

Extension of the Lindhard model, to calculate the "quenching factor" at low energies

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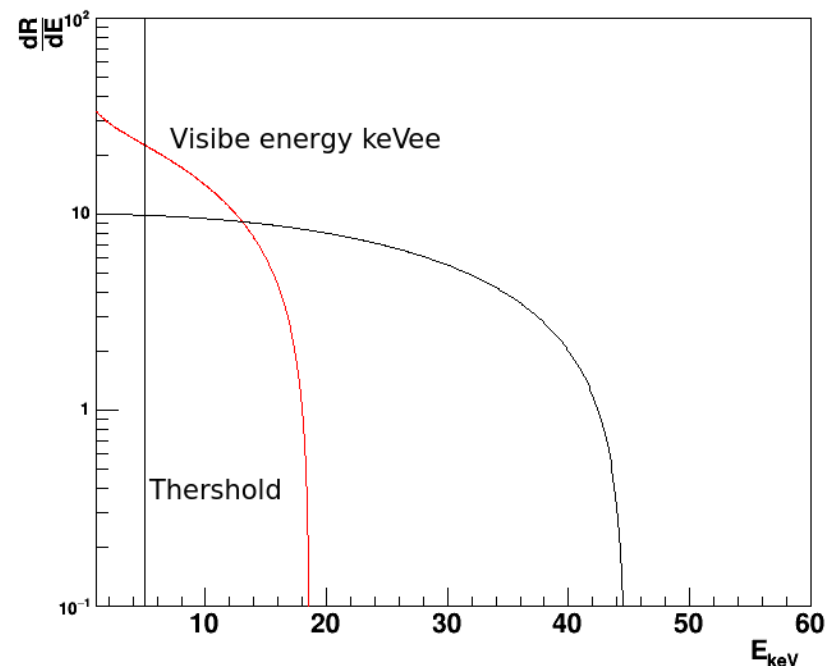
Quenchig Factor

- The amount of electronic excitation produced by a recoiling atom is smaller than that produced by a recoiling electron of the same energy.
- The Q.F is the fraction of energy that is transformed into ionization energy in a crystal.

$$f_n = \frac{\eta}{\varepsilon}$$

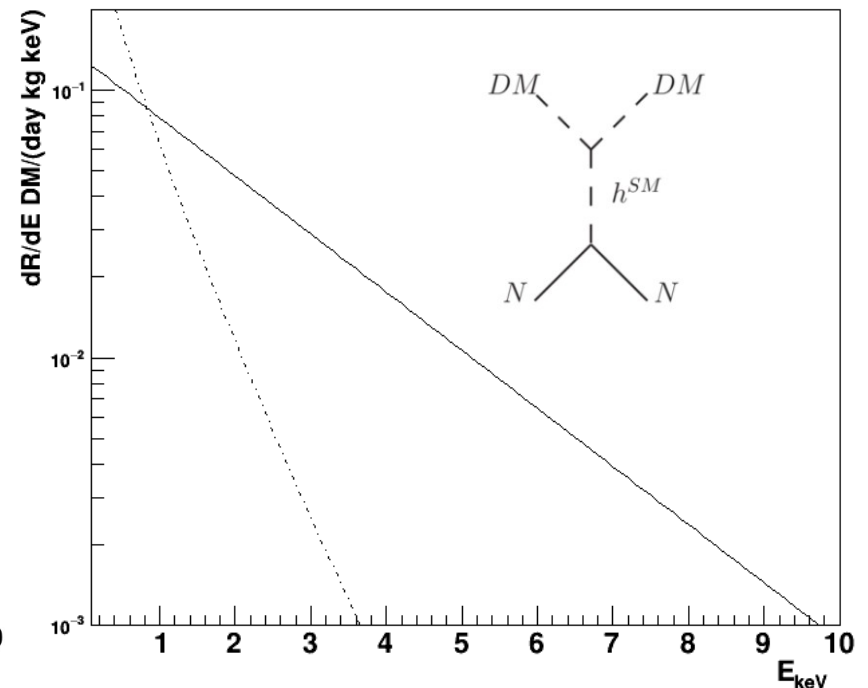
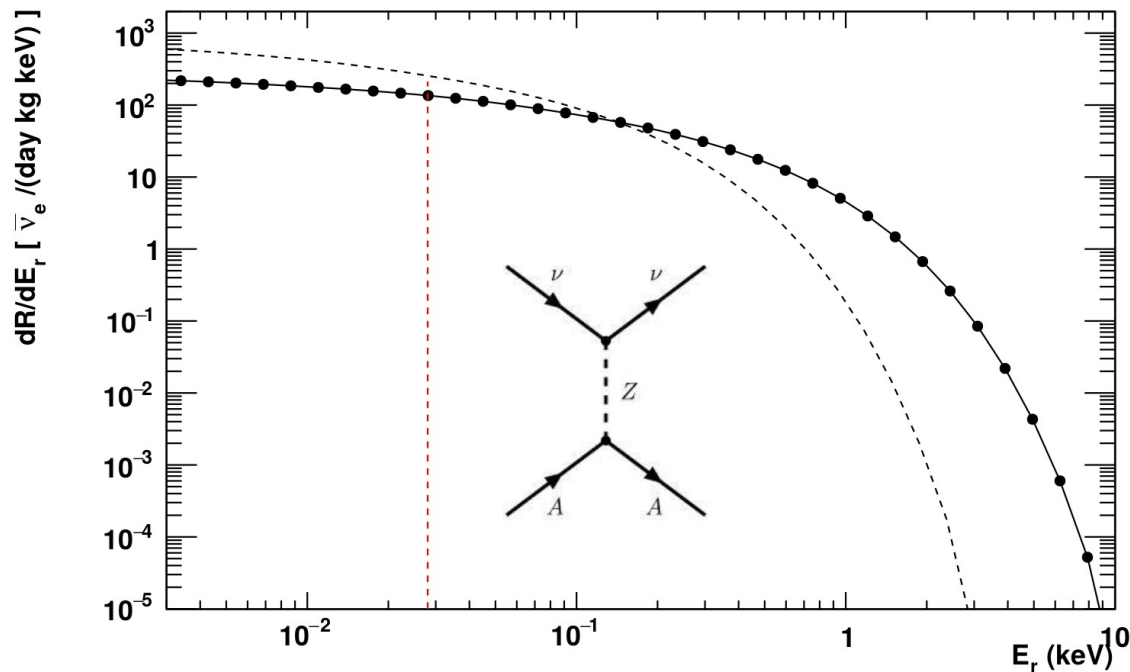
The Q.F shift the spectrum to lower energies.

$$\frac{dR}{dE_v} = \frac{dR}{dE} \frac{dE}{dE_v} = \frac{dR}{dE} \left(f_n + E \frac{df_n}{dE} \right)^{-1}$$



Quenching factor

- In the direct dark matter searches experiments or in the measurements of the coherent elastic neutrino-nucleus scattering, the signal usually consists in the detection of the low energy (~ 1 keV) of the recoiling target atoms.
- As the experiments have reduced their detection thresholds well below $1 \sim \text{keV}$, the understanding of the quenching at low energies has become increasingly important.



φ can be η or ν

Lindhard Model

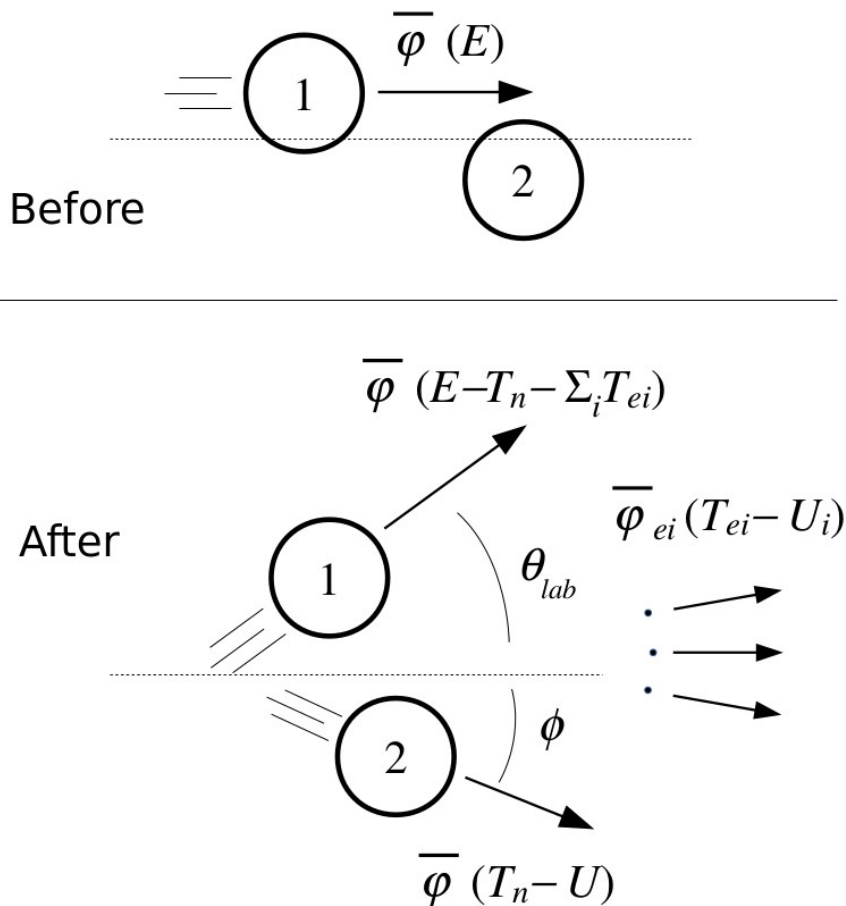
$$\int d\sigma_{n,e} [\underbrace{\bar{\varphi}(E - T_n - \Sigma T_{ei})}_{\text{Energy due to electronic stripping process}} + \underbrace{\bar{\varphi}(T_n - U)}_{\text{Energy due to recoil process}} - \underbrace{\varphi(E)}_{\text{projectile atom before collision}}] = 0$$

Target atom after collision projectile atom after collision projectile atom before collision

Energy is divided in nuclear and electronic average losses (ν , atomic movement, η ionization losses).

$$E = \bar{\nu}(E) + \bar{\eta}(E).$$

- The physical effect (φ) of the incident particle equals the sum of the physical effect of the scattered particle and that of the struck atom
- When writing the above equation Lindhard took into account the energy conservation in the scattering of two ions. To deal with the equation he made four basic approximations: i) electrons do not contribute to the atomic movement process, ii) the atomic-binding energy U can be neglected, iii) the electronic stripping process is expanded up to first order, iv) nuclear and electronic effects are separable.



$$k\varepsilon^{1/2}\bar{v}'(\varepsilon) = \int_0^{\varepsilon^2} \frac{dt}{2t^{3/2}} f(t^{1/2}) \times \left\{ \bar{v}\left(\varepsilon - \frac{t}{\varepsilon}\right) - \bar{v}(\varepsilon) + \bar{v}\left(\frac{t}{\varepsilon}\right) \right\}$$

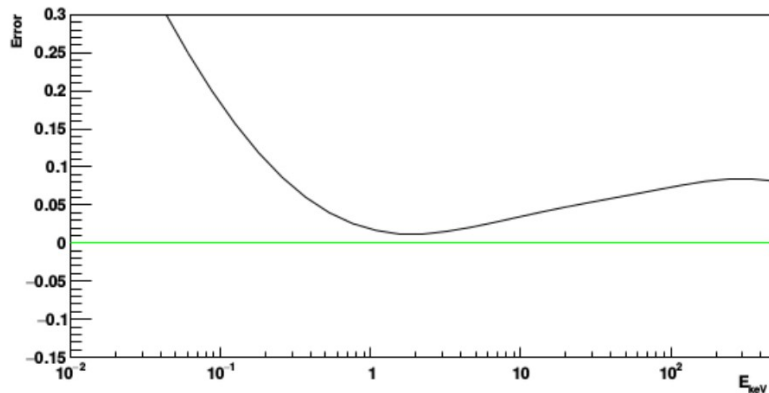
Diagram illustrating the components of the Lindhard equation:

- reduced energy** (points to ε)
- nuclear energy loss** (points to $f(t^{1/2})$)
- electronic energy loss** (points to $k\varepsilon^{1/2}\bar{v}'(\varepsilon)$)
- target atom after collision** (points to $\bar{v}(\varepsilon - \frac{t}{\varepsilon})$)
- projectile atom before collision** (points to $-\bar{v}(\varepsilon)$)
- projectile atom after collision** (points to $\bar{v}(\frac{t}{\varepsilon})$)

ε : dimensionless reduced energy.

- Lindhard after these approximations can deduce this equation .
- Where $t=\varepsilon^2\sin^2(\theta/2)$, $f(t)$ is related to the nuclear cross section scattering.
- Lindhard equation has a problem: the boundary condition at $E \rightarrow 0$, is $v(E) \rightarrow E$, so $v' \rightarrow 1$ in this limit. But the Integral equation predict $v' \rightarrow 0$.
- So the Integral equation can only be solve for energies far from zero.

- Therefore, the equation can be solved making to the Integral equation more raw approximations.
- He found a “reasonable” parametrization from this equation for energies below $\epsilon < 1$.
- This allowed Lindhard to solve numerically his equation for energies $\epsilon > 1$.
- So this approach doesn't allow us to give a good prediction in low energy regions.



The parametrization doesn't solve the Equation at low energies.

$$\bar{\nu}_l(\epsilon) = \frac{\epsilon}{1 + kg(\epsilon)}$$

$$g(\epsilon) = 3\epsilon^{0.15} + 0.7\epsilon^{0.6} + \epsilon$$

Standar high
energy
parametrization



DASK
computer
45 kW
3.5 Tons

Extension of Lindhard model

- We introduce the atomic-binding energy, considering the changing the term

$$\bar{\nu}(t/\varepsilon) \rightarrow \bar{\nu}(t/\varepsilon - u)$$

- Expand the 1st term in the integral equation $\varphi(\varepsilon - t/\varepsilon - \Sigma T_{ei})$ to 2nd order in ΣT_{ei} , (Probability of electronic stripping process term).
- The lower limit of integration (over dt) can no longer be zero. The min t is now εu . (predicted also by model of scattering of two semi-hard spheres).
- In this model we are going to consider u constant.

- So with this modification we can deduce a new Integral modified equation with atomic-binding energy.

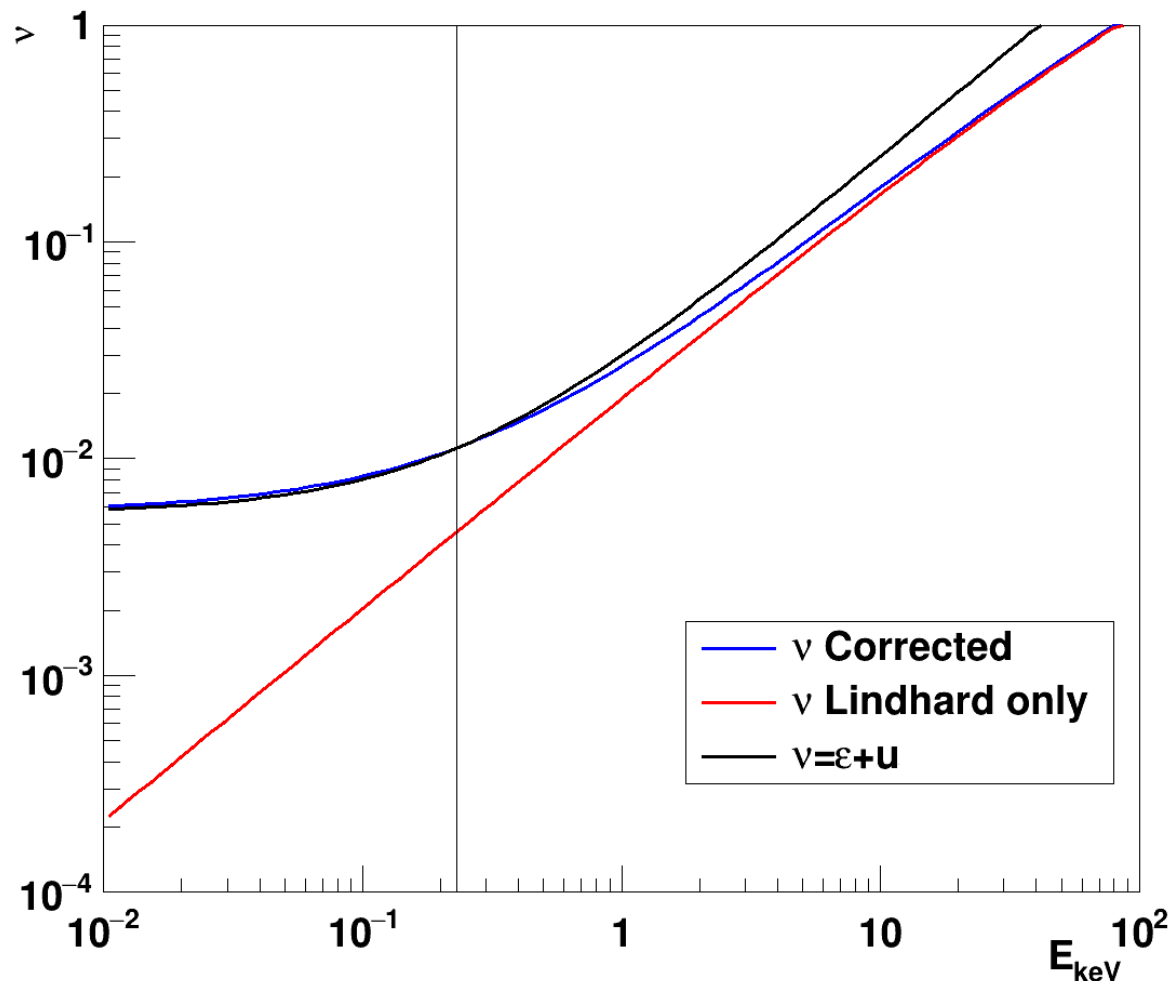
$$-\frac{1}{2}k\varepsilon^{3/2}\bar{v}''(\varepsilon) + k\varepsilon^{1/2}\bar{v}'(\varepsilon) = \int_{\varepsilon u}^{\varepsilon^2} \frac{dt f(t^{1/2})}{2t^{3/2}} [\bar{v}(\varepsilon - t/\varepsilon) + \bar{v}(t/\varepsilon - u) - \bar{v}(\varepsilon)]$$

- This equation ***can be solve numerically from $\varepsilon \geq u$*** . So the equation predict a threshold energy of u.
- The equation requires the following behavior of the solution.

$$\bar{v}_u(\varepsilon) = \begin{cases} \varepsilon + u & \varepsilon < u \\ \varepsilon + u - g(\varepsilon) & \varepsilon \geq u \end{cases} \quad f_n = \frac{\bar{\eta}}{\varepsilon + \bar{u}} = 1 - \frac{\bar{v}}{\varepsilon + \bar{u}}.$$

- The equation admits a solution featuring a "kink" at $\varepsilon=u$ (discontinuous 1st derivative).
- We solve the equation by the shooting method pursuing the boundary condition at high energies that $v'' \rightarrow 0$ and for $\varepsilon=u$

Alternative good parametrization



- We can implement a good analytical approximation to solve the Integral equation.
- The idea is to lessen the ionization contribution, subtracting a fraction of the electronic stopping power.
- $\eta \rightarrow \eta - C' - C\epsilon^{1/2}$

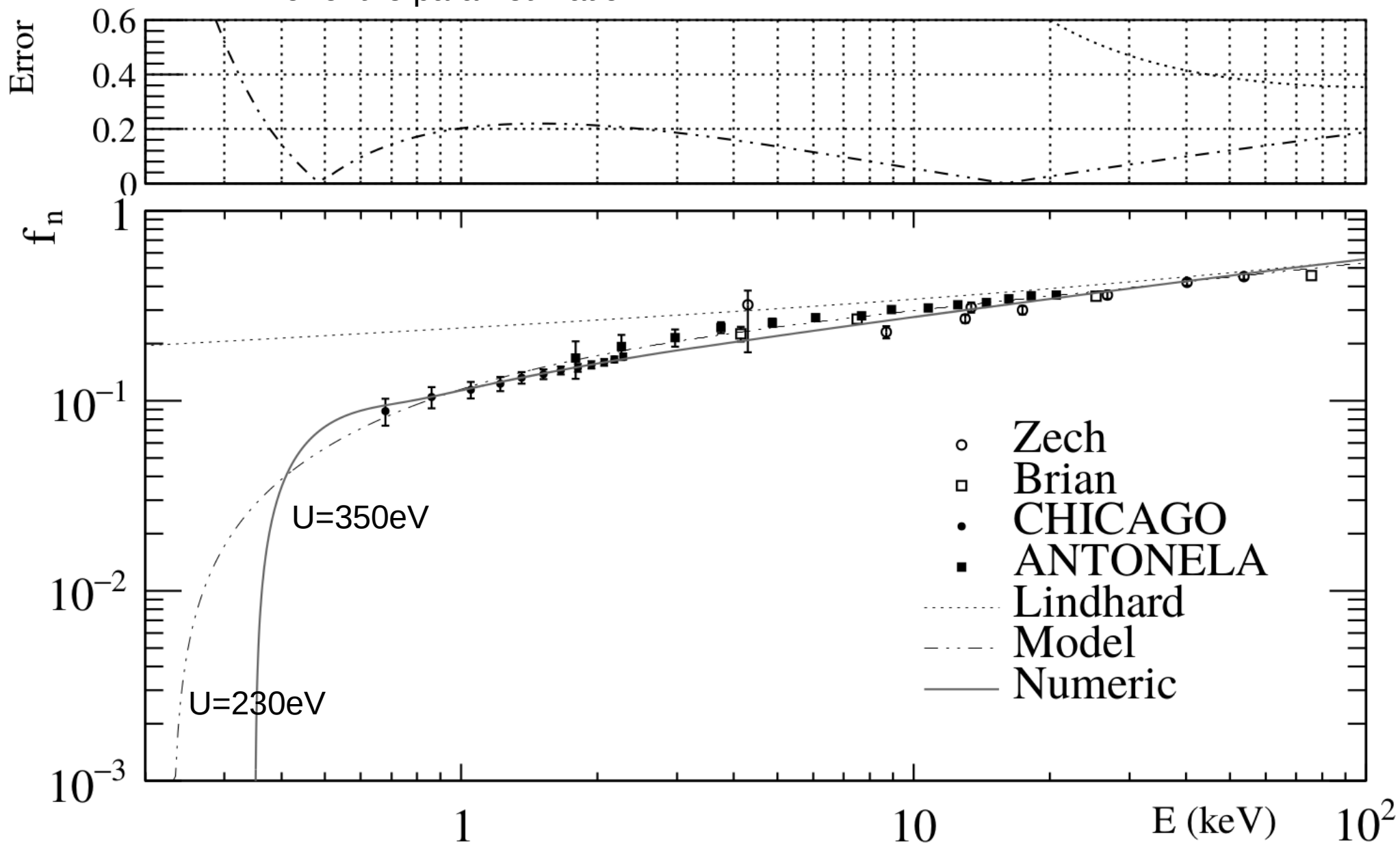
Where C , C' and u are estimated from a fit to the available data.

Results for Si, Ge and Xe.

- We assumed as a first simple model a constant atomic-binding energy (average).
- And also we assume that the k factor, of the electronic stopping power is constant too.

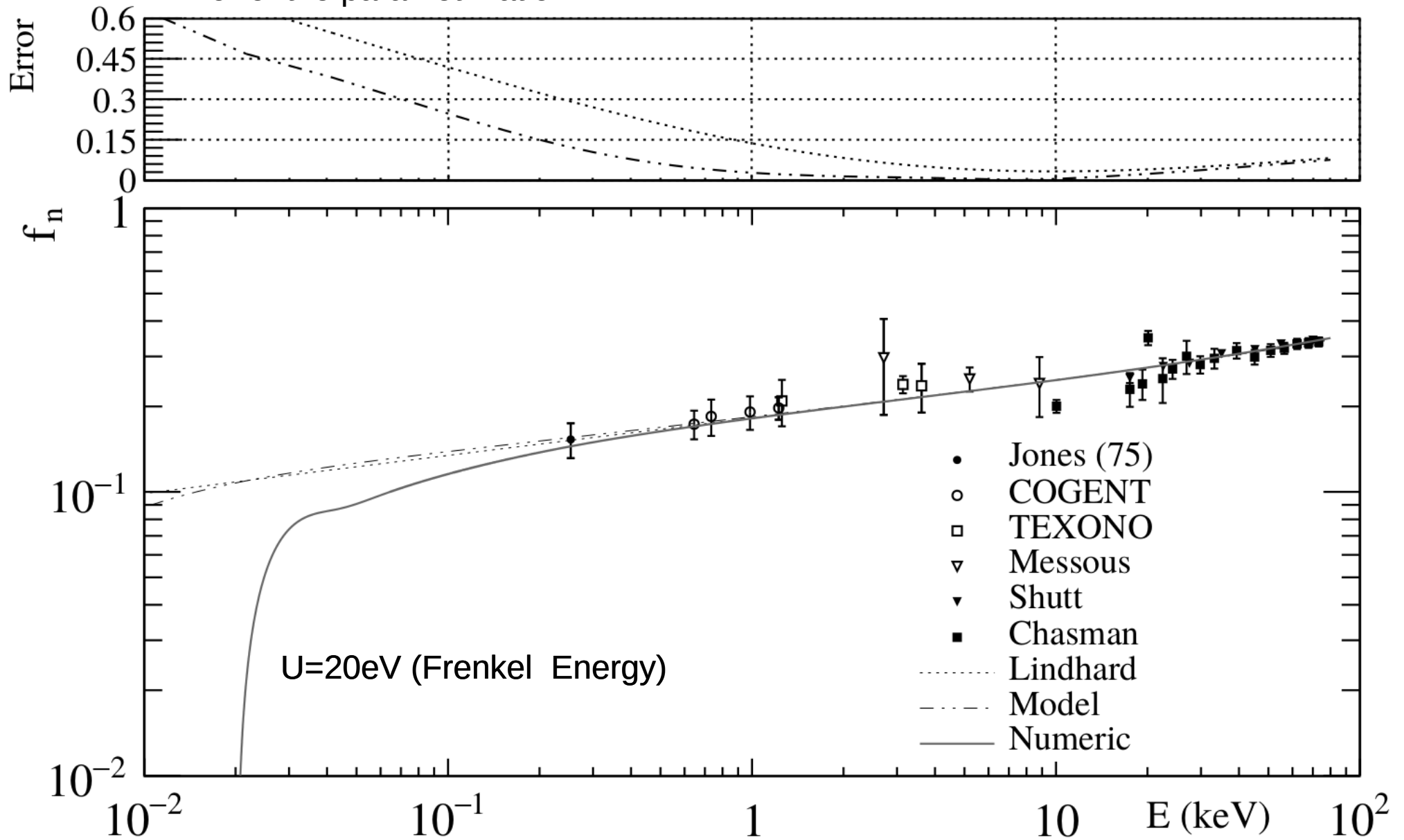
Si

Error of the parametrization



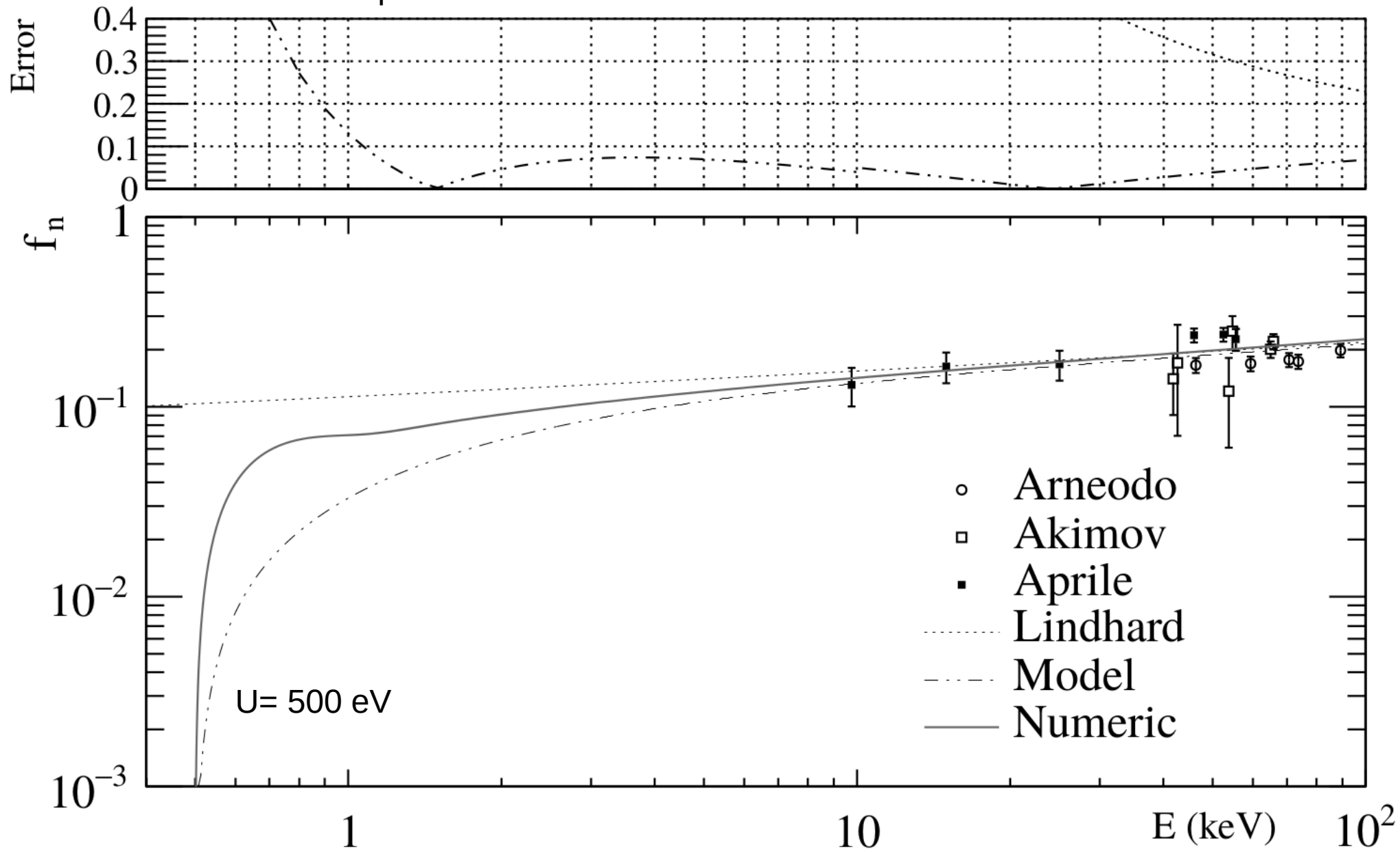
Ge

Error of the parametrization



Xe

Error of the parametrization



Conclusions

- We have implemented an extension to Lindhard model which predicts a cut off at the binding energy.
- We study this new equation and deduce a numeric algorithm that solve the problem with no need of simplifications.
- Also the model now only required the specifications of k and u to have a QF.
- We think that this QF model can have an important impact in the new upcoming results from dark matter and neutrino experiments.
- Other approach (Sorensen) tried to solve this, but in an incomplete way, that doesn't describe the data available.
- We are going to publish this work very soon.

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