Dark Matter scattering in low threshold detectors

Simon Knapen

LBNL



http://dingercatadventures.blogspot.com/2012/08/

SK, J. Kozaczuk, T. Lin: arXiv 2104.12786, 2101.08275, 2011.09496 B. Campbell-Deem, SK, T. Lin, E. Villarama: arXiv 2205.02250



Light dark matter direct detection

What do we need?

Experiment:*

- 1. Ultra-low threshold calorimeters (Spice/HeRALD, ...)
- 2. Single electron detectors (SENSEI, DAMIC, superCDMS, ...)



* Gross oversimplification, but focusing on theory for this talk More on experiment & references in back-up slides

Plot from arXiv 2211.09978

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- 1. Ultra-low threshold calorimeters (Spice/HeRALD, ...)
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Theory:

- 1. Models (constraints are complicated)
- 2. Rate calculations (Collective effects important)
- 3. Background predictions



* Gross oversimplification, but focusing on theory for this talk More on experiment & references in back-up slides

Plot from arXiv 2211.09978

The need for theory



For light DM deBroglie wavelength >> interatomic spacing

"Billiard ball" nuclear recoil not applicable!

Calculations needed



Electronic signals



SPICE prototype detector R. Anthony-Petersen et. al. arXiv 2208.02790

L. Barak et. al. arXiv 2004.11378

SENSEI detector

Moriond EW 2023

Lawrence Berkeley National Lab

Nuclear recoil

$$\omega = \frac{q^2}{2m_N}$$



SK, T. Lin, M. Pyle, K. Zurek: arXiv 1712.06598 S. Griffin, SK, T. Lin, M. Pyle, K. Zurek: arXiv 1807.10291

Nuclear recoil

$$\omega = \frac{q^2}{2m_N}$$



Phonon regime

$$q \ll \sqrt{2m_N\omega}$$

Momentum exchange is a good expansion parameter (phonons are goldstones, similar to chiral perturbation theory)



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Leading Order



 $\mathcal{O}(q), \ \mathcal{O}(q^2) \text{ or } \mathcal{O}(q^4)$

(Depends on DM model & phonon branch)

SK, T. Lin, M. Pyle, K. Zurek: arXiv 1712.06598 S. Griffin, SK, T. Lin, M. Pyle, K. Zurek: arXiv 1807.10291





 $\omega = \frac{q^2}{2m_N}$



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B. Campbell-Deem, SK, T. Lin, E. Villarama: arXiv 2205.02250

(Depends on DM model & phonon branch)



B. Campbell-Deem, P. Cox, SK, T. Lin, T. Melia : 1911.03482

<u>NnLO</u>



<u>N∞LO = nuclear recoil</u>



 $\sim \delta \left(\omega - \frac{q^2}{2m_N} \right)$

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Brian Campbell-Deem (UCSD)



Ethan Villarama (UCSD)

Inclusive all orders result

d: labels atoms (e.g. Ga and As) n: number of phonons

$$\frac{d\sigma}{d^3\mathbf{q}d\omega} \sim \sum_d^{\mathfrak{n}} A_d^2 e^{-2W_d(\mathbf{q})} \sum_n \left(\frac{q^2}{2m_d}\right)^n \frac{1}{n!} \left(\prod_{i=1}^n \int d\omega_i \frac{D_d(\omega_i)}{\omega_i}\right) \delta\left(\sum_j \omega_j - \omega\right).$$
Partial density of states



B. Campbell-Deem, SK, T. Lin, E. Villarama: arXiv 2205.02250

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Partial density of states
$$q \gg \sqrt{2\omega m_{d}}$$

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B. Campbell-Deem, SK, T. Lin, E. Villarama: arXiv 2205.02250

Inclusive all orders result

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B. Campbell-Deem, SK, T. Lin, E. Villarama: arXiv 2205.02250



First calculation for all m_{χ} and arbitrary energy threshold!

(Just an example plot, have results for most relevant models and target materials)

B. Campbell-Deem, SK, T. Lin, E. Villarama: arXiv 2205.02250

Calculations needed

Phonon signals



Electronic signals



SPICE prototype detector R. Anthony-Petersen et. al. arXiv 2208.02790 SENSEI detector L. Barak et. al. arXiv 2004.11378

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Electrons are complicated





- e- are not at rest
- e- are not localized
- e- are not alone
 - → screening

Problem: Calculate wave functions & stick them into matrix element calculation Essig et. al. arXiv 1509.01598



Essig et. al. arXiv 1509.01598

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Essig et. al. arXiv 1509.01598

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Problem: Calculate wave functions & stick them into matrix element calculation Essig et. al. arXiv 1509.01598

Equivalent problem: Calculate rate of energy dissipation in the crystal

SK, J. Kozaczuk, T. Lin: arXiv 2101.08275 Y. Hochberg et. al.: arXiv 2101.08263

Schematic argument

Coulomb potential in a dielectric:

$$H = eQ_{\chi} \int \frac{d^3 \mathbf{k}}{(2\pi)^2} \frac{1}{\epsilon(\mathbf{k},\omega)} \frac{e^{i\mathbf{k}\cdot\mathbf{x}}}{k^2}$$

In QFT language:

$$\sim \sqrt{\frac{1}{\epsilon(\mathbf{k},\omega)}} \frac{1}{k^2}$$

(Non-relativistic limit)

(Exact derivation in the back-up slides)

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Coulomb potential in a dielectric:

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In QFT language:

$$\sim \sqrt{2} \sim \sqrt{\frac{1}{\epsilon(\mathbf{k},\omega)}} \frac{1}{k^2}$$

(Non-relativistic limit)

We are interested in energy dissipation:

$$\sim \sim \lim \left[\frac{-1}{\epsilon(\mathbf{k}, \omega)} \right]$$

"Energy Loss Function" (ELF)

(Exact derivation in the back-up slides)

Calculating the ELF

Simple



(Details in back-up slides)

SK, J. Kozaczuk, T. Lin: arXiv 2101.08275, 2104.12786

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Sophisticated

Calculating the ELF

Simple



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SK, J. Kozaczuk, T. Lin: arXiv 2101.08275, 2104.12786

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Sophisticated

Screening has O(1) effect on integrated rate



Applicable to any mediator that couples to e⁻ density (e.g. scalar mediator and dark photon mediator yield *identical* scattering rate)

SK, J. Kozaczuk, T. Lin: arXiv 2101.08275, 2104.12786

Bonus: The Migdal effect

A hard nuclear recoil can shake some electrons into the conduction band



(Analogous to internal conversion process in collider physics)

Had been calculated for isolated atoms, but not yet for realistic crystals Ibe et. al. arXiv 1707.07258

In a real material, spectator ions must be accounted for!

SK, J. Kozaczuk, T. Lin: arXiv 2011.09496 Liang et.al. : arXiv 2011.13352

Bonus: The Migdal effect

Result:



Bonus: The Migdal effect





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DarkELF



You'd like to quickly calculate DM scattering rates, but don't want to learn Density Functional Theory?

Then DarkELF is the answer for you!

https://github.com/tongylin/DarkELF

- Python 3 package with lots of example Jupyter notebooks
- No dependencies other than numpy, scipy and Vegas
- All DFT results included as look-up tables, no DFT code necessary
- Library of most common materials (Si, Ge, GaAs, diamond, sapphire, etc etc)



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Summary

We calculated:

- DM-phonon scattering to *all orders* in the multiphonon expansion
- DM-electron scattering, *including screening*
- The Migdal effect in semi-conductors

All calculations publicly available in our DarkELF package

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Some future / ongoing work:

- Background processes, e.g. Frenkel pair recombination (ongoing)
- Include core and semi-core electrons in our DFT calculations (ongoing)
- Quenching factor calculation for low energy recoils (very hard)
- Going beyond the isotropic limit (future)

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Extra slides

Electron Detectors

Sensitivity to single e- excitations has already been demonstrated



SENSEI already has 50g-day exposure in shallow underground site

Phonon Detectors

SPICE conceptual design



Qualitative progress in background mitigation





SPICE / HeRALD



detector hanging on aluminium wire bonds

R. Anthony-Petersen et. al. arXiv 2208.02790

Approximations

1. Incoherent approximation



Interference is critical (e.g. Bragg diffraction)



Interference can be neglected

Approximations

1. Incoherent approximation



Interference is critical (e.g. Bragg diffraction)



Interference can be neglected

2. Anharmonic approximation



3. Isotropic approximation

Multiphonon results

<u>Numerics</u>



B. Campbell-Deem, SK, T. Lin, E. Villarama: arXiv 2205.02250

Need to use linear response theory, essentially non-relativistic QFT

Susceptibility: how does the crystal respond to a density perturbation?

$$\chi(\omega, \mathbf{k}) = -\frac{i}{V} \int_{0}^{\infty} dt \, e^{i\omega t} \langle [n_{\mathbf{k}}(t), n_{-\mathbf{k}}(0)] \rangle$$

$$\downarrow$$
Crystal volume
Electron number density operator

This is the non-relativistic, retarded Green's function (fully dressed)

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This is the non-relativistic, retarded Green's function (fully dressed)

Now we use the fluctuation-dissipation theorem

$$\operatorname{Im}\chi(\omega,\mathbf{k}) = -\frac{1}{2}(1 - e^{-\beta\omega})S(\omega,\mathbf{k}) \qquad \beta \equiv \frac{1}{k_B T}$$

With the dynamical structure factor defined as

$$S(\omega, \mathbf{k}) \equiv \frac{2\pi}{V} \sum_{i, f} \frac{e^{-\beta E_i}}{Z} |\langle f | n_{-\mathbf{k}} | i \rangle|^2 \delta(\omega + E_i - E_f)$$

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Fermi's golden rule

Now consider the response to an external electromagnetic perturbation. The induced electron number density is

$$\begin{split} \langle \delta n(\mathbf{k},\omega) \rangle &= \langle n(\mathbf{k},\omega) H_{coul} \rangle \\ &= -\frac{e}{k^2} \chi(\mathbf{k},\omega) \rho_{ext}(\mathbf{k},\omega) \end{split} \qquad \text{with} \qquad H_{coul} = -e \int \frac{d^3 \mathbf{k}}{(2\pi)^2} \frac{e^{i\mathbf{k}\cdot\mathbf{x}}}{k^2} n(-\mathbf{k},\omega) \rho_{ext}(\mathbf{k},\omega) \end{split}$$

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$$= -\frac{e}{k^2} \chi(\mathbf{k},\omega) \rho_{ext}(\mathbf{k},\omega)$$

Using Maxwell's equations

$$i\mathbf{k} \cdot \mathbf{D}(\mathbf{k}, \omega) = 4\pi \rho_{ext}(\mathbf{k}, \omega)$$
$$i\mathbf{k} \cdot \mathbf{E}(\mathbf{k}, \omega) = 4\pi \rho_{ext}(\mathbf{k}, \omega) - 4\pi e \langle \delta n(\mathbf{k}, \omega) \rangle$$

with
$$D(\mathbf{k},\omega) = \epsilon(\mathbf{k},\omega)\mathbf{E}(\mathbf{k},\omega)$$

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$$\mathbf{D}(\mathbf{k},\omega) = \epsilon(\mathbf{k},\omega)\mathbf{E}(\mathbf{k},\omega)$$

Which results in the relation

$$\frac{1}{\epsilon(\omega, \mathbf{k})} = 1 + \frac{4\pi\alpha_{em}}{k^2}\chi(\omega, \mathbf{k}),$$

Now plugging this into the fluctuation-dissipation theorem

$$S(\omega, \mathbf{k}) = \frac{k^2}{2\pi\alpha_{em}} \frac{1}{1 - e^{-\beta\omega}} \operatorname{Im}\left[\frac{-1}{\epsilon(\omega, \mathbf{k})}\right]$$

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Energy Loss Function (ELF)

DM-electron scattering rate

Full formula



Advantages of using the ELF:

- Screening included automatically
- ELF has been measured and calculated extensively in the condensed matter literature

Applicable to any mediator that couples to e⁻ density (e.g. scalar mediator and dark photon mediator yield *identical* scattering rate)

SK, J. Kozaczuk, T. Lin: arXiv 2101.08275, 2104.12786

Lindhard model





Homogenous, free electron gas:

$$\epsilon_{\rm Lin}(\omega,k) = 1 + \frac{3\omega_p^2}{k^2 v_F^2} \lim_{\eta \to 0} \left[f\left(\frac{\omega + i\eta}{k v_F}, \frac{k}{2m_e v_F}\right) \right]$$

with

$$v_F = \left(\frac{3\pi\omega_p^2}{4\alpha m_e^2}\right)^{1/3} \frac{1}{2} \text{Plasmon frequency}$$

$$f(u, z) = \frac{1}{2} + \frac{1}{8z} \left[g(z - u) + g(z + u)\right]$$

$$g(x) = (1 - x^2) \log\left(\frac{1 + x}{1 - x}\right)$$

Features:

- Pauli blocking
- e-h pair continuum
- Plasmon width
- Low k region
- Bandgap

Mermin model

 $\begin{bmatrix} \mathbf{Mermin} \\ \mathbf{$

Homogenous, free electron gas with dissipation (Γ)

$$\epsilon_{\mathrm{Mer}}(\omega,k) = 1 + \frac{(1+i\frac{\Gamma}{\omega})(\epsilon_{\mathrm{Lin}}(\omega+i\Gamma,k)-1)}{1+(i\frac{\Gamma}{\omega})\frac{\epsilon_{\mathrm{Lin}}(\omega+i\Gamma,k)-1}{\epsilon_{\mathrm{Lin}}(0,k)-1}}.$$

Fit a linear combination of Mermin oscillators to optical data:

$$=\sum_{i} A_{i}(k) \operatorname{Im}\left[\frac{-1}{\epsilon_{\operatorname{Mer}}(\omega,k;\omega_{p,i},\Gamma_{i})}\right]$$



M. Vos, P. Grande: chapidif package Data from Y. Sun et. al. Chinese Journal of Chemical Physics 9, 663 (2016)

Features:

 $\epsilon(\omega,k)$

- Pauli blocking
- e-h pair continuum
- Plasmon width
- **M** Low k region
- Bandgap

GPAW method



Compute the ELF from first principles with timedependent Density Functional Theory methods (TD-DFT)

Puts atoms on periodic lattice and model interacting e- as non-interacting e- + effective external potential (Kohn-Sham method)

Inner shell e- are treated as part of the ion (frozen core approximation)



GPAW: https://wiki.fysik.dtu.dk/gpaw/

Features:

- Pauli blocking
- e-h pair continuum
- Plasmon width
- **M** Low k region
- ☑ Bandgap

Differential rate



Mermin & GPAW in very good agreement except:

- Single ionization e- region (background dominated)
- High energy region (subdominant)

(Agreement is less good in massive mediator case; work in progress)

Differential rate



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Integrated rate: Other materials

Using the *Mermin* method we can easily scan over many possible targets:



So far only GPAW results for Ge and Si, other materials are work in progress