega, \sigma) d\omega < \infty$$

$$2) \quad \int S(\omega) d\omega < \infty$$

$$3) \quad K(\omega, \sigma) \neq \delta(\omega - \sigma)$$

Conditions required N.1:

$$1) \quad \Phi(\sigma) = \int S(\omega) K(\omega, \sigma) d\omega < \infty$$

This is required in order to change \int

remember ...

$$S(\omega) = \sum_n |\langle n | \Theta | 0 \rangle|^2 \delta(\omega - E_n + E_0)$$

$$\Phi(\sigma) = \int S(\omega) K(\omega, \sigma) d\omega =$$

1) exchange \int and \sum

$$\Phi(\sigma) = \sum_n \int d\omega \delta(\omega - E_n + E_0) K(\omega, \sigma) |\langle 0 | \Theta | n \rangle|^2$$

2) integrate the δ -function

$$\begin{aligned} \Phi(\sigma) &= \sum_n K(E_n - E_0, \sigma) \langle 0 | \Theta^\dagger | n \rangle \langle n | \Theta | 0 \rangle = \\ &= \sum_n \langle 0 | \Theta^\dagger K(H - E_0, \sigma) | n \rangle \langle n | \Theta | 0 \rangle \end{aligned}$$

Condition required N.2:

$$2) \quad \int S(\omega) d\omega < \infty$$

$$\text{since} \quad \int S(\omega) d\omega = \langle 0 | \Theta^+ \Theta | 0 \rangle$$

It means that $\Theta | 0 \rangle$ has finite norm

and therefore has bound state asymptotic behaviour

Therefore it is allowed to insert
a complete **bound basis**

$$\langle 0 | \Theta^+ \mathbb{K}(H - E_0, \sigma) \Theta | 0 \rangle = \Phi(\sigma) =$$

$$\sum_{\mu\nu} \langle 0 | \Theta^+ |\mu\rangle \langle \mu | \mathbb{K}(H_{\mu\nu} - E_0, \sigma) |\nu\rangle \langle \nu | \Theta | 0 \rangle$$

After diagonalizing $H_{\mu\nu}$ the transform would be simply

$$\sum_{\lambda} \mathbb{K}(\varepsilon_{\lambda} - E_0, \sigma) |\langle \lambda | \Theta | 0 \rangle|^2 = \Phi(\sigma)$$

Condition required N.3:

3) $K(\omega, \sigma)$ not $\delta(\omega - \sigma)$!!!

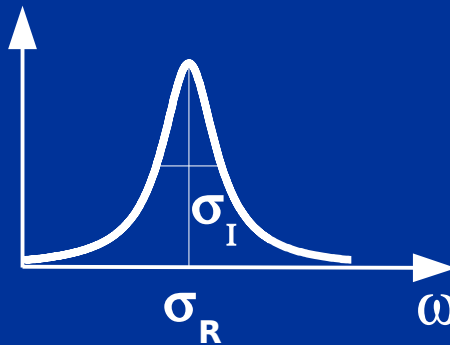
$$\Phi(\sigma) = \int S(\omega) \delta(\omega - \sigma) d\omega = S(\sigma)$$

$$\langle 0 | \Theta^\dagger \delta(H - E_0 - \sigma) \Theta | 0 \rangle =$$

$$-1/\pi \operatorname{Im} \left[\langle 0 | \Theta^\dagger (H - \sigma - E_0 + i\varepsilon)^{-1} \Theta | 0 \rangle \right]$$

*In this case the insertion of a complete bound basis would correspond to **Continuum discretization approximation** !!!*

The LIT Kernel satisfies the condition N.1



$$\Phi(\sigma_R, \sigma_I) = \sigma_I / \pi \int [(\omega - \sigma_R)^2 + \sigma_I^2]^{-1} S(\omega) d\omega < \infty$$

The LIT Kernel satisfies the conditions....

Moreover it is not a δ -function, but
one of **its representations**.

Therefore makes the inversion easier!!

ε infinitesimal!

$$S(\omega) = -1/\pi \operatorname{Im} \left[\langle 0 | \Theta^+ (H - E_0 - \omega + i\varepsilon)^{-1} \Theta | 0 \rangle \right]$$

$$\Phi(\sigma) = -\Gamma/\pi \operatorname{Im} \left[\langle 0 | \Theta^+ (H - E_0 - \sigma_R + i\sigma_I)^{-1} \Theta | 0 \rangle \right]$$

σ_I finite!

Illustration of the problem:

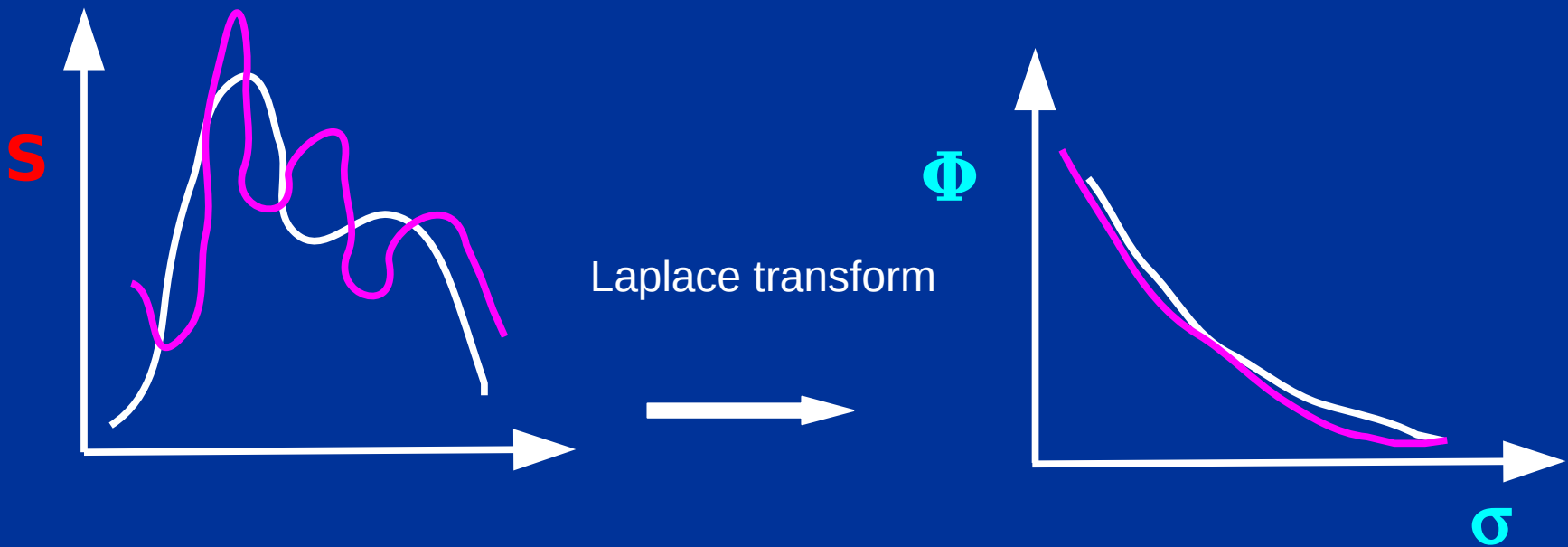


Illustration of the problem:

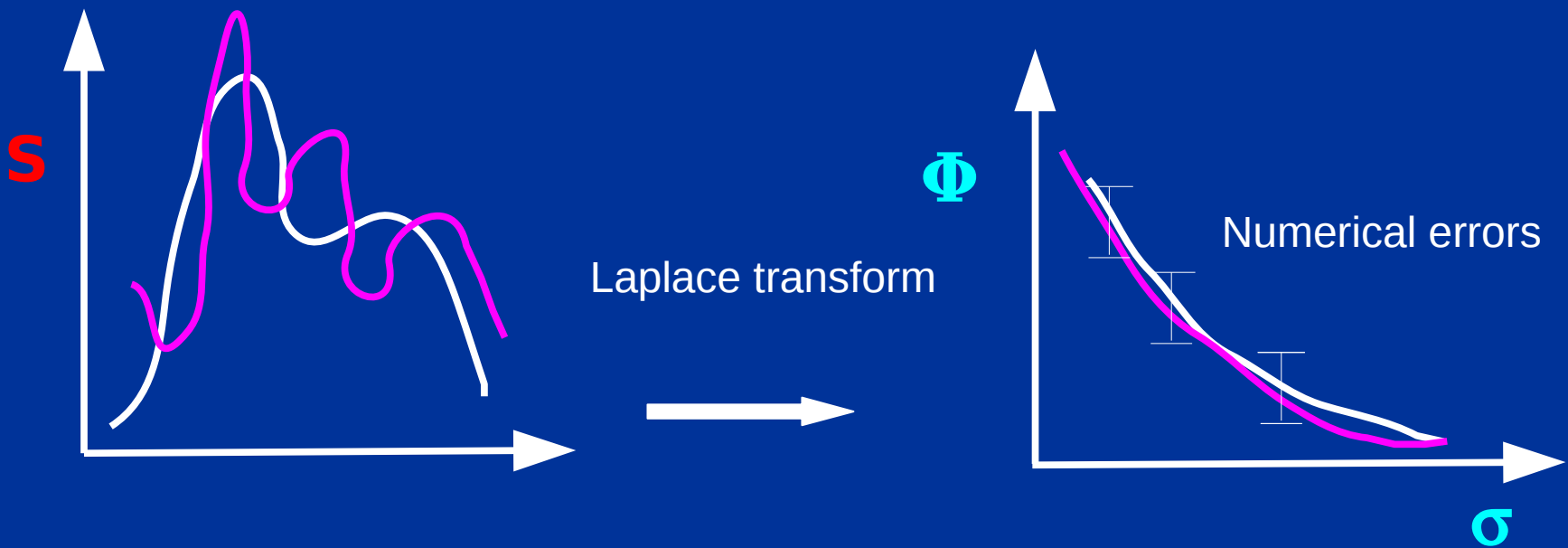
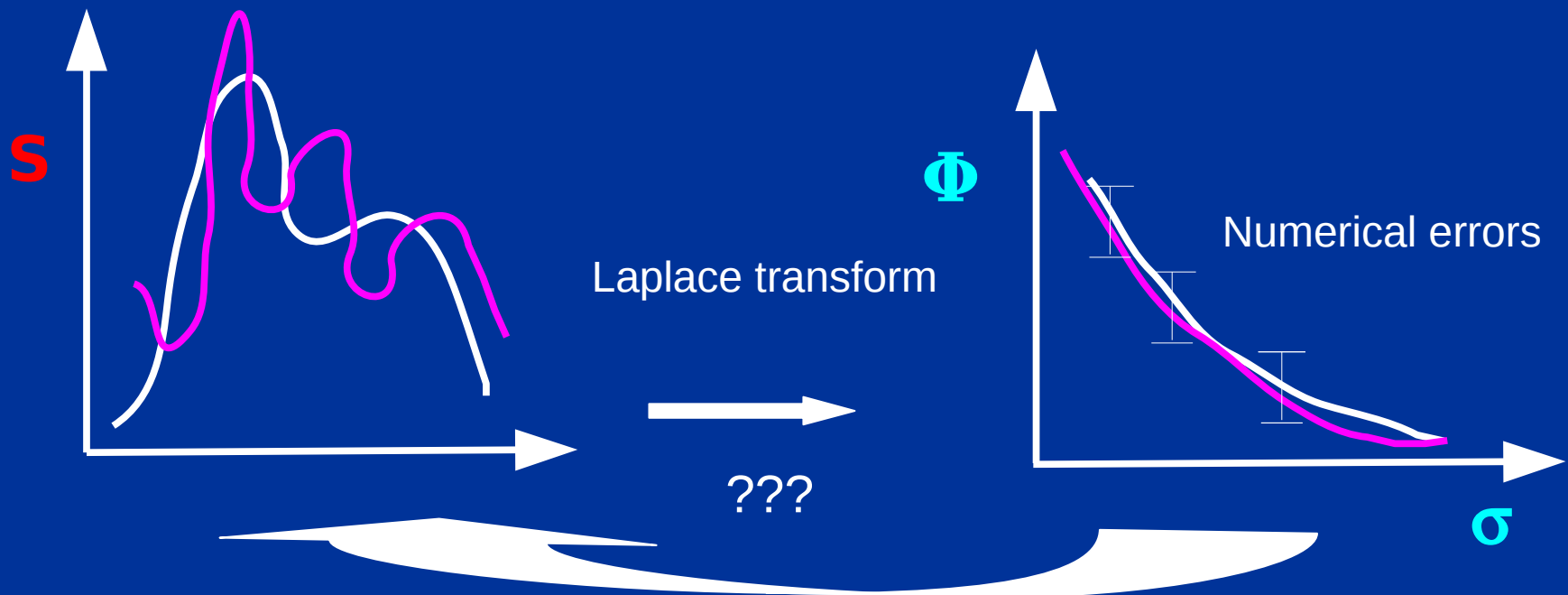
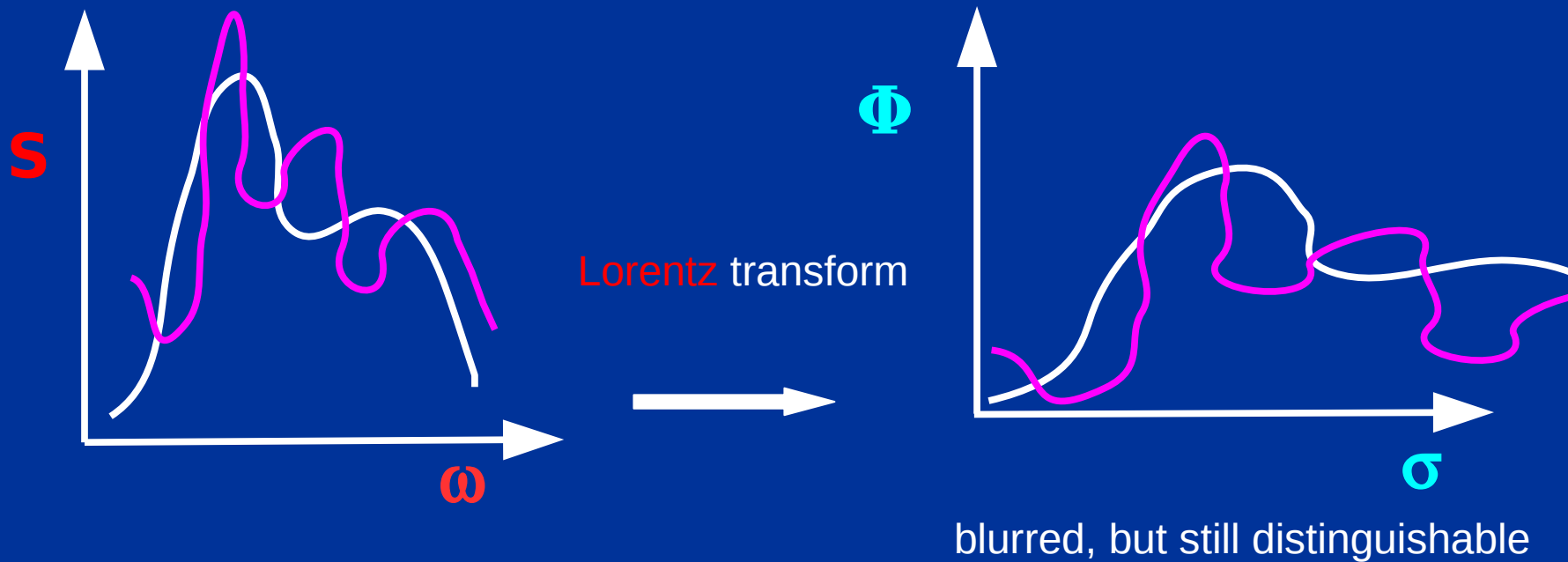


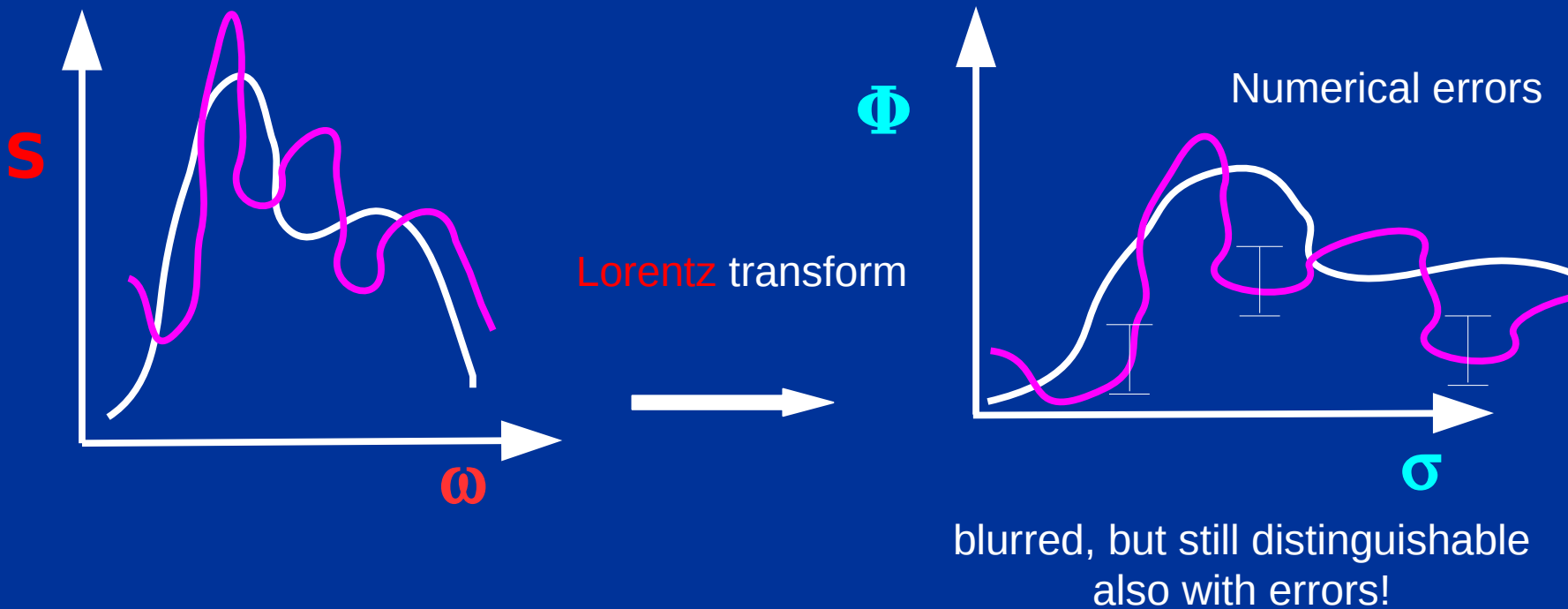
Illustration of the problem:



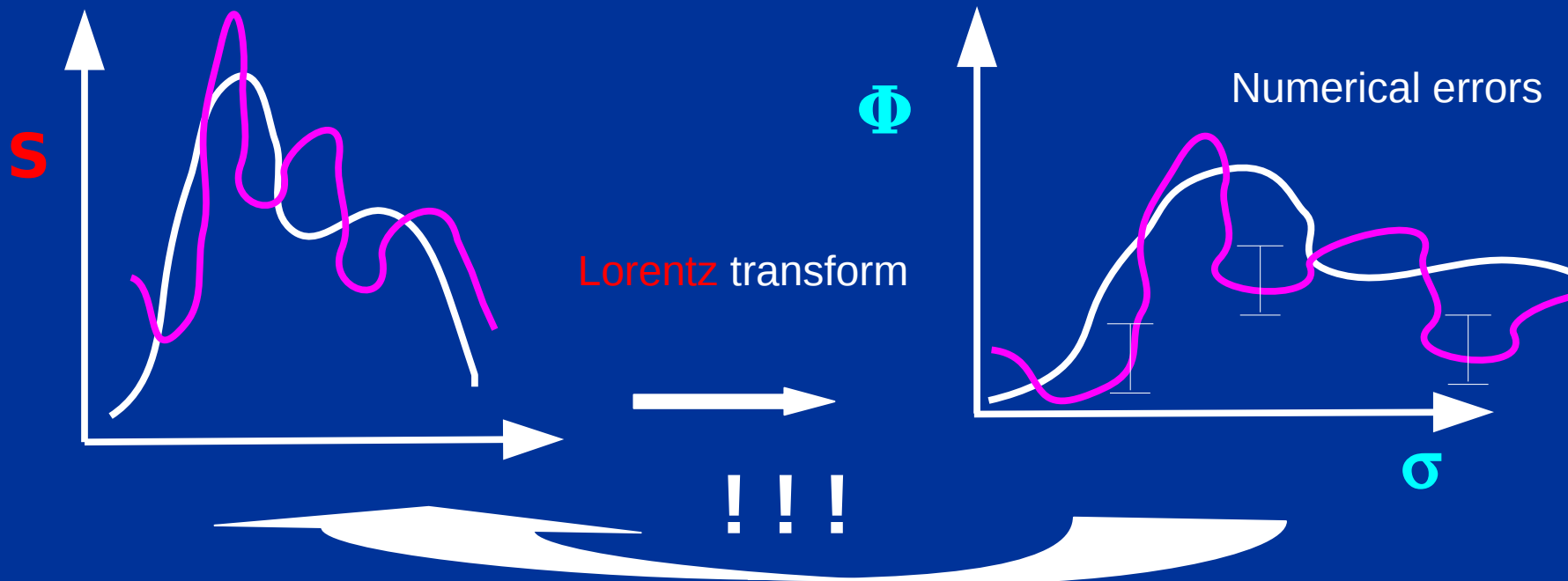
Advantages of a delta function representation:



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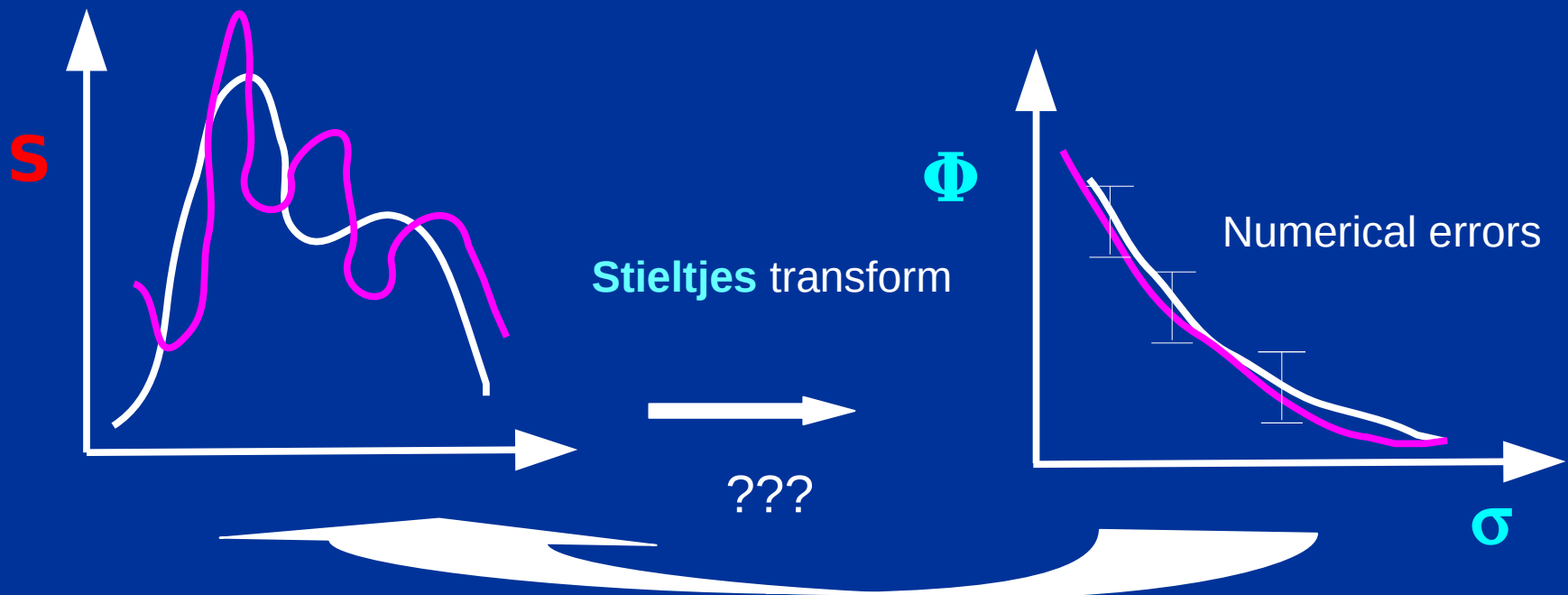
Moreover,
if the width of the Lorentzian were
the same as the experimental resolution
one would not need to invert the transform!

Other kernels?

The Stieltjes Kernel:

$$K(\omega, \sigma) = (\omega + \sigma)^{-1}$$

Illustration of the problem:
Same as Laplace!



**However, it may be useful
for another purpose:**

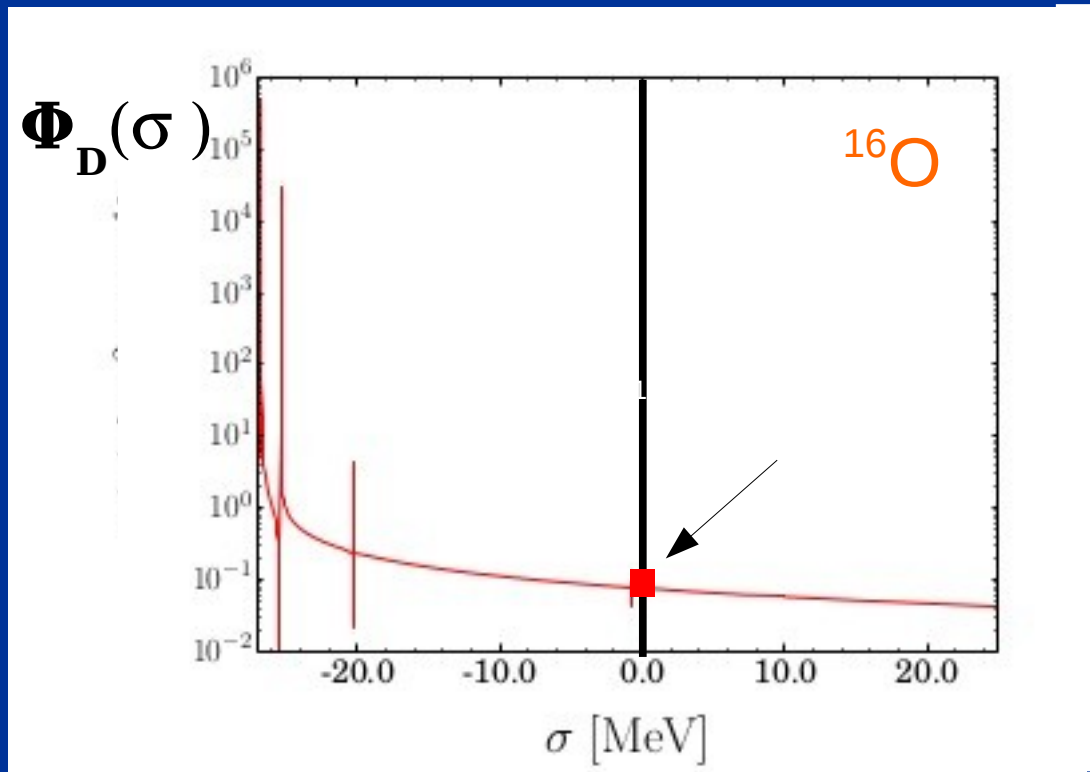
In fact:

$$\lim_{\sigma \rightarrow 0} \Phi(\sigma) = \int S(\omega) \omega^{-1} d\omega = \alpha_{\Theta}$$

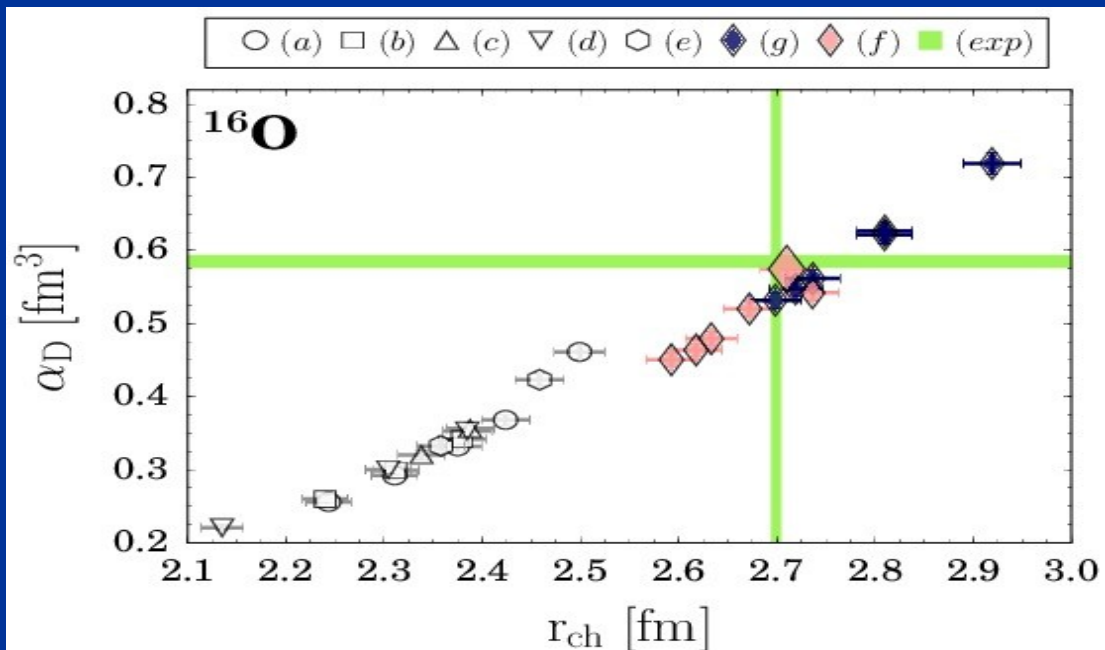
“generalized polarizability”
e.g. electric polarizability, magnetic susceptibility,
compressibility etc... depending on Θ

Recent results
on α_{Θ} with $\Theta = D$
(El. Dipole Polarizability)

Electric Dipole Polarizability as limit of the Stieltjes transform for $\sigma \rightarrow 0$

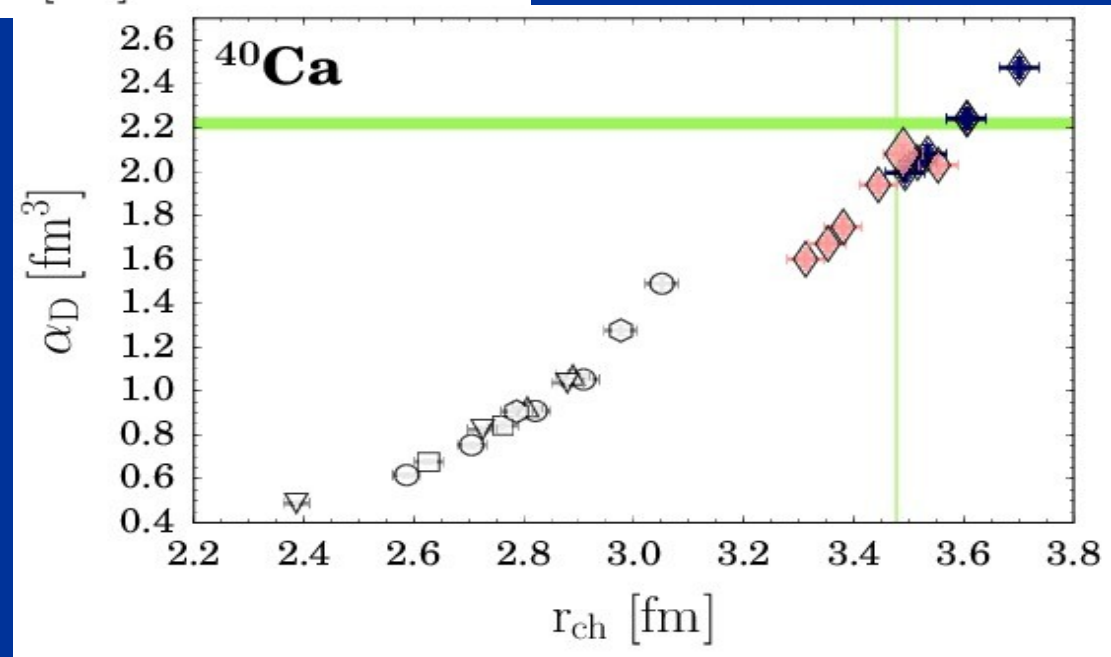


M.Miorelli et al. Phys. Rev. C 94, 034317 (2016)
b.s. expansion: Coupled Cluster



Interesting correlation
with the proton charge radius

Role of 3b-force



G. Hagen et al.
Nature Phys. 2016

A Transform with a kernel suitable for Monte Carlo methods:

[A.Roggero, F. Pederiva, G.O. Phys. Rev. B 88, 115138 (2013)]

combination of Sumudu kernels:

$$K(\omega, \sigma, \mathbf{P}) = N \sigma \left(\frac{e^{-\mu \omega/\sigma}}{\sigma} - \frac{e^{-\nu \omega/\sigma}}{\sigma} \right)^{\mathbf{P}}$$

$$\nu/\mu = b/a \quad \nu - \mu = \frac{\ln [b] - \ln [a]}{b - a} \quad b > a > 0 \text{ integer}$$

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$$K(\omega, \sigma, \mathbf{P}) \xrightarrow{\quad} \delta(\omega - \sigma)$$

$\mathbf{P} \xrightarrow{\quad} \infty$

A Transform with a kernel suitable for Monte Carlo methods:

[A.Roggero, F. Pederiva, G.O. Phys. Rev. B 88, 115138 (2013)]

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$$K(\omega, \sigma, \mathbf{P}) = N \sigma \left(\frac{e^{-\mu \omega/\sigma}}{\sigma} - \frac{e^{-\nu \omega/\sigma}}{\sigma} \right)^{\mathbf{P}}$$
$$= N \sum_k^{\mathbf{P}} (-1)^k \binom{k}{\mathbf{P}} e^{-\tau(\mathbf{P}, k, \sigma) \omega}$$

Finite sum of Laplace Kernels!

A Transform with a kernel suitable for Monte Carlo methods:

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$$= N \sum_k^{\mathbf{P}} (-1)^k \binom{k}{\mathbf{P}} e^{-\tau(\mathbf{P}, k, \sigma) \omega}$$

$$\tau(\mathbf{P}, k, \sigma) = \log(b/a) [\mathbf{P} a/(b - a) + k] / \sigma$$

Small width ---> large \mathbf{P} ---> **large** imaginary time

Applied only to bosonic system: Liquid Helium

The transform is calculated with
AFDMC and then inverted with MEM

See A.Roggero, F. Pederiva, G.O. Phys. Rev. B 88, 115138 (2013)]

Let's remember:

$$\Phi(\sigma) = \int d\omega K(\omega, \sigma) S(\omega)$$


In order to obtain $S(\omega)$ one needs to invert the transform

Problem:

Sometimes the “inversion” of $\Phi(\sigma)$ may be problematic

New Kernels?

What about “wavelets”?

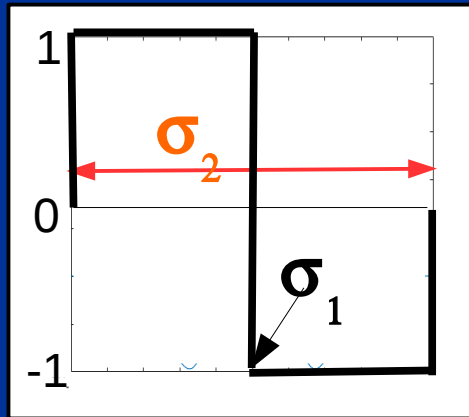
A wavelet Kernel is an oscillating function but with a "window".

It has 2 parameters:

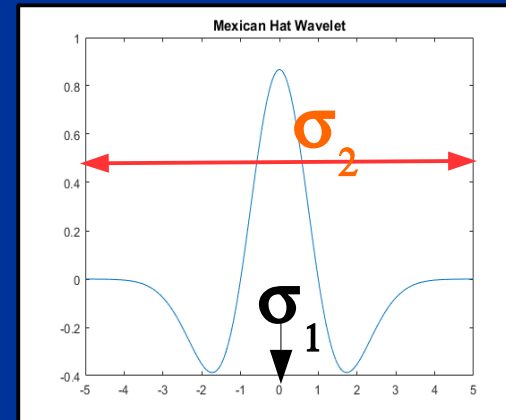
σ_2 drives the frequency of the oscillation

σ_1 drives the position of the window over the ω range

discrete



continuous



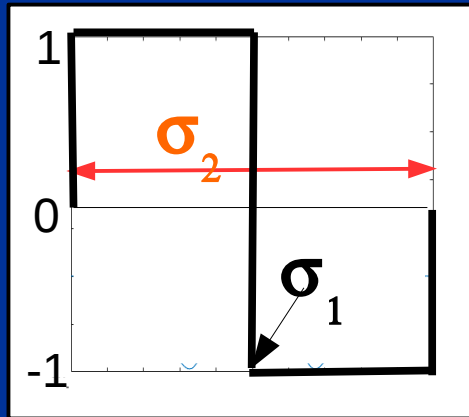
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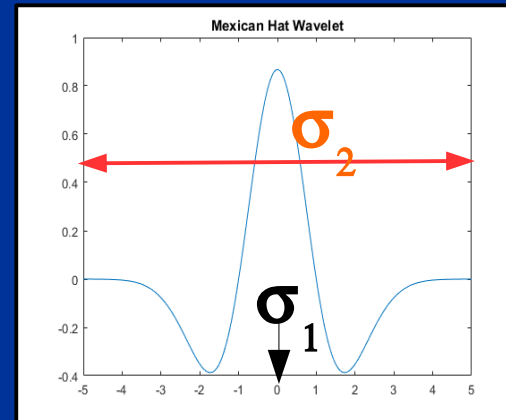
σ_2 drives the frequency of the oscillation

σ_1 drives the position of the window over the ω range

discrete



continuous



*They combine the power of the **Fourier Kernel** (in detecting frequencies of oscillations) and the **Lorentz Kernel** (in picking the information around specific ω ranges)*

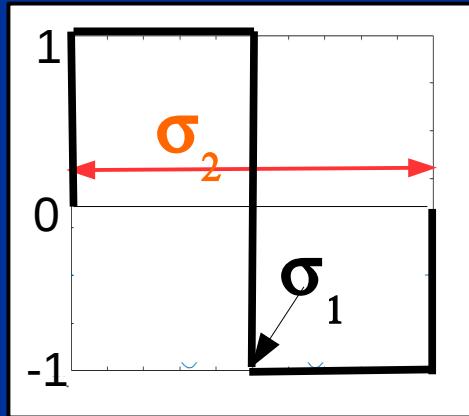
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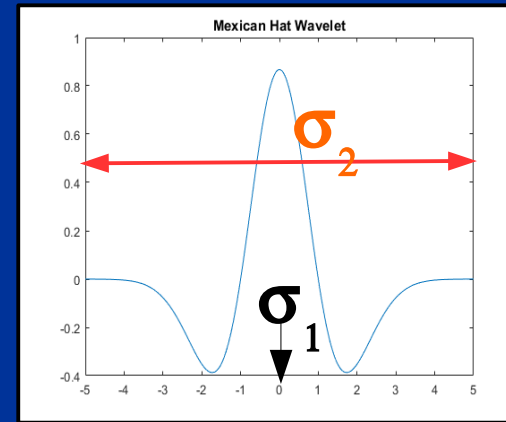
σ_1 drives the frequency of the oscillation

σ_2 drives the position of the window over the ω range

discrete



continuous



Since wavelets are *orthonormal* functions in principle
their inversion is straightforward !

[linear combination of $\Phi(\sigma_1, \sigma_2)$]

Integral transform

$$\Phi(\sigma) = \int d\omega K(\omega, \sigma) S(\omega)$$

If $K(\omega, \sigma) \equiv \mathbf{K}_\sigma(\omega)$ are the elements of an orthogonal basis

$\Phi(\sigma) = \Phi_\sigma$ are the coefficients of the expansion of $S(\omega)$ on that basis

then
$$S(\omega) = \sum_\sigma \Phi_\sigma \mathbf{K}_\sigma(\omega)$$

Integral transform

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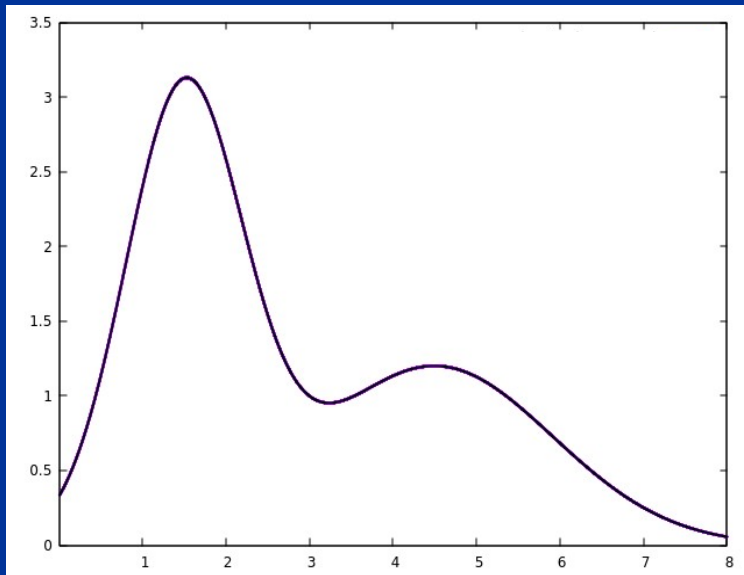
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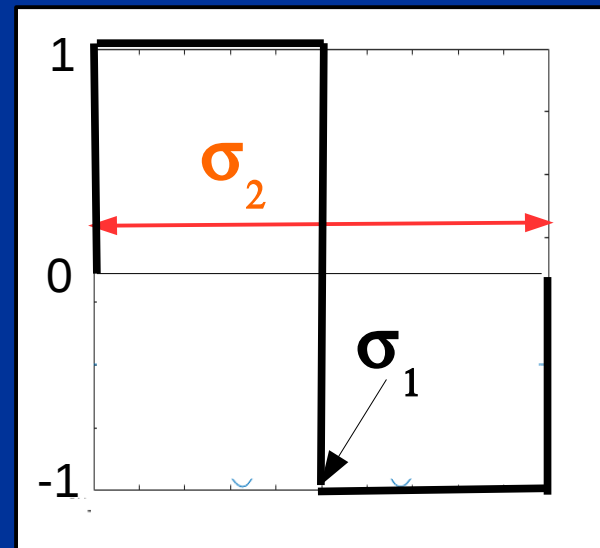
Namely the inversion is straightforward!

A model study (discrete wavelets)

Our model $S(\omega)$



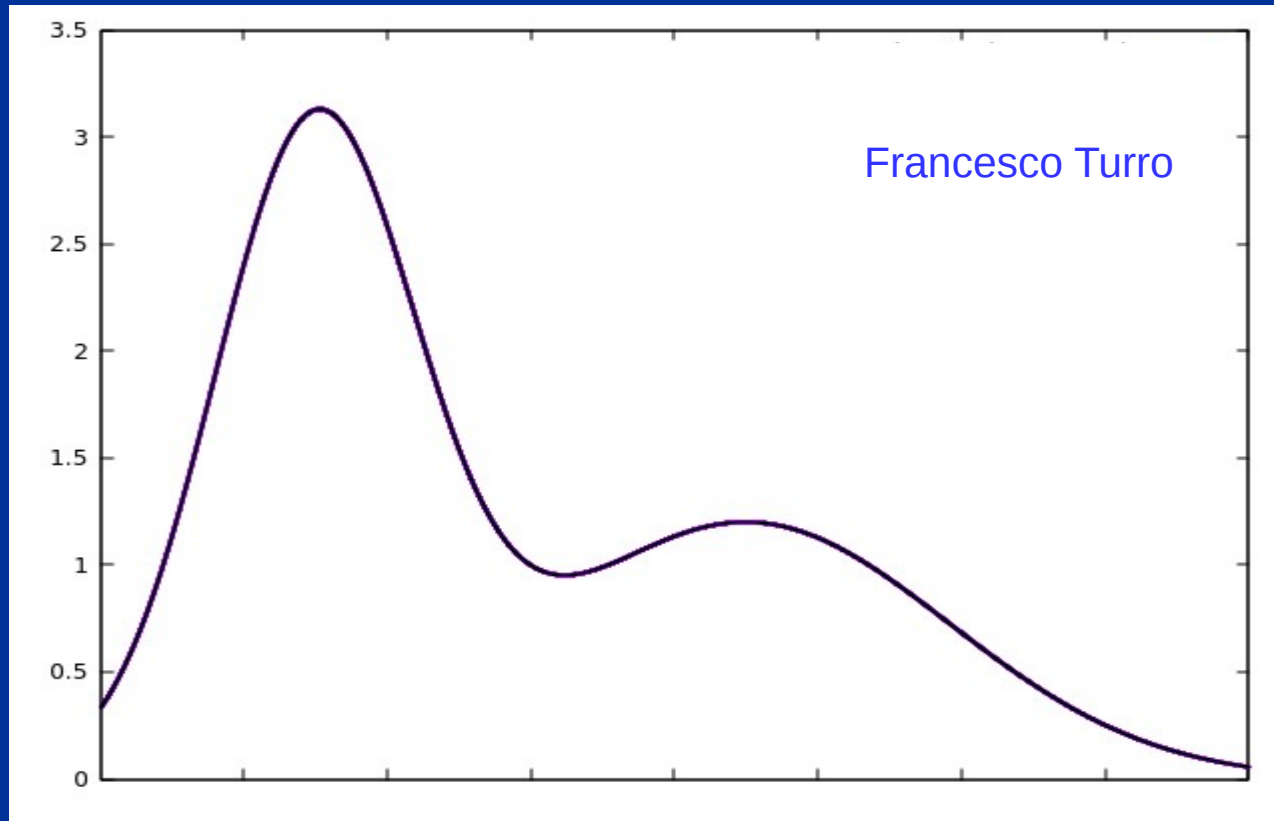
A wavelet kernel
(Haar wavelets)



$$K(\omega, \sigma_1, \sigma_2)$$

Model $S(\omega)$ and reconstructed from wavelet transform:

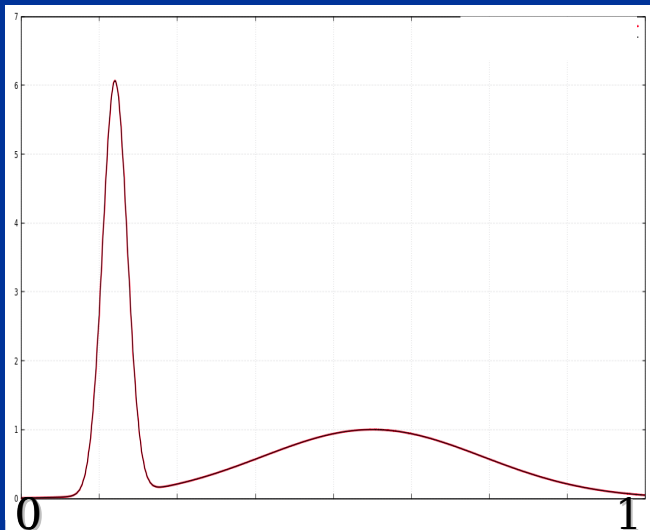
identical!



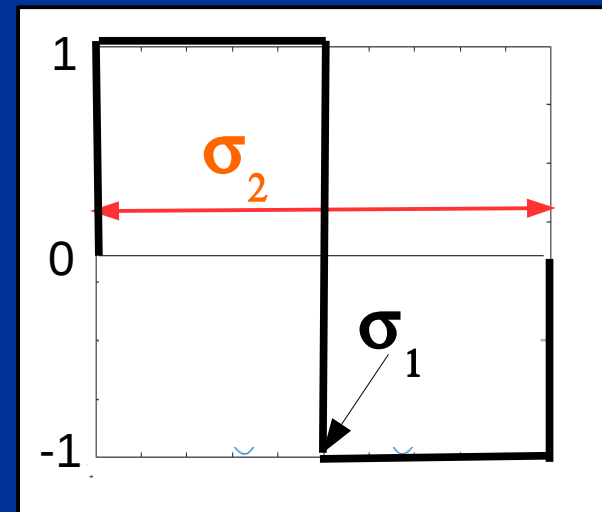
Another model study

(narrow resonance, discrete wavelets)

Our model $S(\omega)$



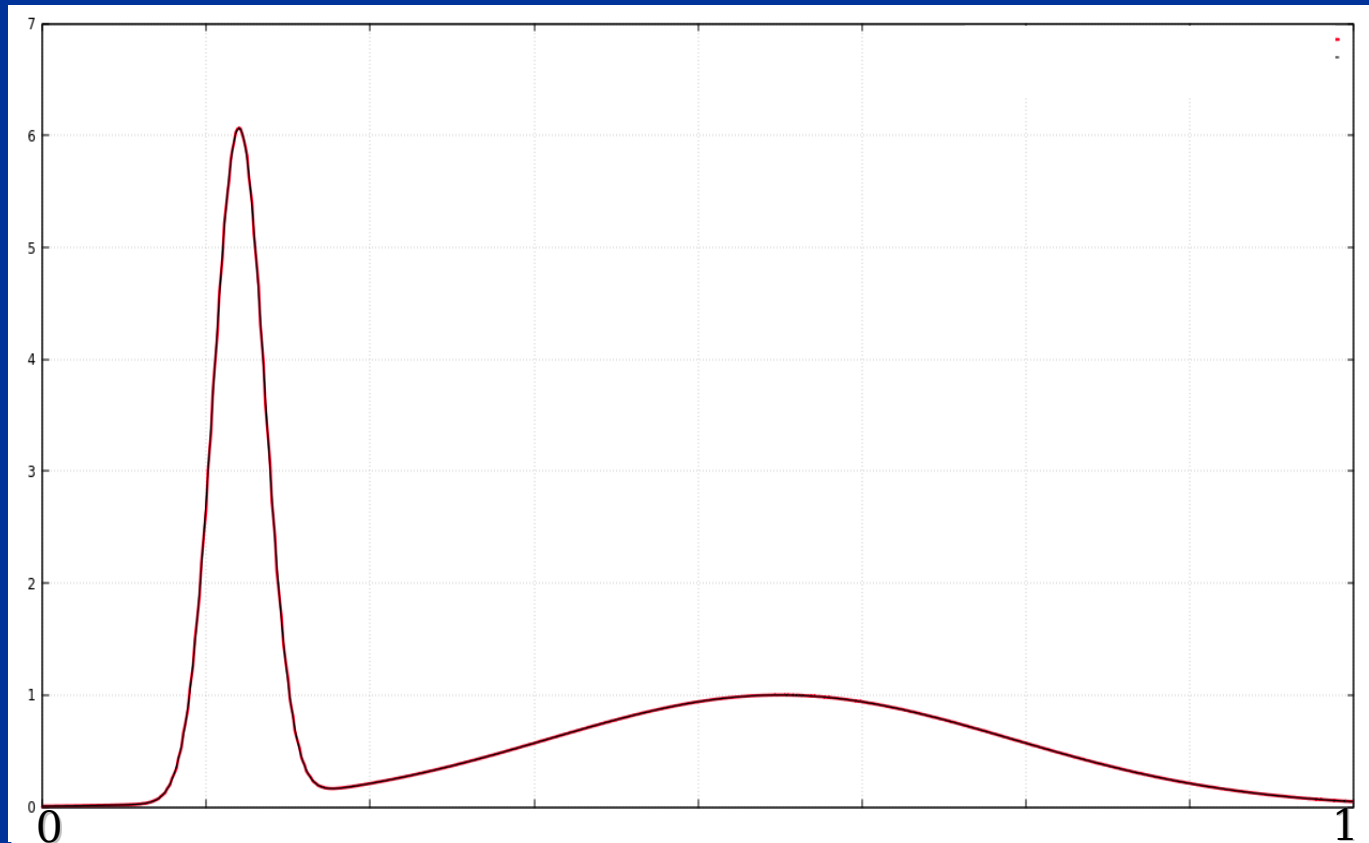
A wavelet kernel



$$K(\omega, \sigma_1, \sigma_2)$$

Model $S(\omega)$ and reconstructed from wavelet transform:

again identical!

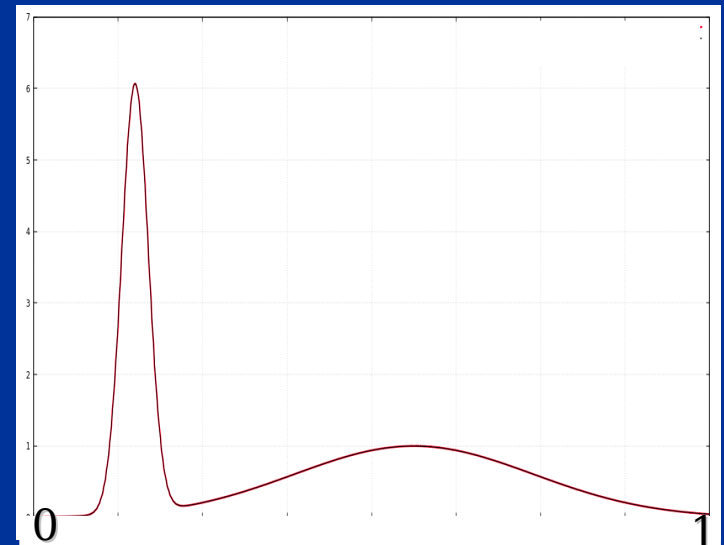
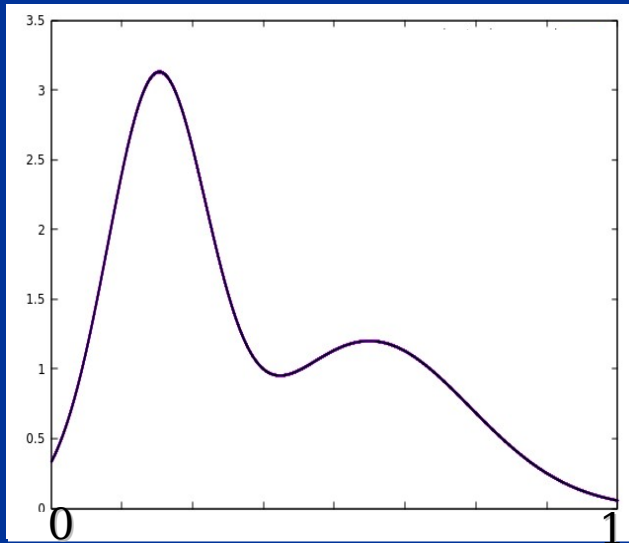


Which information has been used to reconstruct $S(\omega)$???

Which information has been used to reconstruct $S(\omega)$???

values of $K(\omega, \sigma_1, \sigma_2)$ with different widths

$$\sigma_2 = 1/2^J, \quad J=1-5$$



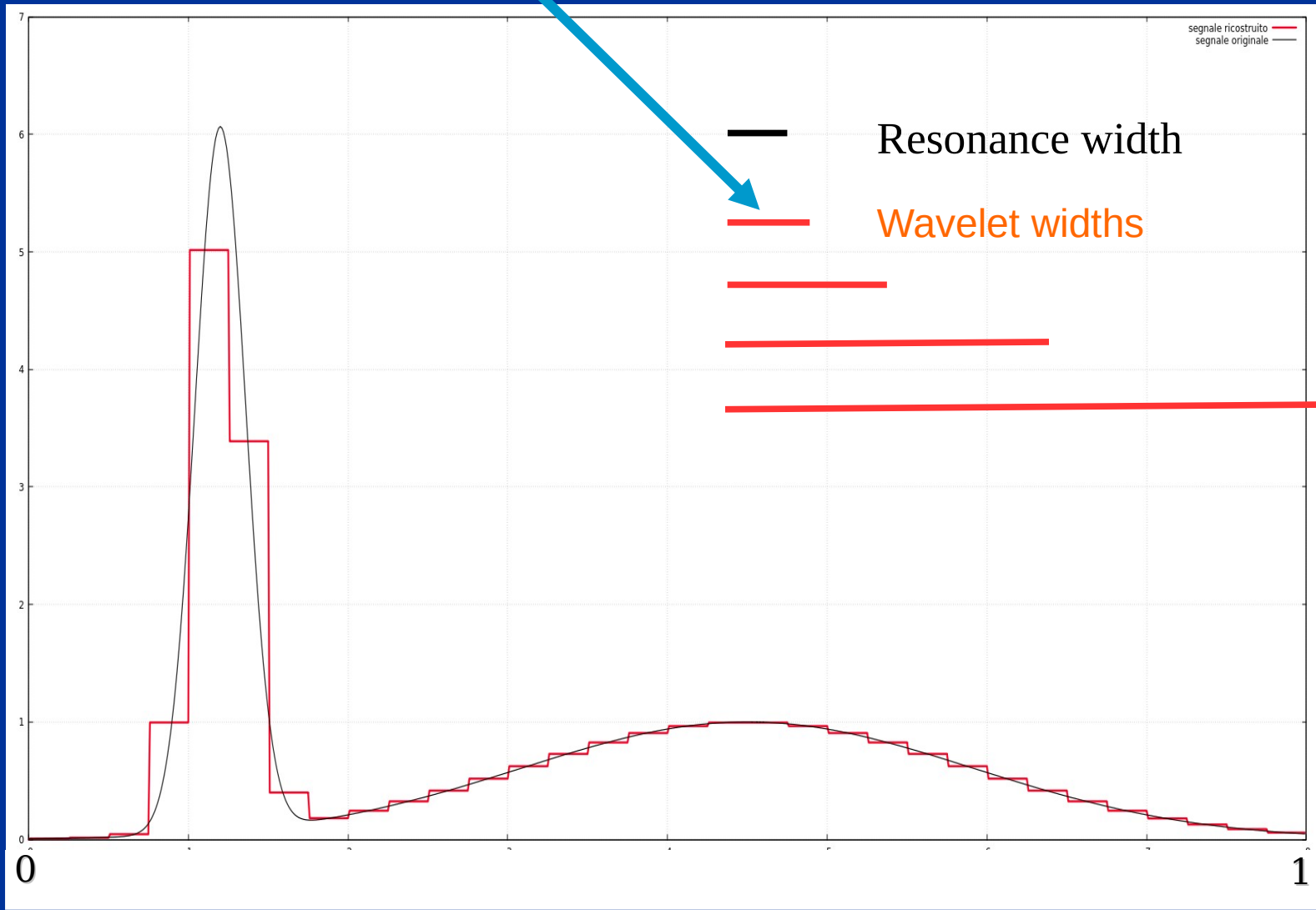
namely a lot of different resolutions up to $\sigma_2 = 0.03$!!!

**This may not be possible
with diagonalization in realistic
cases!**

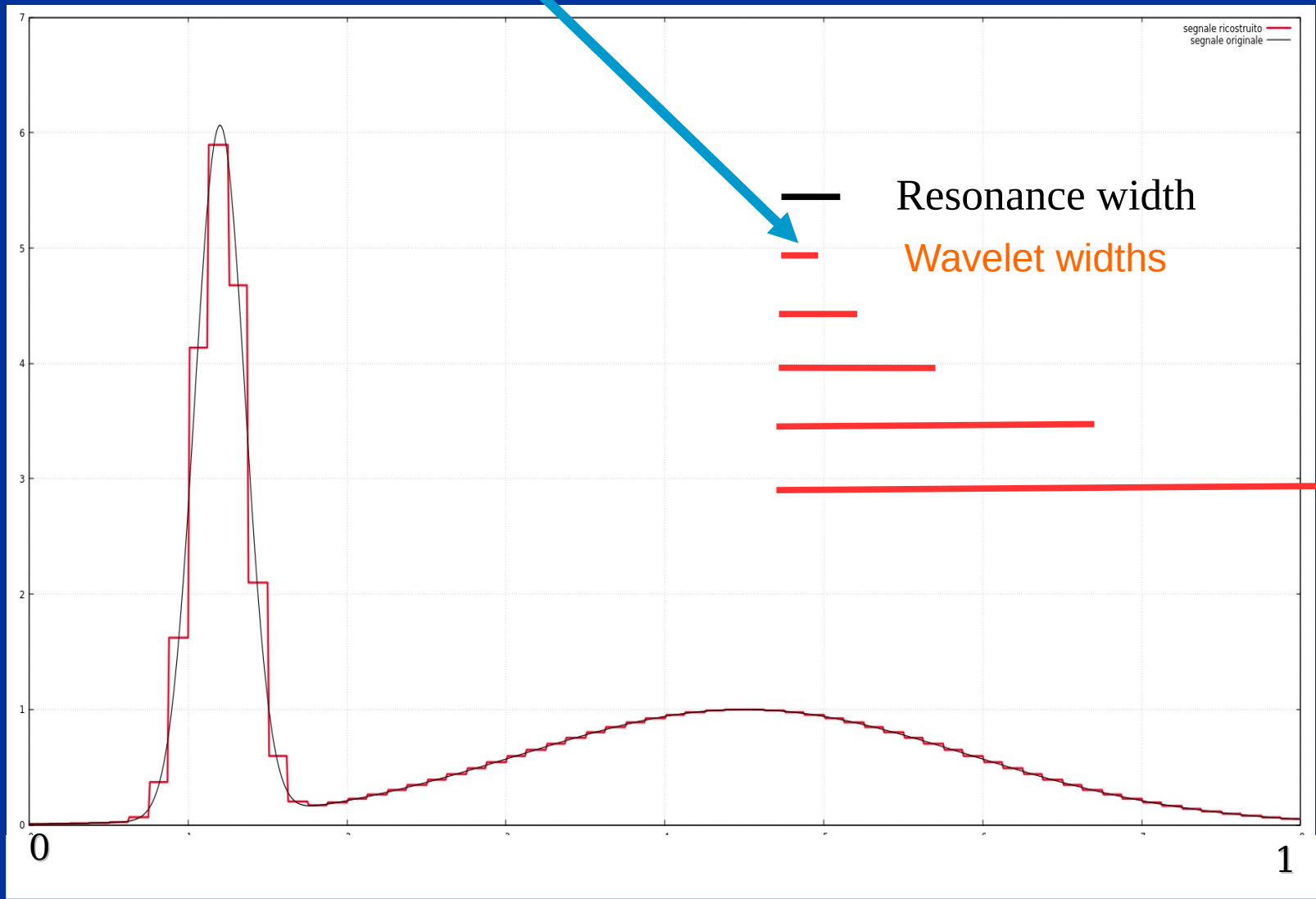
Remember Winfried's talk:

***<<one must have enough Hamiltonian eigenvalues
In the energy range of interest!>>***

Hp. on smallest “resolution” (low density of ϵ_λ):



Hp. on smallest “resolution” (higher density of ϵ_λ):



Summary:

- IT methods are alternative approaches to overcome the many-body scattering problem
- They are suitable for perturbative (inclusive and exclusive) as well as non-perturbative reactions
- New kernels have to be explored
- Big potentialities not yet exploited

Acknowledgements

to all people who have taken part in the
IT adventure over 20 years

- **Victor Efros**
- Winfried Leidemann
- Nir Barnea
- *Sonia Bacca*
- *Sofia Quaglioni*
- Ed Tomusiak
- The CC people (Gaute Hagen, Thomas Papenbrock, *Mirko Miorelli...*)
- The MC people (Francesco Pederiva, *Alessandro Roggero*)
- ...

**Thanks to the organizers
for the invitation!!**