

INTRO TO NON-EQUILIBRIUM 2PI EFFECTIVE ACTION TECHNIQUES

Gert Aarts

Physics Department, Swansea University



Swansea University
Prifysgol Abertawe

INTRODUCTION

nonequilibrium quantum field theory:

- framework with many applications
- in early universe:
inflation, baryon asymmetry, phase transitions, ...
- in relativistic heavy ion collisions
probing strongly interacting matter/extreme QCD
- in atomic physics, BEC, plasma physics, ...

INTRODUCTION

in this lecture:

- emphasis on methods
- relativistic quantum fields
- a few illustrations

REFERENCES

- I'll discuss work of many (relativistic) people, not properly inserting references throughout

Berges, Cox (2000)

Aarts, Berges, + Ahrensmeier, Baier, Serreau

Berges + Borsanyi, Serreau, Wetterich, + Reinosa

Cooper, Dawson, Mihaila

Juchem, Cassing, Greiner

Müller, Lindner

Arrizabalaga, Smit, Tranberg

Rajantie, Tranberg

Aarts + Bonini, Wetterich, + Martinez Resco, + Tranberg

Jeon, Yaffe

Calzetta, Hu

Carrington *et al*

...

OUTLINE

- what is nonequilibrium field theory?
- mean field theory
- 2PI effective action
- a few selected applications
- transport

QUANTUM DYNAMICS

GENERAL FORMULATION

well-defined problem:

- initial conditions: density matrix ρ_D
- time evolution: Heisenberg e.o.m. $\mathcal{O}(t) = e^{iHt} \mathcal{O} e^{-iHt}$
- observables:

$$\langle \mathcal{O}(t) \rangle = \text{Tr } \rho_D \mathcal{O}(t) \quad \langle \mathcal{O}(t) \mathcal{O}(t') \rangle = \text{Tr } \rho_D \mathcal{O}(t) \mathcal{O}(t') \quad \text{etc.}$$

- in equilibrium: $\rho_D \sim e^{-H/T}$, commutes with the evolution operator
- time translation invariance:

$$\langle \mathcal{O}(t) \rangle = \langle \mathcal{O}(0) \rangle \quad \langle \mathcal{O}(t) \mathcal{O}(t') \rangle = G(t - t')$$

QUANTUM DYNAMICS

GENERAL FORMULATION

out of equilibrium:

$$\langle \mathcal{O}(t) \rangle = \text{Tr } \rho_D \mathcal{O}(t) \quad \langle \mathcal{O}(t) \mathcal{O}(t') \rangle = \text{Tr } \rho_D \mathcal{O}(t) \mathcal{O}(t') \quad \text{etc.}$$

- density matrix ρ_D arbitrary ($[H, \rho_D] \neq 0$)
- initial value problem: start at $t = t_0$
- time translation invariance is broken:

$$\langle \mathcal{O}(t) \rangle = G(t - t_0) \quad \langle \mathcal{O}(t) \mathcal{O}(t') \rangle = G(t - t_0, t' - t_0)$$

- relation to the initial conditions: memory
- effective independence of t_0 as $t \rightarrow \infty$?

NONEQUILIBRIUM DYNAMICS

MAIN OBSTRUCTION

no exact solution method available

- use approximation methods
- language of unequal-time correlation functions
- n -point functions: hierarchy of coupled equations

approximation: truncate hierarchy

- problem not specific for quantum dynamics
- fluctuations: quantum and/or statistical

⇒ consider also classical statistical field theory

NONEQUILIBRIUM DYNAMICS

CORRELATION FUNCTIONS

- quantum field theory

$$\langle \phi(x)\phi(y) \rangle = \underbrace{\int d\phi' d\phi'' \langle \phi' | \rho_D | \phi'' \rangle}_{\text{initial cond.}} \underbrace{\int \mathcal{D}\phi e^{iS} \phi(x)\phi(y)}_{\text{quantum evolution (path integral)}}$$

- classical statistical field theory

$$\langle \phi(x)\phi(y) \rangle_{\text{cl}} = \int \underbrace{D\pi D\phi \rho_{\text{cl}}[\pi, \phi]}_{\text{initial prob. distribution and phase space integral}} \underbrace{\phi(x)\phi(y) + \text{equations of motion}}_{\text{classical evolution}}$$

NONEQUILIBRIUM DYNAMICS

ASIDE

in classical statistical field theory:

- “exact” evolution can be found numerically

$$\langle \phi(x)\phi(y) \rangle_{\text{cl}} = \int D\pi D\phi \rho_{\text{cl}}[\pi, \phi] \phi(x)\phi(y) + \text{e.o.m.}$$

- sample initial conditions from $\rho_{\text{cl}}[\pi, \phi]$
- solve e.o.m. for each of them
- average over initial conditions

note:

- classical thermal statistics: $n_{\text{cl}}(\omega) = T/\omega$
- Rayleigh-Jeans divergence
- careful with interpretation at late times

MEAN FIELD APPROXIMATIONS

SIMPLEST ATTEMPT

- equation of motion: $(\square + m^2)\phi = -\frac{\lambda}{6}\phi^3$
- expectation values:

$$\begin{aligned} \langle \phi(x) \rangle & \quad \text{coupled to} \quad \langle \phi^3(x) \rangle \\ G(x, y) = \langle \phi(x)\phi(y) \rangle & \quad \text{coupled to} \quad \langle \phi^3(x)\phi(y) \rangle \\ & \quad \text{etc.} \end{aligned}$$

- mean field/Gaussian/Hartree approximation: replace

$$\phi^3 \rightarrow 3\langle \phi^2 \rangle \phi \quad (\langle \phi \rangle = 0 \text{ for simplicity})$$

- self-consistent equation for two-point function

$$\left[\square + m^2 + \frac{\lambda}{2} G(x, x) \right] G(x, y) = 0$$

MEAN FIELD APPROXIMATIONS

SIMPLEST ATTEMPT

- successfully truncated hierarchy of correlation functions
- Gaussian approximation for $G(x, y) = \langle \phi(x)\phi(y) \rangle$
- same in quantum and classical theory

alas:

- approximation has a nonthermal fixed point
- best seen using equal-time correlation functions

$$G_{\phi\phi}(x - y, t) = \langle \phi(x, t)\phi(y, t) \rangle$$

$$G_{\pi\pi}(x - y, t) = \langle \pi(x, t)\pi(y, t) \rangle$$

$$G_{\pi\phi}(x - y, t) = \frac{1}{2} \langle \pi(x, t)\phi(y, t) + \phi(x, t)\pi(y, t) \rangle$$

MEAN FIELD APPROXIMATIONS

SIMPLEST ATTEMPT

- Gaussian approximation:

$$\partial_t G_{\phi\phi}(p, t) = 2G_{\pi\phi}(p, t)$$

$$\partial_t G_{\pi\phi}(p, t) = -\bar{\omega}_p^2 G_{\phi\phi}(p, t) + G_{\pi\pi}(p, t)$$

$$\partial_t G_{\pi\pi}(p, t) = -2\bar{\omega}_p^2 G_{\pi\phi}(p, t)$$

$$\text{with } \bar{\omega}_p^2 = p^2 + m^2 + \frac{\lambda}{2} \langle \phi^2 \rangle$$

- conserved quantity for every momentum mode p

$$C^2(p) = G_{\phi\phi}(p, t)G_{\pi\pi}(p, t) - G_{\pi\phi}^2(p, t)$$

$$\partial_t C(p) = 0$$

MEAN FIELD APPROXIMATIONS

SIMPLEST ATTEMPT

- nonthermal fixed point:

$$G_{\pi\pi}^*(p) = \bar{\omega}_p^2 G_{\phi\phi}^*(p)$$

$$G_{\pi\phi}^*(p) = 0$$

$$C^2(p) = G_{\phi\phi}^*(p) G_{\pi\pi}^*(p) \quad \text{fixed by initial ensemble}$$

$$\bar{\omega}_p^{*2} = p^2 + m^2 + \frac{\lambda}{2} \langle \phi^2 \rangle^*$$

explicit solution:

- $\langle \phi^2 \rangle^*$ determined by gap equation

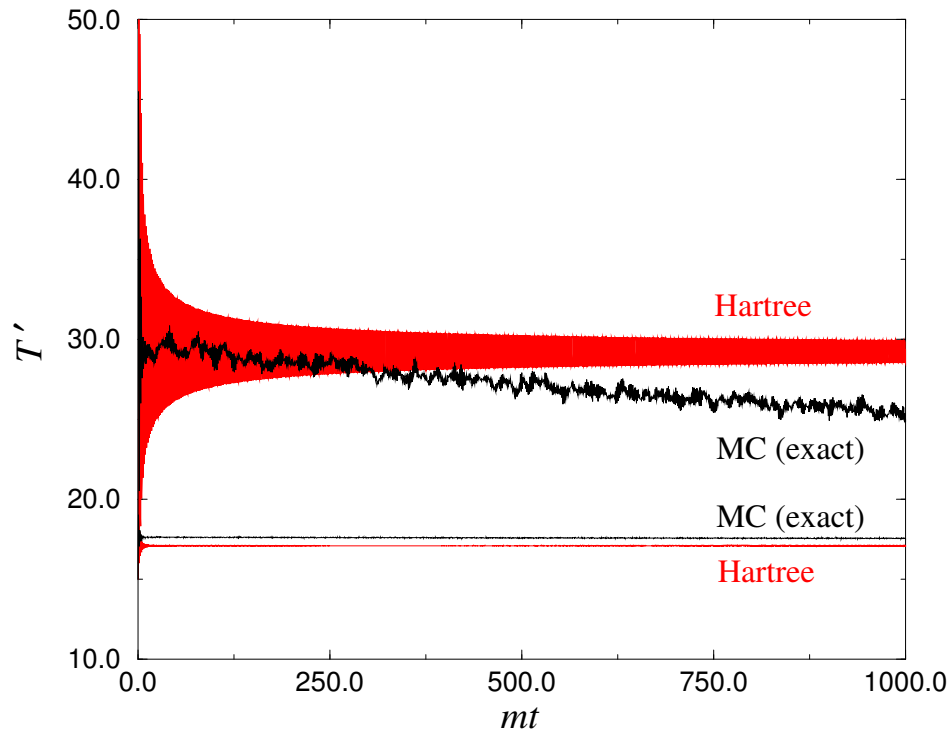
- $G_{\pi\pi}^*(p) = C(p) \bar{\omega}_p^*$ $G_{\phi\phi}^*(p) = C(p) / \bar{\omega}_p^*$

fixed point relevant for actual nonperturbative dynamics?

NONTHERMAL FIXED POINTS

G.A., BONINI AND WETTERICH

- classical test in 1 + 1 dimensions



classical mode
temperature:

$$T(p, t) = G_{\pi\pi}(p, t)$$

in classical thermal
equilibrium:

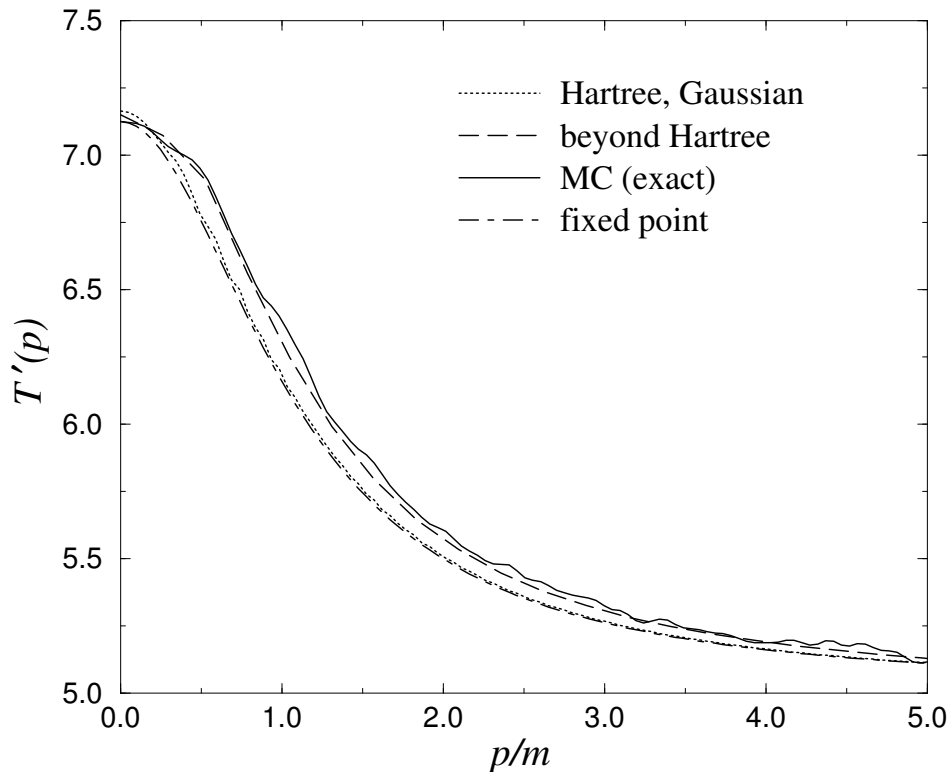
$$T(p, t) = T$$

- Hartree approximation: oscillating around nonthermal fixed point

NONTHERMAL FIXED POINTS

G.A., BONINI AND WETTERICH

- momentum-dependent “temperature” profile



classical mode
temperature:

$$T(p, t) = G_{\pi\pi}(p, t)$$

fixed point:

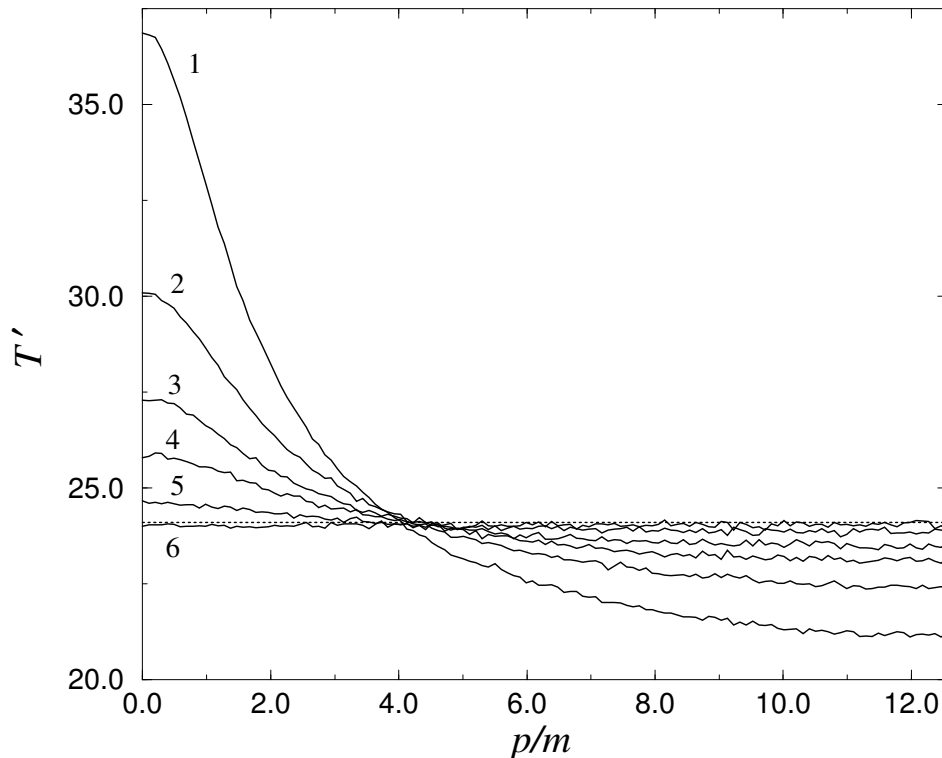
$$T^*(p) = T_0 \left[1 + \frac{\lambda}{2} \frac{\langle \phi^2 \rangle^*}{p^2 + m^2} \right]^{1/2}$$

- initial response determined by nonthermal fixed point, also for exact (MC) evolution

NONTHERMAL FIXED POINTS

G.A., BONINI AND WETTERICH

- momentum-dependent “temperature” profile



classical mode
temperature:

$$T(p, t) = G_{\pi\pi}(p, t) \rightarrow T$$

$$t_1 < t_2 < \dots < t_6$$

- fixed point relevant at early times
- exact (MC) evolution eventually thermalizes:
all modes the same temperature

NONEQUILIBRIUM QUANTUM FIELDS?

WISH LIST

- mean field approximation (dramatically) inadequate
- need to include scattering

want:

- stable time evolution
 - nontrivial due to secularities: many schemes break down when $t \sim 1/(\text{expansion parameter})$
- connection with well-established approaches, e.g. kinetic theory
- dynamics at very late times: conservation laws and hydrodynamics, transport
- ...

KINETIC THEORY

CONNECTION WITH ESTABLISHED METHODS

example:

- Boltzmann equation: $(\partial_t + \mathbf{v}_{\mathbf{p}} \cdot \partial_X) f(\mathbf{p}, X) = C[f]$

$$X = (t, \mathbf{x}) \quad \mathbf{v}_{\mathbf{p}} = \mathbf{p}/E_{\mathbf{p}} \quad p^0 = E_{\mathbf{p}} \quad (\text{onshell})$$

- real particles undergo isolated collisions
- collision kernel for two-to-two scattering processes:

$$C[f] = \frac{1}{2} \int_{\mathbf{p}'\mathbf{k}\mathbf{k}'} |\mathcal{M}|^2 (2\pi)^4 \delta^4(p + p' - k - k') \\ [(1 \pm f_{\mathbf{p}}) (1 \pm f_{\mathbf{p}'}) f_{\mathbf{k}} f_{\mathbf{k}'} - f_{\mathbf{p}} f_{\mathbf{p}'} (1 \pm f_{\mathbf{k}}) (1 \pm f_{\mathbf{k}'})]$$

- stationary solution: $f(\mathbf{p}, X) \rightarrow n(E_{\mathbf{p}}) = 1/[e^{E_{\mathbf{p}}/T} \mp 1]$

KINETIC THEORY

BEYOND KINETIC THEORY?

assumptions:

- onshell particles: phase space distribution
- isolated collisions, well separated in space and time
- ‘slowly varying’, gradient expansion

relax these assumptions:

quantum field theory

⇒ dynamics of correlation functions, in particular two-point functions

KINETIC THEORY

TWO-POINT FUNCTIONS

- Wightman functions:

$$G^>(x, y) = \langle \phi(x)\phi(y) \rangle = G^<(y, x)$$

- spectral function:

$$\rho(x, y) = i\langle [\phi(x), \phi(y)] \rangle = i(G^>(x, y) - G^<(x, y))$$

- statistical function:

$$F(x, y) = \frac{1}{2}\langle [\phi(x), \phi(y)]_+ \rangle = \frac{1}{2}(G^>(x, y) + G^<(x, y))$$

two-point functions closely related to particle distribution functions, after series of manipulations

KINETIC THEORY

TWO-POINT FUNCTIONS

- separation of slow and fast variables: Wigner transform

$$X = \frac{1}{2}(x+y) \quad (x-y) \rightarrow p \quad \Rightarrow \quad G^>(x,y) \rightarrow G^>(p,X)$$

- in equilibrium: Kubo-Martin-Schwinger (KMS) condition
periodicity of the trace (X independent)

$$G^>(x,y) \sim \text{Tr} e^{-H/T} \phi(x)\phi(y) \quad \Rightarrow \quad G^>(\omega, \mathbf{p}) = e^{\omega/T} G^<(\omega, \mathbf{p})$$

- all 2-point functions related to the spectral density

$$G^>(\omega, \mathbf{p}) = [n_B(\omega) + 1] \rho(\omega, \mathbf{p}) \quad G^<(\omega, \mathbf{p}) = n_B(\omega) \rho(\omega, \mathbf{p})$$

- noneq. distr. function: $G^<(p, X) = f(p, X) \rho(p, X)$

- onshell approximation $f(\mathbf{p}, X)$, with $p^0 = E_{\mathbf{p}}(X)$

2PI EFFECTIVE ACTION

FIELD THEORY APPROACH

therefore:

- two-point function important role
- obeys Dyson equation: $G^{-1} = G_0^{-1} - \Sigma$
- what is self energy Σ ?
- formalize: action principle

two-particle irreducible effective action

or

Φ -derivable approach

Luttinger/Ward, Baym, Cornwall/Jackiw/Tomboulis,

2PI EFFECTIVE ACTION

FIELD THEORY APPROACH

- generating functional with local and bilocal sources

$$Z[J, K] = e^{iW[J, K]} = \int \mathcal{D}\varphi e^{i(S[\varphi] + J_i \varphi^i + \frac{1}{2} \varphi^i K_{ij} \varphi^j)}$$

- Legendre transform: $\frac{\delta W}{\delta J_i} = \phi^i$, $\frac{\delta W}{\delta K_{ij}} = \phi^i \phi^j + G^{ij}$

$$\Gamma[\phi, G] = W[J, K] - J_i \phi^i - \frac{1}{2} K_{ij} (\phi^i \phi^j + G^{ij})$$

- effective action can be written as

$$\Gamma[\phi, G] = S[\phi] + \frac{i}{2} \text{Tr} \ln G^{-1} + \frac{i}{2} \text{Tr} G_0^{-1} (G - G_0) + \Gamma_2[\phi, G]$$

- variational principle (in absence of sources)

$$\frac{\delta \Gamma}{\delta \phi} = 0, \quad \frac{\delta \Gamma}{\delta G} = 0 \quad \Rightarrow \quad G^{-1} = G_0^{-1} - \Sigma[G], \quad \Sigma = 2i \frac{\delta \Gamma_2}{\delta G}$$

2PI EFFECTIVE ACTION

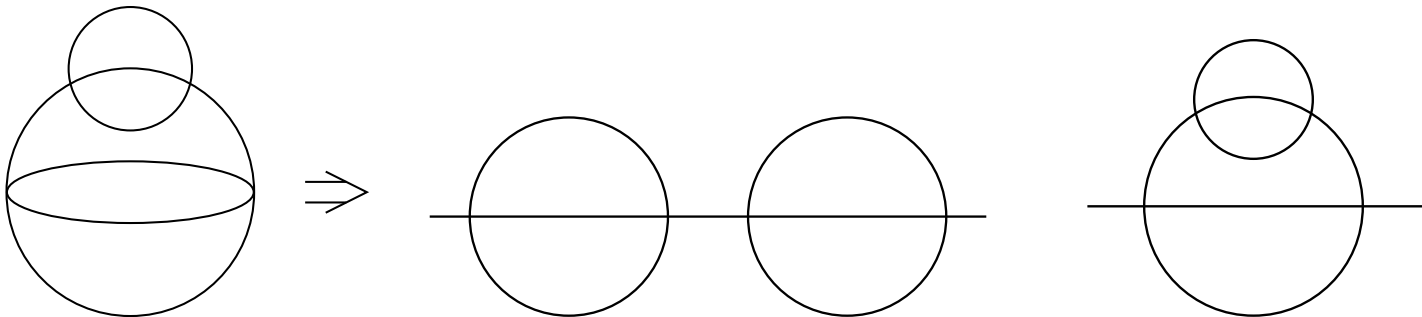
FIELD THEORY APPROACH

- action principle, at the extremum

$$(\square + V'[\phi]) \phi + \frac{\delta\Gamma_2[\phi, G]}{\delta\phi} = 0 \quad G^{-1} = G_0^{-1}[\phi] - \Sigma[\phi, G]$$

- prescription for the self energy $\Sigma = 2i\delta\Gamma_2/\delta G$
- Γ_2 is 2PI $\Leftrightarrow \Sigma$ is 1PI, depends on full G

example:

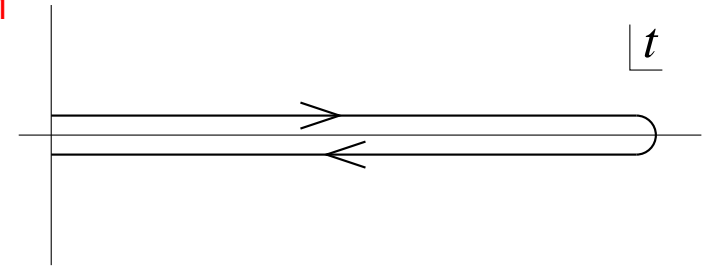


- avoid overcounting

NONEQUILIBRIUM DYNAMICS

INITIAL VALUE PROBLEM

solve equations in real time:



$$\langle \phi(x)\phi(y) \rangle = \underbrace{\int d\phi' d\phi'' \langle \phi' | \rho_D | \phi'' \rangle}_{\text{initial cond.}} \underbrace{\int \mathcal{D}\phi e^{iS} \phi(x)\phi(y)}_{\text{quantum evolution (path integral)}}$$

initial cond.

quantum evolution
(path integral)

use Schwinger-Keldysh contour for initial value problems

$$i (\square_x + m^2) G(x, y) = \int_{\mathcal{C}} dz \Sigma(x, z) G(z, y) + \delta_{\mathcal{C}}(x - y)$$

action principle along complex-time path \mathcal{C}

NONEQUILIBRIUM DYNAMICS

INITIAL VALUE PROBLEM

- Green functions: $G^>$, $G^<$ etc.
- minimal choice
- decompose contour propagator in real and imaginary parts:

$$G(x, y) = F(x, y) - \frac{i}{2} \text{sign}(x^0 - y^0) \rho(x, y)$$

statistical function
even, anti-commutator

spectral function
odd, commutator

- spectral function is a commutator:

$$\rho(x, y) \Big|_{x^0=y^0} = 0, \quad \partial_{x^0} \rho(x, y) \Big|_{x^0=y^0} = \delta(\mathbf{x} - \mathbf{y})$$

NONEQUILIBRIUM DYNAMICS

INITIAL VALUE PROBLEM

manifestly real and causal equations

$$\begin{aligned} [\square_x + m^2] F(x, y) &= - \int_0^{x^0} dz^0 \int d\mathbf{z} \Sigma_\rho(x, z) F(z, y) \\ &\quad + \int_0^{y^0} dz^0 \int d\mathbf{z} \Sigma_F(x, z) \rho(z, y) \\ [\square_x + m^2] \rho(x, y) &= - \int_{y^0}^{x^0} dz^0 \int d\mathbf{z} \Sigma_\rho(x, z) \rho(z, y) \end{aligned}$$

with $\Sigma_{F,\rho}$ given in terms of F and ρ

predicting the future = remembering the past

NONEQUILIBRIUM DYNAMICS

INITIAL VALUE PROBLEM

- action principle
- conserved energy ($\langle\phi\rangle = 0$):

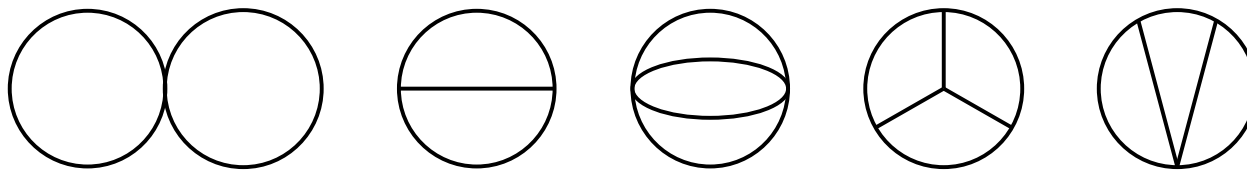
$$E = \int d^3x \frac{1}{2} [\partial_{x^0} \partial_{y^0} + \partial_{x^i} \partial_{y^i} + m^2] F(x, y) \Big|_{x=y} \\ + \frac{1}{4} \int d^3x \int_0^{x^0} dz^0 \int d^3z [\Sigma_\rho(x, z) F(z, x) - \Sigma_F(x, z) \rho(z, x)]$$

- conserved for every truncation

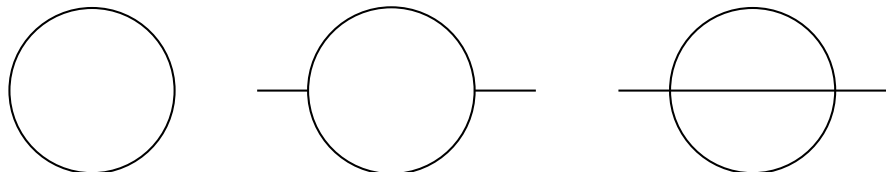
2PI TRUNCATIONS

LOOP AND $1/N$ EXPANSIONS TO NEXT-TO-LEADING ORDER

- so far exact, approximation enters via truncation of Γ_2
- systematic, in practice loop and $1/N$ expansions
- three-loop expansion ($\langle\phi\rangle = 0$)



- diagrams 1, 2, 3 well-studied (no internal vertices)
- self energies:

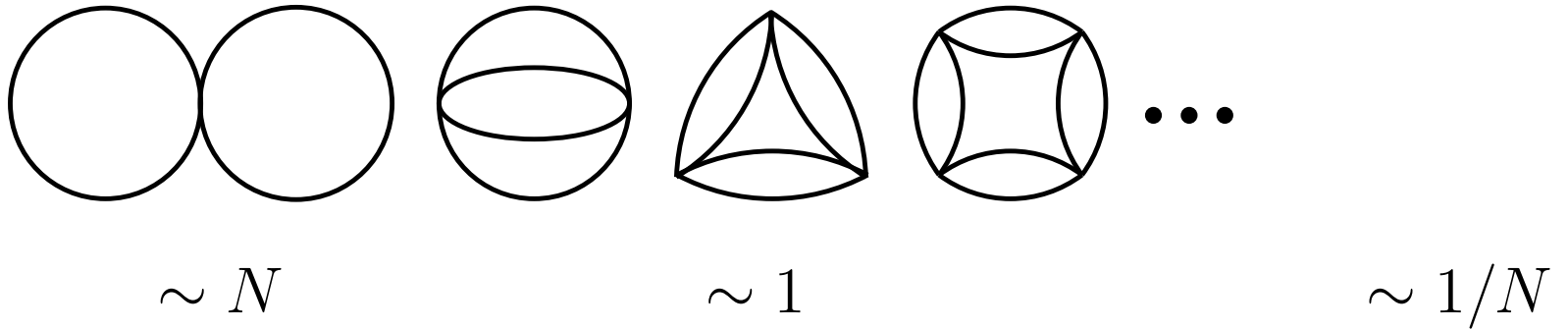


2PI TRUNCATIONS

LOOP AND $1/N$ EXPANSIONS TO NEXT-TO-LEADING ORDER

large N expansion:

- $O(N)$ model, vertex $\sim 1/N$ (with $\langle \phi \rangle = 0$ for simplicity)



2PI TRUNCATIONS

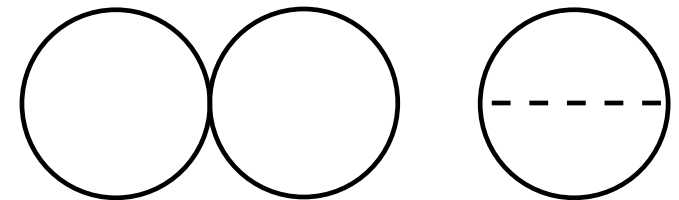
LOOP AND $1/N$ EXPANSIONS TO NEXT-TO-LEADING ORDER

large N expansion:

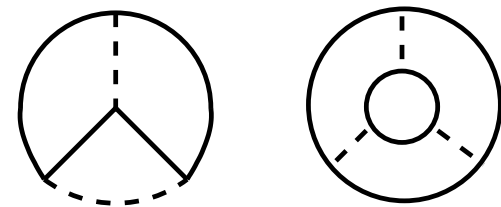
- $O(N)$ model, vertex $\sim 1/N$ (with $\langle \phi \rangle = 0$ for simplicity)
- efficient formulation: use chain of bubbles

$$\text{---} \langle \text{---} \rangle \text{---} = \text{---} \times \text{---} + \text{---} \circ \text{---} \text{---}$$

\Rightarrow effective two-loop approximation



- NNLO contribution ($\sim 1/N$):

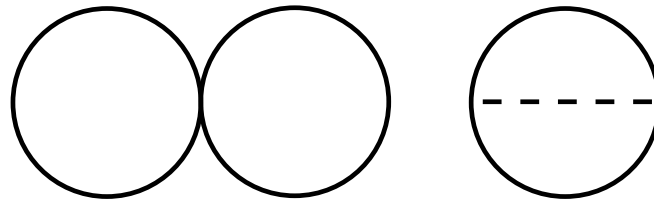


2PI TRUNCATIONS

LOOP AND $1/N$ EXPANSIONS TO NEXT-TO-LEADING ORDER

large N expansion:

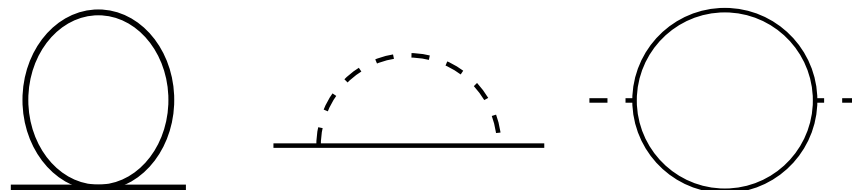
- $O(N)$ model, vertex $\sim 1/N$ (with $\langle \phi \rangle = 0$ for simplicity)



- dressed propagators:

$$G^{-1} = G_0^{-1} - \Sigma \quad D^{-1} = D_0^{-1} - \Pi$$

- self energies:



2PI TRUNCATIONS

LOOP AND $1/N$ EXPANSIONS TO NEXT-TO-LEADING ORDER

- closed set of self-consistent equations:

$$\begin{aligned} [\square_x + M^2(x)] F(x, y) = & - \int_0^{x^0} dz^0 \int d\mathbf{z} \Sigma_\rho(x, z) F(z, y) \\ & + \int_0^{y^0} dz^0 \int d\mathbf{z} \Sigma_F(x, z) \rho(z, y) \end{aligned}$$

$$[\square_x + M^2(x)] \rho(x, y) = - \int_{y^0}^{x^0} dz^0 \int d\mathbf{z} \Sigma_\rho(x, z) \rho(z, y)$$

with

$$\Sigma_F(x, y) = -\frac{\lambda}{3N} \left[F(x, y) D_F(x, y) - \frac{1}{4} \rho(x, y) D_\rho(x, y) \right]$$

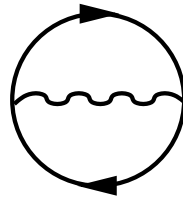
$$\Sigma_\rho(x, y) = -\frac{\lambda}{3N} [\rho(x, y) D_F(x, y) + F(x, y) D_\rho(x, y)]$$

2PI TRUNCATIONS

LOOP AND $1/N$ EXPANSIONS TO NEXT-TO-LEADING ORDER

large N expansion:

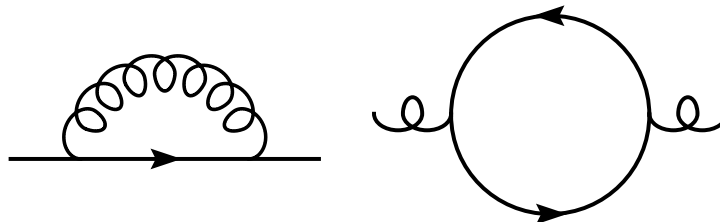
- large N_f gauge theory, vertex $e^2 \sim 1/N$



- dressed propagators:

$$G^{-1} = G_0^{-1} - \Sigma \quad D^{-1} = D_0^{-1} - \Pi$$

- self energies:



SOLUTIONS

NUMERICAL

- solve integro-differential equations on a spacetime lattice
- straightforward discretization, no further approximation
- expensive numerically due to “memory kernel”

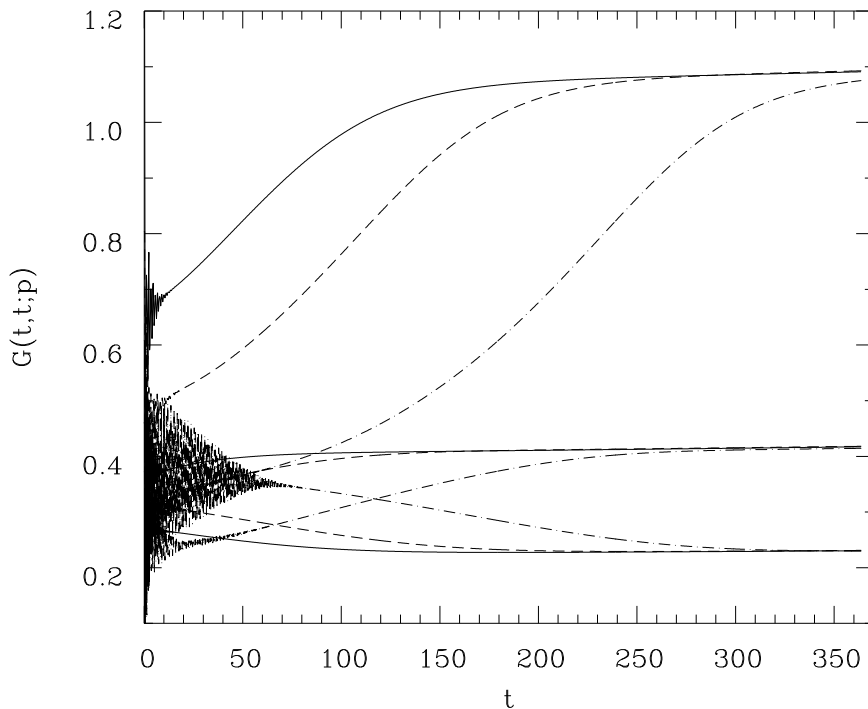
some applications

LOSS OF MEMORY

THERMALIZATION

first results by Berges and Cox (2000):

- take different initial conditions (or density matrices) with the total energy density identical
- independence of initial conditions at late times



3-loop expansion in $\lambda\phi^4$
in 1 + 1 dimensions

time evolution of different
momentum modes $F(t, t; p)$

PRECISION TESTS

CLASSICAL 2PI APPROXIMATION

- 2PI approach in classical statistical field theory
- possibility to compare with “exact” solution
- sampling of initial conditions + numerical integration of classical equation of motion

example of classical limit: three-loop approximation

$$\Sigma_{\rho}(x, z) = -\frac{\lambda^2}{2} \rho(x, z) \left[F^2(x, z) - \frac{1}{12} \rho^2(x, z) \right],$$
$$\Sigma_F(x, z) = -\frac{\lambda^2}{6} F(x, z) \left[F^2(x, z) - \frac{3}{4} \rho^2(x, z) \right]$$

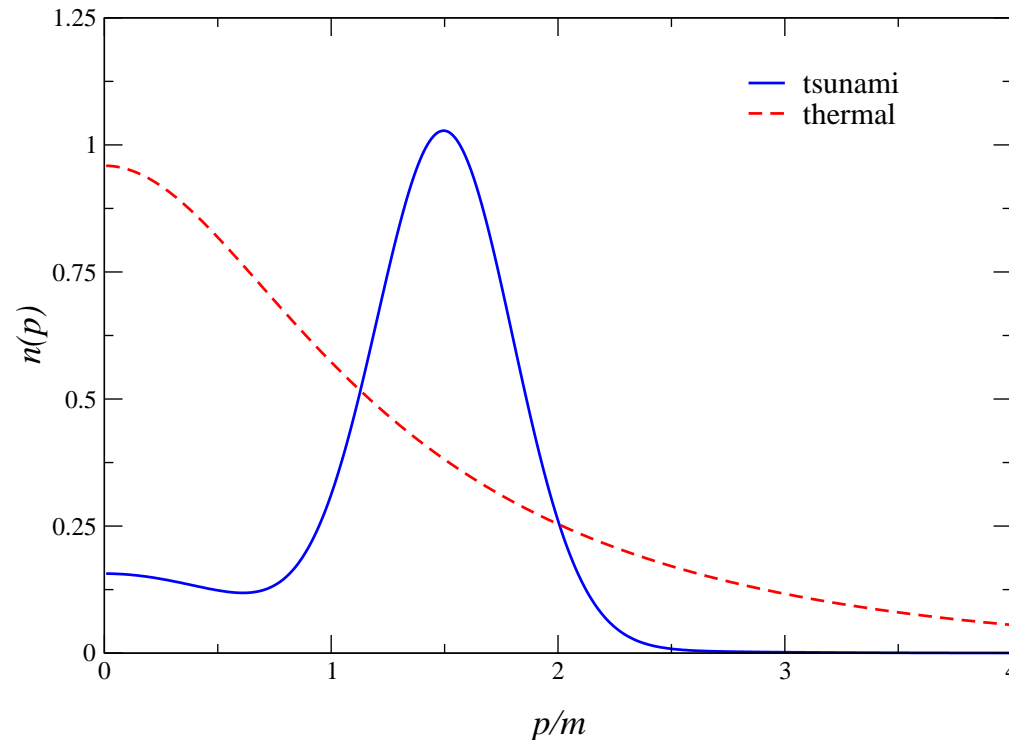
classically:

$$\Sigma_{\rho}^{\text{cl}}(x, z) = -\frac{\lambda^2}{2} \rho(x, z) F^2(x, z) \quad \Sigma_F^{\text{cl}}(x, z) = -\frac{\lambda^2}{6} F^3(x, z)$$

NONEQUILIBRIUM INITIAL CONDITIONS

TSUNAMI

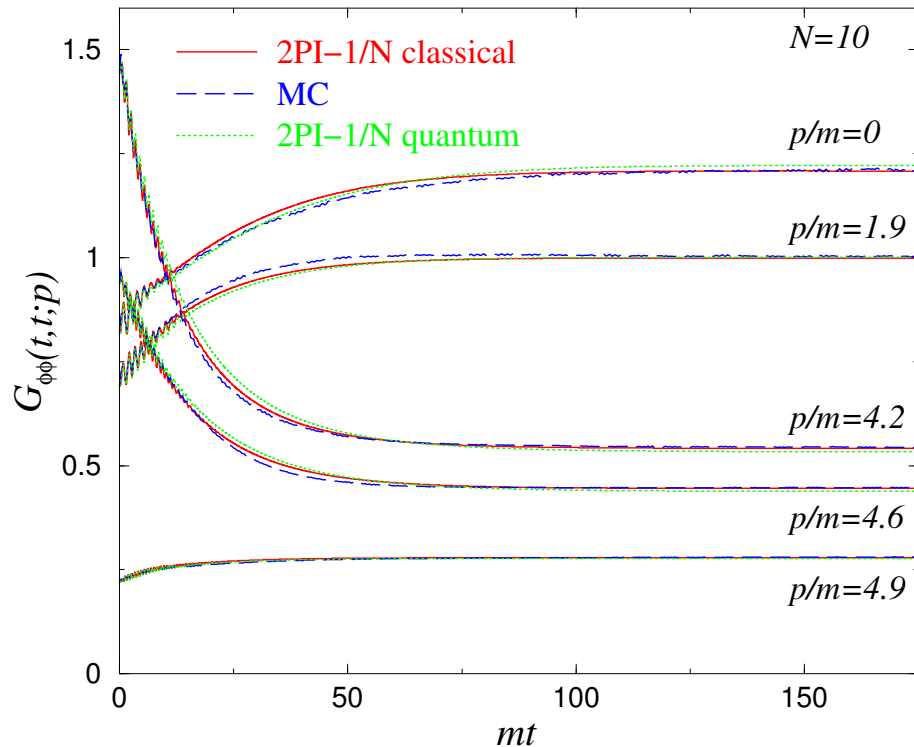
- Gaussian initial conditions far from equilibrium
- specify $F(t, t'; \mathbf{p})$, $\partial_t F(t, t'; \mathbf{p})$, $\partial_t \partial_{t'} F(t, t'; \mathbf{p})$ at $t = t' = 0$ in terms of initial particle number $n(\mathbf{p})$



- easily implemented in exact and 2PI dynamics

PRECISION TESTS

G.A. & BERGES

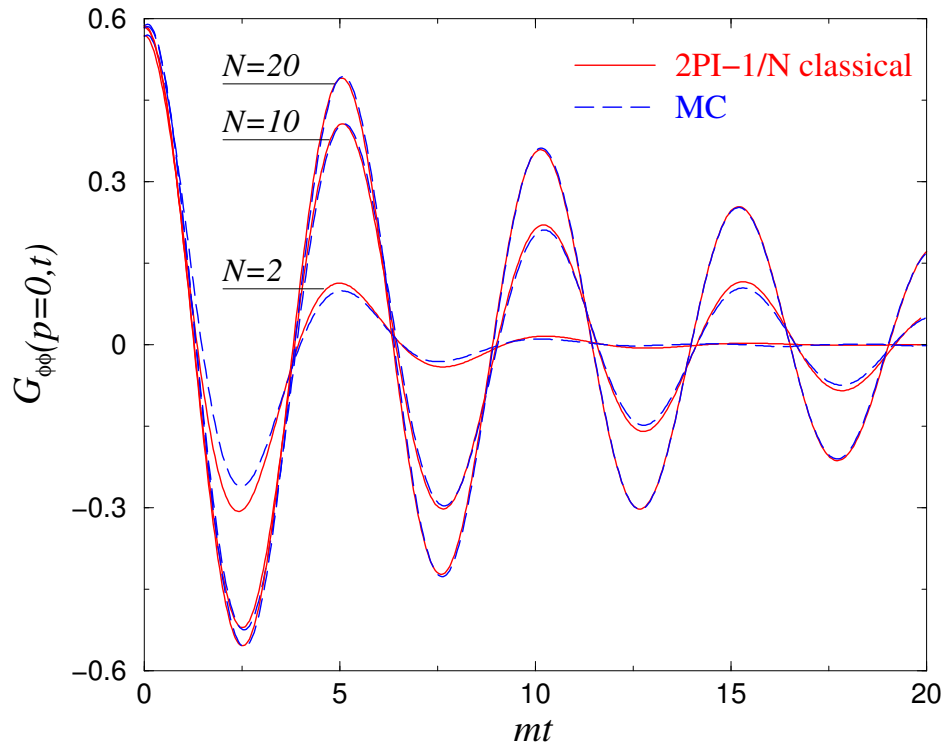


- tsunami initial conditions
- equal-time correlation function: ‘particle number’
- high energy density: compare quantum and classical evolution

- evolution from $2\text{PI}-1/N$ expansion in agreement with ‘exact’ evolution, also for late times.
- reliable description of both early and late times
- capable of describing equilibration

PRECISION TESTS

G.A. & BERGES



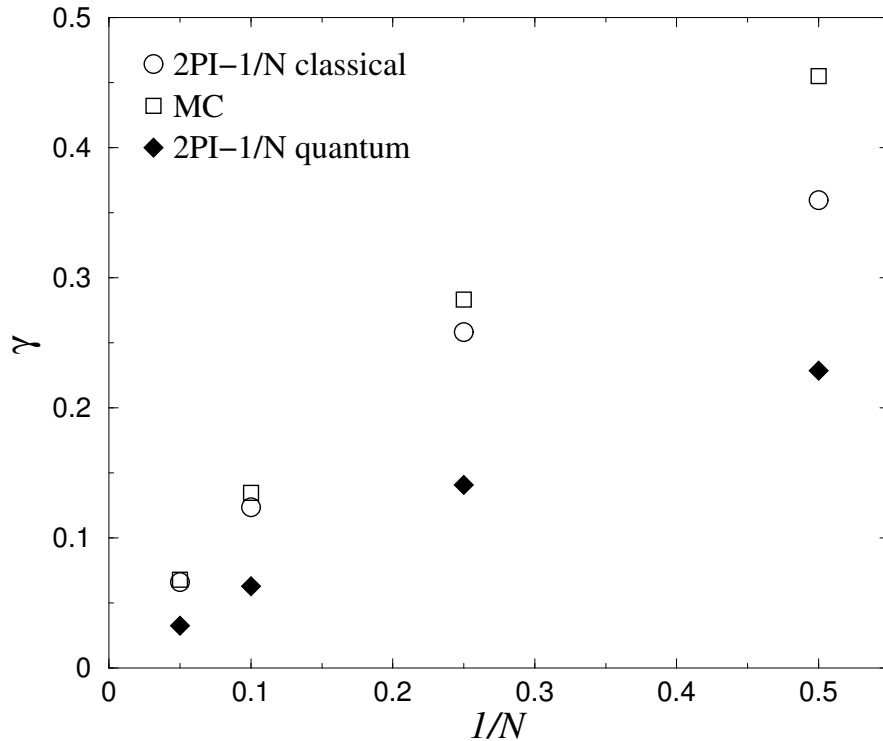
2PI-1/ N expansion

unequal-time correlation
function

- Monte Carlo: sample of 80.000 initial conditions
- 2PI-1/ N : one (expensive) numerical solution
- quantitative agreement for larger N

PRECISION TESTS

G.A. & BERGES



2PI-1/ N expansion

assume ansatz

$$G(t, 0; \mathbf{p}) \sim e^{-\gamma t} \cos mt$$

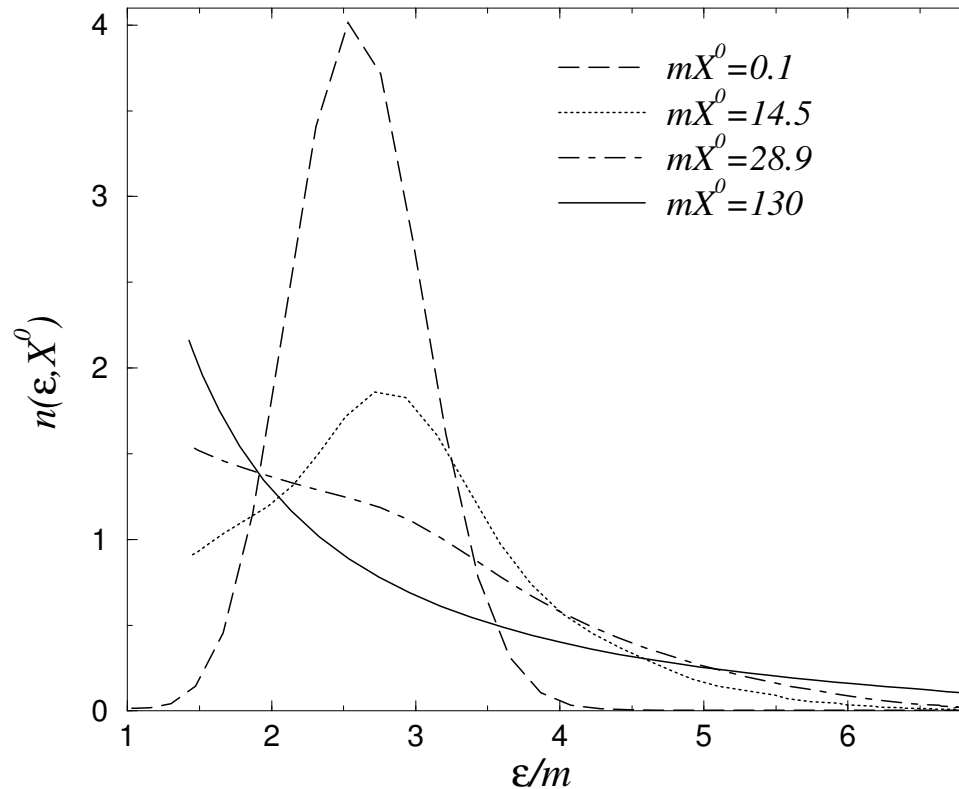
fit γ and m

- compare classical 2PI with classical exact
- quantitative agreement for larger N
- compare classical 2PI with quantum 2PI
- quantum \neq classical!

(NOT) KINETIC THEORY

G.A. & BERGES

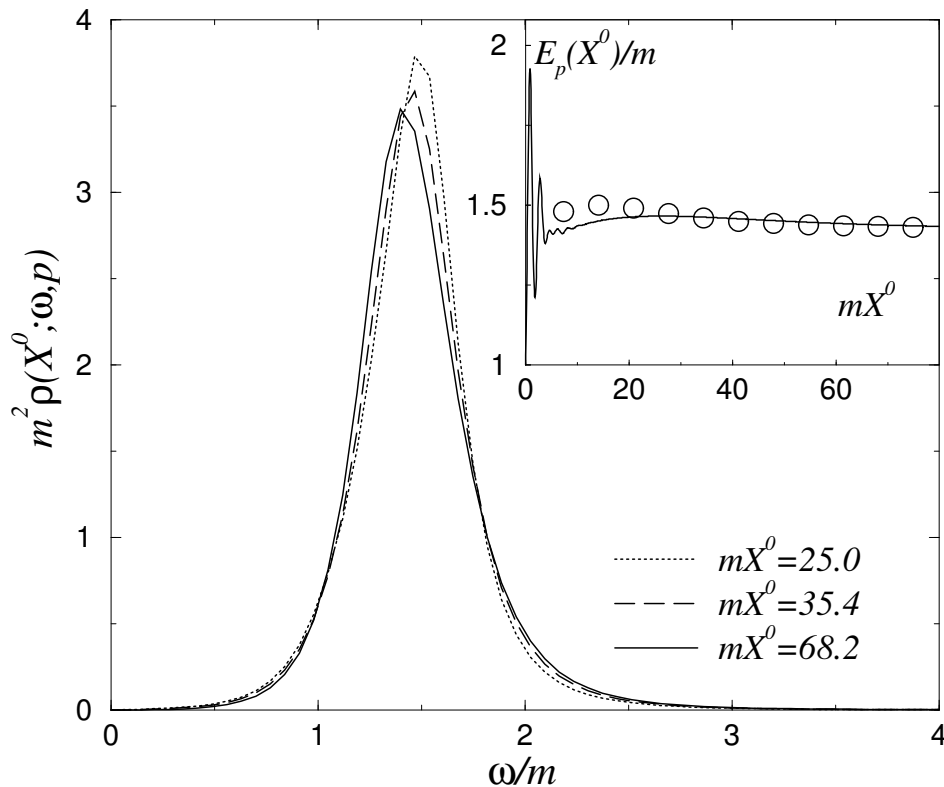
- separation of fast and slow variables
- effective particle number distribution is evolving fast and wildly



(NOT) KINETIC THEORY

G.A. & BERGES

- self-consistent evolution of the spectral function $\rho(t, t'; \mathbf{p})$
- no quasiparticle approximation
- Wigner transform: $\rho(t, t'; \mathbf{p}) \rightarrow \rho(\omega, \mathbf{p}; X^0)$



$$X^0 = (t + t')/2$$

- quasiparticle peak
- non-zero width
- slowly evolving

ET CETERA

much more work has been done:

- quick establishment of equation of state (prethermalization)
- fermions
- momentum anisotropy
- (tachyonic) preheating
- warm inflation
- renormalization
- ...

- cold atoms

TRANSPORT

FINAL STAGES

unified picture:

- dynamics far from equilibrium with 2PI truncations
- system will (eventually) equilibrate and thermalize

precise question:

- which scattering processes are certainly included?
- which scattering processes are certainly not included?

TRANSPORT

FINAL STAGES

- final stages of evolution
- dynamics of nearly conserved quantities
- hydrodynamic modes are slowest

- energy-momentum
- charges
- ...

evolve according to “low-energy effective field theory”

=

hydrodynamics

NEAR EQUILIBRIUM: TRANSPORT COEFFICIENTS

KUBO RELATIONS AND LINEAR RESPONSE

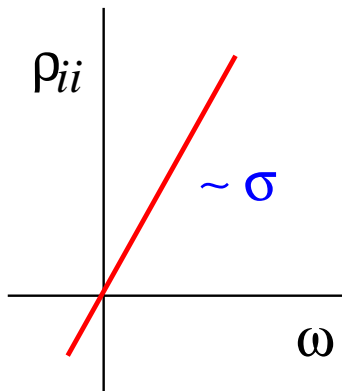
electrical conductivity:

$$\sigma = \frac{1}{6} \frac{\partial}{\partial \omega} \rho^{ii}(\omega, \mathbf{0}) \Big|_{\omega=0}$$

shear viscosity:

$$\eta = \frac{1}{20} \frac{\partial}{\partial \omega} \rho_{\pi\pi}(\omega, \mathbf{0}) \Big|_{\omega=0}$$

spectral densities:



$$\rho^{\mu\nu}(\omega, \mathbf{p}) = \int d^4x e^{ipx} \langle [j^\mu(x), j^\nu(0)] \rangle_{\text{eq}}$$

$$\rho_{\pi\pi}(\omega, \mathbf{p}) = \int d^4x e^{ipx} \langle [\pi_{ij}(x), \pi_{ij}(0)] \rangle_{\text{eq}}$$

with $j^\mu = \bar{\psi} \gamma^\mu \psi$, $\pi_{ij} = T_{ij} - \frac{1}{3} \delta_{ij} T_k^k = \partial_i \phi \partial_j \phi - \frac{1}{3} \delta_{ij} \partial_k \phi \partial_k \phi$

transport coefficients

\sim

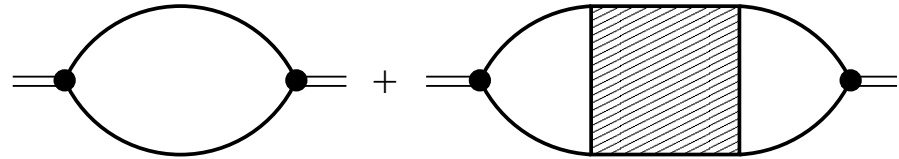
slope of current-current
spectral functions at $\omega = 0$

NEAR EQUILIBRIUM: TRANSPORT COEFFICIENTS

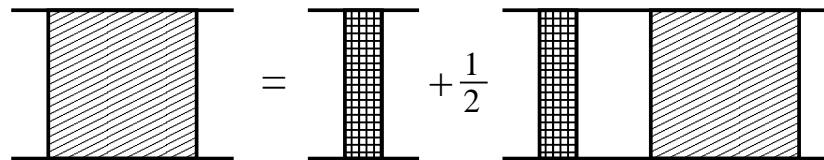
G. A. AND J. M. MARTINEZ RESCO

imaginary part of correlators of bilocal operators

2PI effective action
as generating functional:



generates ladder diagrams:



with kernel or rung:

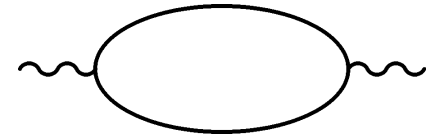
$$\Lambda_{ij;kl} = 4i \frac{\delta^2 \Gamma_2}{\delta G^{ij} \delta G^{kl}} = 2 \frac{\delta \Sigma_{ij}}{\delta G^{kl}}$$

kernel (and self energy) determined by Γ_2

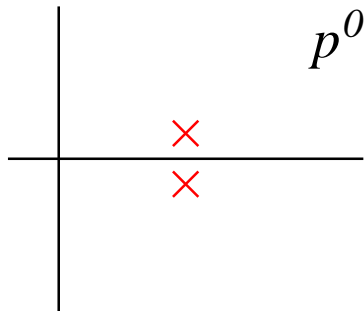
NEAR EQUILIBRIUM: TRANSPORT COEFFICIENTS

DRESSED PROPAGATORS

- pinching poles:



$$\lim_{\omega \rightarrow 0} \rho^{ii}(\omega, \mathbf{0}) = 4e^2 \omega \int \frac{d^4 p}{(2\pi)^4} n'_F(p^0) G_R(p) G_A(p)$$



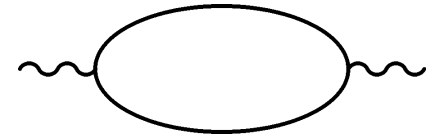
propagators in the loop carry the same momentum, product of retarded (R) and advanced (A) propagators

- with bare propagators ill-defined

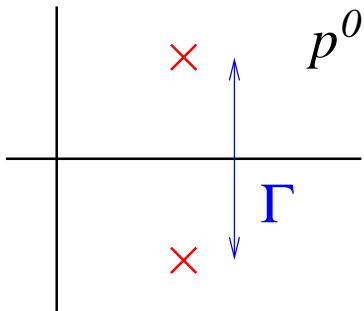
NEAR EQUILIBRIUM: TRANSPORT COEFFICIENTS

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propagators in the loop carry the same momentum, product of retarded (R) and advanced (A) propagators

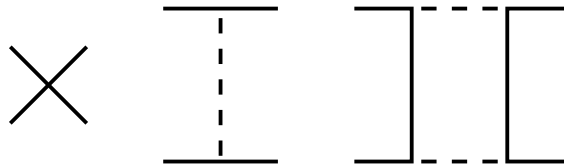
- inclusion of thermal width $\Gamma \sim 1/N$ required

\Rightarrow finite collision time/mean free path in a medium resummed nonperturbatively

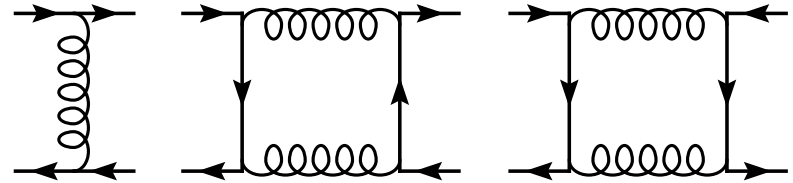
NEAR EQUILIBRIUM: TRANSPORT COEFFICIENTS

RUNGS AND LADDER DIAGRAMS

● $O(N)$ model



● large N_f QED/QCD

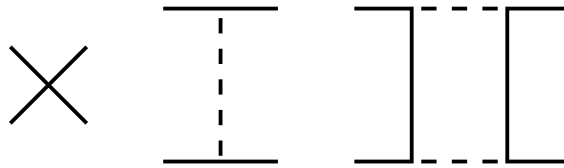


+ subleading terms in the $1/N$ expansion

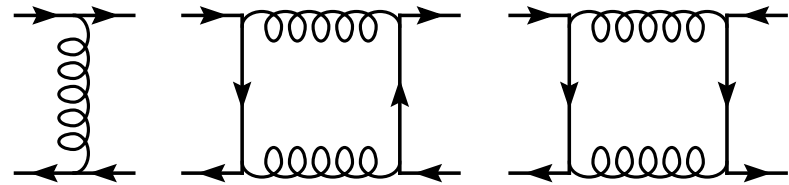
NEAR EQUILIBRIUM: TRANSPORT COEFFICIENTS

RUNGS AND LADDER DIAGRAMS

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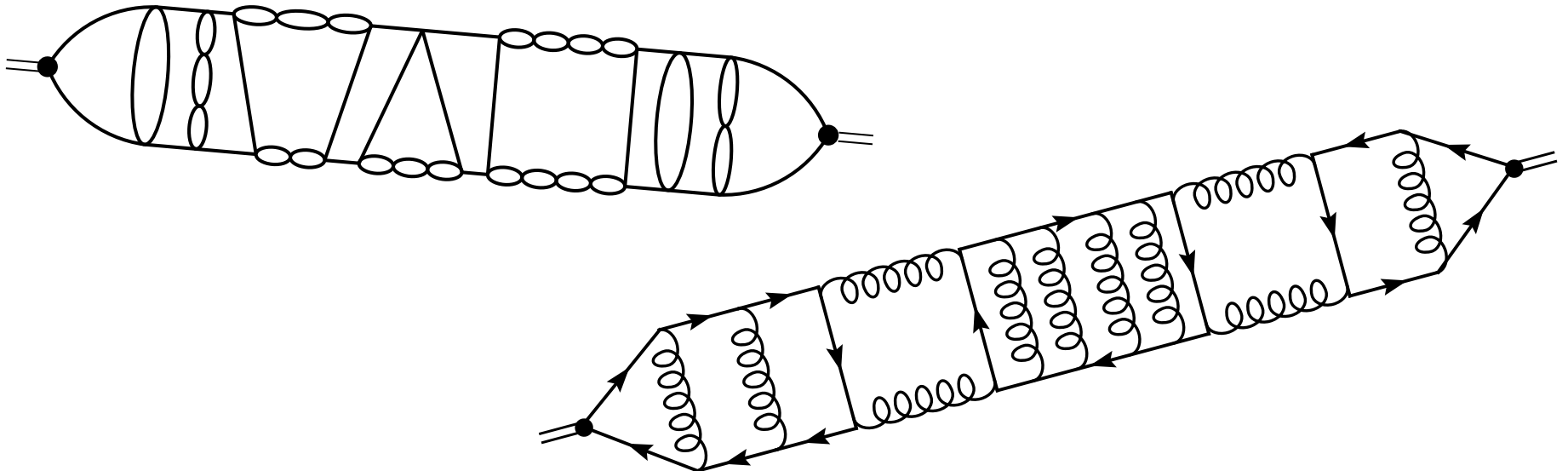


● large N_f QED/QCD



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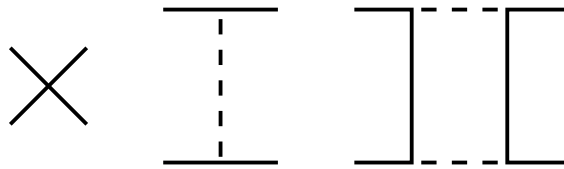
typical ladder diagrams:



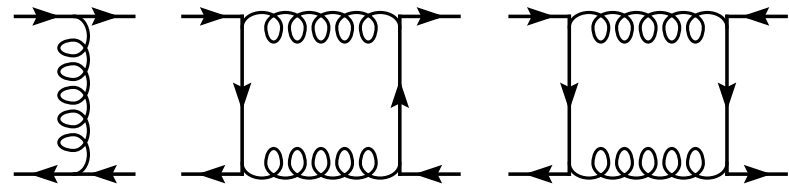
NEAR EQUILIBRIUM: TRANSPORT COEFFICIENTS

RUNGS AND LADDER DIAGRAMS

● $O(N)$ model



● large N_f QED/QCD



+ subleading terms in the $1/N$ expansion

● power counting

- positive powers of N : closed scalar or fermion loops and pairs of propagators with pinching poles
- negative powers of N : vertices

all contributions to LO in $1/N$ expansion

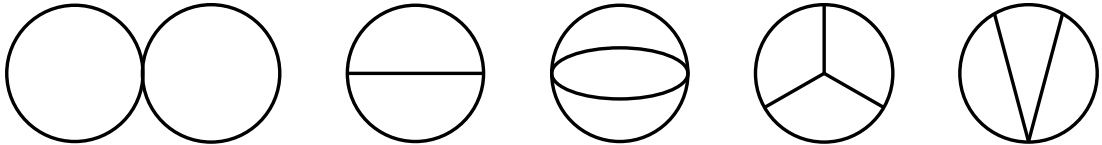
- subleading terms: cannot be neglected for self-consistent dynamics far from equilibrium

TRANSPORT

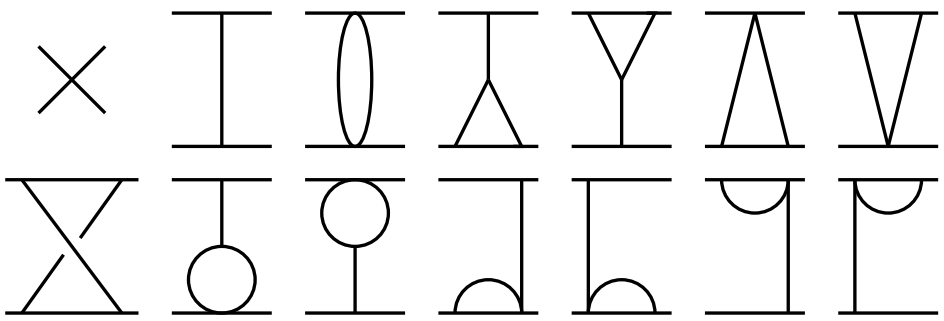
PRECISE QUESTION

which scattering processes are (not) included?

- 3-loop expansion in $g\phi^3 + \lambda\phi^4$ theory



- kernel

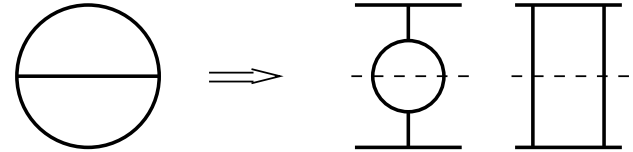


- a lot of scattering processes

TRANSPORT

PRECISE QUESTION

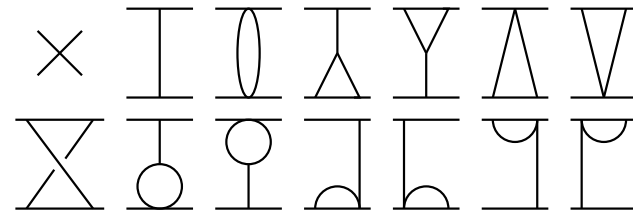
- 2-loop approximation:
(iterated)



- sum of squares of 2 \rightarrow 2 scattering processes

$$|\mathcal{M}|^2 \sim g^4 [|G(s)|^2 + |G(t)|^2 + |G(u)|^2]$$

- 3-loop approximation:



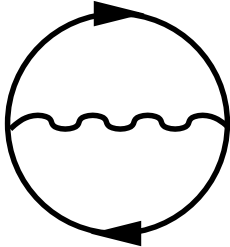
- square of sum of 2 \rightarrow 2 scattering processes
(+ subleading vertex corrections)

$$|\mathcal{M}|^2 \sim \left| \lambda + g^2 [G(s) + G(t) + G(u)] \right|^2$$

- interference included

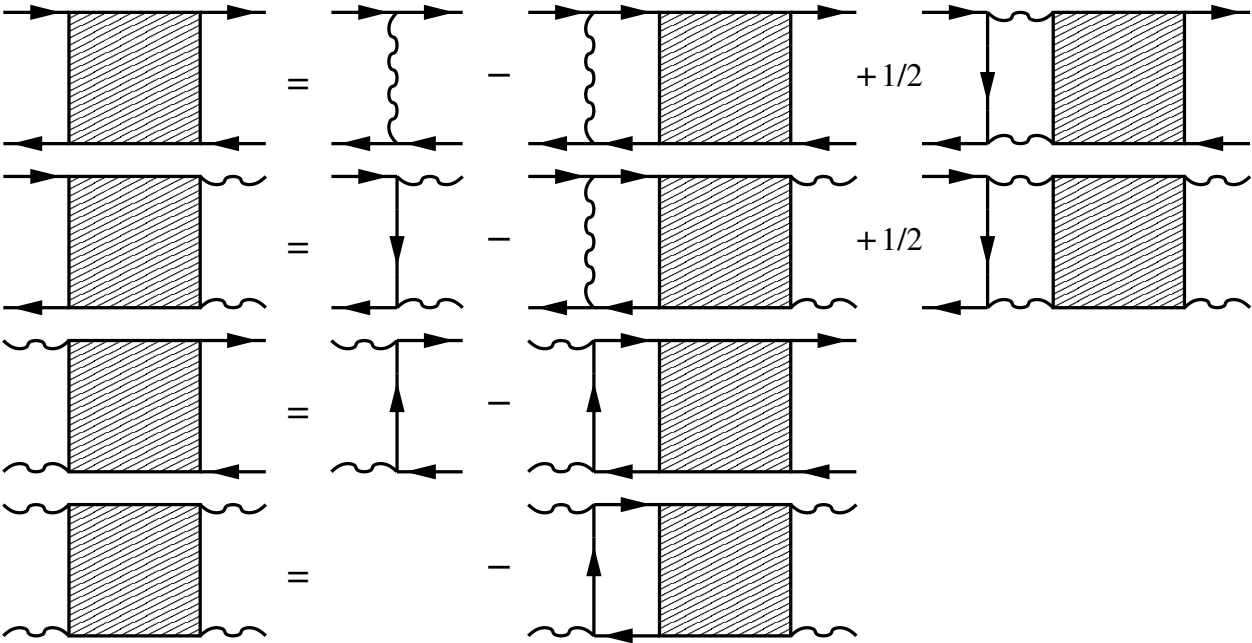
TRANSPORT

PRECISE QUESTION



- 2-loop or large N_f expansion in QED

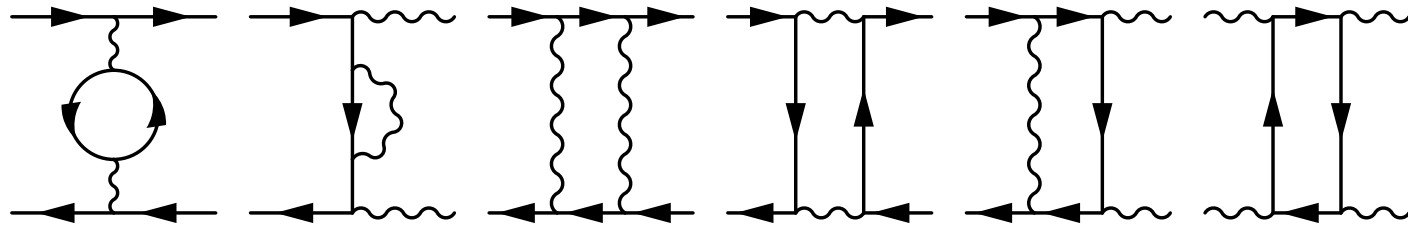
- coupled integral equations:



TRANSPORT

PRECISE QUESTION

- scattering kernel



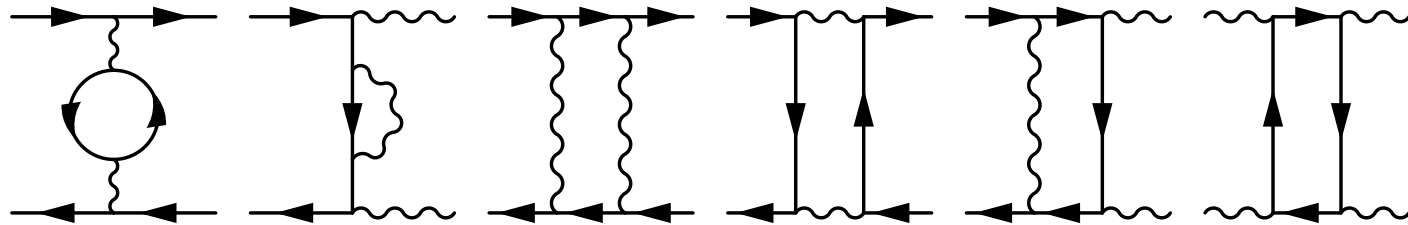
- weak coupling in the leading log approximation
(transport coefficient $\sim 1/e^4 \ln 1/e$)

- rung 1 \Rightarrow t -channel Coulomb scattering
- rung 2 \Rightarrow Compton scattering, pair annihilation

TRANSPORT

PRECISE QUESTION

- scattering kernel



- weak coupling in the leading log approximation (transport coefficient $\sim 1/e^4 \ln 1/e$)
 - rung 1 \Rightarrow t -channel Coulomb scattering
 - rung 2 \Rightarrow Compton scattering, pair annihilation
- leading order large N_f QED:
 - rung 1 and 4 \Rightarrow Coulomb scattering in all channels (no interference)

TRANSPORT

SUMMARY

scalars/fermions (with current truncations):

- most transport coefficients correct to LO
- notable exception: bulk viscosity Calzetta and Hu

gauge theories (two loop truncations):

- correct to leading log
- correct at leading order in large N_f
- full leading order requires use of 3PI effective action

(Carrington et al)

OUTLOOK

done:

- scalars/fermions: most formal aspects studied
- some applications
- gauge theories: formal developments in progress

to do:

- more applications for scalars/fermions possible
- gauge theories: more formal developments
- gauge theories: numerical implementation and tests