


Atomic and molecular transitions for neutrino and LDM detection

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in collaboration with Martin Bauer and Javier Perez-Soler

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 **QuMAP**
Quantum Materials and
Astroparticle Physics



UNIVERSITY OF
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Contents

- Motivation: CvB and DM
- Generalised atomic transitions
- Pair absorption
- The future: molecules and ML

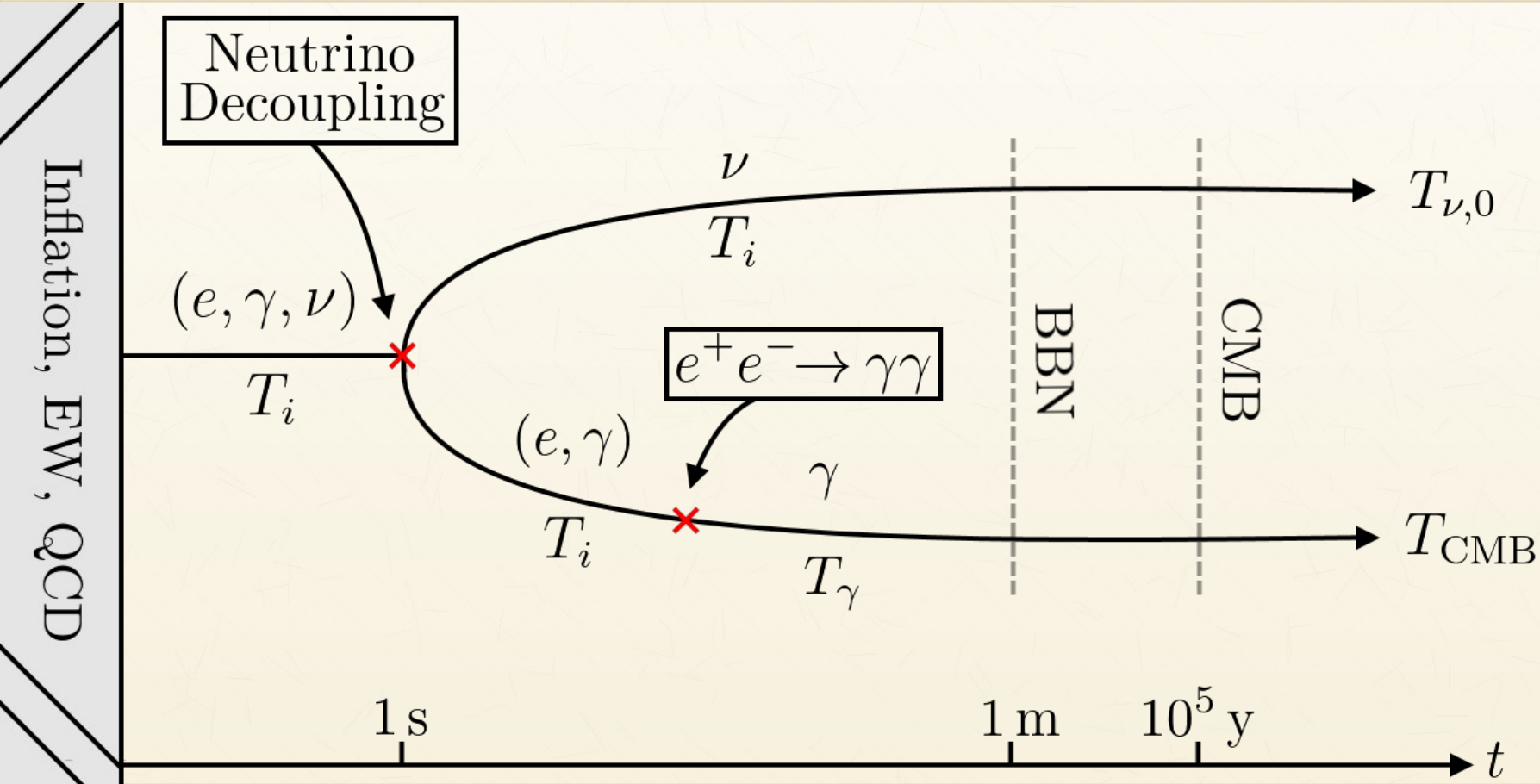


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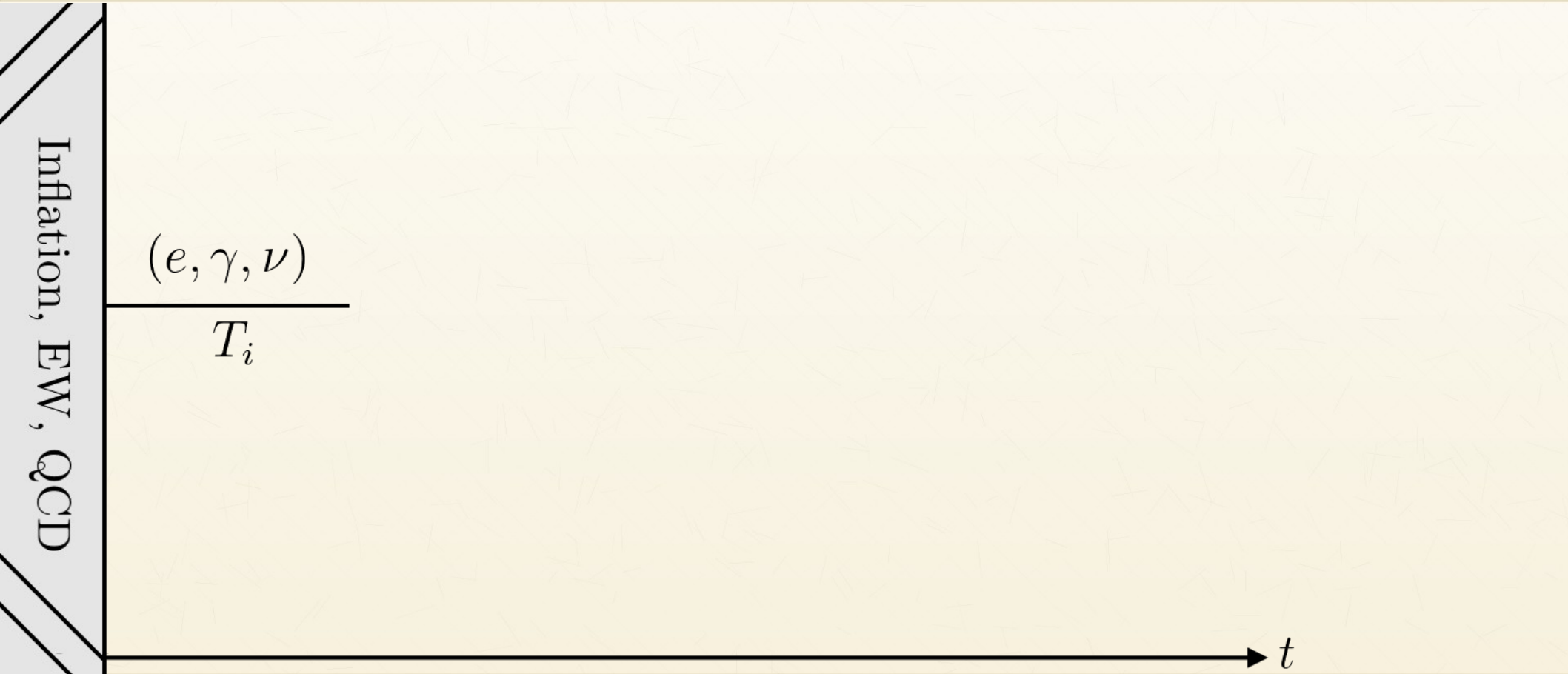
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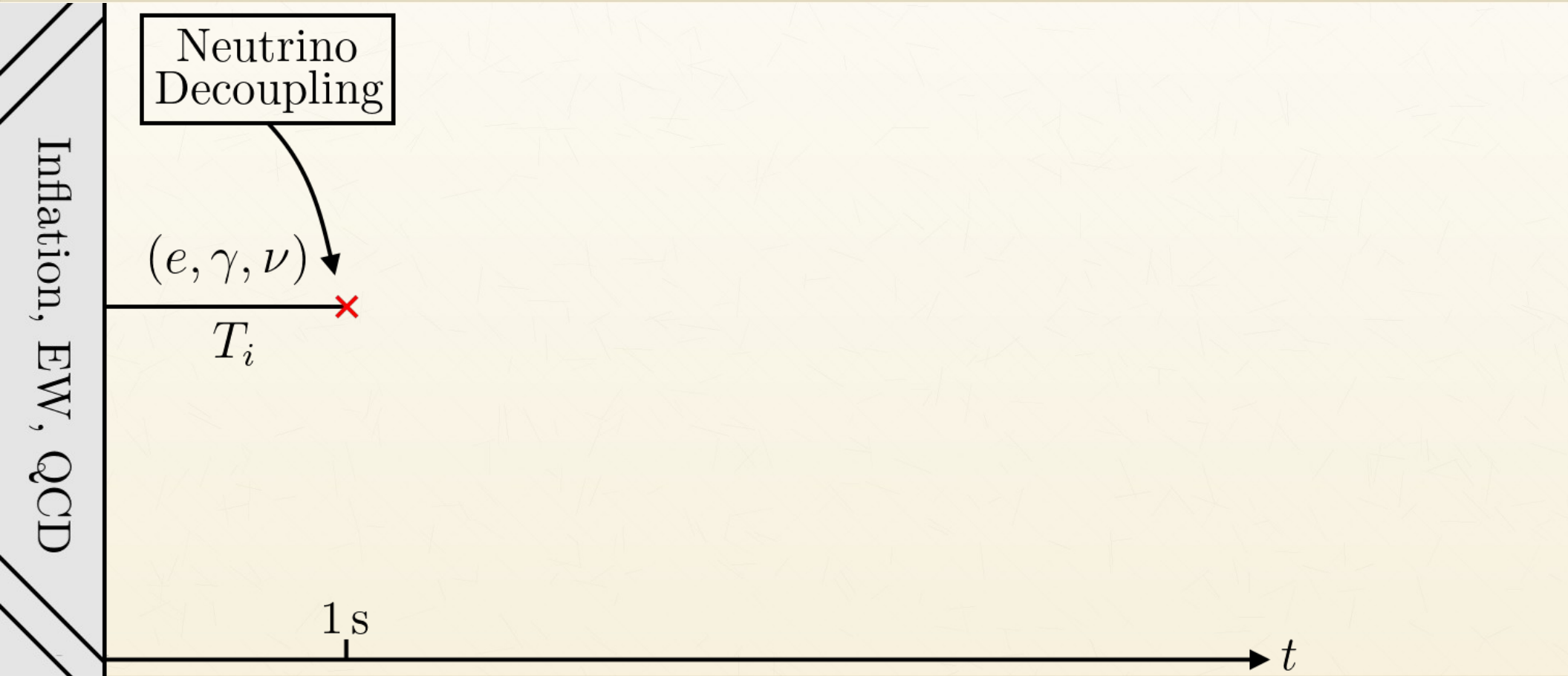
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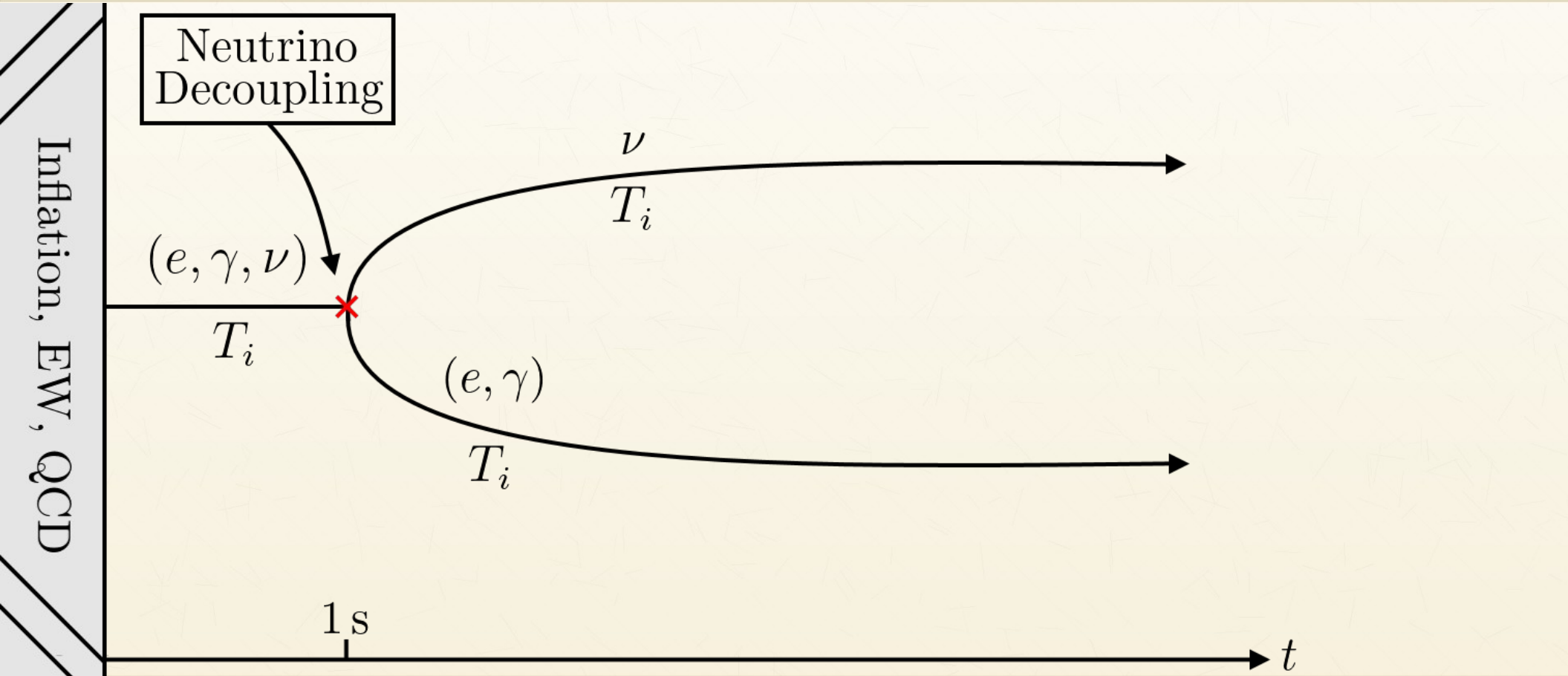
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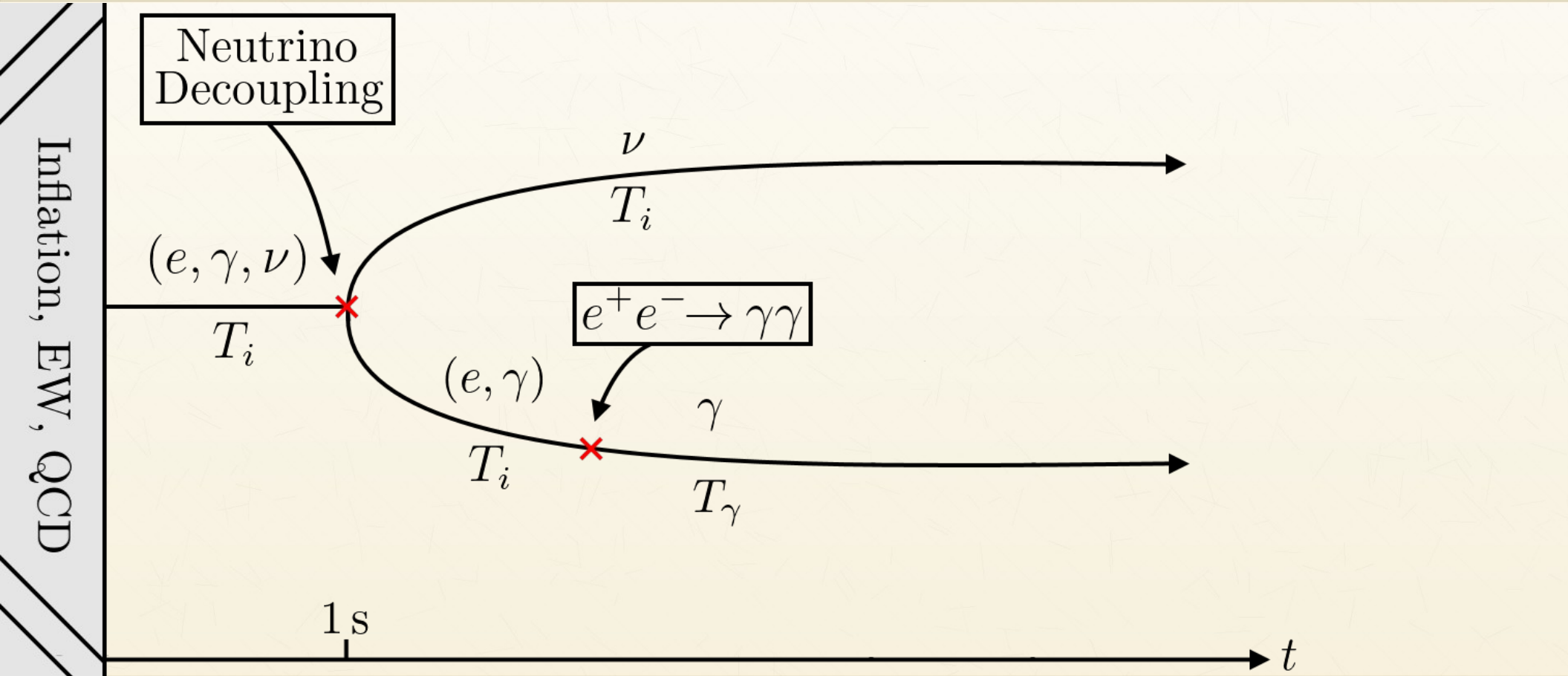
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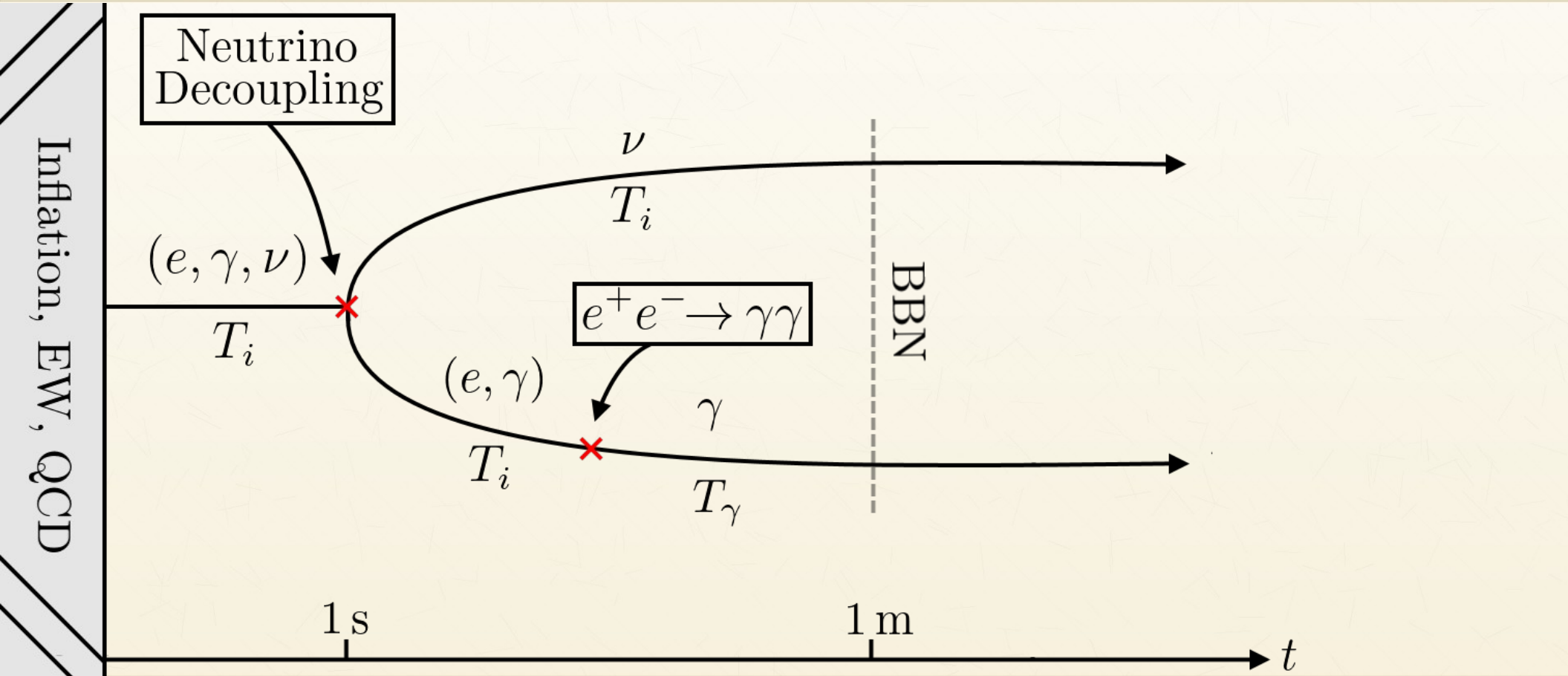
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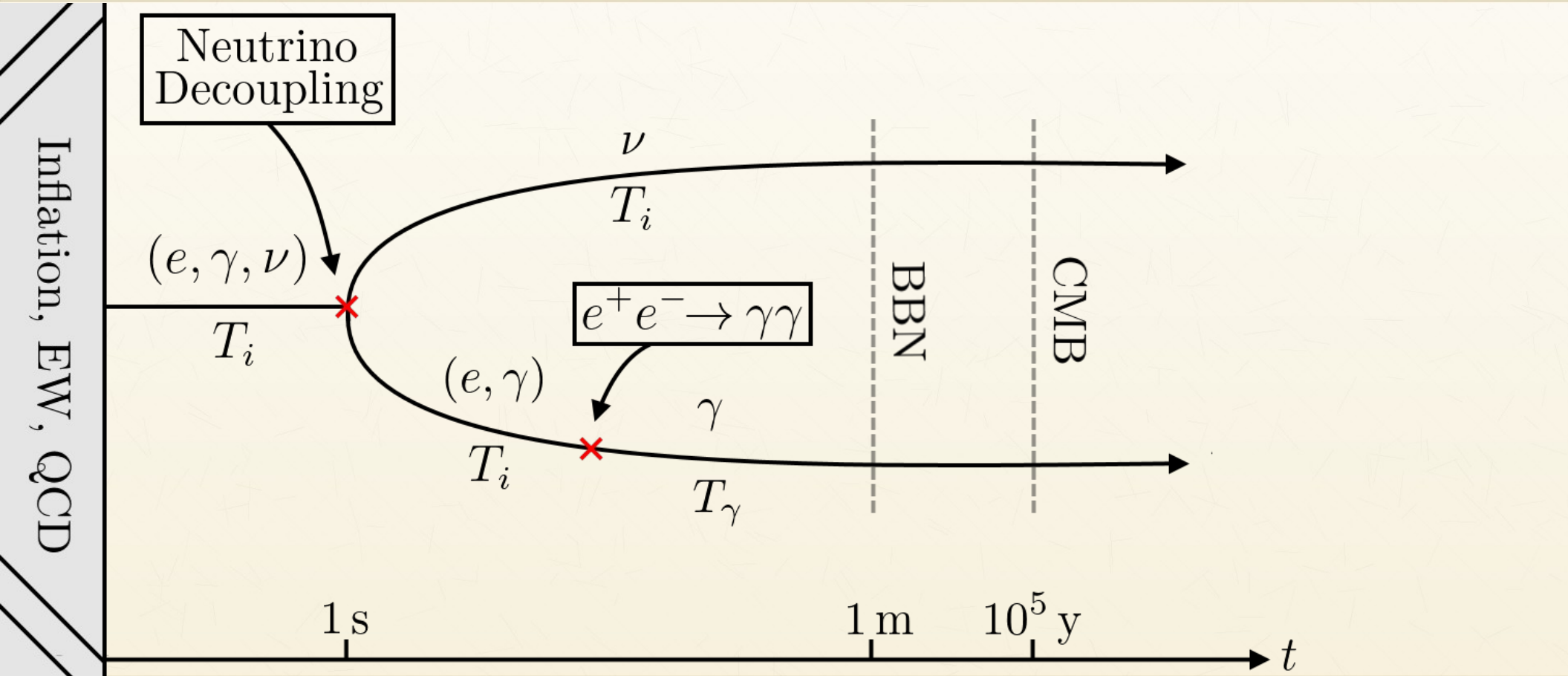
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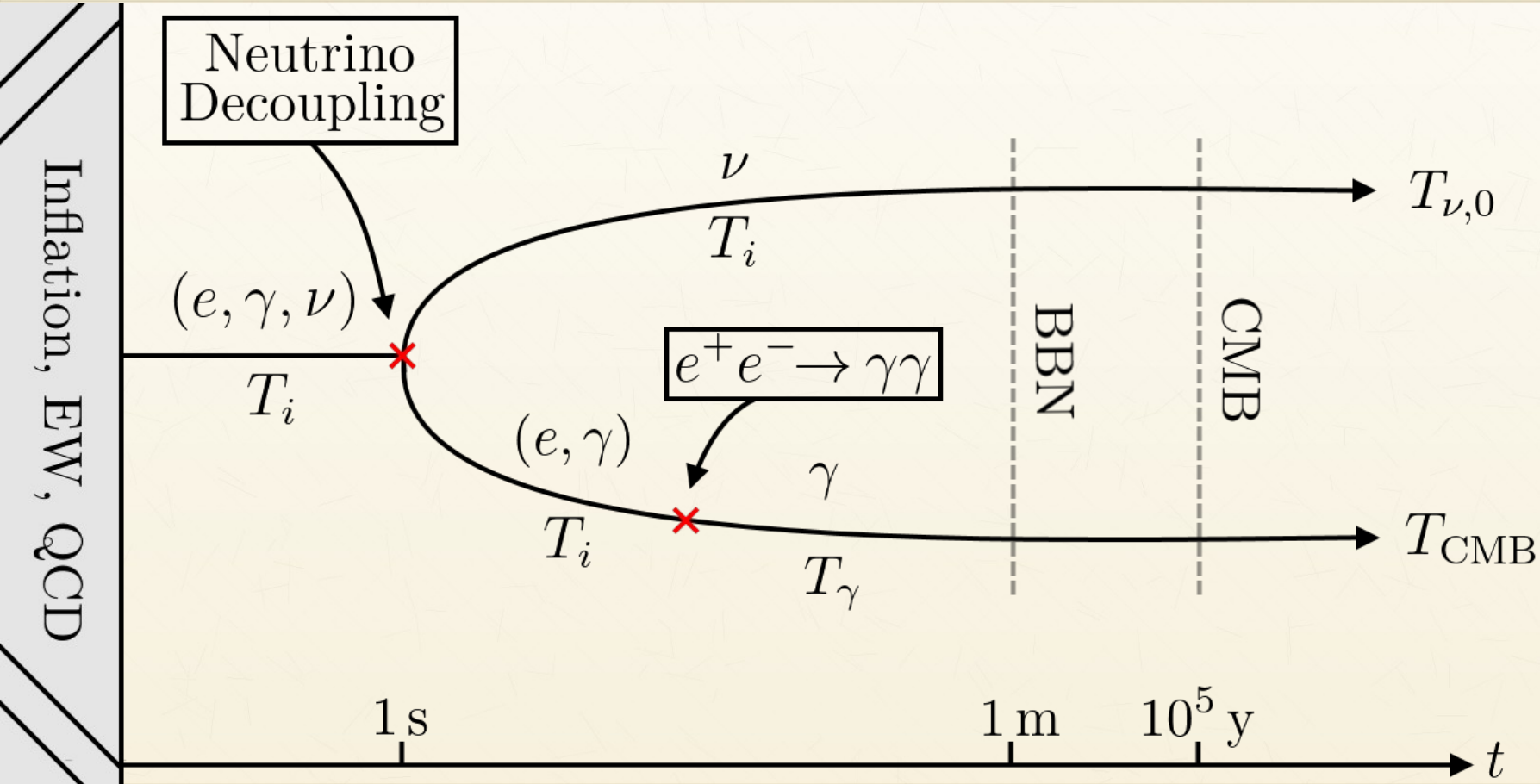
Motivation: The CvB



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Motivation: The CvB

- Redshifted to temperature:

$$T_{\nu,0} = \left(\frac{4}{11} \right)^{\frac{1}{3}} T_{\text{CMB}}$$

Motivation: The CvB

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$$T_{\nu,0} = 0.168 \text{ meV}$$

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- Following a Fermi-Dirac distribution with:

Motivation: The CvB

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$$T_{\nu,0} = 0.168 \text{ meV}$$

- Following a Fermi-Dirac distribution with:

$$n_{\nu} = 56 \text{ cm}^{-3}$$

Motivation: The CvB

- Why do we care?

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- Why do we care?



Cosmology

Motivation: The CvB

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Neutrino physics

Motivation: The CvB

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- Verify Λ CDM

Neutrino physics

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- NR neutrinos
- Dirac vs Majorana

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Neutrino physics

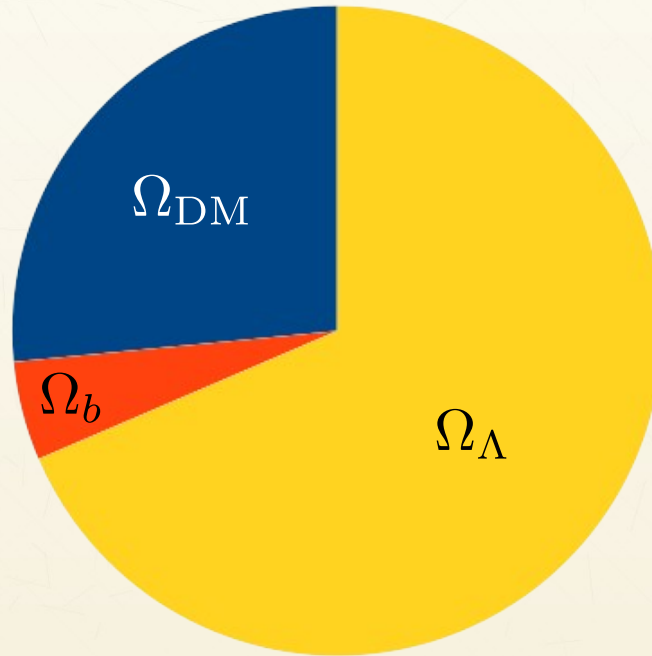
- NR neutrinos
- Dirac vs Majorana
- Absolute neutrino mass

Motivation: DM

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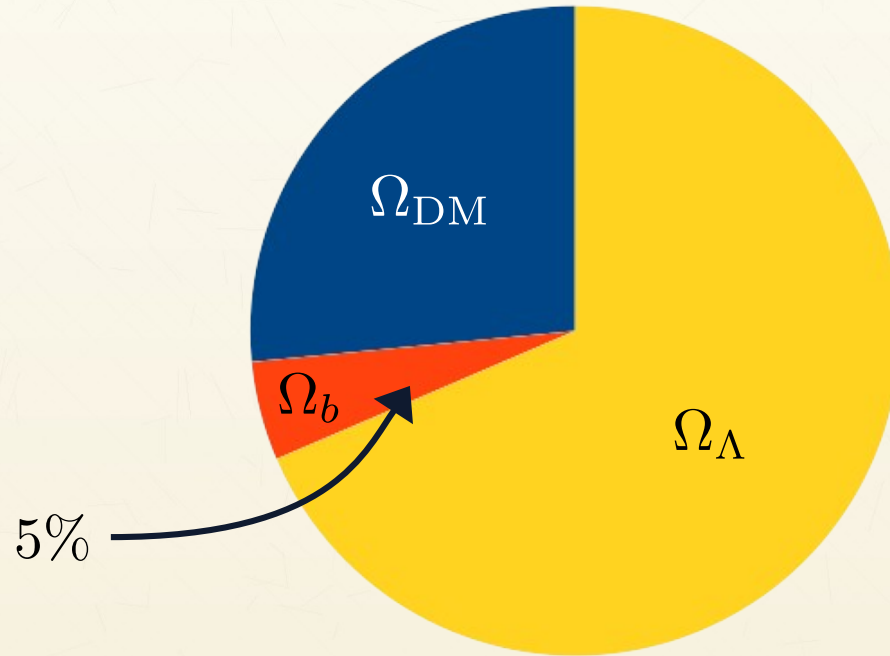
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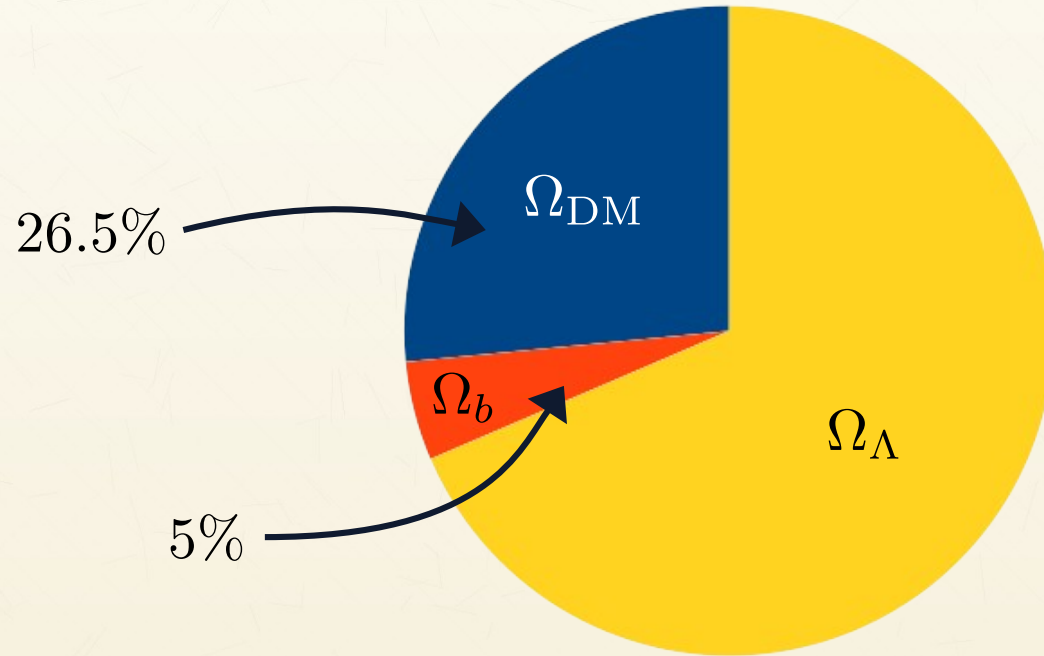
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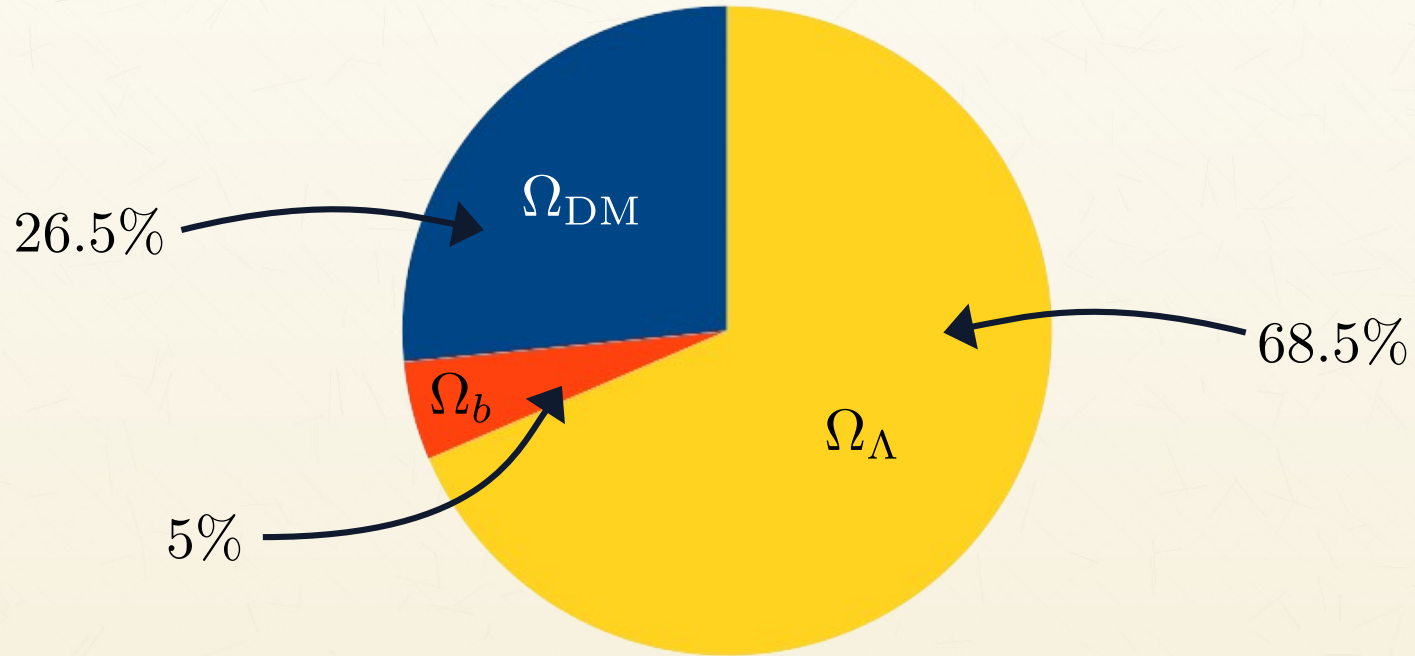
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Motivation: Detection

- How do we detect light, cold relics?

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Elastic

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$$R = \frac{n}{\Lambda^a} E_k^b$$

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Density

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Density

UV scale

Inelastic

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Diagram illustrating the relationship between variables in the Elastic detection equation:

- n is linked to Density.
- Λ^a is linked to UV scale.
- E_k^b is linked to Kinetic energy.

Inelastic

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Density

UV scale

Kinetic energy

Inelastic

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UV scale → Λ^a Kinetic energy → E_k^b Density → n

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$$R = \frac{n}{\Lambda^a} E_{\text{thresh}}^b$$

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Threshold energy → E_{thresh}^b

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$$R_\nu = G_F^2 n_\nu E_{k,\nu}^2$$

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Fermi
constant



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$$R_\nu = G_F^2 n_\nu m_\nu^2$$

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- How do we detect light, cold relics?

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$$R_\nu = 3 \cdot 10^{-22} \text{ y}^{-1}$$

at $m_\nu = 10 \text{ meV}$

Inelastic

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$$R_\nu = 3 \cdot 10^{-22} \text{ y}^{-1}$$

at $m_\nu = 10 \text{ meV}$

Inelastic

- Threshold

$$R_\nu = 2 \cdot 10^{-16} \text{ y}^{-1}$$

at $m_\nu = 10 \text{ meV}$

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$$R_\varphi = 3 \cdot 10^{-20} \text{ y}^{-1}$$

at $m_\varphi = 10 \text{ meV}$

Inelastic

- Threshold

$$R_\varphi = 1 \cdot 10^{-7} \text{ y}^{-1}$$

at $m_\varphi = 10 \text{ meV}$

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$$R_\varphi = 1 \cdot 10^{-7} \text{ y}^{-1}$$

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Focus of today's talk



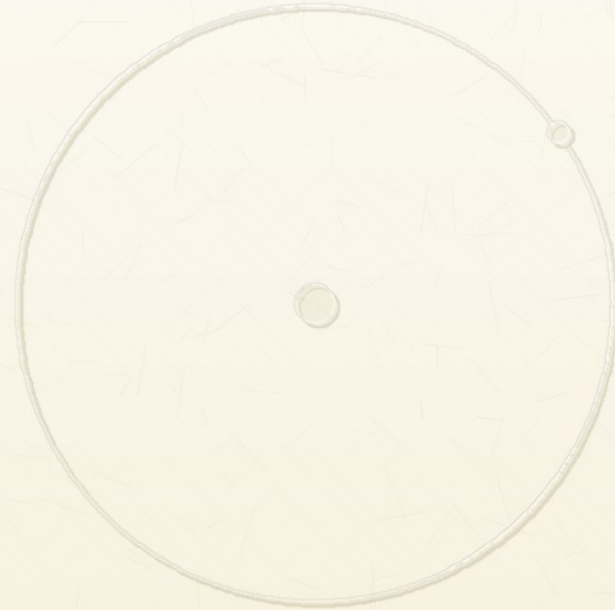
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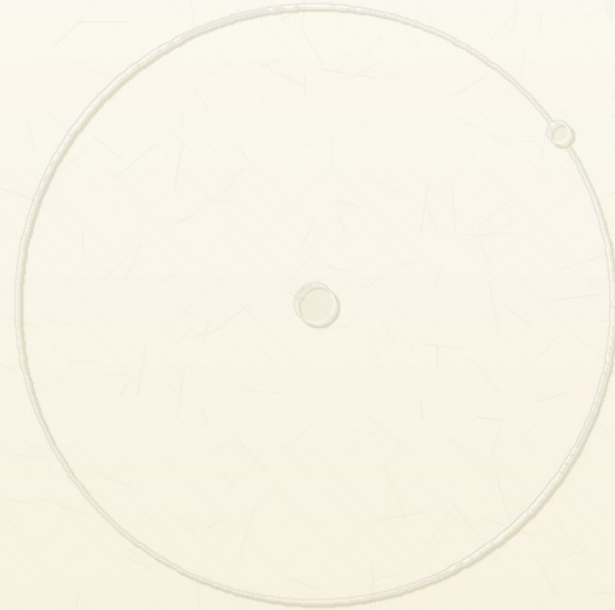
Generalised atomic transitions

- We need a system with:



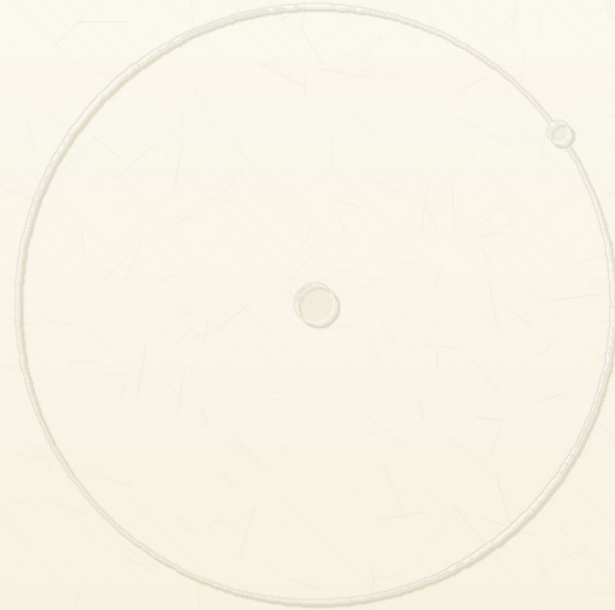
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- We need a system with:
 - Tiny threshold, $\mathcal{O}(\text{eV})$



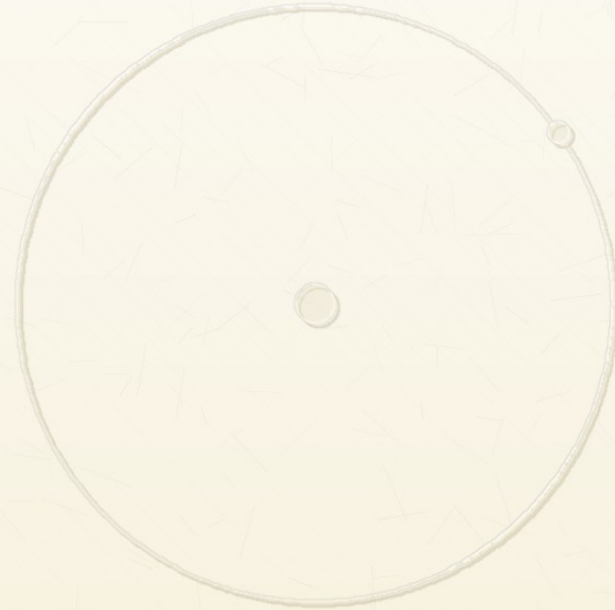
Generalised atomic transitions

- We need a system with:
 - Tiny threshold, $\mathcal{O}(\text{eV})$
 - Clean signal



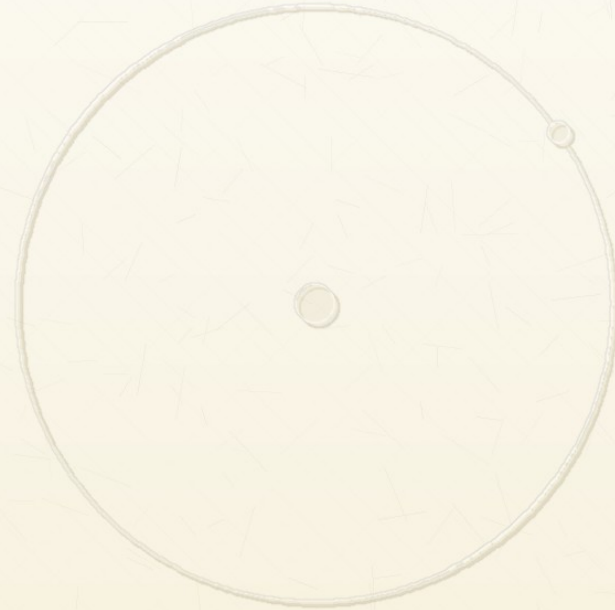
Generalised atomic transitions

- We need a system with:
 - Tiny threshold, $\mathcal{O}(\text{eV})$
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- Atomic systems!



Generalised atomic transitions

- We need a system with:
 - Tiny threshold, $\mathcal{O}(\text{eV})$
 - Clean signal
- Atomic systems!
- But what are we looking for?



Generalised atomic transitions

- First option, new long range forces:

Generalised atomic transitions

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$$U(r) = \frac{Z\alpha_{\text{EM}}}{r}$$

Generalised atomic transitions

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$$U(r) = \frac{Z\alpha_{\text{EM}}}{r} \longrightarrow U(r) = \frac{Z\alpha_{\text{EM}}}{r} + \beta\mathcal{R}(r)$$

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- Where, in general:

Generalised atomic transitions

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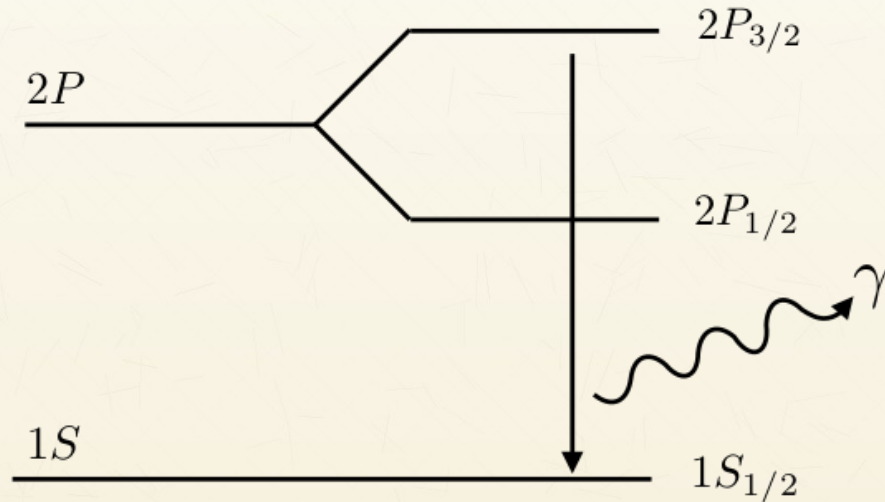
$$\mathcal{R}(r) = \frac{1}{r^n} e^{-mr}$$

Generalised atomic transitions

- Second option, new transitions:

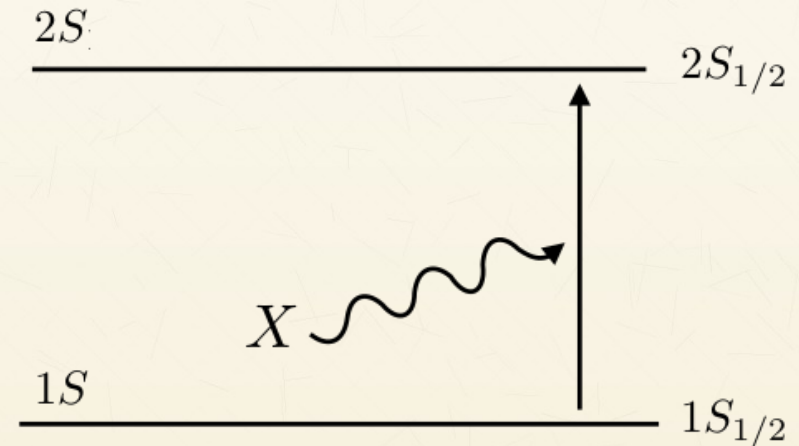
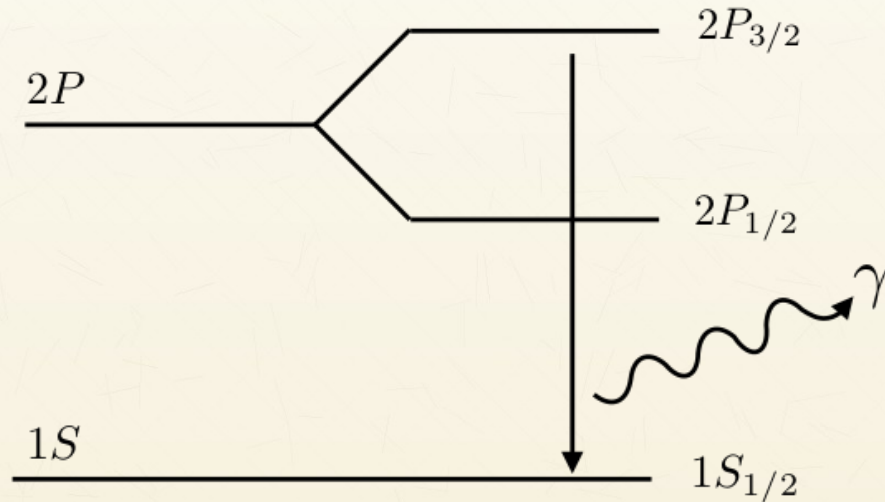
Generalised atomic transitions

- Second option, new transitions:



Generalised atomic transitions

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Generalised atomic transitions

- How do we compute these rates?

$$d\Gamma = 2\pi |\mathcal{M}_{fi}|^2 \delta(\dots) d\rho$$

Generalised atomic transitions

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Amplitude



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Amplitude

Energy conservation

Generalised atomic transitions

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Amplitude

Energy conservation

Phase space

Generalised atomic transitions

- How do we compute these rates?

$$d\Gamma = 2\pi |\mathcal{M}_{fi}|^2 \delta(\dots) d\rho$$

Amplitude

Phase space

Energy conservation

$$\mathcal{M}_{fi} = \langle f | H_{\text{int}}(\psi_e, A_\mu, \varphi, \dots) | i \rangle$$

Generalised atomic transitions

- We work with relativistic wavefunctions, but why?

Generalised atomic transitions

- We work with relativistic wavefunctions, but why?
- Because it's easier...lets us use the language of QFT:

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$$A_\mu \bar{\psi}_e \gamma^\mu \psi_e \not\Rightarrow \vec{p}_e \cdot \vec{A}$$

Generalised atomic transitions

- We work with relativistic wavefunctions, but why?
- Because it's easier...lets us use the language of QFT:

$$A_\mu \bar{\psi}_e \gamma^\mu \psi_e \not\Rightarrow \vec{p}_e \cdot \vec{A}$$

- More precise at large Z , and more transitions

Generalised atomic transitions

- What's the catch?

Generalised atomic transitions

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- Very computationally expensive:

Generalised atomic transitions

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 - Four component wavefunctions!

Generalised atomic transitions

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 - Lorentz indices!

Generalised atomic transitions

- What's the catch?
- Very computationally expensive:
 - Four component wavefunctions!
 - Lorentz indices!
- Classical labelling (e.g. M1, E1, ...) obscured

Generalised atomic transitions

- Let's count the terms: $N_{\text{terms}} = 1$

Generalised atomic transitions

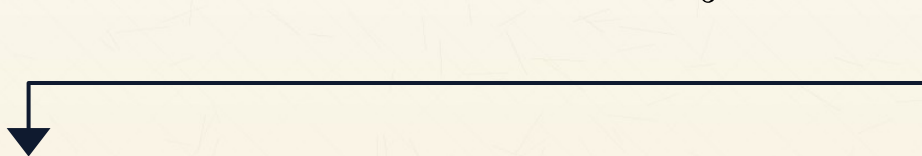
- Let's count the terms: $N_{\text{terms}} = 1$

$$\mathcal{M}_{fi} = \langle f | \int d^3x \mathcal{H}_{\text{int}}(x) | i \rangle$$

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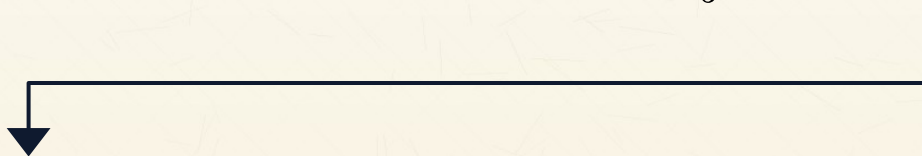
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$$\mathcal{H}_{\text{int}} = \sum_L (\bar{\psi}_e \Gamma_L^{\{\mu\}} \psi_e) \mathcal{O}_{L, \{\mu\}}$$

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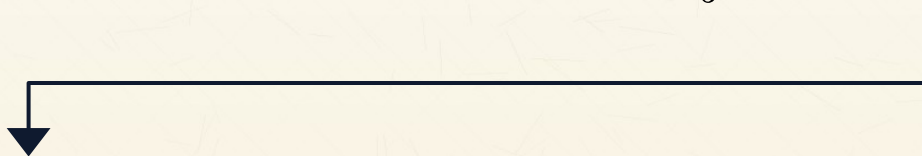
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$$\Gamma_L^{\{\mu\}} \in \{1, \gamma^5, \gamma^\mu, \gamma^\mu \gamma^5, \sigma^{\mu\nu}\}$$

Generalised atomic transitions

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$$\psi_e(\vec{x}) \sim \begin{pmatrix} F(r)\Omega(\theta, \phi) \\ G(r)\Omega'(\theta, \phi) \end{pmatrix}$$

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$$\sigma_i \Omega_q = \sum_{q'} c_{q'} \Omega_{q'}$$

Generalised atomic transitions

- Let's count the terms: $N_{\text{terms}} = N_L \times N_\mu \times N_{q'}$

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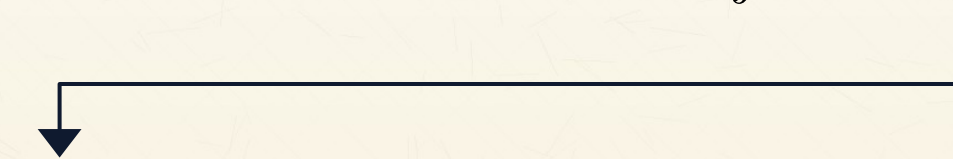
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Generalised atomic transitions

- Let's count the terms: $N_{\text{terms}} = N_L \times N_\mu \times N_{q'} \times 2$

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- Example, QED:

Generalised atomic transitions

- Let's count the terms: $N_{\text{terms}} = 3 \times N_{\mu} \times N_{q'} \times 2$

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- Example, QED:

$$\mathcal{O}_{L, \{\mu\}} \in \{eA_{\mu}, \mu_B F_{\mu\nu}, \mu_E \tilde{F}_{\mu\nu}\}$$

Generalised atomic transitions

- Let's count the terms: $N_{\text{terms}} = 3 \times 5 \times N_{q'} \times 2$

$$\mathcal{M}_{fi} = \langle f | \int d^3x \mathcal{H}_{\text{int}}(x) | i \rangle$$

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Generalised atomic transitions

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- We really want the squared amplitude:

Generalised atomic transitions

- Let's count the terms: $N_{\text{terms}} = (\text{a lot})^2$

$$\mathcal{M}_{fi} = \langle f | \int d^3x \mathcal{H}_{\text{int}}(x) | i \rangle$$

- We really want the squared amplitude:

$$|\mathcal{M}_{fi}|^2 = \left| \langle f | \int d^3x \mathcal{H}_{\text{int}}(x) | i \rangle \right|^2$$

Generalised atomic transitions

- Let's count the terms: $N_{\text{terms}} = \sum (\text{a lot})^2$

$$\mathcal{M}_{fi} = \langle f | \int d^3x \mathcal{H}_{\text{int}}(x) | i \rangle$$

- And what if we want to sum over polarisations?

$$|\mathcal{M}_{fi}|^2 = \sum_{m,m'} |\langle f | \int d^3x \mathcal{H}_{\text{int}}(x) | i \rangle|^2$$

Generalised atomic transitions

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$$\mathcal{M}_{fi} = \langle f | \int d^3x \mathcal{H}_{\text{int}}(x) | i \rangle$$

- Factorise:

$$|\mathcal{M}_{fi}|^2 = \left| \sum \mathcal{I}_r \cdot \mathcal{I}_\Omega \right|^2$$

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Radial integral: sets the scale



Generalised atomic transitions

- Let's count the terms: $N_{\text{terms}} = \sum (\text{a lot})^2$

$$\mathcal{M}_{fi} = \langle f | \int d^3x \mathcal{H}_{\text{int}}(x) | i \rangle$$

- Factorise:

$$|\mathcal{M}_{fi}|^2 = \left| \sum \mathcal{I}_r \cdot \mathcal{I}_\Omega \right|^2$$

Radial integral: sets the scale

Angular integral: selection rules

Generalised atomic transitions



<https://gitlab.com/JShergold/cinco>

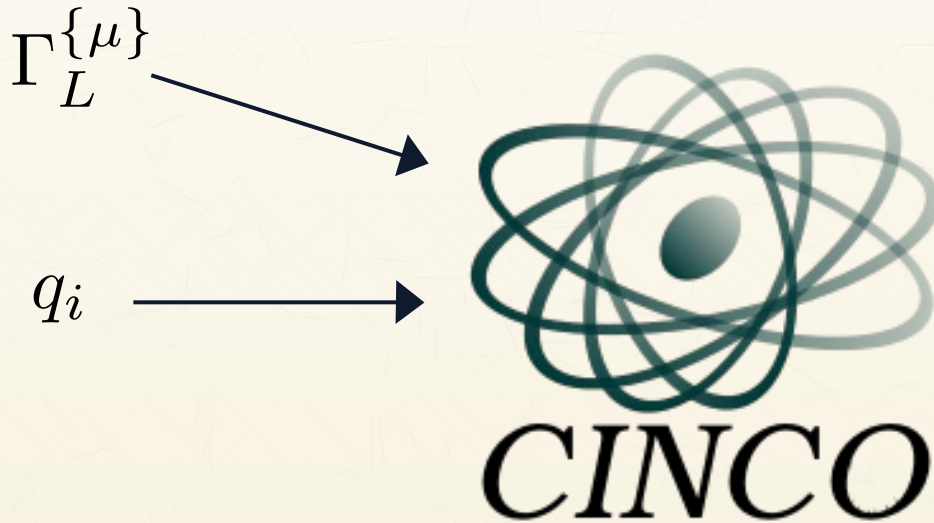
Generalised atomic transitions

$\Gamma_L^{\{\mu\}}$



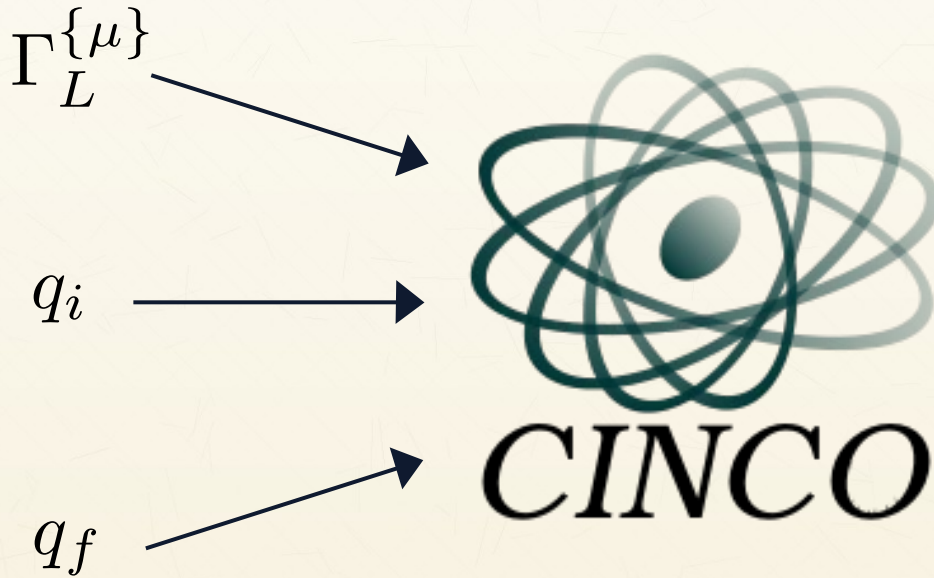
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Generalised atomic transitions



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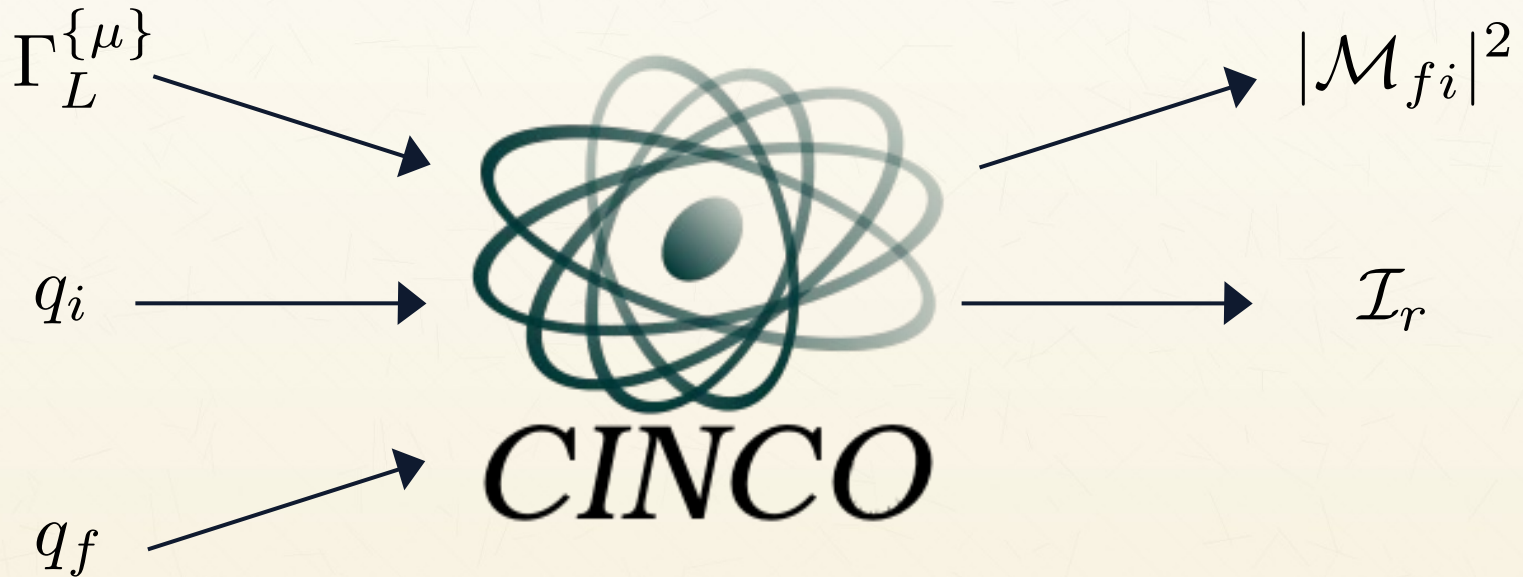
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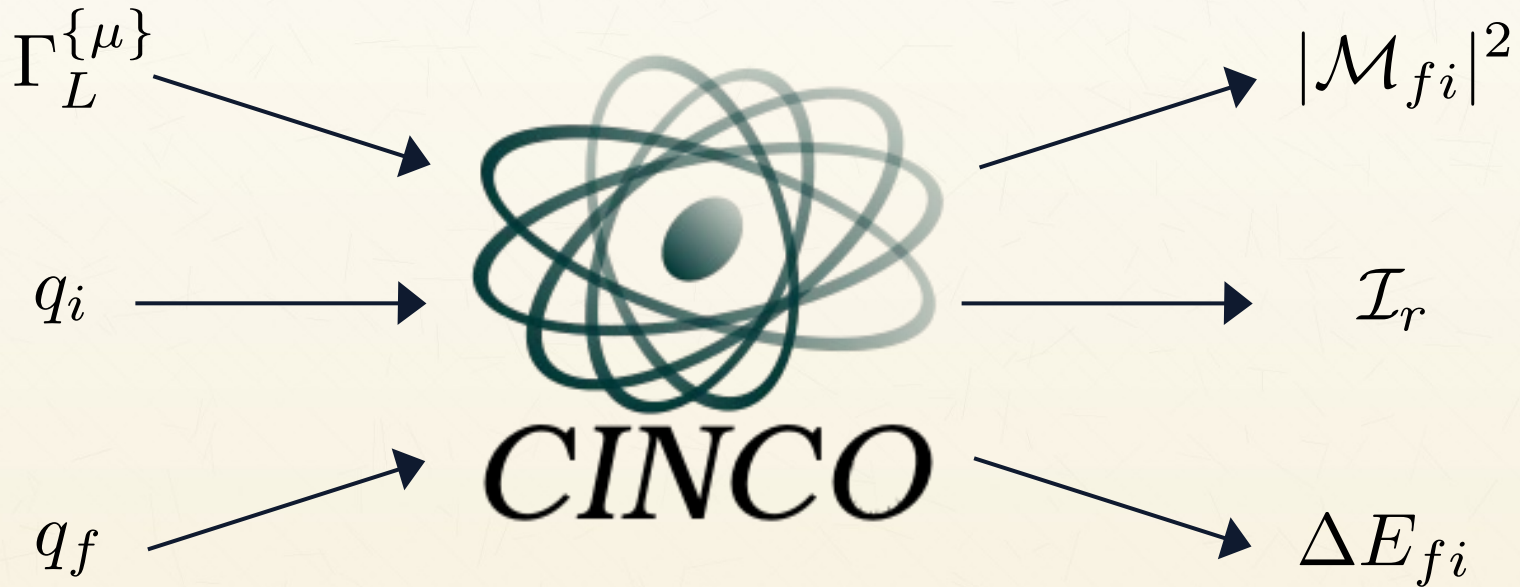
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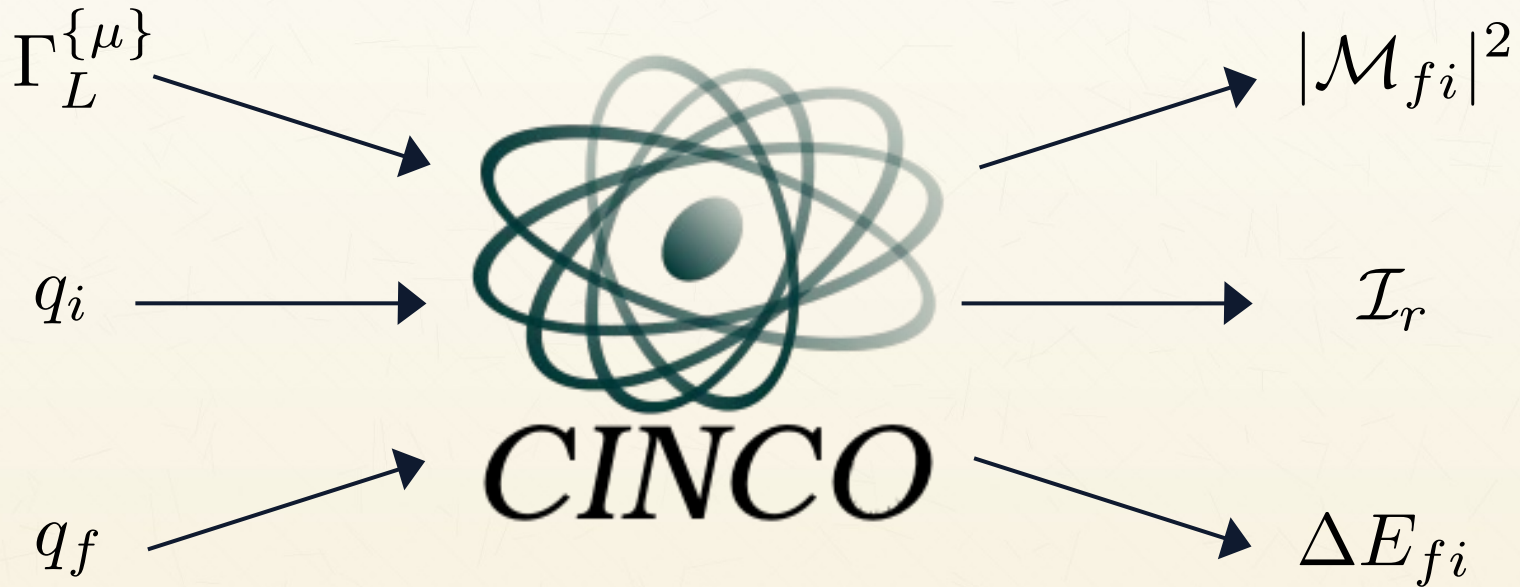
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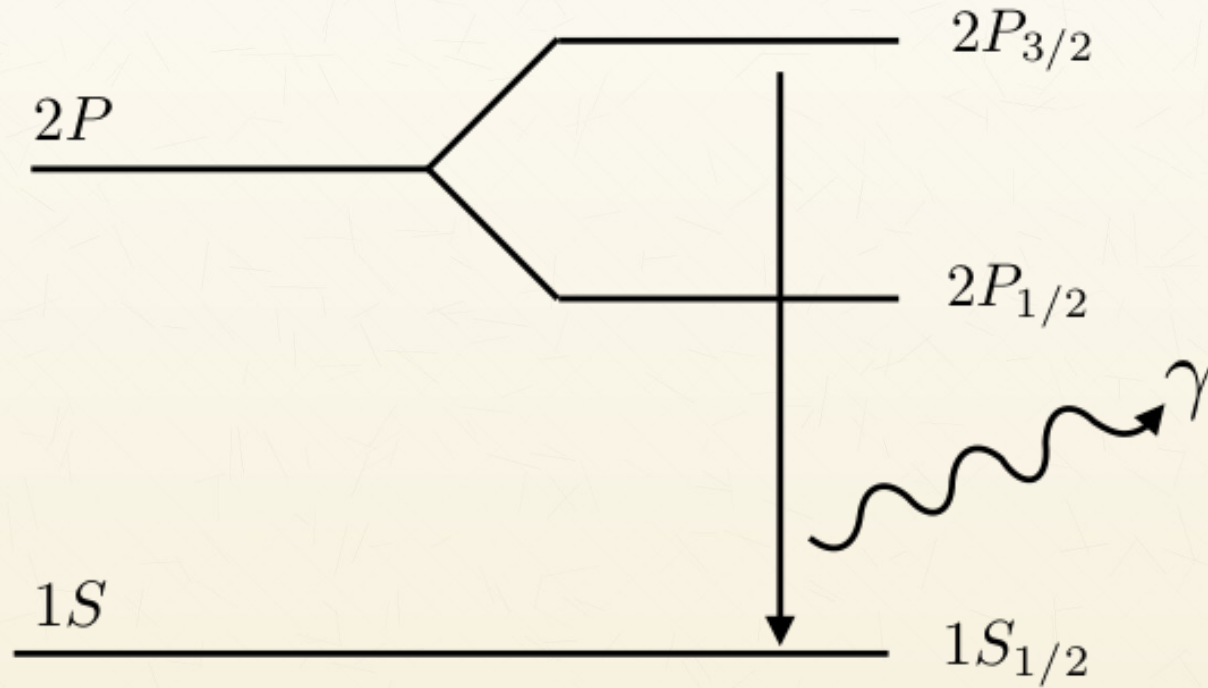
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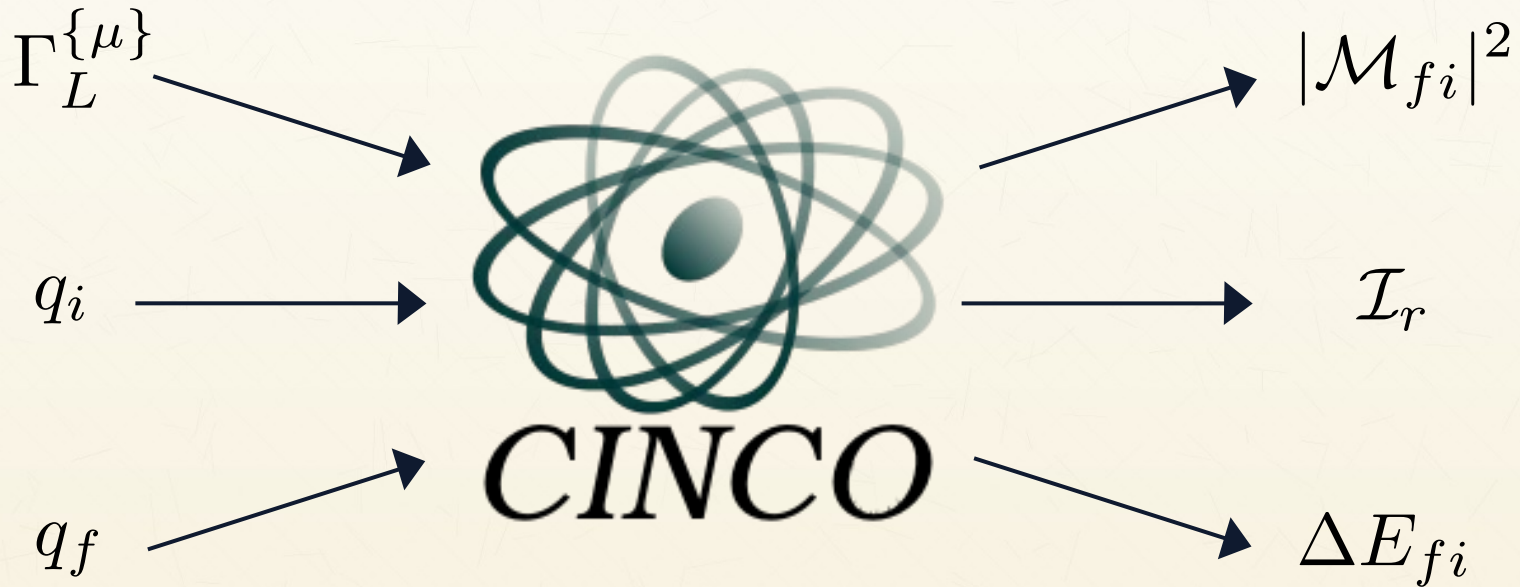


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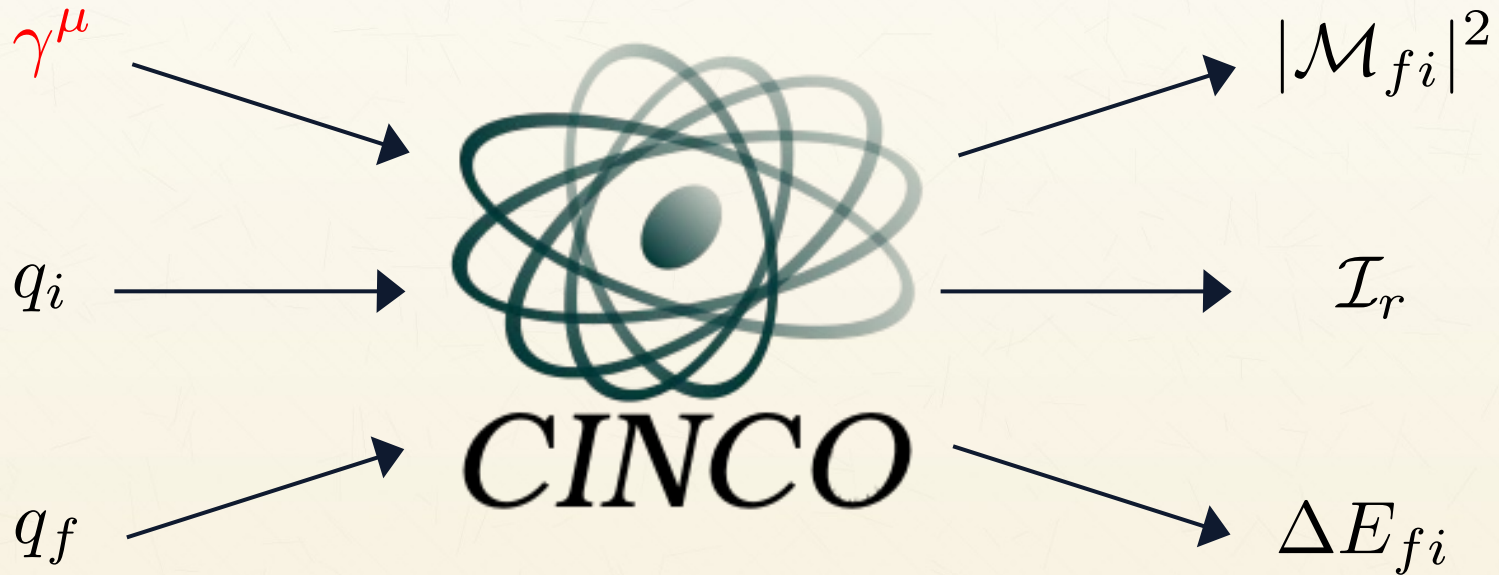


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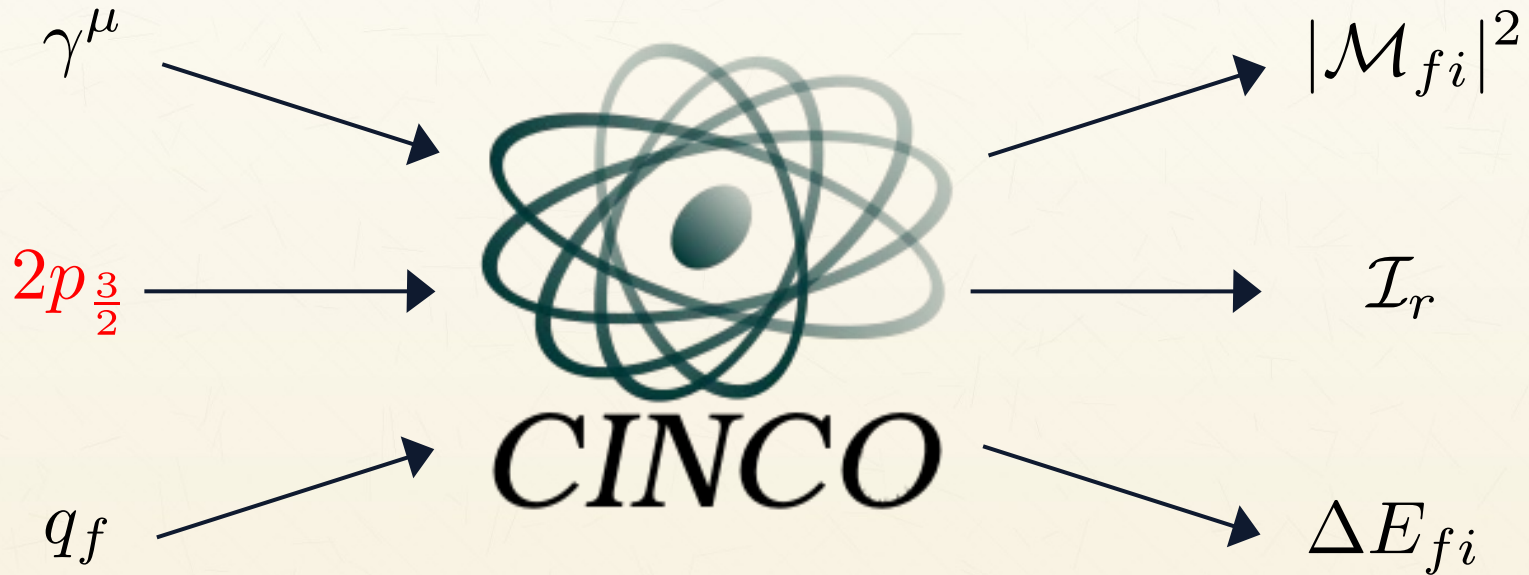
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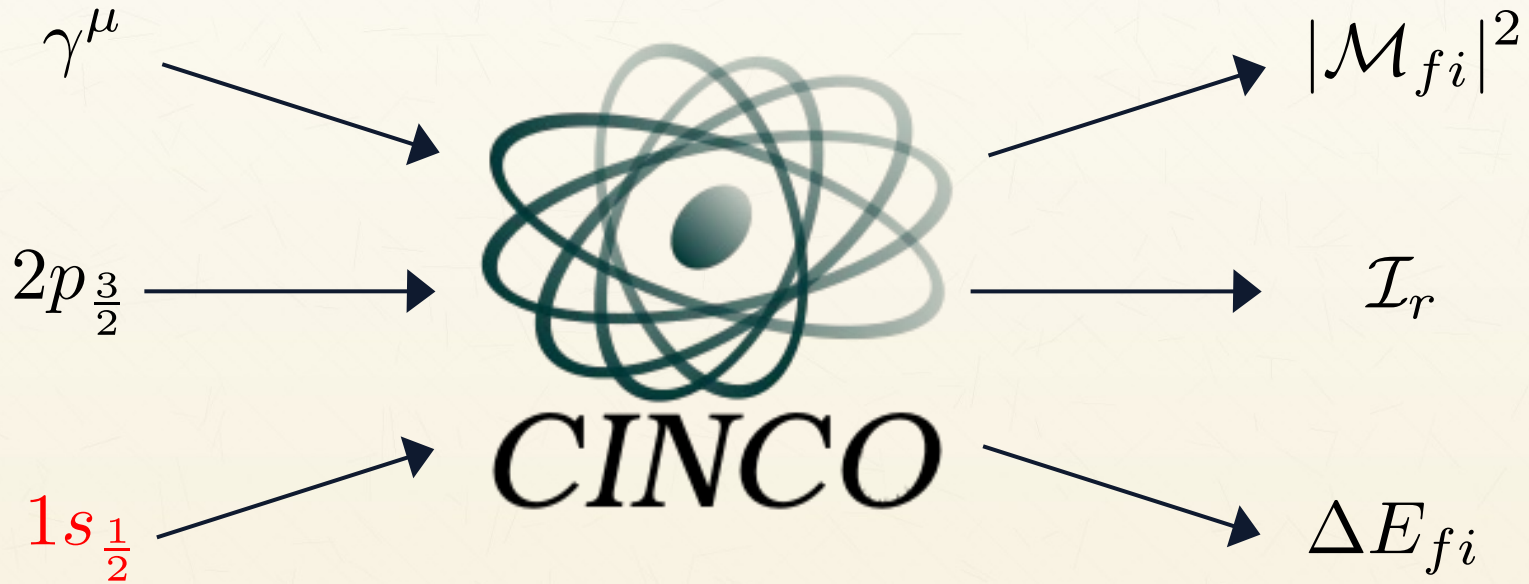
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Generalised atomic transitions

- Transition amplitude:

<https://gitlab.com/JShergold/cinco>

Generalised atomic transitions

- Transition amplitude:

$$\langle |\mathcal{M}_{fi}|^2 \rangle = \frac{4}{9} \left(\langle \vec{\mathcal{O}}_V \rangle^* \cdot \langle \vec{\mathcal{O}}_V \rangle \right) \mathcal{I}_{gf}^2$$

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<https://gitlab.com/JShergold/cinco>

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- The expectation value (not done by CINCO):

$$\langle \mathcal{O}_{V,\mu} \rangle = \langle \gamma | A_\mu | 0 \rangle = \frac{e}{\sqrt{2E_\gamma}} \epsilon_\mu^*$$

<https://gitlab.com/JShergold/cinco>

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- Transition amplitude:

$$\langle |\mathcal{M}_{fi}|^2 \rangle = \frac{8\pi\alpha_{\text{EM}}}{9E_\gamma} (\vec{\epsilon} \cdot \vec{\epsilon}^*) \mathcal{I}_{gf}^2$$

<https://gitlab.com/JShergold/cinco>

Generalised atomic transitions

- Phase space integral:

$$\Gamma_{2p_{3/2} \rightarrow 1s_{1/2}} = \frac{16\pi\alpha_{EM}}{9} \Delta E_{fi} \mathcal{I}_{gf}^2$$

<https://gitlab.com/JShergold/cinco>

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$$\Gamma_{2p_{3/2} \rightarrow 1s_{1/2}} = 6.2650 \cdot 10^8 \text{ s}^{-1}$$

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- Motivation: CvB and DM
- Generalised atomic transitions
- **Pair absorption**
- The future: molecules and ML

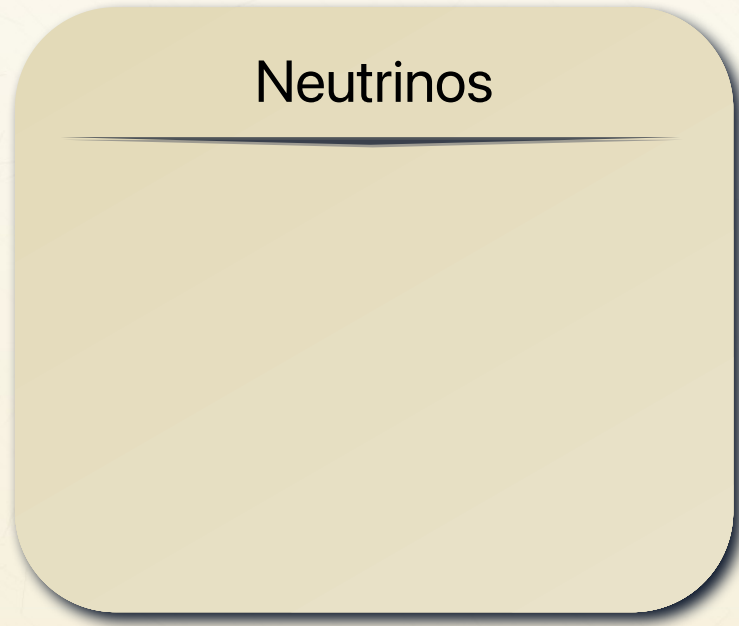


Pair absorption

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Dark Matter

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Effective dim-6 coupling



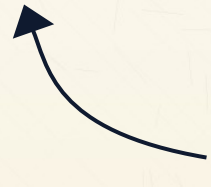
Pair absorption

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$$R_{\text{one}} = G_{\text{eff}}^2 E_{\text{thresh}}^2 n_{\varphi}$$

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$$R_{\text{pair}} > R_{\text{one}}$$

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$$n_\varphi > E_{\text{thresh}}^3$$

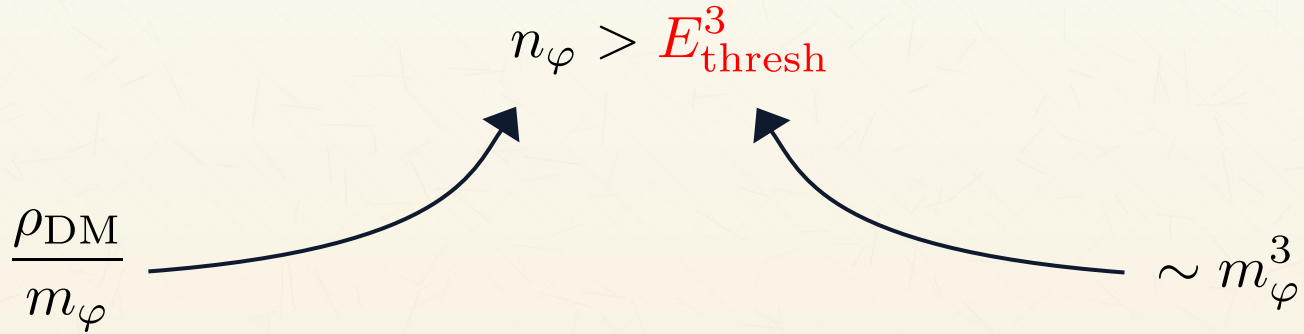
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$$\frac{\rho_{\text{DM}}}{m_\varphi} \rightarrow n_\varphi > E_{\text{thresh}}^3$$

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$0.3 - 0.4 \text{ GeV cm}^{-3}$



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- Precisely where atomic transitions sit!


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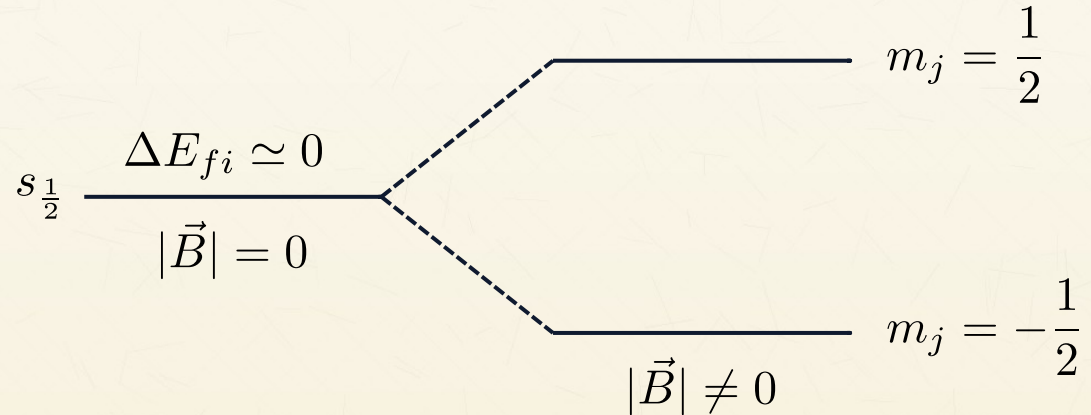
$$s_{\frac{1}{2}} \frac{\Delta E_{fi} \simeq 0}{|\vec{B}| = 0}$$

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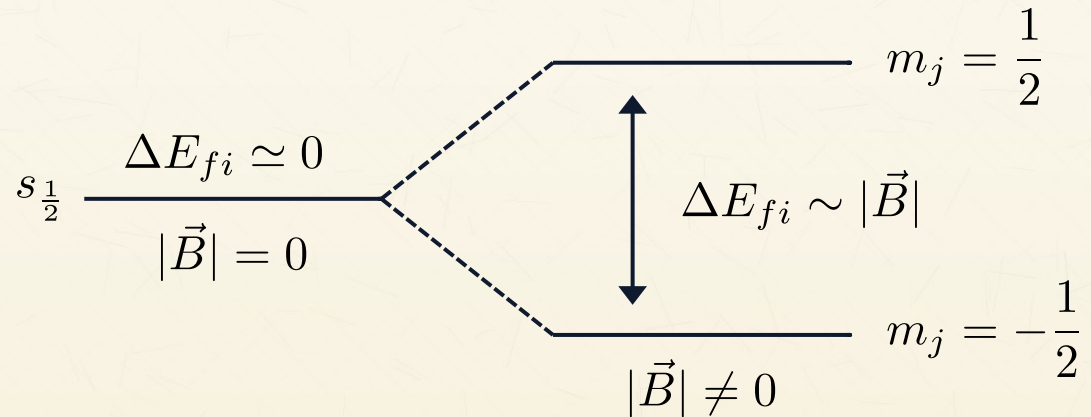


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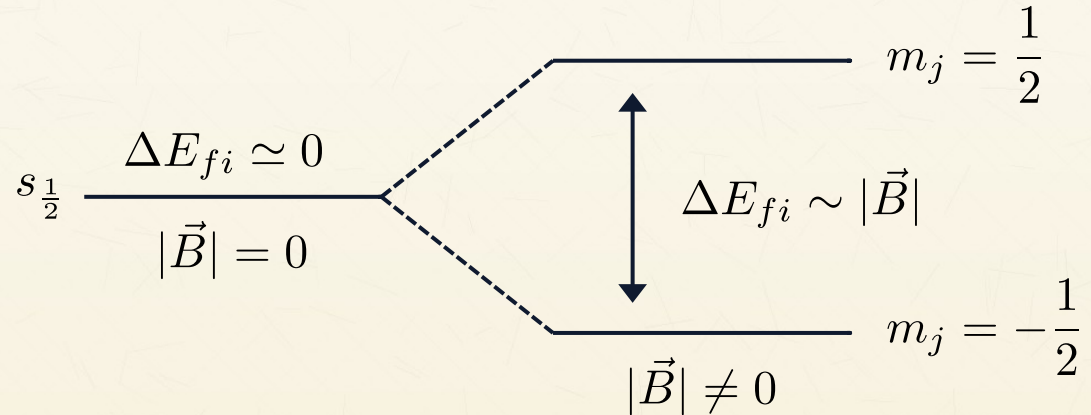
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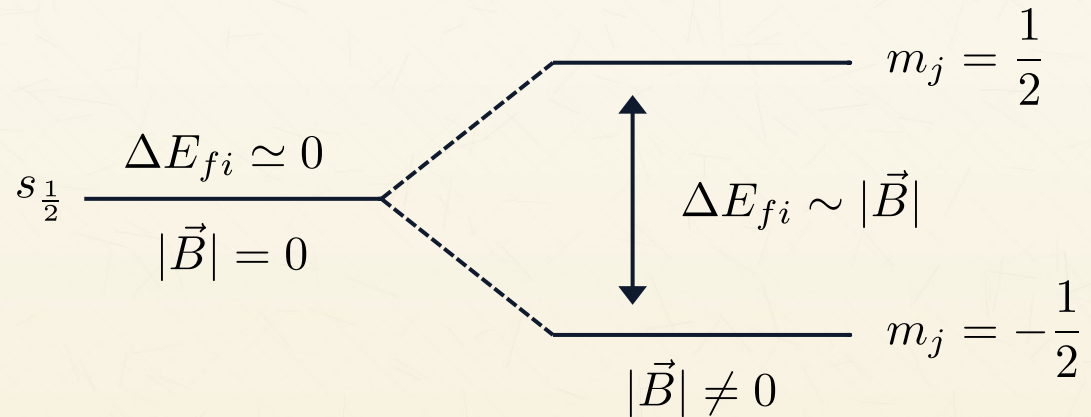
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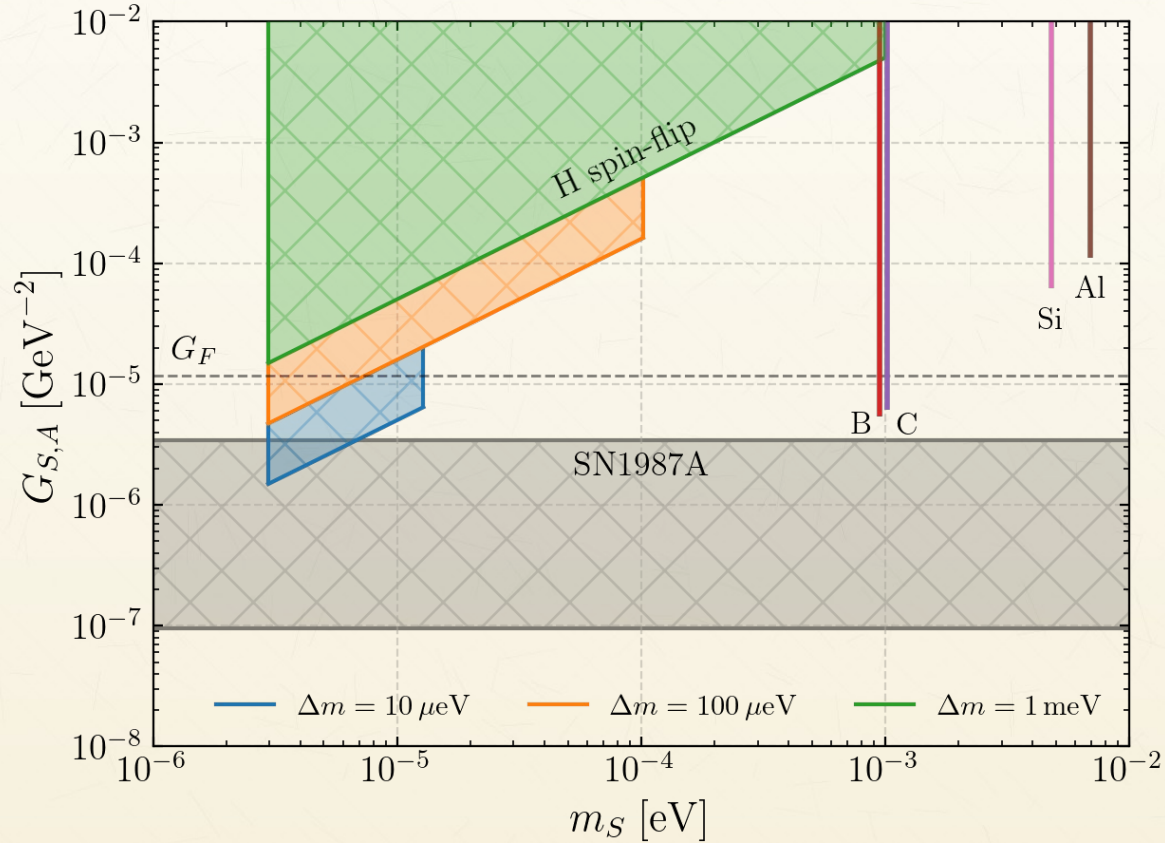
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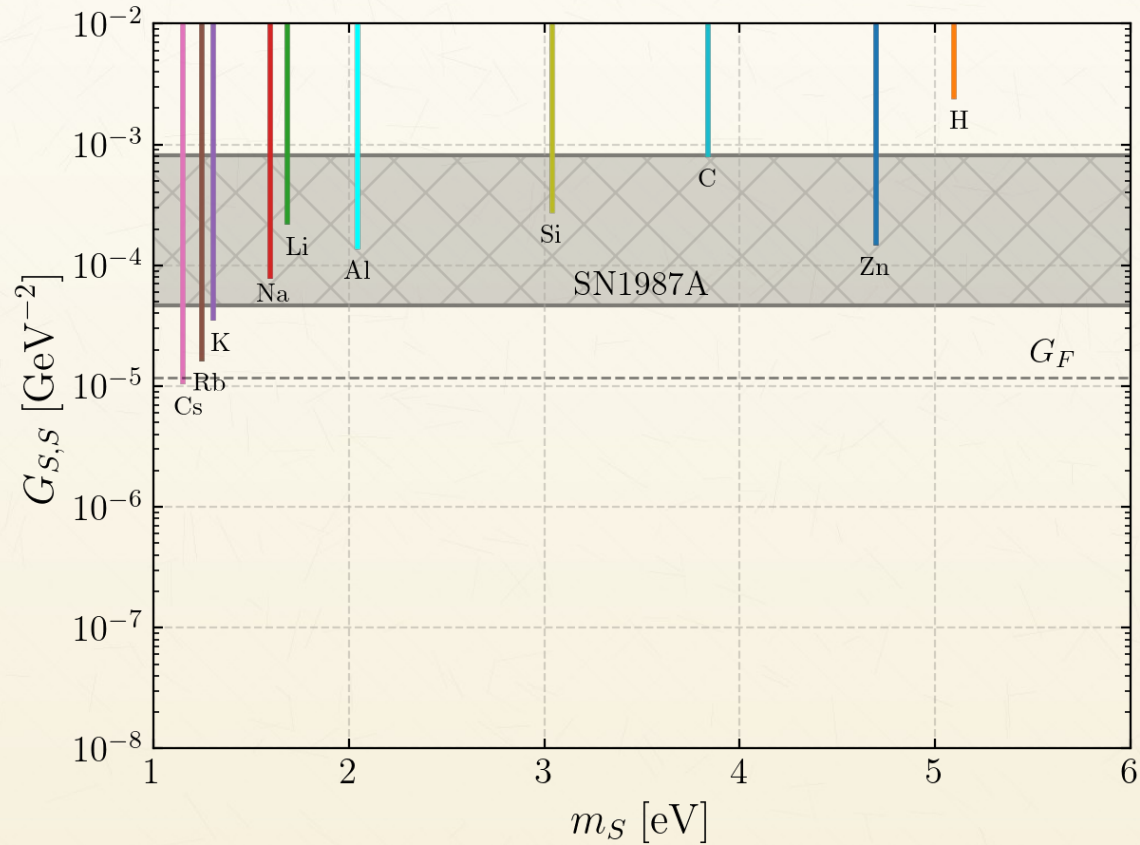
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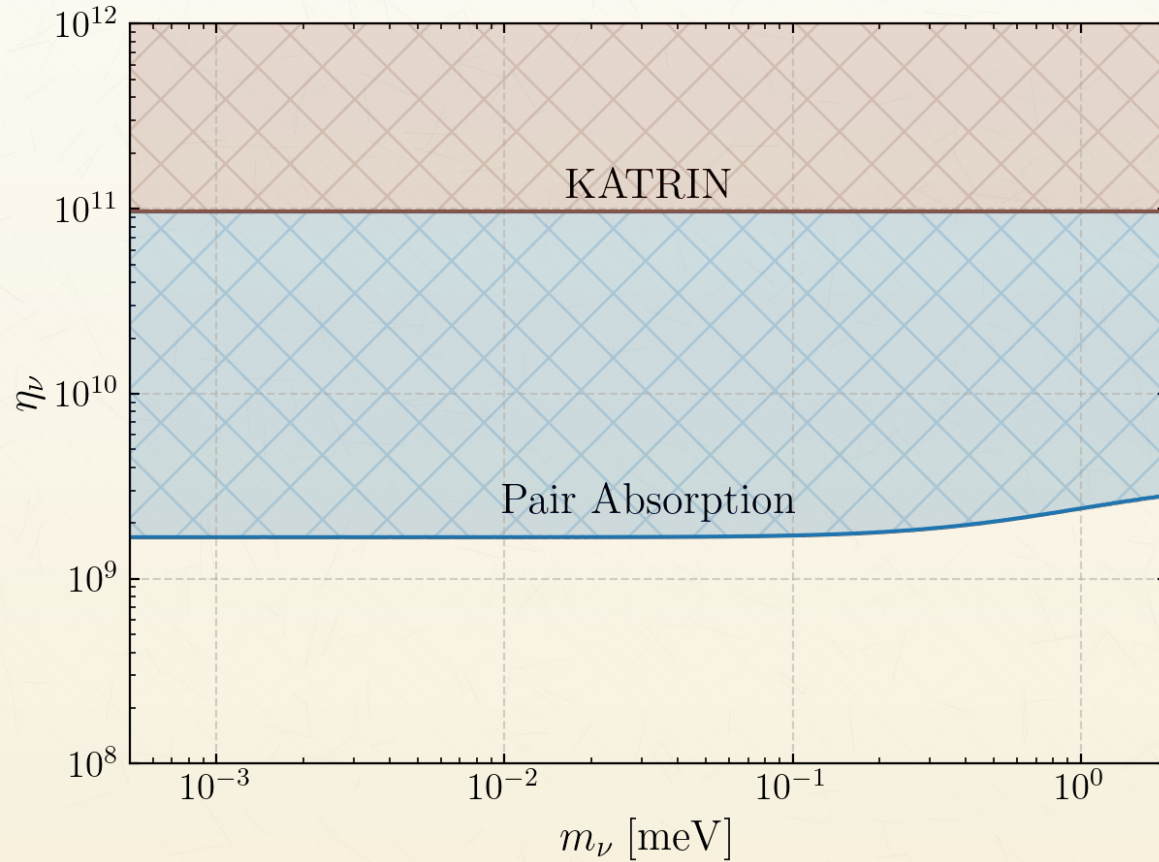
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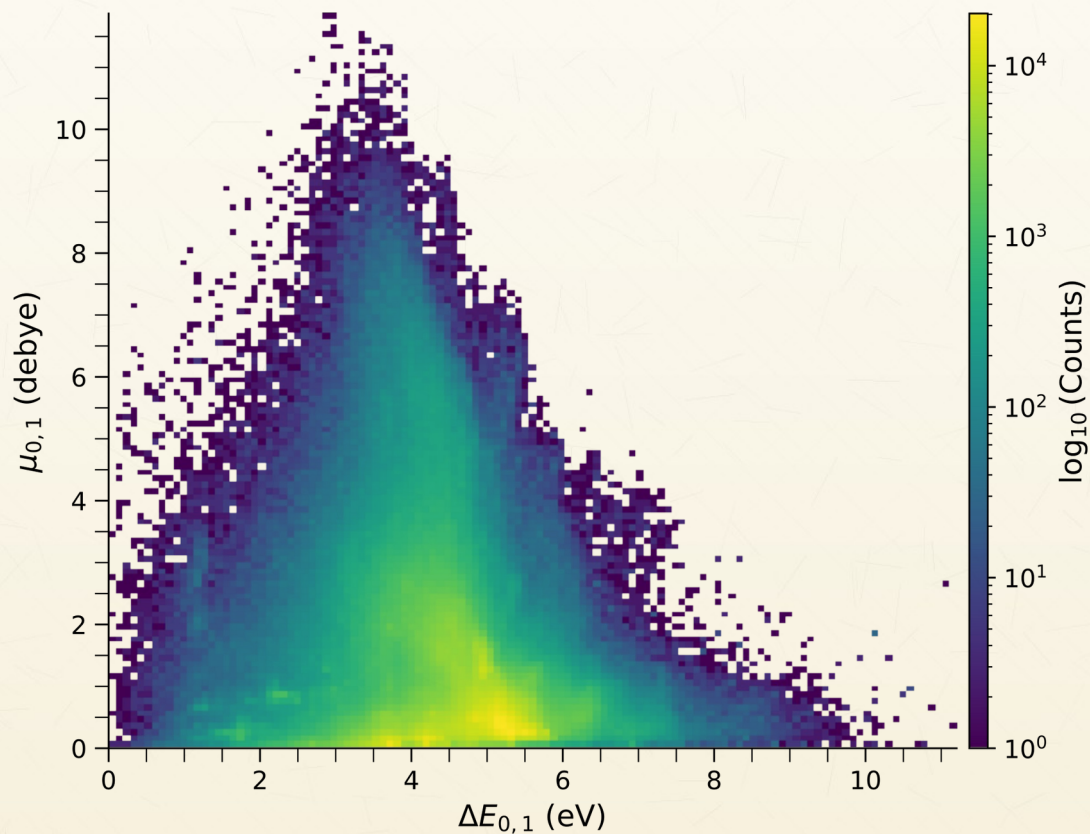
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The future: molecules and ML



The future: molecules and ML

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- Watch this space!

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- Molecules and crystals, the future?

**“I really don’t like the beige.” – My
fiancée**

Thank you
Any questions?