

Modelling dark matter-electron interactions in materials

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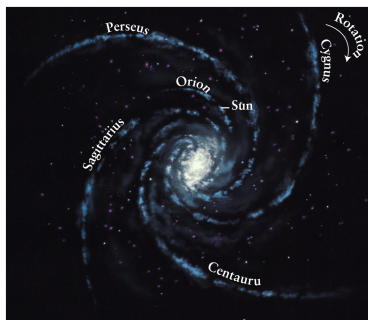
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*Knut och Alice
Wallenbergs
Stiftelse*

Why?

Basic principles of DM direct detection

- Face-on view of our galaxy:



- The sun's orbital motion induces a flux of DM particles through our planet

- When a DM particle crosses a terrestrial detector can deposit energy by interaction with its constituents
- DM direct detection experiments search for such *rare, energy depositions*
- Expected rate of DM “signal events”

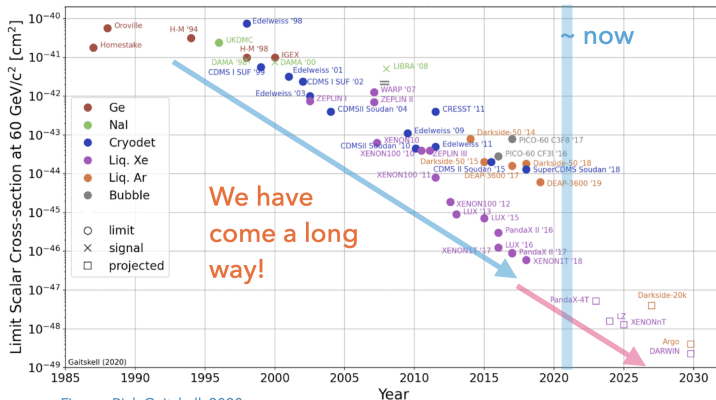
$$d\mathcal{R} = \frac{\rho_\chi}{m_\chi} \int d\mathbf{v} |\mathbf{v}| f_\chi(\mathbf{v} + \mathbf{v}_\oplus) d\sigma$$

Astrophysics

Particle physics,
Nuclear physics,
Solid state physics

History of DM direct detection

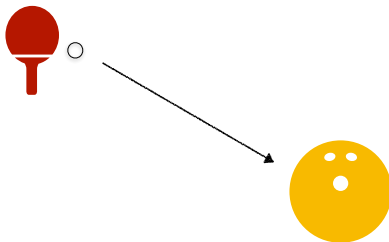
- Exclusion limits and projected sensitivities on the strength of DM-nucleon interactions



- Why has DM so far escaped a direct detection?

A possible explanation for the lack of DM detection

- A simple explanation for the lack of DM detection is that it is lighter than nucleons (<1 GeV), and therefore too light to cause an observable nuclear recoil



- Not unlike a light pingpong ball being too light to move a heavy bowling ball ...
- If this is true, DM should be searched for in the recoils of a lighter target:
the electron

How?

- Leveraging on previous results on the scattering of DM by nucleons bound in nuclei:

J. Fan, M. Reece and L. T. Wang,
JCAP **11** (2010), 042

A. L. Fitzpatrick, W. Haxton, E. Katz, N. Lubbers and Y. Xu,
JCAP **02** (2013), 004

- We developed an effective theory to model DM-electron interactions in materials:

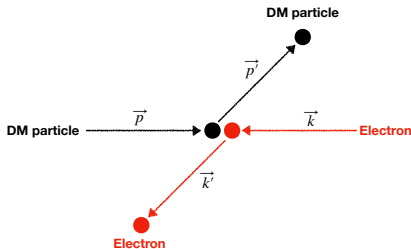
R. Catena, D. Cole, T. Emken, M. Matas, N. Spaldin, W. Tarantino and E. Urdshals,
JCAP **03** (2023), 052

R. Catena, T. Emken, M. Matas, N. A. Spaldin and E. Urdshals,
Phys. Rev. Res. **3** (2021) no.3, 033149

R. Catena, T. Emken, N. A. Spaldin and W. Tarantino,
Phys. Rev. Res. **2** (2020) no.3, 033195

An effective theory approach / assumptions

- Consider the scattering of a DM particle of mass m_χ by a free electron of mass m_e ,



- In the non-relativistic limit, this process is characterised by a double separation of scales:

$$|\mathbf{q}|/m_e \ll 1,$$

$$|v| \ll 1,$$

$$\mathbf{q} = \mathbf{p} - \mathbf{p}'$$

$$v = p/m_\chi$$

- Its amplitude $\mathcal{M}_{\chi e}$ is invariant under Galilean transformations, translations and rotations

An effective theory approach / amplitude

- What is the predicted form for $\mathcal{M}_{\chi e}$ in our non-relativistic effective theory?
We find:

$$\mathcal{M}_{\chi e}(\mathbf{q}, \mathbf{v}^\perp) = \sum_i c_i \langle \mathcal{O}_i \rangle$$

Diagram illustrating the components of the effective theory amplitude $\mathcal{M}_{\chi e}(\mathbf{q}, \mathbf{v}^\perp)$:

- Matrix elements**: Points to $\langle \mathcal{O}_i \rangle$.
- Rotationally invariant operators in the DM-electron spin space**: Points to \mathcal{O}_i .
- Unknown coupling constants**: Points to c_i .
- Sum over operator type**: Points to the summation index i .
- Out of the four momenta \vec{p} , \vec{p}' , \vec{k} and \vec{k}' only two are independent: \vec{q} and \vec{v}^\perp** : Points to the arguments \mathbf{q} and \mathbf{v}^\perp of the amplitude.

Examples of \mathcal{O}_i operators:

$$\mathcal{O}_1 = \mathbb{1}_\chi \mathbb{1}_e, \quad \mathcal{O}_4 = \mathbf{S}_\chi \cdot \mathbf{S}_e, \quad \mathcal{O}_7 = \mathbf{S}_\chi \cdot \mathbf{v}^\perp, \quad \mathcal{O}_{11} = i\mathbf{S}_\chi \cdot \mathbf{q}/m_e, \quad \dots$$

Electron wave function overlap integrals

- We use $\mathcal{M}_{\chi e}$ to calculate the rate of transitions from the electronic state "1" to "2"

$$d\mathcal{R}_{1\rightarrow 2} \propto \left| \int \frac{d^3\mathbf{k}}{(2\pi)^3} \psi_2^*(\mathbf{k} + \mathbf{q}) \mathcal{M}_{\chi e}(\mathbf{v}^\perp, \mathbf{q}) \psi_1(\mathbf{k}) \right|^2$$

where $\mathbf{v}^\perp = \mathbf{v} - \mathbf{q}/(2\mu_{\chi e}) - \mathbf{k}/m_e$

- For $\mathcal{M}_{\chi e}(\mathbf{v}^\perp, \mathbf{q}) \neq \mathcal{M}_{\chi e}(\mathbf{q})$, $\mathcal{M}_{\chi e}$ cannot be moved outside the integral sign
- $\mathcal{M}_{\chi e}$ depends on \mathbf{v}^\perp , and thus on \mathbf{k} in the case of anapole and magnetic dipole interactions
- It also depends on \mathbf{v}^\perp in a number of simplified models with vector mediators

R. Catena, D. Cole, T. Emken, M. Matas, N. Spaldin, W. Tarantino and E. Urdshals,
JCAP **03** (2023), 052

DM-induced electronic transition rate

- Our total rate formula:

$$\mathcal{R}_{\text{theory}} = \frac{n_{\chi}}{128\pi m_{\chi}^2 m_e^2} \int d(\ln \Delta E) \int dq q \hat{n}(q, \Delta E) \sum_{\ell=1}^r \Re \left[\mathcal{R}_{\ell}^*(q, v) \overline{\mathcal{W}}_{\ell}(q, \Delta E) \right]$$

Velocity integral

Sum over up to
 $r=7$
response functions

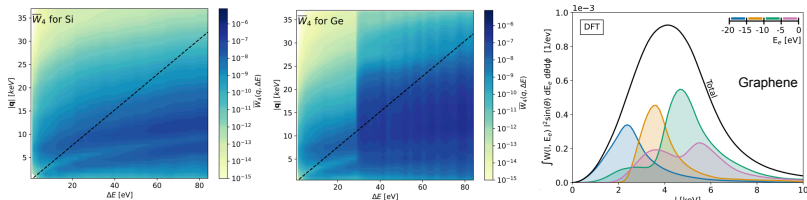
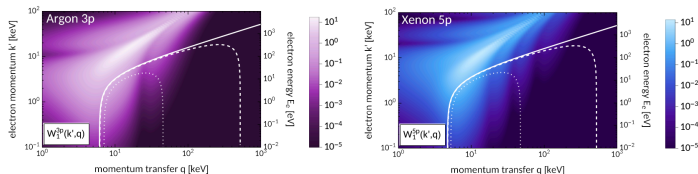
Free electron physics:
couplings and kinematics

Response function:
electron wave function
overlap integral

- It predicts a factorisation between the **free electron physics** encoded in \mathcal{R}_{ℓ} and the **material physics** encoded in the response functions \mathcal{W}_{ℓ}

Applications

Response function evaluation

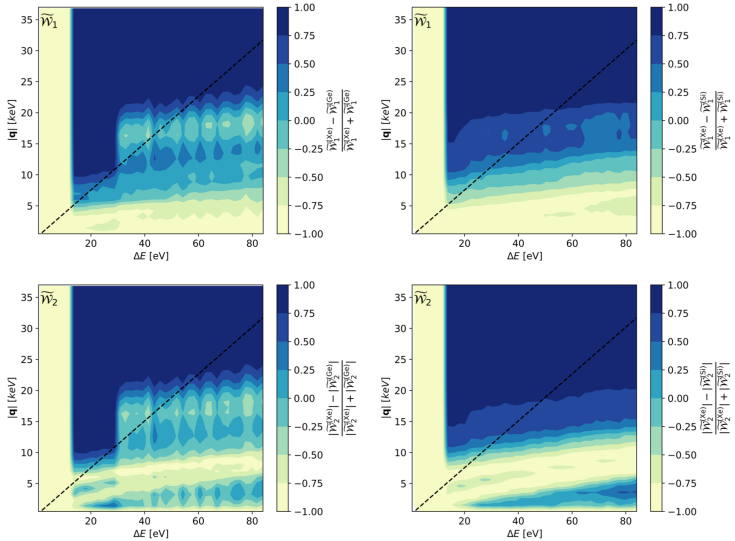


Xenon and Argon: R. Catena, T. Emken, N. A. Spaldin and W. Tarantino,
Phys. Rev. Res. 2 (2020) no.3, 033195

Germanium and Silicon: R. Catena, T. Emken, M. Matas, N. A. Spaldin and E. Urdshals,
Phys. Rev. Res. 3 (2021) no.3, 033149

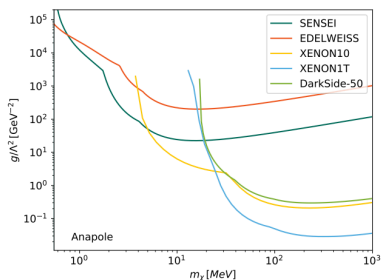
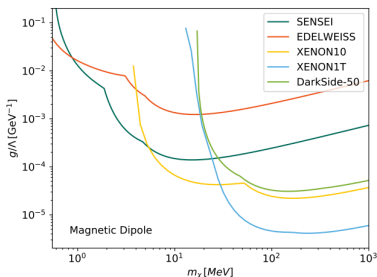
Graphene: R. Catena, T. Emken, M. Matas, N. A. Spaldin and E. Urdshals,
arXiv:2303.15497 [hep-ph]

Response function comparison



Exclusion limits

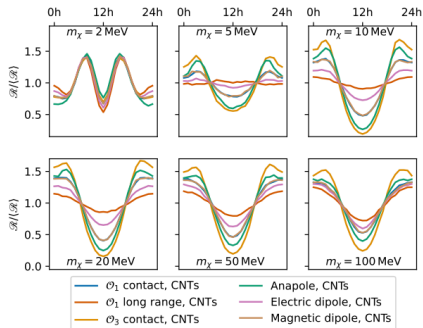
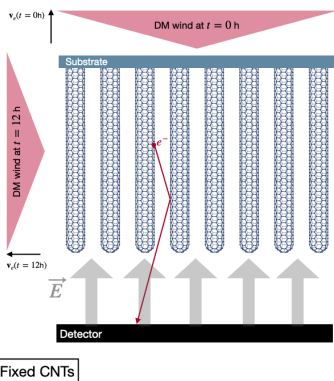
- The formalism allows us to perform calculations within models, e.g. anapole and magnetic dipole DM, which were not tractable before (lacking the required \mathcal{W}_ℓ 's)



R. Catena, T. Emken, M. Matas, N. A. Spaldin and E. Urdshals,
arXiv:2303.15509 [hep-ph]

General predictions for new detector materials

- It also allows us to make predictions for new direct detection materials, e.g. graphene, for general DM-electron interactions



Summary

- We developed a non-relativistic effective theory to model general DM-electron interactions in materials
- Our formalism predicts a factorisation between the **free electron physics** and the **material physics** encoded in the response functions \mathcal{W}_ℓ
- It allows us to perform calculations within models, e.g. anapole and magnetic dipole DM, which were not tractable before (lacking the required \mathcal{W}_ℓ 's)
- Furthermore, it enables us to assess the potential of new direct detection materials (graphene) for a general form of the underlying DM-electron interaction