

# **Non-Dipolarity of Channeling Radiation at GeV Beam Energies**

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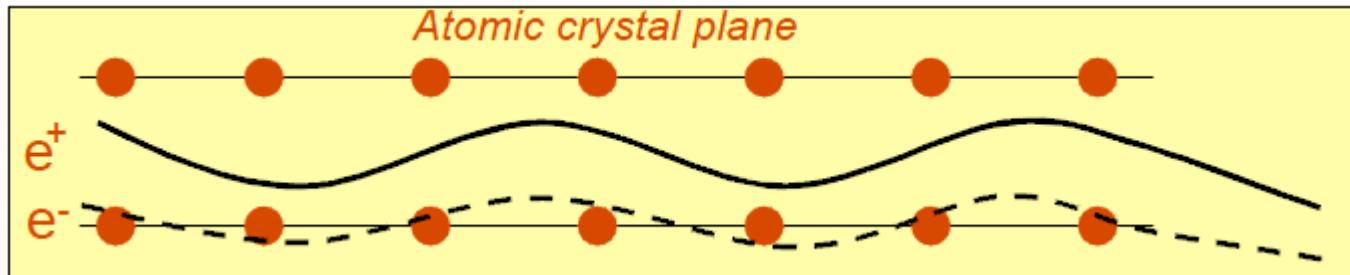
**26.09.2016, del Garda Italy**

## **Outline**

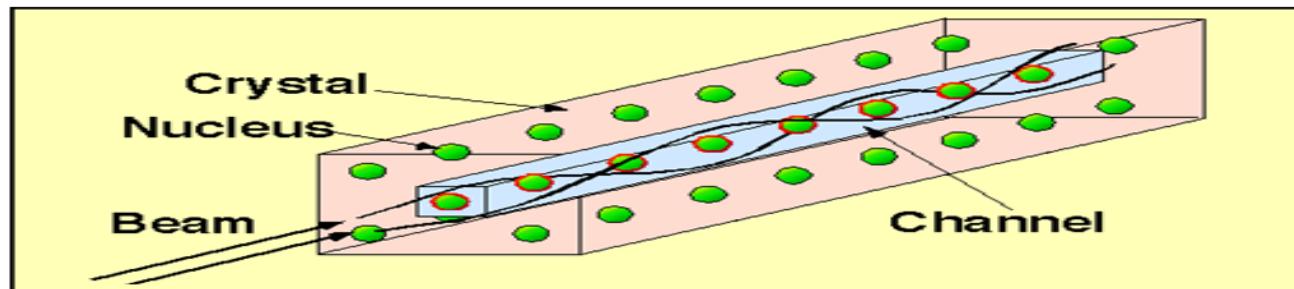
- 1. Introduction**
- 2. Continuum potential**
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- 6. Comparison of non-dipole with dipole approximation**
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## 1. Introduction

Planar channeling: one-dimensional problem



Axial channeling: two-dimensional problem



## 2. Continuum potential

### Planar channeling

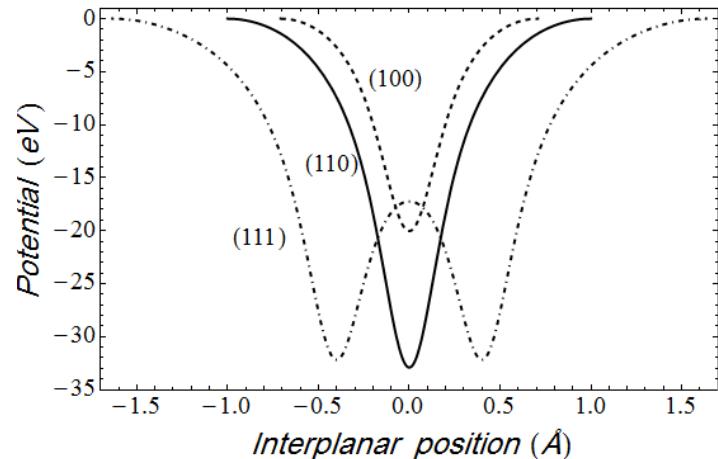
$$V(x) = \sum_n v_n e^{inx}$$

$$v_n = -\frac{2\pi}{V_c} a_0^2 (e^2 / a_0) \sum_j e^{-M_j(\vec{g})} e^{-i\vec{g} \cdot \vec{r}_j} \sum_{i=1}^4 a_i e^{\left(-\frac{1}{4} \left(\frac{b_i}{4\pi^2}\right) (ng)^2\right)}$$

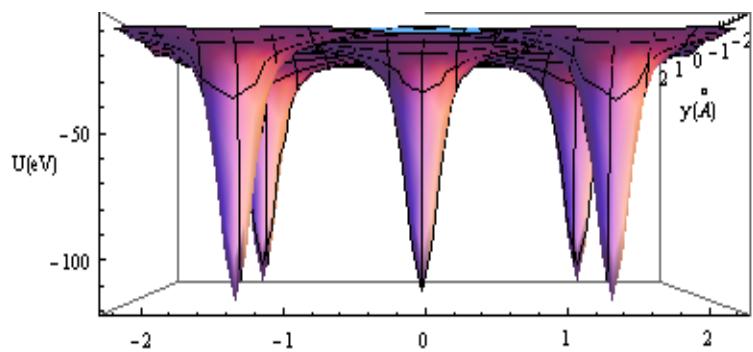
### Axial channeling

$$V(x, y) = \sum_{\vec{g}_m} v_{\vec{g}_m} e^{i\vec{g}_m \cdot \vec{r}_\perp}$$

$$v_{\vec{g}_m} = -\frac{2\pi}{V_c} a_0^2 (e^2 / a_0) \sum_j e^{-i\vec{g} \cdot \vec{r}_j} \sum_{i=1}^4 a_i e^{\left(-\frac{1}{4} \left(\frac{b_i}{4\pi^2} + 2\langle u_j^2 \rangle\right) |\vec{g}_m|^2\right)}$$



The planar continuum potentials of diamond for electrons



The <100> axial continuum potential of germanium for electrons

### 3. Theory of planar channeling radiation (Quantum)

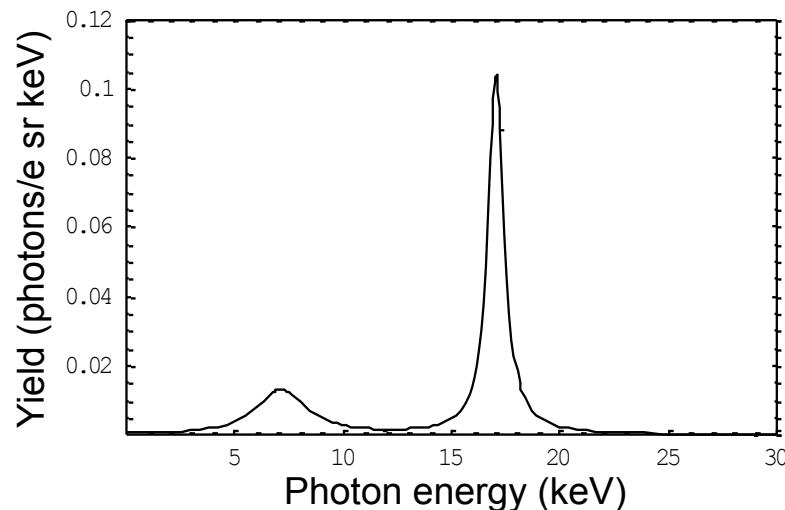
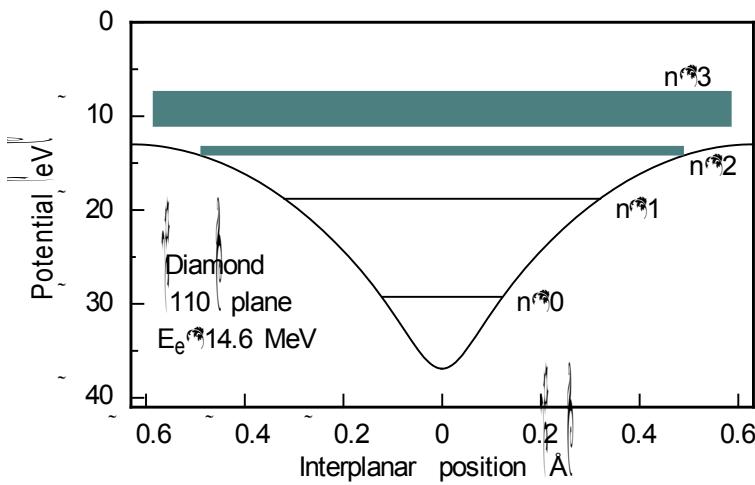
Quantum mechanical model

$E_e < 100 \text{ MeV}$

$$-\frac{\hbar^2}{2m_e\gamma} \frac{d^2\psi(x)}{dx^2} + V(x)\psi(x) = E\psi(x) \longrightarrow \text{Wave functions } \psi_i(x) \text{ and eigenvalues } E_i$$

$$E_0 = 2\gamma^2(E_i - E_f)$$

$$\frac{d^2N_{CR}(i \rightarrow f)}{d\Omega_\gamma dE_\gamma} = \frac{\alpha\lambda_c^2}{\pi\hbar c} 2\gamma^2(E_i - E_f) \left| \left\langle \psi_f(x) \left| \frac{d}{dx} \right| \psi_i(x) \right\rangle \right|^2 \int_0^z dz P_i(z) \times \frac{\Gamma_{\text{tot}}/2}{(E_\gamma - E_0)^2 + 0.25\Gamma_{\text{tot}}^2}$$



A Mathematica package for calculation of planar channeling radiation spectra of relativistic electrons channeled in a diamond-structure single crystal (quantum approach)

## 4. Theory of planar channeling radiation (classical) dipole approximation

Classical model       $E_e > 100\text{MeV}$

$$\text{Planar : } \gamma m \ddot{x}(t) = F = -\frac{\partial V(x)}{\partial x}$$

$$\text{Angular-energy distribution: } \frac{d^2 E}{d\omega d\Omega} = \frac{e^2}{4\pi^2 c} \left| \int_0^\tau e^{i(\omega t - \vec{k} \cdot \vec{r})} \frac{\vec{n} \times ((\vec{n} - \vec{\beta}) \times \vec{\beta})}{(1 - \vec{\beta} \cdot \vec{n})^2} dt \right|^2$$

$$c\vec{\beta} = \dot{\vec{r}}(t) \quad \quad \quad \vec{k} = \omega \vec{n} / c$$

$$\vec{r}(t) = x(t)\vec{e}_x + ct\vec{e}_z$$

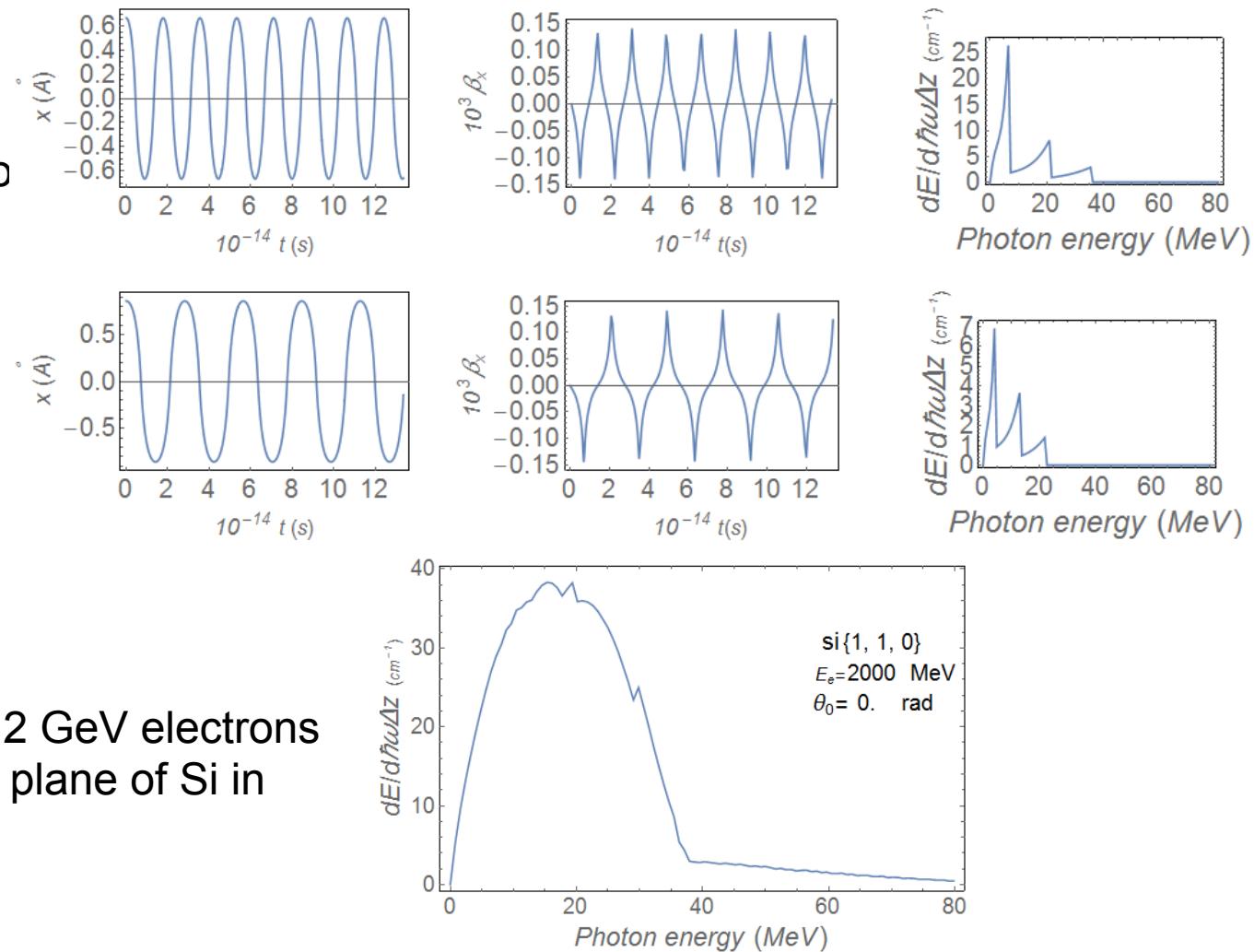
Total radiated energy  
in thin crystal:

$$\frac{dE}{d\omega \Delta z} = \frac{e^2}{c^4 T^2} \sum_{n=1}^{\infty} \Theta[1 - \eta_n] (\eta_n^2 - \eta_n + \frac{1}{2}) \cdot |\dot{x}_{\tilde{\omega}}|^2$$

$$\eta_n = \frac{T\omega}{4\pi\gamma^2 n}; \quad \tilde{\omega} = \frac{2\pi n}{T}; \quad \dot{x}_{\tilde{\omega}} = \int_0^T \dot{x} e^{i\tilde{\omega}t} dt$$

## 4. Dipol approximation

Trajectories, velocities and CR spectra for two different incidence points of 2 GeV electrons to (110) plane of a Si crystal.



Radiation spectrum of 2 GeV electrons channeled along (110) plane of Si in dipol approximation.

Simulation of planar channeling-radiation spectra of relativistic electrons and positrons channeled in a diamond-structure or tungsten single crystal (classical approach)

Paper: Nucl. Instrum. Methods B 342 (2015) 144

Program: Classical Planar Channeling Radiation Package <http://prof.bsu.ac.ir/~azadegan/>

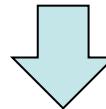
## 5. Non-dipole approximation

$$m\gamma_0 \ddot{x} = F_x = -\frac{\partial U(x)}{\partial x} \quad \gamma_0 = 1/\sqrt{1 - v^2/c^2}$$

At relativistic energies the longitudinal velocity component is coupled with the transverse component through  $\implies \gamma = \sqrt{1 - (v_x^2 + v_z^2)/c^2}$

conservation law for the longitudinal momentum component

$$d/dt(\gamma m v_z) = 0$$



$$v_z(t) = v_z(0) \sqrt{\frac{1 - v_x^2(t)/c^2}{1 - v_x^2(0)/c^2}} \approx v_z(0) \left[ 1 - \frac{1}{2c^2} (v_x^2(t) - v_x^2(0)) \right].$$

longitudinal component:

$$z(t) = \int_0^t v_z(t') dt' = v_z(0) \left( 1 + \frac{1}{2} \frac{v_x^2(0)}{c^2} \right) t - \frac{1}{2} \frac{v_z(0)}{c^2} \int_0^t v_x^2(t') dt'$$

## 5. Non-dipole approximation

$$\langle \dot{z} \rangle = \bar{\beta}_z c = \frac{1}{T} \int_0^T \dot{z}(t) dt$$

$$\vec{r}(t) = x(t)\vec{e}_x + (\bar{\beta}_z ct + z(t))\vec{e}_z \quad \begin{matrix} x(t) \text{ and } z(t) \text{ are periodic with period } T \\ \text{and } T/2, \text{ respectively} \end{matrix}$$

$$\frac{d^2 E}{d\omega d\Omega} = \frac{e^2 \omega^2}{4\pi^2 c} \left| \int_0^\tau \vec{n} \times \vec{\beta} e^{i(\omega t - \vec{k} \cdot \vec{r})} dt \right|^2$$

$$\frac{d^2 \bar{E}}{d\omega d\Omega} = \frac{e^2 \omega^2}{2\pi c^2} \sum_{n=1}^{\infty} \frac{I_n}{(1 - \bar{\beta}_z \cos(\vartheta))} \delta\left(\omega - \frac{\omega_n}{(1 - \bar{\beta}_z \cos(\vartheta))}\right) \quad \omega_n = 2\pi n/T$$

$$I_n = |a_n|^2 - |\vec{n} \cdot \vec{a}_n|$$

$$\vec{a}_n = \frac{1}{T} \int_0^T \vec{\beta}(t) \exp\left(-i \frac{\omega}{c} [x(t) \cos(\varphi) \sin(\vartheta) + z_p(t) \cos(\vartheta)]\right) e^{i\omega_n t} dt$$

## 5. Non-dipole approximation

The frequency spectrum is obtained by integration over all emission angles  $\vartheta$  and  $\varphi$ . Integration over angle  $\varphi$  can be taken easily.

$$\begin{aligned}\vec{a}_n &= \frac{1}{T} \int_0^T \vec{\beta}(t) e^{i \frac{2\pi n t}{T}} \int_0^{2\pi} \exp \left( -i \frac{\omega}{c} [x(t) \cos(\varphi) \sin(\vartheta) \right. \\ &\quad \left. + z_p(t) \cos(\vartheta)] \right) d\varphi dt \\ &= \frac{2\pi}{T} \int_0^T \vec{\beta}(t) J_0 \left( \frac{\omega x(t) \sin(\vartheta)}{c} \right) \exp \left[ i \left( \omega_n t - \frac{\omega z(t) \cos(\theta)}{c} \right) \right] dt\end{aligned}$$

$J_0(x)$  is the zero order Bessel function and integration over time must be done numerically.

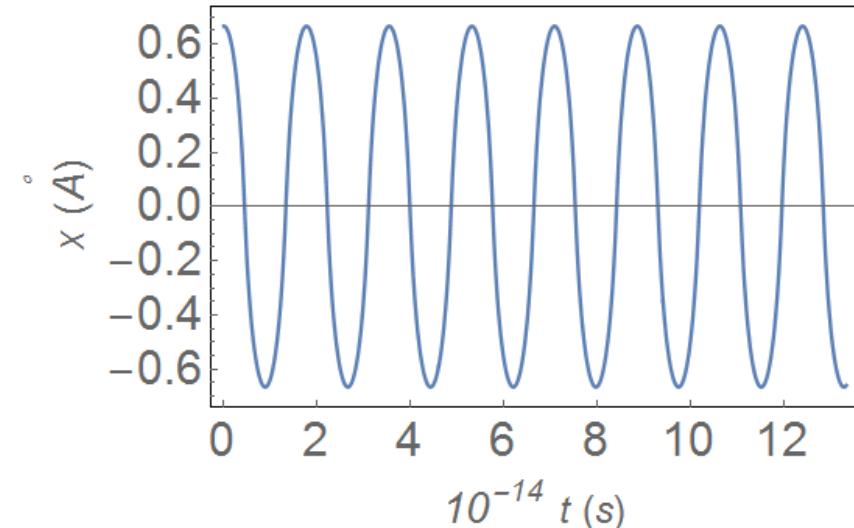
Due to  $\delta$ -function under the integral over angle  $\vartheta$ , the spectrum is restricted by two limits  $\omega_{\min} = \omega_n / (1 + \beta)$  and  $\omega_{\max} = \omega_n / (1 - \beta)$ ,

$$\begin{aligned}\frac{d^2 \bar{E}}{d\omega^2} &= \frac{e^2 \omega^2}{2\pi c^2} \sum_{n=1}^{\infty} \int_{-1}^1 \frac{I_n(\vartheta)}{\left(1 - \bar{\beta}_z \cos(\vartheta)\right)} \delta\left(\omega - \frac{\omega_n}{\left(1 - \bar{\beta}_z \cos(\vartheta)\right)}\right) d(\cos(\vartheta)) \\ &= \frac{e^2 \omega}{2\pi c^2} \sum_{n=1}^{\infty} I_n(\theta_m) \quad \theta_m = \cos^{-1}\left(\frac{1 - \omega/\omega_n}{\bar{\beta}_z}\right) \quad \omega_{\min} \leq \omega \leq \omega_{\max}\end{aligned}$$

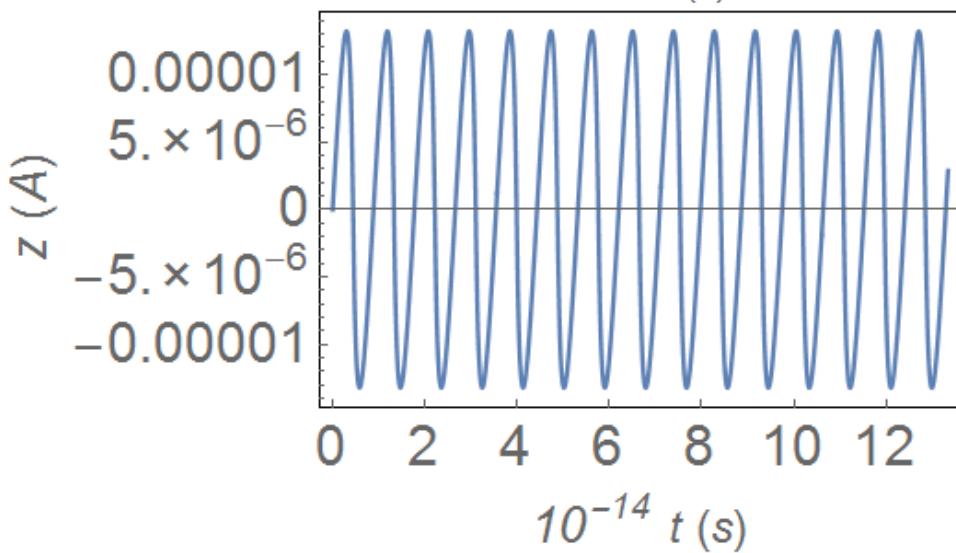
## 5. Non-dipole approximation

2 GeV electron  
(110) plane of Si

Transversal component

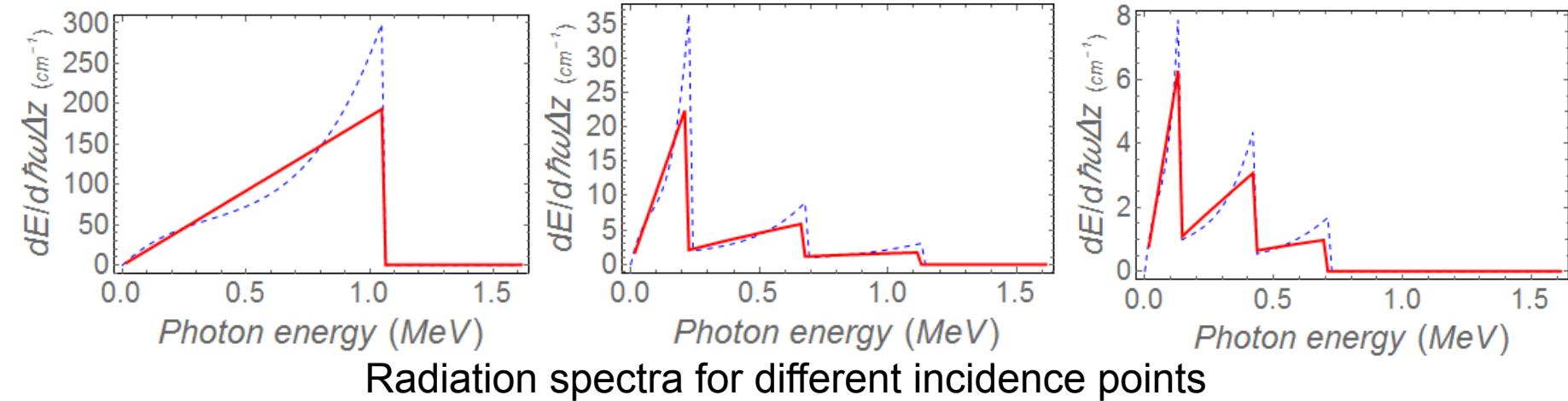


Longitudinal component



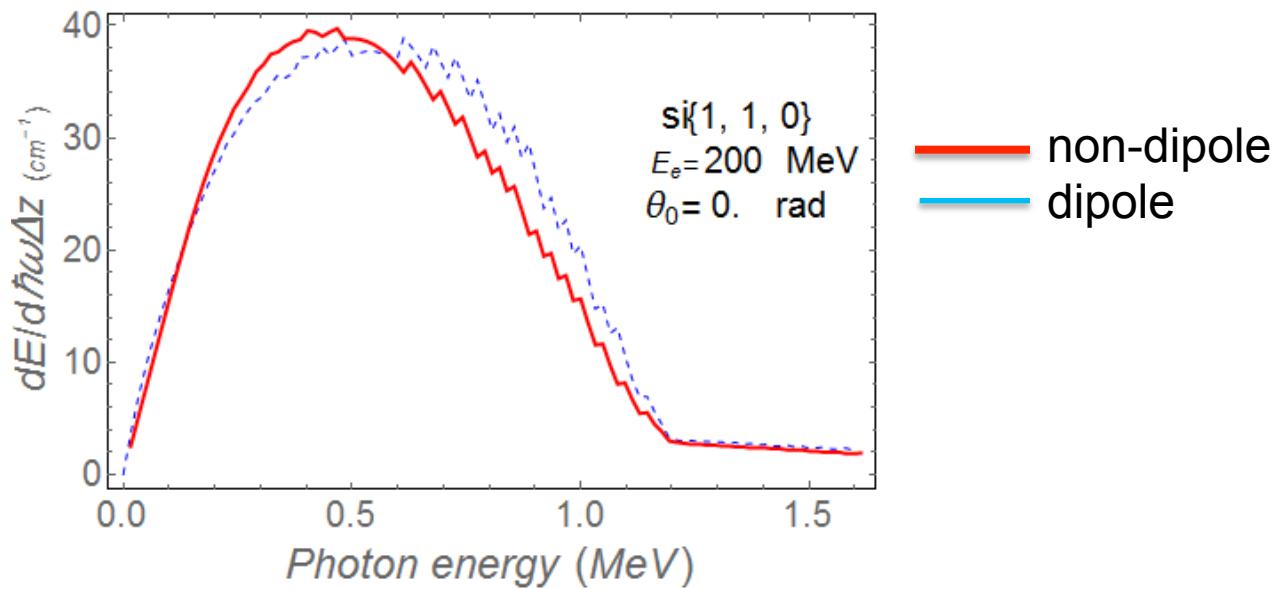
## 5. Comparison of non-dipole with dipol approximation

$E_e = 200 \text{ MeV}$

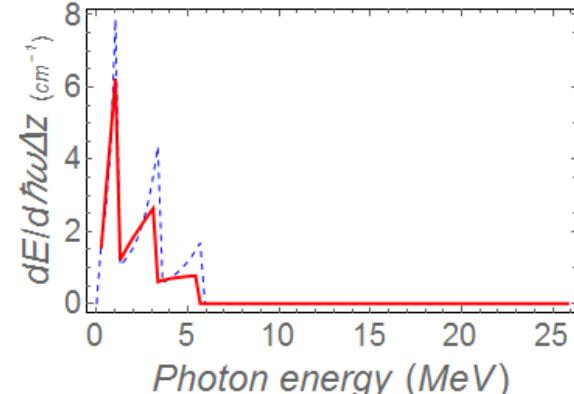
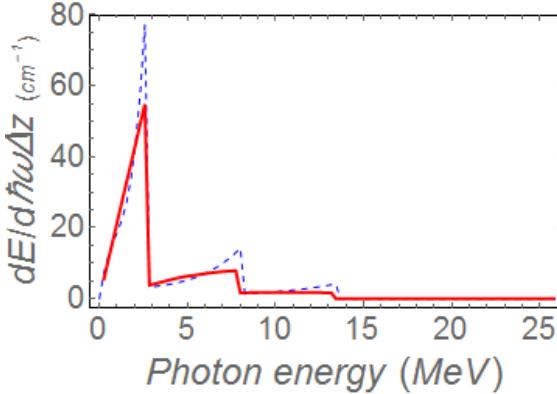
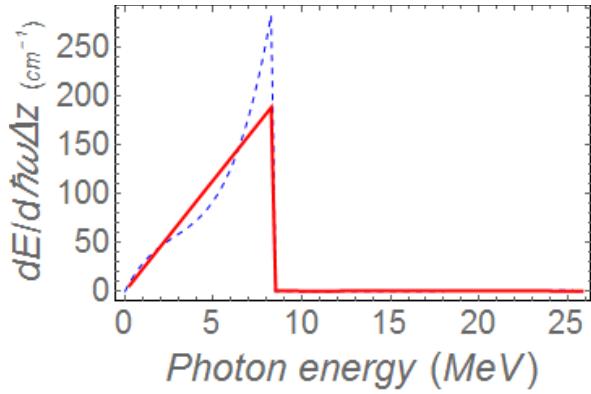


Radiation spectra for different incidence points

Total radiation spectra  
for 200 MeV electrons  
Si (110) plane

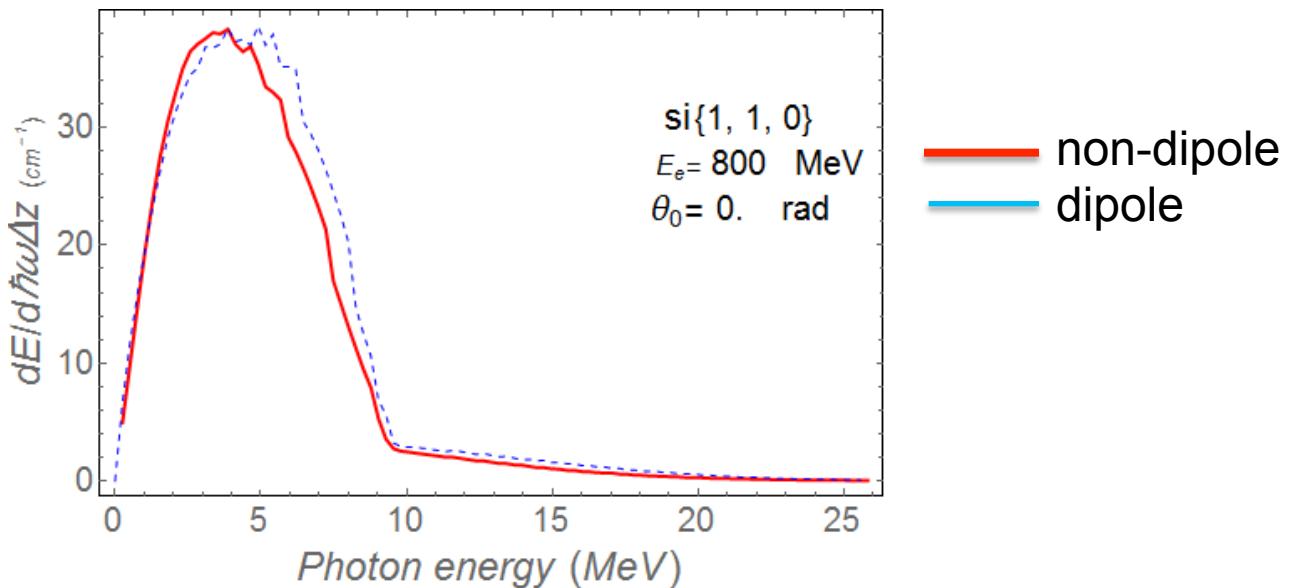


## 5. Comparison of non-dipole with dipol approximation $E_e=800$ MeV

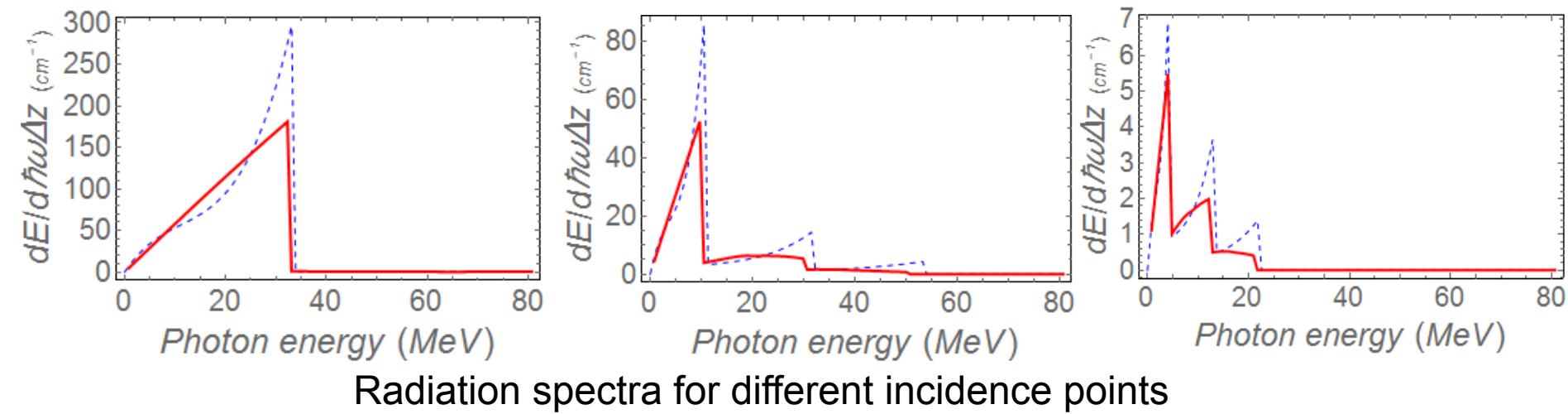


Radiation spectra for different incidence points

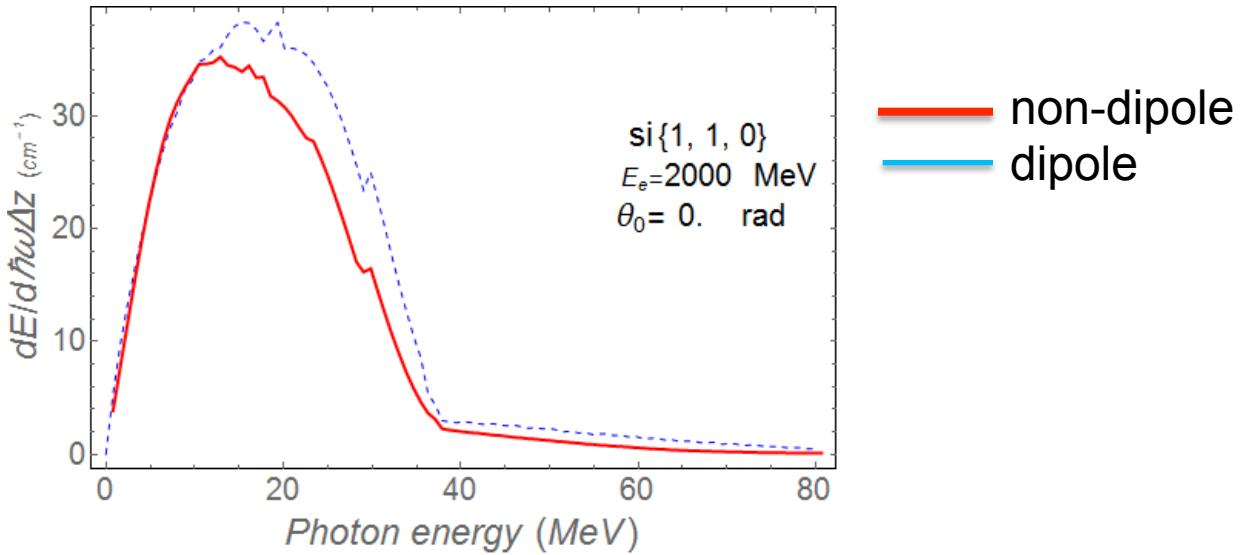
Total radiation spectra  
for 800 MeV electrons  
Si (110) plane



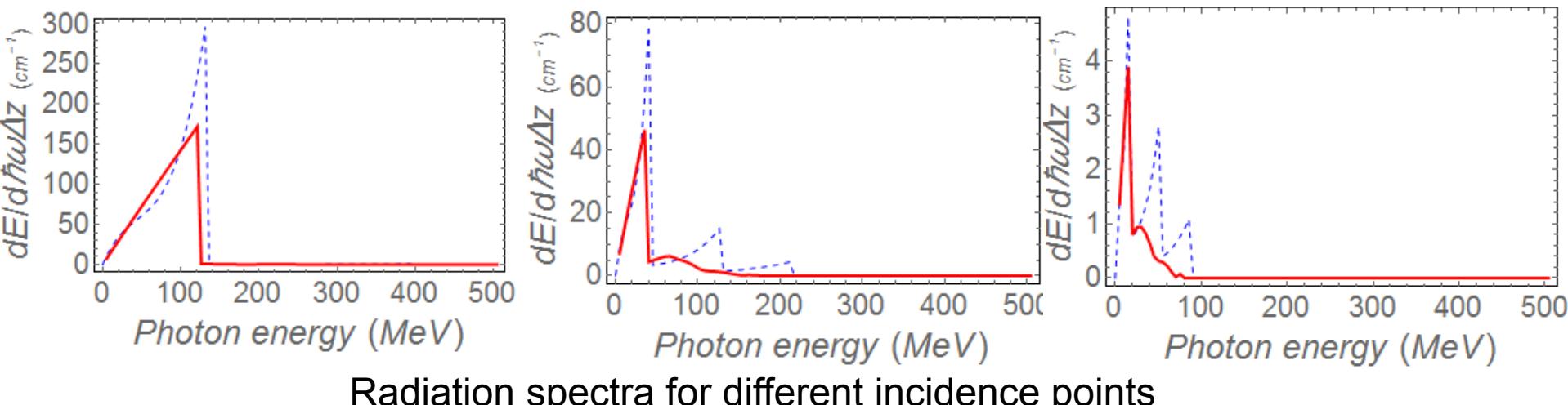
## 6. Comparison of non-dipole with dipol approximation E<sub>e</sub>=2 GeV



Total radiation spectra  
for 2 GeV electrons  
Si (110) plane

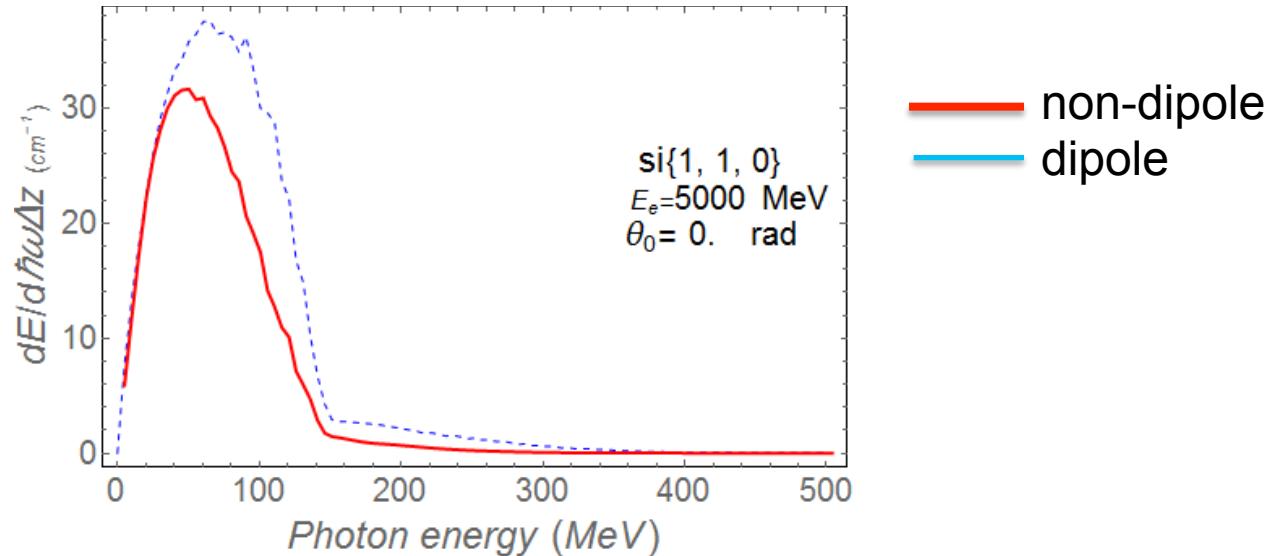


## 6. Comparison of non-dipole with dipol approximation $E_e=5 \text{ GeV}$



Radiation spectra for different incidence points

Total radiation spectra  
for 5 GeV electrons  
Si (110) plane



## 7. Summary

- We treated planar as well as axial channeling radiation at different energies and developed several software codes (*Mathematica*) appropriate for quantum as well as for classical calculations. Users can download the codes from the internet.
- We investigated the influence of non-dipolarity of channeling radiation. This effect can not be neglected at beam energies larger than about 1 GeV.
- This effect is also important for the simulation of positron production by means of channeling radiation.

# **Thank you**