# Vacuum polarization in 1+1 spacetime dimensions

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### **Abstract**

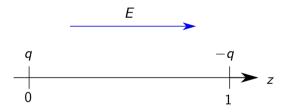
In my M.Sc. project, I study the effects of the backreaction of a charged Klein-Gordon field coupled to an external electric field, in 1+1 dimensional spacetime. In this talk, I summarize the statement of the problem I intend to solve, and present preliminary results and observations from my project.

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- ► Preliminaries
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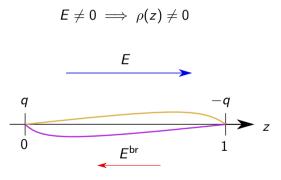
# The set-up

### 1 Preliminaries



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#### 1 Preliminaries



### Semi classical Klein-Gordon-Maxwell (KGM) equations

The goal is to solve the system of coupled differential equations

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With the gauge covariant derivative

$$D_{\mu} = \partial_{\mu} + ieA_{\mu},$$

and the boundary conditions

$$\begin{cases} \phi \mid_{z=0,1} = 0 \\ F^{\mu\nu} \mid_{z=0,1} = \lambda \end{cases}$$

#### 1 Preliminaries

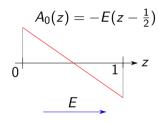


Figure: The background electric potential

1. Constant electric field of strength  $\lambda$  pointing towards positive z.

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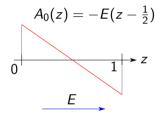


Figure: The background electric potential

- 1. Constant electric field of strength  $\lambda$  pointing towards positive z.
- 2. Under the Coulomb gauge

$$A_0(t,z) = -Ez, A_1(t,z) = 0$$

up to additive constant

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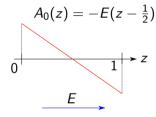


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$$A_0(t,z) = -Ez, A_1(t,z) = 0$$

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3. To ensure anti-symmetric solutions

$$A_0(t,z)=-E(z-\frac{1}{2})$$

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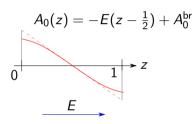


Figure: The full electric potential (background+backreaction)

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$$A_0(t,z) = -Ez, A_1(t,z) = 0$$

up to additive constant

3. To ensure anti-symmetric solutions

$$A_0(t,z) = -E(z - \frac{1}{2}) + A_0^{br}(z)$$

#### 1 Preliminaries

### Klein-Gordon equation

With the chosen gauge the KGM equations turn into

$$((\partial_t + ieA_0)^2 - \partial_z^2 + m^2)\phi(t, z) = 0,$$
 
$$\partial_z^2 A_0^{br} = -\langle \rho(z) \rangle_{\phi},$$

with

$$A_0(z) = -E\left(z - \frac{1}{2}\right) + A_0^{\text{br}}(z),$$
  
 $\rho(z) = ie\left((D_0\phi)^*\phi - \phi^*D_0\phi\right)$ 

# Time independent Klein-Gordon equation

1 Preliminaries

### **Mode equation**

$$\phi(t,z) = \phi_n(z)e^{-i\omega_n t} \implies$$

# Time independent Klein-Gordon equation

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### Mode equation

$$\phi(t,z) = \phi_n(z)e^{-i\omega_n t} \implies \left( \left[\omega_n - eA_0(z)\right]^2 + \frac{d^2}{dz^2} - m^2 \right)\phi_n = 0$$

### Time independent Klein-Gordon equation

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Ignoring the backreaction of the scalar field (External field approximation)

$$\left(\left[\omega_n + \lambda\left(z - \frac{1}{2}\right)\right]^2 + \frac{d^2}{dz^2} - m^2\right)\phi_n = 0, \ \lambda = eE$$

### **Analytic solutions**

#### 2 External field approximation

When  $A_0(z) = -\lambda(z - \frac{1}{2})$ , the KG equation can be solved analytically

$$\phi_n(z) = a_n D_{i rac{m^2}{2\lambda} - rac{1}{2}} \left( rac{1+i}{\sqrt{\lambda}} \left( \omega_n + \lambda \left( z - rac{1}{2} 
ight) 
ight) 
ight) \ + b_n D_{-i rac{m^2}{2\lambda} - rac{1}{2}} \left( rac{i-1}{\sqrt{\lambda}} \left( \omega_n + \lambda \left( z - rac{1}{2} 
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ight)$$

with  $D_{\nu}(z)$  the parabolic cylinder functions.

## Instabilities of the external field approximation

#### 2 External field approximation

The external field approximation yields instabilities for critical  $\lambda$  values [AW83]:

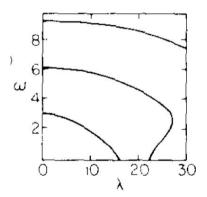


Figure: Energy of the first three modes as the background field strength  $\lambda$  increases, without considering the backreaction of the field. Courtesy of [AW83].

### Backreaction avoids those instabilities?

#### 2 External field approximation

The main claim in [AW83] is that considering the backreaction of the scalar field raises the energy levels, avoiding instabilities

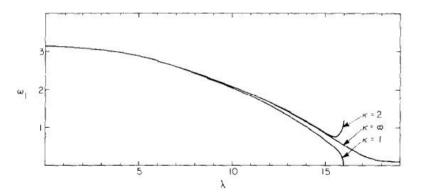


Figure: Positive energy levels for increasing  $\lambda$  with backreaction considered. [AW83]

### Vacuum polarization

#### 3 Quantisation and renormalisation

Vacuum polarization is calculated as the zeroth component of the charge density current

$$\rho(z) = ie((D_0\phi)^* \phi - \phi^* D_0\phi).$$

This operator is non-linear on the fields  $\implies$  ill-defined expectation value.

# Mode expansion of the vacuum polarization

3 Quantisation and renormalisation

In [AW83] the vacuum polarization is directly calculated as

$$\langle 0 | \rho(z) | 0 \rangle = ie \langle 0 | \phi^* D_0 \phi - \phi (D_0 \phi)^* | 0 \rangle =$$

$$e \left( \sum_{n>0} (\omega_n - eA_0) |\phi_n|^2 + \sum_{n<0} (\omega_n - eA_0) |\phi_n|^2 \right)$$

### **Hadamard states**

#### 3 Quantisation and renormalisation

Assume the two-point function w(x, x') is of Hadamard form

$$\langle 0 | \phi(x) \phi^*(x') | 0 \rangle = w(x, x') = \underbrace{H(x, x')}_{\text{Divergent}} + \underbrace{R(x, x')}_{\text{Smooth}}$$

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w(x, x') is a divergent distribution as  $x' \to x$ , but so is H(x, x'). In a way similar to the normal ordering prescription of RQFT, define

$$\langle 0|\, D_{\alpha}\phi(x)\, (D_{\beta}\phi(x))^*\, |0\rangle := \lim_{\substack{v'\to x\\ }} \left[D_{\alpha}D_{\beta'}^{\prime*}\left(w(x,x')-H(x,x')\right)\right].$$

## Renormalized vacuum polarization

3 Quantisation and renormalisation

### Hadamard Point-splitting[WZ20]

$$\langle 0 | \rho(z) | 0 \rangle = ie \langle 0 | \phi^* D_0 \phi - \phi (D_0 \phi)^* | 0 \rangle =$$

$$e \lim_{\tau \to 0} \left( \sum_{n > 0} (\omega_n - eA_0) |\phi_n|^2 e^{i\omega_n(\tau + i\epsilon)} + \sum_{n < 0} (\omega_n - eA_0) |\phi_n|^2 e^{-i\omega_n(\tau + i\epsilon)} \right) + \frac{e^2}{\pi} A_0$$

### Renormalized vacuum polarization

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### Mode sum formula [AW83]

$$\langle 0 | \rho(z) | 0 \rangle = ie \langle 0 | \phi^* D_0 \phi - \phi (D_0 \phi)^* | 0 \rangle =$$

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# Renormalized vacuum polarization

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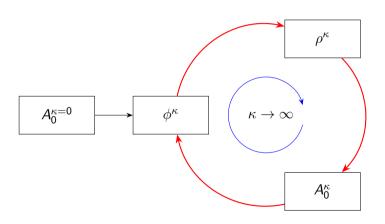
$$\langle 0| \, 
ho(z) \, |0\rangle = e \sum_{n=-N, n \neq =0}^{N} (\omega_n - eA_0) |\phi_n|^2 + \frac{e^2}{\pi} A_0$$

### Mode sum formula [AW83]

$$\langle 0 | \rho(z) | 0 \rangle = e \sum_{n=-1, n\neq =0}^{1} (\omega_n - eA_0) |\phi_n|^2$$

# **Closing the loop**

### 3 Quantisation and renormalisation



### Closing the loop

#### 3 Quantisation and renormalisation

$$\left(\left[\omega_n^{\kappa+1}-eA_0^{\kappa}(z)\right]^2+rac{d^2}{dz^2}-m^2
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ho^{\kappa}(z'')dz''dz'.$$

### Closing the loop

#### 3 Quantisation and renormalisation

$$\left(\left[\omega_n^{\kappa+1} - eA_0^{\kappa}(z)\right]^2 + \frac{d^2}{dz^2} - m^2\right)\phi_n^{\kappa+1} = 0$$
  $A_0^{\kappa}(z) = -\lambda\left(z - \frac{1}{2}\right) - \int_{\frac{1}{2}}^z \int_0^{z'} \rho^{\kappa}(z'')dz''dz'.$ 

### As a fixed point problem

Reminiscent of fixed point problems

$$A_0^{\kappa+1}=f(A_0^{\kappa})$$

4 Fixed point problems

#### **Definition**

For X a metric space, a fixed point  $x \in X$  of a function  $f : X \to X$  is defined as

$$x = f(x)$$
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$$A_0 \longrightarrow (\omega_n, \phi_n) \longrightarrow \rho \longrightarrow -\int \int \rho \longrightarrow A_{\text{background}} + A_{\text{induced}}(A_0)$$

# One dimensional fixed point problems

4 Fixed point problems

### Fixed point

Find, if it exists

$$\lim_{n\to\infty} f^n(x), f^n := \underbrace{f \circ \ldots \circ f}_{\text{n times}}$$

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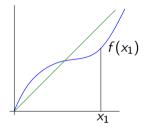


Figure: Iterations in fixed point problems

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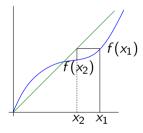


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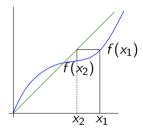


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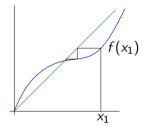


Figure: Iterations in fixed point problems

## The relaxing update rule

4 Fixed point problems

To avoid numerical instabilities due to  $\Delta \lambda$  being too big,

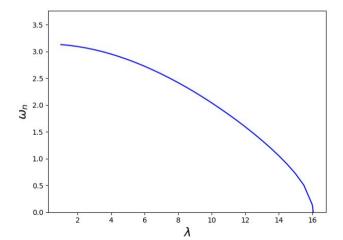
$$A_0^{\kappa+1} = cA_0^{\kappa} + (1-c)(A_{\mathsf{background}} + A_{\mathsf{induced}}^{\kappa}), 0 < c \lesssim 1$$

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# The energy of the modes for each $\lambda$

### 5 Results

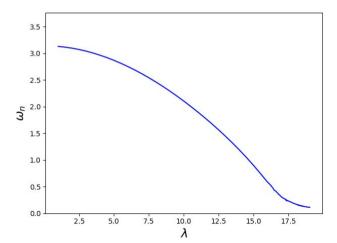
Figure:  $\omega_1$  as a function of  $\lambda$  for the massless case in the **external field approximation**.



# The energy of the modes for each $\lambda$

### 5 Results

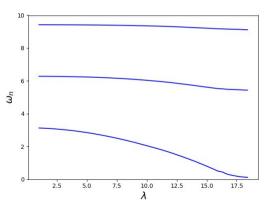
Figure:  $\omega_1$  as a function of  $\lambda$  in the mode sum prescription with N=1.



# The energy of the modes for each $\lambda$

### 5 Results

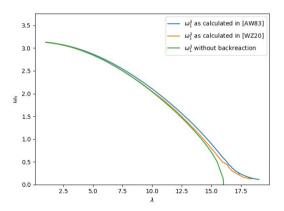
Figure: The energy of the first three modes of the scalar field in the Hadamard point-splitting prescription with cutoff N=12 as  $\lambda$  increases.



# The energy of the modes at each $\lambda$

#### 5 Results

Figure:  $\omega_1$  as a function of  $\lambda$  for the mode sum formula prescription, Hadamard point-splitting prescription and external field approximation.



# Self consistent vacuum polarizations for low $\lambda$

#### 5 Results

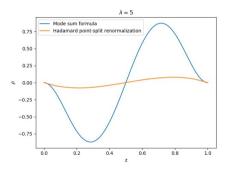


Figure:  $\lambda = 5$  comparison of the self consistent vacuum polarization.

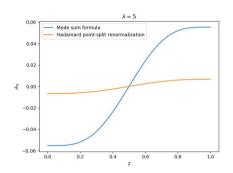


Figure:  $\lambda = 5$  comparison of the self consistent induced  $A_0$ 

# Self consistent vacuum polarizations for high $\lambda$

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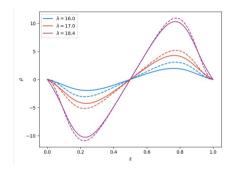
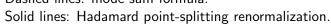


Figure: Higher values of  $\lambda$  comparison of the self consistent vacuum polarization.

Dashed lines: mode sum formula.



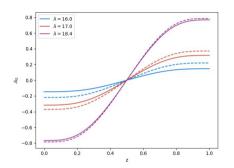


Figure: Higher values of  $\lambda$  comparison of the self consistent vacuum polarization.

# Vacuum polarization for different $\lambda$

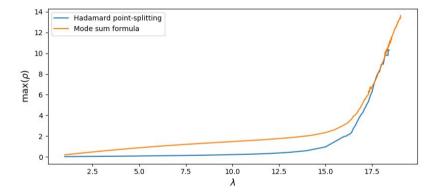


Figure: The maximum of the vacuum polarization for different values of the background electric field strength.

5 Results

# Asymptotic behaviour?

6 Standing questions

It is not yet safe to say that backreaction completely avoids instabilities of the solutions for each background electric field strength  $\lambda$ .

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6 Standing questions

It is not yet safe to say that backreaction completely avoids instabilities of the solutions for each background electric field strength  $\lambda$ .

There are two main problems that can arise for a given configuration

- 1. The root finding algorithm for the boundary conditions finds no  $\omega_1$
- 2. There is no convergence

These two issues can be solved by a smaller  $\Delta\lambda$ , or a relaxing parameter c closer to 1. It can get however to the order of  $\Delta\lambda\sim 10^{-7}$ .

## **Good convergence**

### 6 Standing questions

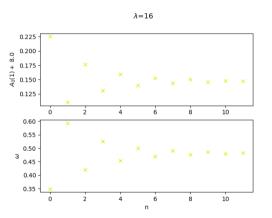


Figure: The iterations in the potential and the energy of the first mode in a case of good convergence.

## No convergence

### 6 Standing questions

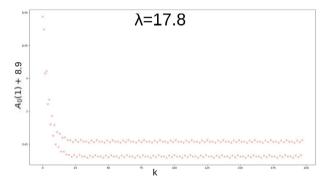


Figure: The values of the potential for constant  $\lambda=17.8$ , as  $\kappa\to\infty$ .

# Periodic points in the context of fixed points

6 Standing questions

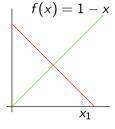


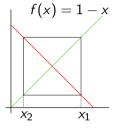
Figure: The periodic points of the function f(x) = 1 - x

Eventhough  $x = \frac{1}{2}$  is a fixed point of f(x), it cannot be found by taking the limit  $\lim_{n\to\infty} f^n(x_1)$ . Indeed,

$$x_{2n} := f^{2n}(x_1) = x_1$$
  
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# Periodic points in the context of fixed points

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# **Dynamic relaxation**

7 Possible fixes

In the 1-D case, no convergence appeared when f'(x) = -1.

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In the 1-D case, no convergence appeared when f'(x) = -1. Recall the relaxing update law

$$A_0^{\kappa+1} = cA_0^{\kappa} + (1-c)(A_{\mathsf{background}} + A_{\mathsf{induced}}^{\kappa}), 0 < c \lesssim 1$$

# **Dynamic relaxation**

#### 7 Possible fixes

In the 1-D case, no convergence appeared when f'(x) = -1. Recall the relaxing update law

$$A_0^{\kappa+1} = cA_0^{\kappa} + (1-c)(A_{\mathsf{background}} + A_{\mathsf{induced}}^{\kappa}), 0 < c \lesssim 1$$

Choose c at every mesh point  $z_n$ , so that convergence is fastest, e.g. The convergence for the function  $f(x) = \frac{1}{2}$  is immediate.

## **Extrapolation**

7 Possible fixes

Since the problem arises from starting too far away from the next self consistent solution, try to predict by extrapolation.

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#### 7 Possible fixes

Since the problem arises from starting too far away from the next self consistent solution, try to predict by extrapolation.

## Point-wise extrapolation

 $A_0^{\lambda_n}(z), A_0^{\lambda_{n-1}}(z)$  the self consistent potentials for  $\lambda_n, \lambda_{n-1}$ , respectively. Guess the next self consistent potential (and use it as a starting  $A_0$ ) by

$$A_0^{\lambda_{n+1}}(z) = rac{A_0^{\lambda_n}(z) - A_0^{\lambda_{n-1}}(z)}{\lambda_n - \lambda_{n-1}}(\lambda_{n+1} - \lambda_n) + A_0^{\lambda_n}(z)$$
 (1)

- ▶ Preliminaries
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- We used the proper prescription for the vacuum polarization  $\langle \rho(z) \rangle$  calculated in [WZ20], to study the effect the backreaction of the Klein-Gordon field has on a background constant electric field.
- We find that the backreaction raises the energy of the modes enough so as to avoid instabilities found when ignoring backreaction.
- However, convergence is still not fast enough to be able to correctly study the whole of the  $\lambda$  parameter space.

- ▶ Preliminaries
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- 1. Redoing the calculations for Neumann boundary conditions,
- 2. Studying the effect of the size of the interval has for the different boundary conditions,
- 3. Redoing the calculations for mixed (Robin) boundary conditions.

- ▶ Preliminaries
- ▶ Results
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```
[AW83] Jan Ambjørn and Stephen Wolfram. "Properties of the vacuum. 2. Electrodynamic". ln: Annals of Physics 147.1 (1983), pp. 33-56. ISSN: 0003-4916. DOI: https://doi.org/10.1016/0003-4916(83)90066-0. URL: https://www.sciencedirect.com/science/article/pii/0003491683900660.
```

[WZ20] Jonathan Wernersson and Jochen Zahn. "Vacuum polarization near boundaries". In: *Phys. Rev. D 103, 016012 (2021)* (Oct. 12, 2020). DOI: 10.1103/PhysRevD.103.016012. arXiv: 2010.05499v2 [hep-th].

# Vacuum polarization in 1+1 spacetime dimensions